

Overview of Comments

| | |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----|
| ■ Frank wanted to discuss the title (M) | 1 |
| ■ If we introduce this abbreviation, should we replace the uses of “control-flow graph” with “CFG” in the paper? (M) | 1 |
| ■ The IDE paper already states that each IFDS problem can be solved with IDE. I wonder if this sounds like the equivalence transformation was our contribution. We just use the equivalence transformation to explain the CC transformation, and for the proofs. (M) | 2 |
| ■ Do we need to mention the imprecise map from the CC-IDE result to the IFDS result? It doesn’t really seem relevant, and could be confusing. (O) | 2 |
| ■ would it make sense to add an edge in Figure 1 from the IFDS problem to the IFDS result? (F) | 2 |
| ■ This section is current under revision. Please do not edit.. (F) | 3 |
| ■ The figure that we deleted was constructed so that there would be a realizable and a non-realizable path. Now that we deleted that figure, let’s delete this example as well. (M) | 6 |
| ■ The four references take away space in the References section. Remove some of them? (M) | 7 |
| ■ I’m not saying any more that IFDS is a version of Sharir&Pnueli’s functional approach. (M) | 7 |
| ■ I also removed the comparison of IFDS with other algorithms. I’m not pointing out the benefits of IFDS any more, so we might fail to show the importance of the algorithm. (M) | 7 |
| ■ Provide reference explaining context sensitivity? (M) | 7 |
| ■ Tell the reader here to ignore the labels on the edges of the graph? (M) | 8 |
| ■ We never use the interpretation function, so I deleted it, too (M) | 9 |
| ■ I think the rest of this section will not make sense to anyone who does not already understand it. Should I try to replace it with one paragraph that tries to give an intuitive understanding of what the IDE algorithm is doing? (M) | 11 |
| ■ In Figure 3, should we say that ψ corresponds to the function’s return value? (M) | 12 |
| ■ we agreed to move this figure to the initial example section (O) | 12 |
| ■ Move Figure 3 up to where it’s mentioned first? If yes, the caption won’t make sense. (M) | 12 |
| ■ The following definitions are very general, and not related to the transformation. Can we move them somewhere earlier? (O) | 12 |
| ■ incorporate into previous section (O) | 14 |

| | |
|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----|
| ■ Here, you defined the output of a transformation as just $G\#$ and an edge function. Later, you also include the lattice. Should we include the lattice here? (O) | 16 |
| ■ What if r points to null? (O) | 16 |
| ■ define $n_1(e)$ globally somewhere early on, and consider using a more decriptive name like src instead of n_1 (O) | 17 |
| ■ move material from following section here (O) | 18 |
| ■ Prove this lemma in the appendix. The parts of the proof can be found in section 5. (O) | 19 |
| ■ The normalization function \mathcal{N} was only needed in composition. I therefore “inlined” it into composition to save space. This makes composition a bit harder to intuitively verify, but saves complexity here. The verification can be in the proof in the supplementary appendix. (O) | 19 |
| ■ Prove this lemma in the appendix. The parts of the proof can be found in section 5. (O) | 19 |
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| ■ Sagiv et al. define a notion of an “efficient” implementation of micro-functions. If the micro-functions are “efficient”, then the asymptotic complexity bounds proven by Sagiv hold. Our implementation is “efficient” if we assume that the number of correlated receivers is a constant independent of the size of the program. That seems a stretch. If it’s linear in the size of the program, then we would need to add a factor to their complexity bounds. Their bound is $O(ED^3)$, while ours would be $O(ED^3R)$. But then we can’t directly use their complexity proof; we would have to adapt it. I’m leaning towards leaving out any mention of “efficient” implementation. (O) | 19 |
| ■ Should we say \mathcal{R}_{IDE} instead of \mathcal{R} everywhere, to be consistent with \mathcal{R}_{IFDS} ? (O) | 19 |
| ■ should this be a section or subsection? should it be moved somewhere else (or maybe not included at all)? (O) | 20 |
| ■ We should incorporate the four functions into the explanation of IFDS in the background. Then we don’t need them here. (O) | 21 |
| ■ We need to decide what exactly we want to include in the evaluation. The numbers of correlated calls are interesting. But I think we want to leave out all mention of the taint analysis. Perhaps we should have no empirical results at all (and only the theory/proofs)? (O) | 26 |
| ■ We should convert this section into two or three sentences to be added to the conclusion. They should focus especially on the interprocedurally-correlated receivers. (O) | 28 |

Data Flow Analysis in the Presence of Correlated Calls

Marianna Rapoport¹, Ondřej Lhoták¹, and Frank Tip²

¹ University of Waterloo
{mrapoport, olhotak}@uwaterloo.ca
² Samsung Research America
ftip@samsung.com

Frank wanted to discuss the title (M)

Abstract. We present a technique to improve the precision of data-flow analyses on object-oriented programs in the presence of *correlated calls*. Two method calls are correlated if they are polymorphic and are invoked on the same object. Correlated calls are problematic because they can make existing data-flow analyses consider certain infeasible data-flow paths as valid. This leads to loss in precision of the analysis solution. We show how infeasible paths can be eliminated for *Inter-procedural Finite Distributive Subset* (IFDS) problems, a large class of data-flow analysis problems. We show how the precision of IFDS problems can be improved in the presence of correlated calls, by using the *Inter-procedural Distributive Environment* (IDE) algorithm to eliminate infeasible paths. Using IDE, we eliminate the infeasible paths and obtain a more precise result for the original IFDS problem. Our analysis is implemented in Scala, using the WALA framework for static program analysis on Java bytecode.

1 Introduction

Data-flow analysis computes an approximation of how values may flow through a program, and has applications in compiler optimization, programming tools, and computer security, and many other areas. Data-flow analyses operate on *control-flow graphs* (CFGs) that model the order in which the instructions of a program are executed. This is typically done by associating *flow functions* that represent how data is propagated with the edges of the control-flow graph. A *confluence operator* specifies how the data facts that have been computed along different paths should be merged when the paths join.

Since a control-flow graph is an over-approximation of the possible flows of control in concrete executions of a program, it may contain *infeasible* paths that cannot occur at runtime. The precision of a data-flow analysis algorithm depends critically on its ability to detect and disregard such infeasible paths. The popular *Interprocedural Finite Distributive Subset* (IFDS) algorithm by Reps, Horwitz, and Sagiv [17] is a general data-flow analysis algorithm for computing solutions to finite distributive data-flow problems such as reaching definitions, available

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expressions, and taint analysis. A distinguishing characteristic of IFDS is that it avoids infeasible interprocedural paths in which calls and returns to/from functions are not properly matched. Sagiv, Reps, and Horwitz also presented the *Interprocedural Distributive Environment* (IDE) algorithm [20] that similarly only considers properly matched call/return edges, but that supports a broader range of dataflow problems by expanding the domain of flow functions to *environments* that go beyond the data-flow facts considered by IFDS.

This paper presents an approach to dataflow analysis that avoids a type of infeasible path that arises in object-oriented programs when two or more methods are dynamically dispatched on the same receiver object. In such cases, if the method calls are polymorphic (i.e., if they dispatch to different method definitions depending on the type of the receiver expression at run time), then their dispatch behaviors will be correlated. A recent paper [?] identified this problem but did not present a concrete solution or algorithm, and we are not aware of any existing dataflow analysis that is capable of avoiding infeasible paths that arise in the presence of correlated method calls.

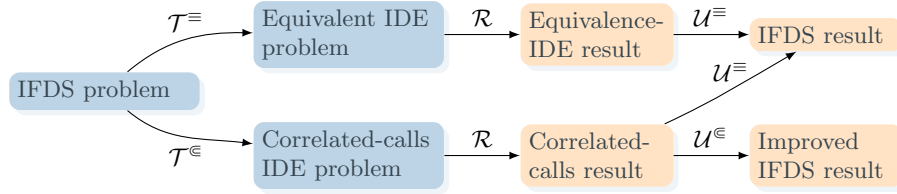


Fig. 1: Transformations between IFDS and IDE problems and their results

The approach taken in our work is to transform a standard IFDS problem into an IDE problem that precisely accounts for infeasible paths due to correlated calls. The results of this IDE problem can be mapped back to the dataflow domain of the original IFDS problem.

We present a formalization of the transformation and prove its correctness as follows. First, we derive an “Equivalence IDE problem” from the original IFDS problem by associating the identity environment functions with all edges, and show that the solution to this Equivalence IDE problem can be mapped back to the solution of the original IFDS problem. Then, we derive a “Correlated Calls IDE Problem” from the original IFDS problem and show how a solution to this problem can be mapped to the solution to the original IFDS problem, but also how a more precise IFDS result can be derived from it. We also show that the correlated-calls analysis is sound, i.e., that it never considers concrete execution paths as infeasible. This is illustrated schematically in Figure 1³.

³ The labels \mathcal{R} , $\mathcal{U}^=$, $\mathcal{U}^€$, $\mathcal{T}^=$, and $\mathcal{T}^€$ on edges in Figure 1 reflect a number of mappings and projections that will be defined in Section ▶ref◀ and that can be ignored here.

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We implemented the correlated-calls transformation and the IDE algorithm in Scala, on top of the WALA framework for static analysis of JVM bytecode [6]. We also report on preliminary experiments in which our correlated-calls transformation is applied to an IFDS formulation of a simple taint analysis. Our results show that solving the resulting IDE problem avoids infeasible paths due to correlated calls as expected.

In summary, the contributions of this paper are as follows:

- We present a general approach for transforming IFDS problems into corresponding IDE problems that avoid infeasible paths due to correlated method calls and prove its correctness.
- We implemented the approach in Scala, on top of the WALA program analysis framework and report on preliminary experiments.

The remainder of this paper is organized as follows. Section 2 presents a motivating example. Section 3 reviews the IFDS and IDE algorithms. Section ?? presents the correlated-calls transformation and a proof of its correctness. The implementation of our approach and preliminary experiments are discussed in Section 5. Related work is discussed in Section ▶ref◀. Finally, conclusions and directions for future work are presented in Section ▶ref◀.

2 Motivating Example

Consider a call site $r.m()$ in an object-oriented programming language, where the variable r is the *receiver* variable of the call site and m is the name of the invoked method⁴. In the rest of the paper, we use the general term *receiver* to mean a receiver variable. At runtime, the actual method that will be invoked by the call site depends on the runtime type of the object referenced by r . If the call site $r.m()$ can be associated with more than one method at compile time, we will say that the call site is *polymorphic*.

For example, in Listing 2, it is not possible to infer statically whether the runtime type of the variable `a` in the `main` method is `A` or `B`. The call `a.foo()` can be dispatched to either `A.foo` or `B.foo`, and `a.bar(v)` can be dispatched to either `A.bar` or `B.bar`. A concrete execution path for the `main` method might therefore go through `A.foo` and `A.bar`, or through `B.foo` and `B.bar`. However, there cannot be an execution path through `A.foo` and `B.bar` or through `B.foo` and `A.bar`.

We call the invocations to methods `foo` and `bar` *correlated*. More generally, correlated calls occur when more than one polymorphic call is invoked on the same receiver variable.

Suppose we wanted to perform a taint analysis on the program in Listing 2. Most dataflow-analysis algorithms, including IFDS, would conservatively assume

⁴ We assume an internal representation of the program in which for each call site $e_r.m()$, the expression e_r has been evaluated to the variable r .

```
class A {  
    String foo {  
        return secret();  
    }  
  
    void bar(String s) {}  
}  
  
class B extends A {  
    String foo {  
        return "not_secret";  
    }  
  
    void bar(String s) {  
        System.out.println(s);  
    }  
}  
  
class Main {  
    public static void main(String[] args) {  
        A a = args == null ? new A() : new B();  
        String v = a.foo();  
        a.bar(v);  
    }  
}
```

Fig. 2: Example program containing correlated calls

that the call `a.bar` could be dispatched to both `A.bar` and `B.bar`, independently of what `a.foo` had been dispatched to in the previous line.

As a result, such an analysis would consider a path through `A.foo` and `B.bar` feasible. This means that the variable `v` would be considered secret. We would conclude that a secret value is passed to `B.bar` and printed to the user. In other words, we would consider the program to leak secret information, which it does not do in any concrete execution.

3 Background

The purpose of the correlated-calls analysis is to solve IFDS problems more precisely than using the standard IFDS algorithm by ruling out some infeasible paths. The correlated-calls analysis works by transforming an IFDS problem to an IDE problem, solving the IDE problem, and transforming the IDE result to a solution to the original IFDS problem. This section describes the general ideas underlying IFDS and IDE.

3.1 Terminology and Notation

We will start by introducing several concepts used by the IFDS and IDE analyses.

A *control-flow graph* is a directed graph in which nodes correspond to instructions and edges represent transfer of control between the instructions during an execution of the program. A control-flow graph has a unique start node, `startmain`, which is the node corresponding to the program entrypoint.

An *intra-procedural* path is a path in a control-flow graph whose nodes are in the same procedure. By contrast, an *inter-procedural path* is one that contains nodes from different procedures.

A *control-flow supergraph* is a control-flow graph in which each procedure p is augmented with an additional *start node* `startp` and *end node* `endp`, and for each call c_q to a procedure q , there is a *call node* `callcq` and subsequent *return node* `returncq`.

A control-flow supergraph allows us to model the control flow in inter-procedural paths. The flow from the caller to the callee is represented using an edge

$$(\text{call}_{c_q}, \text{start}_q).$$

The control flow from the callee back to the caller goes through an edge

$$(\text{end}_q, \text{return}_{c_q}).$$

We will denote the source and end nodes of an edge e as `src(e)` and `end(e)`.

A *flow-sensitive* data-flow analysis is one that takes the order of program instructions into account.

Let each call node in a program be labeled with a distinct opening parenthesis and the corresponding return node with the matching closing parenthesis. For a given path p , let s be the string that is obtained by concatenating the labels

of the nodes in p . Then p is *valid* if s belongs to the language of substrings of balanced parentheses. The set of all inter-procedurally valid paths from the start node to a node n is denoted as $\text{VP}(n)$. The set $\text{VP}(n)$ is a conservative approximation of all concrete execution paths from the start node to n .

A *context-sensitive* data-flow analysis is an analysis that considers only inter-procedurally valid paths.

Example 1. In the supergraph in Figure ??, let us assign $\{, \}$ parentheses to $\text{call}_{A.f}$ and $\text{return}_{A.f}$, and \langle, \rangle parentheses to call_f and return_f . Then the string corresponding to the path

$$p_1 = [\text{call}_{A.f}, \text{start}_f, \text{if } (s == \text{null}), \text{return } s, \text{end}_f, \text{return}_{A.f}]$$

is $\{\}$, which indicates that p_1 is valid. Every prefix of p_1 is also a valid path.

However, the graph also contains an inter-procedurally invalid path

$$p_2 = [\text{call}_f, \text{start}_f, \text{if } (s == \text{null}), \text{return } s, \text{end}_f, \text{return}_{A.f}]$$

with corresponding string $\langle \rangle$.

A *lattice* is a partially ordered set in which each subset has a least upper bound and a greatest lower bound.

A *meet semilattice* $L = (S, \sqcap)$ is defined by a set S and a meet operation \sqcap that is associative, commutative, and idempotent. The meet operation induces a partial order (S, \sqsubseteq) where every subset contains a greatest lower bound: For all $x, y \in S$, $x \sqsubseteq y$ if $x \sqcap y = x$. The greatest lower bound, or top element, of the semilattice is denoted as \top . If k is the length of the longest chains of elements in the semilattice, then the *height* of the semilattice is $k - 1$.

In this work, we will denote a map from a set of keys K to values from set V as

$$\{(k, v) \mid k \in K, v \in V\}. \quad (1)$$

For an arbitrary map m , $m(x)$ is the value to which x is mapped in m . We denote by $m[x \rightarrow y]$ a map identical to m , except that the element x is mapped to y . To avoid excessive parentheses, we write $(m[x_1 \rightarrow y_1])[x_2 \rightarrow y_2]$ as $m[x_1 \rightarrow y_1][x_2 \rightarrow y_2]$.

We will denote the identity function $\lambda x. x$ by id . We will use a typed version of this function in various contexts, where the type of x will vary with the context.

Finally, we introduce the notion of *distributivity*. Given a set D , a function $f : 2^D \rightarrow 2^D$ is distributive if $\forall x_1, x_2 \in 2^D$,

$$f(x_1 \cup x_2) = f(x_1) \cup f(x_2). \quad (2)$$

The figure that we deleted was constructed so that there would be a realizable and a non-realizable path. Now that we deleted that figure, let's delete this example as well. (M)

3.2 IFDS

The IFDS framework is a precise and efficient algorithm for data-flow analysis. IFDS was developed in 1995 by T. Reps, S. Horwitz, and M. Sagiv at the University of Wisconsin and has been used to solve a variety of data-flow analysis problems [4,13,10,24].

The IFDS algorithm is applicable to problems which can be expressed with data-flow functions that satisfy certain restrictions. *Inter-procedural* flow functions specify how data flows from the invocation of a procedure to its start, and from the procedure’s end back to its call site. *Distributive* flow functions are those that distribute over the confluence operator. In the context of IFDS, the confluence operator is called *meet*, and it can be either union or intersection. The data-flow facts on which the analysis operates must be a *finite* set D . Each flow function operates on a *subset* of D (for example, the set of variables in the program) which makes the domain of the flow functions the power set of D .

Given a data-flow problem that satisfies the restrictions of IFDS, the algorithm provides a *context-sensitive* solution in polynomial time. Compared to IFDS, most data-flow analyses are either general but do not run in polynomial time [8,21] or handle a very specific set of problems [9].

Data-Flow Problems Suitable for IFDS Informally, an IFDS analysis can only solve decision problems. An IFDS analysis answers questions of the following kind: “is property X true at program point Y ?”. For example, a taint-analysis problem asks, for each variable v in the program, “is v secret at a given program point?”.

Formally, a data-flow analysis problem is suitable for an IFDS analysis if it can be encoded as an IFDS problem

$$(G^*, D, F, M_F, \sqcap), \quad (3)$$

where $G^* = (N^*, E^*)$ is the supergraph of the input program with nodes N^* and edges E^* , D is a finite set of *data-flow facts*, F is a set of distributive dataflow functions of type $2^D \rightarrow 2^D$, $M_F : E^* \rightarrow F$ is a function that maps supergraph edges to dataflow functions, and M_F is extended to paths by composition⁵. The *meet operator* \sqcap is either union or intersection.

Without loss of generality, we will take *meet* to denote union. It can be shown that any problem where *meet* is defined as intersection can be reformulated into an equivalent one where *meet* is defined as union [17].

Overview of the IFDS Algorithm Given an IFDS problem, for each node $n \in N^*$ the IFDS algorithm computes the *meet-over-all-valid-paths* solution

$$\text{MVP}_F(n) = \bigsqcap_{q \in \text{VP}(n)} M_F(q)(\emptyset). \quad (4)$$

⁵ Let A be a set and $f : E^* \rightarrow (A \rightarrow A)$ a function from supergraph edges to functions on A . We say that f is extended to paths by composition to denote that for a path q consisting of the edges e_1, \dots, e_k , $f(q) = f(e_k) \circ \dots \circ f(e_1) \circ \text{id}$.

The four references take away space in the References section. Remove some of them? (M)

I’m not saying any more that IFDS is a version of Sharir&Pnueli’s functional approach. (M)

I also removed the comparison of IFDS with other algorithms. I’m not pointing out the benefits of IFDS any more, so we might fail to show the importance of the algorithm. (M)

Provide reference explaining context sensitivity? (M)

To compute the meet-over-all-valid-paths solution, each node in the control-flow supergraph is paired with a *fact* $d \in D \cup \{\mathbf{0}\}$, $\mathbf{0} \notin D$, yielding the nodes $N^\#$ of the *exploded supergraph* $G^\# = (N^\#, E^\#)$. Roughly, for each node in the program, a fact denotes a binary property whose value (true or false) we want to find out. The start node of the exploded supergraph is the node $(\text{start}_{\text{main}}, \mathbf{0})$.

The flow functions F define the edges of the exploded supergraph. Using the flow functions, the IFDS algorithm computes the inter-procedurally *realizable* paths from the start to the rest of the exploded graph's nodes. A *realizable* path is a valid path in the exploded supergraph that starts with the entry node $\text{start}_{\text{main}}$.

If there is a realizable path from the node $(\text{start}_{\text{main}}, \mathbf{0})$ to a given node (n, d) , $d \neq \mathbf{0}$, then the fact d is considered to hold at node n . A path to a node $(n, \mathbf{0})$ means that in the control-flow supergraph, there is a path from $\text{start}_{\text{main}}$ to n .

In this way, the IFDS algorithm reduces the input data-flow problem to a graph-reachability problem.

Example 2. In a taint analysis, D is the set of variables in the program. If a fact $d \in D$ is reachable at a given node, then the variable is considered secret at that node. Otherwise, it is considered not secret. The question “is d secret at node n ?” becomes “is there a realizable path from $(\text{start}_{\text{main}}, \mathbf{0})$ to (n, d) ?”.

Example 3. The exploded supergraph for Listing 2 is shown in Figure 3. We can see that there is a realizable path from the start node of the exploded graph to the variable \mathbf{s} at the node $\text{print}(\mathbf{s})$ in the `B.bar` method. This means that at that node, \mathbf{s} is considered secret.

The flow functions $F \subseteq 2^D \rightarrow 2^D$ allow us to establish the edges in the exploded supergraph.

Given a control-flow-graph edge $e = (n_1, n_2) \in E^*$ and a distributive dataflow function $f = M(e)$, the *representation relation* $R_f : (D \cup \{\mathbf{0}\}) \times (D \cup \{\mathbf{0}\})$ of f is defined as

$$R_f = \{(\mathbf{0}, \mathbf{0})\} \cup \{(\mathbf{0}, d_j) \mid d_j \in f(\emptyset)\} \cup \{(d_i, d_j) \mid d_j \in f(\{d_i\}), d_j \notin f(\emptyset)\}.$$

Each pair $(d_i, d_j) \in R_f$ corresponds to an edge $((n_1, d_i), (n_2, d_j))$ in the exploded supergraph.

Note that R_f constructs pairs of dataflow facts so that

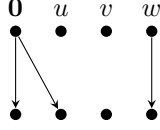
- there is always an edge $(\mathbf{0}, \mathbf{0})$ corresponding to the control-flow-graph edge;
- if there is an edge $(\mathbf{0}, d_j)$, then there is no other edge leading to d_j ; in particular, there is never an edge $(d_i, \mathbf{0})$ where $d_i \neq \mathbf{0}$.

Example 4. The representation relation R_f for a set of data-flow facts $D = \{u, v, w\}$ and dataflow function $f = \lambda S. S \setminus \{v\} \cup \{u\}$ looks as follows:

$$R_f = \{(\mathbf{0}, \mathbf{0}), (\mathbf{0}, u), (w, w)\}.$$

The corresponding exploded-graph edges are shown below.

Tell the reader here to ignore the labels on the edges of the graph? (M)



The representation relation lets us decompose a flow function into functions that operate on each fact individually. This is possible due to distributivity: we can apply the flow function on each single fact and take the union of the results, rather than applying the function to the union of the facts.

The representation relation allows us to compactly represent the composition and meet operations which are required for the IFDS algorithm.

For two representation relations R_{f_1} , R_{f_2} , the composition and meet operations are defined as follows:

$$R_{f_1} \circ R_{f_2} = \{(d_1, d_3) \mid \exists d_2 : (d_1, d_2) \in R_{f_1}, (d_2, d_3) \in R_{f_2}\}. \quad (5)$$

and

$$R_{f_1} \sqcap R_{f_2} = R_{f_1} \cup R_{f_2}. \quad (6)$$

The representation relation distributes over composition and meet:

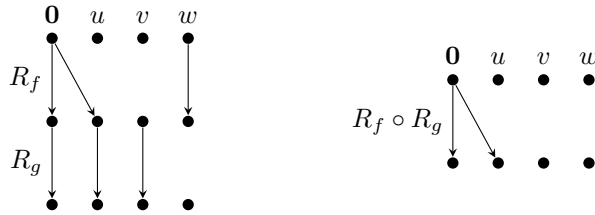
$$R_{f_1} \circ R_{f_2} = R_{f_1 \circ f_2}, \quad \text{and} \quad R_{f_1} \sqcap R_{f_2} = R_{f_1 \sqcap f_2}. \quad (7)$$

On the exploded graph, the composition of two functions is represented by the paths that are formed when the exploded-graph edges are combined.

Example 5. If $g = \lambda S. S \setminus \{w\}$ and f is defined as in example 4, then

$$R_f \circ R_g = \{(\mathbf{0}, \mathbf{0}), (\mathbf{0}, u)\},$$

as illustrated by the corresponding exploded graph edges:



We presented an overview of the IFDS analysis. IFDS problems are transformed into IDE problems by the correlated-calls analysis. The IDE framework is described in the next section.

We never use the interpretation function, so I deleted it, too (M)

3.3 IDE

There exists an entire class of data-flow problems that cannot be formulated as IFDS problems. Informally, the problems cannot be formulated as decision problems. For instance, a constant-propagation problem asks, for each variable v in the program, “if v is a constant at a given program point, what is v ’s value?”. The questions asked by constant propagation are of the form “if property X (*v being a constant*) is true at program point Y , what is the value of some property Z (*the value of the constant*) corresponding to X ?”. It turns out that problems with such questions can often be solved by the IDE algorithm.

The IDE framework is an expressive extension to IFDS that was created by the same authors in 1996. The problems that IDE can solve include, but are not limited to, IFDS problems [17]. Just as the IFDS algorithm, the IDE algorithm is suitable for data-flow analyses that can be encoded with inter-procedural, distributive flow functions. However, in IDE, the domain of the flow functions is not restricted to sets D of data-flow facts. The IDE domain of a flow function consists of *environments* that map data-flow facts from the set D to lattice elements.

As an example, in a constant propagation problem, an IDE environment would map each variable to the (possibly) constant value that it is bound to. To illustrate the distinction between IFDS and IDE we could say that IFDS can find out which variables in a program are constants, whereas IDE can additionally retrieve the values of the constant variables. Instead of just telling us whether a fact holds or not, the IDE analysis can provide us with additional information about facts.

Just as in the IFDS analysis, the IDE algorithm reduces a data-flow problem to a graph-reachability problem. Additionally, for each program point, the algorithm computes an *environment* $\text{Env}(D, L)$, where data-flow facts are mapped to values of a lattice L .

For example, using the IDE analysis, we can encode a restricted version of a constant-propagation analysis⁶. The data-flow facts correspond to program variables, and the lattice incorporates all possible values for constants. If a fact d in the exploded supergraph is reachable at node n , and $\text{Env}(d) \notin \{\perp, \top\}$, it means that the variable associated with d is a constant. Furthermore, the value of the constant can be inferred from the environment for the corresponding node and is equal to $\text{Env}(d)$.

Formally, an IDE problem is defined as a four-tuple

$$(G^*, D, L, M_{\text{Env}}), \quad (8)$$

where G^* is a control-flow supergraph, D is a set of data-flow facts, and L is a meet semilattice with finite height. Finally, $M_{\text{Env}} : E^* \rightarrow (\text{Env}(D, L) \rightarrow \text{Env}(D, L))$ is a function from the edges of the control-flow supergraph to distributive *environment transformers*. M_{Env} is extended to paths by composition.

⁶ In the general case, constant propagation cannot be encoded with distributive flow functions and is therefore not suitable for an IDE analysis [12].

Given an IDE problem, for each node $n \in N^*$ and fact $d \in D$, the IDE algorithm computes the meet-over-all-valid-paths solution

$$\text{MVP}_{\text{Env}}(n, d) = \bigcap_{q \in \text{VP}(n)} M_{\text{Env}}(q)(\Omega)(d), \quad (9)$$

where M_{Env} is extended to paths by composition and

$$\Omega = \lambda d. \top \quad (10)$$

is the top element in the environment lattice $\text{Env}(D, L)$.

Environment Transformers For each node in the control-flow graph, the result of an IDE analysis computes an environment $\text{Env}(D, L)$, which is a map from data-flow facts to lattice elements.

Instead of flow functions that show how to propagate facts, the IDE framework uses distributive environment transformers to propagate environments. For each edge (n_1, n_2) in the control-flow supergraph, an environment transformer indicates how the environment at node n_1 is modified at node n_2 .

From Section 3.2 we know that flow functions can be represented with exploded-graph edges. To represent environment transformers, we will construct *labeled* exploded-graph edges, where each edge is associated with a distributive *micro function*⁷ $f : L \rightarrow L$. A micro function shows how to change a lattice element for a given node and fact.

For instance, if an IDE problem is equivalent to an IFDS problem, the edges of the exploded supergraph are the same for both problems. In the IDE problem, the edges of the exploded supergraph are labeled with identity micro functions.

We extend the meet operator to work on micro functions by defining

$$(f_1 \sqcap f_2)(l) = f_1(l) \sqcap f_2(l) \quad (11)$$

for all $l \in L$.

In IDE problems, the auxiliary fact analogous to $\mathbf{0}$ in IFDS is denoted as Λ .

In this way, each edge in the exploded graph is labeled with a micro function. The mapping from exploded-graph edges to the corresponding micro functions is stored in *edge functions*, denoted as $\text{EdgeFn} : E^\# \rightarrow (L \rightarrow L)$.

Overview of the IDE Algorithm Given a labeled exploded supergraph, the IDE algorithm computes the environments for all nodes in the control-flow graph.

The algorithm first computes the lattice elements $l_{n,d}$ that correspond to each reachable node (n, d) in the exploded supergraph. The union of the exploded nodes (n, d) for a given control-flow node n , mapped to the corresponding lattice elements $l_{n,d}$, form the environment Env_n for that node:

$$\text{Env}_n = \{(d, l_{n,d}) \mid (n, d) \in N^\#\}. \quad (12)$$

The overall idea behind computing the lattice elements $l_{n,d}$ is the following. For each inter-procedurally realizable path

$$p = [(\text{start}_{\text{main}}, A), (n_1, d_1), \dots, (n_k, d_k)]$$

that starts with the entrypoint of the exploded supergraph, we compute the micro function f_p that corresponds to p . The micro function consists of the composition of all individual micro functions with which the edges of p are labeled:

$$f_p = \text{EdgeFn}((n_{k-1}, d_{k-1}), (n_k, d_k)) \circ \dots \circ \text{EdgeFn}(\text{start}_{\text{main}}, A), (n_1, d_1)). \quad (13)$$

Let the lattice element that (n, d) is mapped to according to path p be denoted as $l_{n,d}^p$. As shown in Sagiv et al. [20], the lattice element can be obtained by applying f_p to the bottom element:

$$l_{n,d}^p = f_p(\perp). \quad (14)$$

Let Q be the set of paths that start at the entry point and end at the given node (n, d) . The lattice element $l_{n,d}$ is the meet of the lattice elements corresponding to all the paths in Q :

$$l_{n,d} = \bigcap_{q \in Q} l_{n,d}^q.$$

This is a general outline of the IDE analysis. We use the IDE framework to improve the precision of IFDS problems in the presence of correlated calls. The next section describes how this is done.

In Figure 3, should we say that ψ corresponds to the function's return value? (M)

we agreed to move this figure to the initial example section (O)

Move Figure 3 up to where it's mentioned first? If yes, the caption won't make sense. (M)

The following definitions are very general, and not related to the transformation. Can we move them somewhere earlier? (O)

Edge Functions Let \mathcal{F} be the set of methods in a program with a signature $s_{\mathcal{F}}$.

Definition 1. Let $r.c()$ be a call site on a receiver $r \in R$ with runtime type $t \in T$. Let $s_{\mathcal{F}}$ be the method signature corresponding to the call $c()$. For $s_{\mathcal{F}}$ and t , a lookup function returns the method implementation $f \in \mathcal{F}$ to which the call $r.c()$ is dispatched:

$$\text{lookup}(s_{\mathcal{F}}, t) = f. \quad (15)$$

Definition 2. For a method signature $s_{\mathcal{F}}$ and a method implementation $f \in \mathcal{F}$, the static-type function τ returns the set of types for which the lookup function yields f :

$$\tau(s_{\mathcal{F}}, f) = \{t \mid \text{lookup}(s_{\mathcal{F}}, t) = f\}. \quad (16)$$

In other words, τ computes the set of types for which calls to methods with signatures $s_{\mathcal{F}}$ are dispatched to f .

If there is a supergraph path from a method call with signature $s_{\mathcal{F}}$ to the start of f , then the set $\tau(s_{\mathcal{F}}, f)$ is always non-empty.

⁷ See Sagiv et al. [20] for a formal definition of the representation relation for environment transformers.

I think the rest of this section will not make sense to anyone who does not already understand it. Should I try to replace it with one paragraph that tries to give an intuitive understanding of what the IDE algorithm is doing? (M)

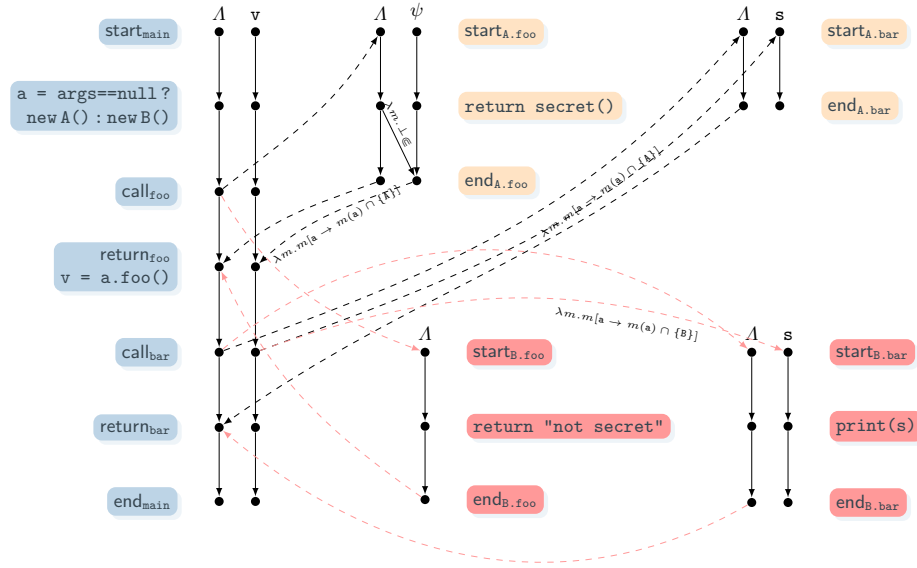


Fig. 3: An example program demonstrating correlated-call edge functions on the Λ -node path for Listing 2. All non-labeled edges are implicitly labeled with identity functions `id`. The variable corresponding to an initial secret value is denoted as ψ .

Definition 3. A call site is called *monomorphic* if it can be dispatched to only one method. If a call site can be dispatched to more than one method it is called *polymorphic*.

Let $r.c()$ be a call on a receiver $r \in R$ with a method signature $s_{\mathcal{F}}$ to a function $f \in \mathcal{F}$. If the call site is monomorphic, then $\tau(s_{\mathcal{F}}, f)$ contains all types $T' \subseteq T$ that are compatible with the static type of r . If the call site is polymorphic, then $\tau(s_{\mathcal{F}}, f) \subset T'$, since some types $t \in T'$ cause dispatch to a method other than f .

incorporate into previous section (O)

3.4 Defining IFDS and IDE Problems

In Section 3, we defined what IFDS and IDE problems are, their applications, and their constraints. In this section, we describe how to create instances of IFDS and IDE problems.

Defining an IFDS Problem Recall that an IFDS problem instance is defined as a five-tuple

$$(G^*, D, F, M_F, \sqcap),$$

where $G^* = (N^*, E^*)$ is the control-flow supergraph of the program, D is the set of dataflow facts, $F \subseteq 2^D \rightarrow 2^D$ is a set of distributive dataflow functions, and the function

$$M_F : E^* \rightarrow (2^D \rightarrow 2^D)$$

maps the supergraph edges to dataflow functions, and is extended to paths by composition.

In practice, an IFDS problem can be defined by providing an exploded supergraph $G^\# = (N^\#, E^\#)$. Each node of $G^\#$ is a pair (n, d) , where $n \in N^*$ is a node in the control-flow supergraph and $d \in (D \cup \{\mathbf{0}\})$, $\mathbf{0} \notin D$, where $\mathbf{0}$ is an auxiliary fact that is necessary for the IFDS algorithm.

The meaning of an edge in the exploded supergraph is the following. Let (n_1, d_1) and (n_2, d_2) be two nodes in the exploded supergraph $G^\#$. Furthermore, assume that if fact d_1 at node n_1 holds, then the fact d_2 at node n_2 also holds. Then there is an edge $(n_1, d_1), (n_2, d_2) \in E^\#$.

Defining an IDE Problem An IDE problem instance is a four-tuple

$$(G^*, D, L, M_{\text{Env}}),$$

where G^* and D are defined in the same way as for IFDS. L is a finite-height lattice that represents the values to which dataflow facts are mapped in an IDE problem. An environment $\text{Env}(D, L)$ maps dataflow facts to lattice elements. Finally, the map

$$M_{\text{Env}} : E^* \rightarrow (\text{Env}(D, L) \rightarrow \text{Env}(D, L))$$

is a function from the control-flow-supergraph edges to environment transformers, extended to paths by composition.

An IDE problem can be defined with a labeled exploded supergraph⁸, in which an edge function

$$\text{EdgeFn} : E^\# \rightarrow (L \rightarrow L) \quad (17)$$

pairs edges with *micro functions*, and is extended to paths by composition.

The set of micro functions of an IDE problem is a subset of $L \rightarrow L$ that is closed under function meet and composition.

The meaning of an edge in the labeled exploded supergraph is the following. Let $e = ((n_1, d_1), (n_2, d_2)) \in E^\#$ be an edge in the exploded supergraph with label $f = \text{EdgeFn}(e)$. Then

- the fact that d_1 holds at node n_1 implies that d_2 holds at n_2 ;
- if at node n_1 the fact d_1 was mapped to a lattice element l_1 by an environment $\text{Env}(D, L)$, then the fact d_2 at node n_2 should be mapped to $f(l_1)$.

As shown in Sagiv et al. [20], the relationship between environment transformers and edge functions can be described with the following equations. For individual edges $(n_1, n_2) \in E^*$,

$$\begin{aligned} M_{\text{Env}}((n_1, n_2))(\text{env})(d) \\ = \text{EdgeFn}((n_1, \perp), (n_2, d))(\top) \sqcap \prod_{d' \in D} \text{EdgeFn}((n_1, d'), (n_2, d))(\text{env}(d')), \end{aligned} \quad (18)$$

where env is an environment $\text{Env}(D, L)$. Informally, for a given control-flow-supergraph edge e and data-flow fact d , the M_{Env} function captures the meet of the edge function applied to all possible exploded-graph edges along e .

For paths p that start with the entry point $\text{start}_{\text{main}}$,

$$M_{\text{Env}}(p)(\Omega)(d) = \prod_{r \in \text{RP}(p, d)} \text{EdgeFn}(r)(\top), \quad (19)$$

where $n \in N^*$, $d \in D$, $p \in \text{VP}(n)$, and RP is the set of all inter-procedurally realizable paths.

To summarize, an IDE problem can be defined by a labeled exploded supergraph

$$(G^\#, L, \text{EdgeFn}), \quad (20)$$

where each edge of the exploded supergraph corresponds to a micro function.

⁸ The exploded supergraph in an IDE problem is defined in the same way as in an IFDS problem. The only difference is that the $\mathbf{0}$ fact is denoted as \perp [17,20].

4 Correlated Calls Analysis

The correlated-calls analysis is defined as a transformation from an arbitrary IFDS problem to a corresponding IDE problem. The solution of the IDE problem is converted to a solution of the original IFDS problem. The converted IFDS result can be more precise than the original IFDS result because it avoids infeasible paths corresponding to correlated calls.

4.1 Transformations from IFDS to IDE

Here, you defined the output of a transformation as just $G^\#$ and an edge function. Later, you also include the lattice. Should we include the lattice here? (O)

Let $G^\#$ be the exploded supergraph of an IFDS problem. A *transformation* $\mathcal{T} : (G^\#) \rightarrow (G^\#, (\top, \perp), \text{EdgeFn})$ converts it into an equivalent IDE problem with a two-point lattice. We consider two transformations:

- an equivalence transformation \mathcal{T}^\equiv (pronounced “t-equiv”) that generates IDE problems with the same precision as the original IFDS problem, and
- a correlated-call transformation \mathcal{T}^\subseteq (pronounced “t-c-c”) that generates IDE problems that exclude infeasible paths.

Both transformations keep the exploded supergraph $G^\#$ the same, and only generate different edge functions.

Equivalence Transformation The lattice for the equivalence transformation \mathcal{T}^\equiv is the two-point lattice $L^\equiv = \{\perp, \top\}$, where \perp means “reachable”, and \top means “not reachable”. The edge functions EdgeFn^\equiv are defined as

$$\text{EdgeFn}^\equiv = \begin{cases} \lambda e. \lambda m. \perp & \text{if } d_1(e) = \Lambda \text{ and } d_2(e) \neq \Lambda; \\ \lambda e. \text{id} & \text{otherwise,} \end{cases} \quad (21)$$

where $d_1(e)$ is the source fact of an edge e and $d_2(e)$ is its target fact. At a “diagonal” edge from a Λ -fact to a non- Λ -fact d , the micro function returns \perp to make the fact d reachable. All other micro-functions are the identity. The equivalence transformation is thus defined as: $\mathcal{T}^\equiv((G^\#)) = (G^\#, L^\equiv, \text{EdgeFn}^\equiv)$.

Correlated-Calls Transformation In the correlated-calls transformation, the lattice elements are maps from receivers to sets of types: $L^\subseteq = \{m : R \rightarrow 2^T\}$, where R is the set of receivers and T is the set of all types. For each receiver r , the map gives an overapproximation of the possible runtime types of r . Sets of types are ordered by the superset relation, and this is lifted to maps from receivers to sets of types, so the bottom element \perp_\subseteq maps every receiver to the set of all types, and the top element \top_\subseteq maps every receiver to the empty set of types. During an actual execution, every receiver r points to an object of some runtime type. Therefore, a data flow fact is unreachable along any feasible path if its corresponding lattice element maps any receiver to the empty set of types.

What if r points to null? (O)

A micro-function $f \in L^\subseteq \rightarrow L^\subseteq$ defines how the map from receivers to types should be updated when an instruction is executed. The micro-function

for most kinds of instructions is the identity. On a call to and return from a specific method m called on receiver r , the micro-function restricts the receiver-to-type map to map r only to types consistent with the polymorphic dispatch to method m . Finally, when an instruction assigns an object of unknown type to a receiver r , the corresponding micro-function updates the map to map r to the set of all types. This is made precise by the following definition:

Definition 4. *Given a previously fixed set $S \subseteq R$ of receivers, the micro-function $\varepsilon_S(e)$ of a supergraph edge e is defined as:*

$$\varepsilon_S(e) = \lambda m . \begin{cases} m[r \rightarrow m(r) \cap \tau(s_{\mathcal{F}}, f)], & \text{if } e \text{ is a call-start edge, } r.c() \text{ is the call site at } \text{src}(e), f \text{ is the called procedure with signature } s_{\mathcal{F}}, \text{ and } r \in S; \\ m[r \rightarrow m(r) \cap \tau(s_{\mathcal{F}}, f)] & \text{if } e \text{ is an end-return edge, } r.c() \text{ is the call corresponding to the return node at } \text{end}(e), f \text{ is the called method with signature } s_{\mathcal{F}}, v_1, \dots, v_k \in S \text{ are the local variables in } f, \text{ and } r \in S; \\ [v_1 \rightarrow \perp_T] \dots [v_k \rightarrow \perp_T], & \\ m[r \rightarrow \perp_T], & \text{if } \text{src}(e) \text{ contains an assignment for } r \in S; \\ m & \text{otherwise.} \end{cases} \quad (22)$$

define $n_1(e)$ globally somewhere early on, and consider using a more descriptive name like src instead of n_1 (\mathbf{O})

In the above definition, the purpose of the set S is to limit the set of considered receivers. We will use S in Section 4.5.

We can now define **EdgeFn**, which assigns a micro-function to each edge in the exploded supergraph. Along a Λ -edge, the micro function is the identity. All other functions can be described with ε_S . On a “diagonal” edge from Λ to a non- Λ fact that corresponds to some data flow fact becoming reachable, $\varepsilon_S(e)$ is applied to the initial map \perp_{\in} that conservatively allows every receiver to point to an object of any type. On all other edges, $\varepsilon_S(e)$ is applied to the existing map before the edge. The is formalized in the following definition.

Definition 5. *For each edge $e = (n_1, d_1) \rightarrow (n_2, d_2)$, $\text{EdgeFn}_S^{\infty}(e)$ is defined as follows:*

$$\text{EdgeFn}_S^{\infty}(e) = \begin{cases} id & \text{if } d_1 = d_2 = \Lambda, \\ \lambda m . \varepsilon_S(e)(\perp_{\in}) & \text{if } d_1 = \Lambda \text{ and } d_2 \neq \Lambda, \\ \lambda m . \varepsilon_S(e)(m) & \text{otherwise.} \end{cases} \quad (23)$$

Example 6. Consider the program from Listing 2, whose exploded supergraph appeared in Figure 3.

Returning a secret value in method **A.foo** creates a “diagonal” edge from the Λ -fact to the secret fact ψ . The diagonal edge is labeled with the micro function $\lambda m . \perp_{\in}$. Thus, at the end node of the method, every receiver is mapped to the set of all types \perp_T .

On the end-return edge from **A.foo** to **main**, the set of types for the receiver **a** is restricted by the micro function $\lambda m . m[\mathbf{a} \rightarrow m(\mathbf{a}) \cap \{\mathbf{A}\}]$ on the edge corresponding to the assignment of the return value ψ to **v**.

Similarly, on the call-start edge from `main` to `B.bar`, the possible types of the receiver `a` are further restricted by the micro-function $\lambda m. m[\mathbf{a} \rightarrow m(\mathbf{a}) \cap \{\mathbf{B}\}]$ on the edge that passes the argument `v` to the parameter `s`.

The composition of these micro functions results in the empty set as the possible types of the receiver `a`, indicating that this data flow path that would result in an information leak is actually infeasible.

Finally, the correlated-calls transformation is defined as $\mathcal{T}_S^{\subseteq}((G^{\#})) = (G^{\#}, L_S^{\subseteq}, \text{EdgeFn}_S^{\subseteq})$.

4.2 Converting IDE Results to IFDS Results

For each program point n , the result of an IFDS analysis gives a set of facts d that may be reached at n . The result of an IDE analysis pairs each such fact d with a lattice element ℓ . Formally, for an IFDS problem P , the result $\mathcal{R}_{\text{IFDS}}$ has type $N \rightarrow D$. Similarly, for an IDE problem Q , the result \mathcal{R}_{IDE} has type $N \rightarrow D \times L$.

Recall that in the equivalence transformation lattice L^{\equiv} , \perp means reachable and \top means unreachable. Therefore, a result ρ of the equivalence IDE analysis is converted to an IFDS result as follows: $\mathcal{U}^{\equiv}(\rho) = \lambda n. \{d \mid \rho(n) = (d, \top)\}$.

In the correlated-calls transformation lattice L^{\subseteq} , a map that maps any receiver to the empty set of possible types means that the corresponding data flow path is infeasible. Therefore, a result ρ of the correlated-calls IDE analysis is converted to an IFDS result as follows: $\mathcal{U}^{\subseteq}(\rho) = \lambda n. \{d \mid \rho(n) = (d, \ell), \forall r. \ell(r) \neq \top\}$.

4.3 Implementation of Correlated Calls Micro-Functions

move material from
following section here
(O)

Conceptually, micro-functions are functions from L to L , where L is the IDE lattice, either L^{\equiv} or L^{\subseteq} in our context. However, the IDE algorithm requires an efficient representation of micro-functions. The chosen representation needs to support the basic micro-functions that we presented in Section 4.1. The representation must also support function application, comparison, and be closed under function composition and meet. We now propose such a representation for the correlated-calls micro-functions.

The representation of a micro-function is a map from receivers to pairs of sets of types $I(r)$ and $U(r)$, where $U(r)$ is required to be a subset of $I(r)$. We use the notation $\langle I, U \rangle$ to represent such a map, and $I(r)$ and $U(r)$ to look up the sets corresponding to a particular receiver r . We define the meaning $\llbracket \langle I, U \rangle \rrbracket$ of a representation $\langle I, U \rangle$ as follows: $\llbracket \langle I, U \rangle \rrbracket = \lambda m. \lambda r. (m(r) \cap I(r)) \cup U(r)$. In words, the micro-function represented by $\langle I, U \rangle$ takes an existing map m from receivers to sets of types, and returns a map that maps each receiver r to the set $m(r)$ given by the original map m intersected with $I(r)$ and unioned with $U(r)$.

All of the basic micro-functions defined in Definition 4 can be expressed in this representation.

The implementation of function application follows directly from the definition of the representation: $\langle I, U \rangle (m) = \lambda r. (m(r) \cap I(r)) \cup U(r)$.

To compare two micro-functions for equality, it suffices to compare the corresponding sets $I(r)$ and $U(r)$ for all receivers r . The following lemma shows that this implementation of comparison corresponds to equality of the represented micro-functions:

Lemma 1. *For any pair of micro-function representations $\langle I, U \rangle, \langle I', U' \rangle$,*

$$\forall r. I(r) = I'(r) \wedge U(r) = U'(r) \iff \llbracket \langle I, U \rangle \rrbracket = \llbracket \langle I', U' \rangle \rrbracket$$

The composition of two micro-function representations is defined as follows: $\langle I, U \rangle \circ \langle I', U' \rangle = \langle \lambda r. (I(r) \cap I'(r)) \cup U(r), \lambda r. (I(r) \cap U'(r)) \cup U(r) \rangle$. The following lemma shows that this implementation corresponds to the composition of the denoted functions:

Lemma 2. *For any pair of micro-function representations $\langle I, U \rangle, \langle I', U' \rangle$,*

$$\llbracket \langle I, U \rangle \rrbracket \circ \llbracket \langle I', U' \rangle \rrbracket = \llbracket \langle I, U \rangle \circ \langle I', U' \rangle \rrbracket$$

The meet of two micro-function representations is defined as follows: $\langle I, U \rangle \sqcap \langle I', U' \rangle = \langle \lambda r. I(r) \cup I'(r), \lambda r. U(r) \cup U'(r) \rangle$. The following lemma shows that this implementation corresponds to the meet of the denoted functions:

Lemma 3. *For any pair of micro-function representations $\langle I, U \rangle, \langle I', U' \rangle$,*

$$\llbracket \langle I, U \rangle \rrbracket \sqcap \llbracket \langle I', U' \rangle \rrbracket = \llbracket \langle I, U \rangle \sqcap \langle I', U' \rangle \rrbracket$$

4.4 Theoretical Results

In the following lemma we show that the result of an IDE problem obtained through a correlated-calls transformation is a subset of the original IFDS result.

Lemma 4 (Precision). *For an IFDS problem P and all $n \in N^*$,*

$$\mathcal{U}^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P)))(n) \subseteq \mathcal{R}_{IFDS}(P)(n). \quad (24)$$

We will next show that our analysis is sound, i.e. that the result of an IDE problem obtained through a correlated-calls transformation removes only facts that occur on infeasible paths.

Lemma 5 (Soundness). *Let $p = [\text{start}_{\text{main}}, \dots, n]$ be a concrete execution path, and let $d \in D$. If $d \in M_F(p)(\emptyset)$, then*

$$d \in \mathcal{U}^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P)))(n). \quad (25)$$

Prove this lemma in the appendix. The parts of the proof can be found in section 5. (O)

The normalization function \mathcal{N} was only needed in composition. I therefore “inlined” it into composition to save space. This makes composition a bit harder to intuitively verify, but saves complexity here. The verification can be in the proof in the supplementary appendix. (O)

Prove this lemma in the appendix. The parts of the proof can be found in section 5. (O)

Prove this lemma in the appendix. The parts of the proof can be found in section 5. (O)

Sagiv et al. define a notion of an “efficient” implementation of micro-functions. If the micro-functions are “efficient”, then the asymptotic complexity bounds proven by Sagiv hold. Our implementation is “efficient” if we assume that the number of correlated receivers is a constant independent of the size of the program. That seems a stretch. If it’s linear in the size of the program, then we would need to add a factor to their complexity bounds. Their bound is $O(ED^3)$, while ours would be $O(ED^3R)$. But then we can’t directly use their complexity proof; we would have to adapt it. I’m leaning towards leaving out any mention of “efficient” implementation. (O)

Should we say \mathcal{R}_{IDE} instead of \mathcal{R} everywhere, to be consistent with \mathcal{R}_{IFDS} ? (O)

4.5 Correlated-Call Receivers

We will now show that in a correlated-calls transformation, it is enough to consider only some of the receivers of set R .

Definition 6. *Let c_1 and c_2 be two call sites on a receiver $r \in R$. If both call sites are polymorphic, then we say that r is a correlated-call receiver.*

should this be a section or subsection? should it be moved somewhere else (or maybe not included at all)? (O)

In other words, a correlated-call receiver is a receiver that has at least two polymorphic call invocations. We will denote the set of correlated-call receivers as R^\subseteq .

We will describe a “reduced” correlated-calls transformation in which we only consider receivers from R^\subseteq and ignore other receivers of R . We will show that IDE problems obtained through ordinary and reduced correlated-calls transformations yield the same results. In other words, we show that if a correlated calls analysis considers only correlated-call receivers, no precision is lost.

Lemma 6. *Let P be an IFDS problem. Then*

$$\mathcal{U}^\subseteq(\mathcal{R}(\mathcal{T}_{R^\subseteq}^\subseteq(P))) = \mathcal{U}^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P))). \quad (26)$$

To summarize, Lemma 5 shows that the result \mathcal{R}_\subseteq of a correlated-calls analysis is sound since it overapproximates the data flow of all possible concrete execution paths. We have also shown in Lemma 4 that the correlated-calls analysis improves the precision of the original IFDS result $\mathcal{R}_{\text{IFDS}}$, because the correlated-calls result \mathcal{R}_\subseteq underapproximates an equivalence-IDE result $\mathcal{R}_\equiv = \mathcal{R}_{\text{IFDS}}$. Finally, we showed in Lemma 6 that a correlated-call transformation to IDE that considers only correlated-call receivers R^\subseteq achieves the same result \mathcal{R}_\subseteq that is obtained when considering all receivers R .

5 Evaluation

This section discusses implementation aspects of the correlated-calls analysis and presents experimental results.

5.1 Implementation of the Analysis

The correlated-calls analysis was implemented in the Scala programming language [16]. We chose Java as the target language for client programs of the analysis. To retrieve information about an input program, such as its control-flow supergraph or the set of receivers and their types, we used the WALA framework for static analysis on Java bytecode [6].

Since WALA currently only contains an implementation of IFDS, we implemented IDE from scratch. Instead of using WALA’s IFDS implementation, to run an IFDS problem, we converted it to an IDE problem and used our own IDE solver.

We should incorporate the four functions into the explanation of IFDS in the background. Then we don't need them here. (O)

IFDS As described in Section 3.4, an IFDS problem is defined in terms of an exploded supergraph. The control-flow supergraph of an input program can be retrieved using WALA. Hence, our implementation of an IFDS problem should be able to convert a control-flow supergraph into an exploded supergraph.

We represent an IFDS problem with a trait, or protocol, that contains declarations of four *flow functions*. Each function has type

$$F : (N \times D \times N) \rightarrow 2^D$$

and defines a set of edges on the exploded graph. Given an edge (n_1, n_2) of the control-flow supergraph and the fact d_1 that corresponds to the source node n_1 , $F(n_1, d_1, n_2)$ returns the set of all facts $d_2 \in D_2$ such that $((n_1, d_1), (n_2, d_2)) \in E^\#$ ⁹. The four functions are:

- **call-start**, for inter-procedural edges from a call node to the start node of the target method;
- **call-return**, for intra-procedural edges from a call node to its return node;
- **end-return**, for inter-procedural edges from the end node of a method to the return node of the callee;
- **default**, for all other intra-procedural edges.

Taint Analysis Using this representation of an IFDS problem, we implemented an IFDS problem instance for taint analysis. We used it as a sample IFDS problem on which to evaluate the correlated-calls-IDE construction.

Let N^* be the control-flow supergraph of a program and D the set of the program variables. Let $\text{encl}(n)$ be a function that returns the enclosing method of a node $n \in N^*$. Finally, let the function $r_m : D \rightarrow 2^D$ be defined as follows:

$$r_m(d) = \begin{cases} \emptyset & \text{if } d \text{ is a local variable in method } m, \\ \{d\} & \text{otherwise.} \end{cases} \quad (27)$$

When defining the flow functions for a taint analysis, we will use r_m to avoid the propagation of local variables, as shown below.

For a fact $d_1 \in D \cup \{\mathbf{0}\}$ and two nodes $n_1, n_2 \in N^*$, the simplified¹⁰ version of flow functions for a taint-analysis looks as follows.

If n_1 is a call node that calls method m , and n_2 is m 's start node,

$$\text{call-start}(n_1, d_1, n_2) = \begin{cases} r_{\text{encl}(n_1)}(d_1) \cup \{v\} & \text{if } a \text{ is the } i\text{th argument of the call,} \\ & d_1 = a, \text{ and } v \text{ is the } i\text{th parameter} \\ & \text{of } m; \\ r_{\text{encl}(n_1)}(d_1) & \text{otherwise.} \end{cases}$$

⁹ In each invocation of a flow function, the fact d_1 is provided by the IDE algorithm.

¹⁰ For simplicity, the shown flow functions do not account for different Java-specific features such as arrays, fields, operations on strings, etc.

If n_1 is a call node with corresponding return node n_2 ,

$$\text{call-return}(n_1, d_1, n_2) = \begin{cases} \{d_1\} & \text{if } d_1 \text{ is a local variable in } \text{encl}(n_1), \\ \emptyset & \text{otherwise.} \end{cases}$$

If c is a call node calling method m , n_1 is m 's end node, and n_2 is c 's return node,

$$\text{end-return}(n_1, d_1, n_2) = \begin{cases} r_{\text{encl}(n_1)}(d_1) \cup \{x\} & \text{if } n_1 \text{ is a return statement} \\ & \text{returning } v, n_2 \text{ is an assignment} \\ & \text{with left-hand side } x, \text{ and } d_1 = v; \\ r_{\text{encl}(n_1)}(d_1) & \text{otherwise.} \end{cases}$$

Otherwise,

$$\text{default}(n_1, d_1, n_2) = \{d_1\}.$$

Example 7. Consider the supergraph in Figure ?? . The call-to-start flow function from method `main` to `f` looks as follows:

$$\begin{aligned} \text{call-start}(\text{call}_{A.f}, a, \text{start}_f) &= r_{\text{main}}(a) \cup \{s\} \\ &= \{s\}. \end{aligned}$$

We can see that correspondingly, the exploded supergraph contains an edge from $(\text{call}_{A.f}, a)$ to (start_f, s) .

IDE The correlated-calls analysis was implemented as an IDE problem instance.

We defined an IDE problem in the same way as an IFDS problem, except that the IDE flow functions are of type

$$(N \times D \times N) \rightarrow 2^{D \times (L \rightarrow L)}.$$

With the new flow functions, we can implement a labeled exploded supergraph, since the new flow functions return a set of facts that are paired with micro functions.

For example, if Q is an IDE problem, then the call-to-start flow function for Q is defined as follows:

$$\begin{aligned} &\text{call-start}^Q(n_1, d_1, n_2) \\ &= \left\{ (d_2, f) \mid d_2 \in D, f \in L^Q \rightarrow L^Q : \text{EdgeFn}^Q((n_1, d_1), (n_2, d_2)) = f \right\}. \end{aligned}$$

The other flow functions are defined analogously.

5.2 Testing

In this section we assess the correctness and effectiveness of the correlated-calls analysis.

Conversion from IFDS to IDE We implemented the equivalence transformation \mathcal{T}^{\equiv} and the correlated-calls transformation $\mathcal{T}_{R^{\epsilon}}^{\subseteq}$ from IFDS to IDE described in Section 4.1. To run an IFDS problem, we converted it to an IDE problem using \mathcal{T}^{\equiv} and $\mathcal{T}_{R^{\epsilon}}^{\subseteq}$ and used our IDE analysis algorithm to run the latter.

Given an IFDS problem described with IFDS flow functions, an equivalence transformation creates an IDE problem described with the following IDE flow functions:

$$\begin{aligned} \text{call-start}^{\equiv}(n_1, d_1, n_2) &= \{(d_2, \epsilon(d_1, d_2)) \mid d_2 \in \text{call-start}(n_1, d_1, n_2)\} \\ \text{call-return}^{\equiv}(n_1, d_1, n_2) &= \{(d_2, \epsilon(d_1, d_2)) \mid d_2 \in \text{call-return}(n_1, d_1, n_2)\} \\ \text{end-return}^{\equiv}(n_1, d_1, n_2) &= \{(d_2, \epsilon(d_1, d_2)) \mid d_2 \in \text{end-return}(n_1, d_1, n_2)\} \\ \text{default}^{\equiv}(n_1, d_1, n_2) &= \{(d_2, \epsilon(d_1, d_2)) \mid d_2 \in \text{default}(n_1, d_1, n_2)\}, \end{aligned}$$

where ϵ is the bottom function on an edge from a Λ -fact to a non- Λ -fact, and the identity function otherwise:

$$\epsilon(d_1, d_2) = \begin{cases} \lambda l. \perp & \text{if } d_1 = \Lambda \text{ and } d_2 \neq \Lambda; \\ \text{id} & \text{otherwise.} \end{cases}$$

We also implemented a correlated-call transformation from IFDS into IDE problems that consider correlated calls. This transformation is described in Section 4.1.

Regression Testing We used regression tests to assess the correctness of the implemented analyses. Each test involves running a certain analysis on one input Java program.

IDE-Implementation Correctness To test the correctness of the IDE algorithm implementation, we implemented a copy-constant-propagation IDE problem [20]. In a copy-constant propagation analysis, a variable is considered constant if it is assigned a constant literal or another variable that is also a constant. For example, in a program

```
int a = 1;
int b = a;
int c = a + b;
int d = a + 2;
```

`a` and `b` are considered constant, but `c` and `d` are not (although `d` would be considered constant in linear-constant propagation).

We tested the propagation of constants on different intra- and inter-procedural data-flow paths, in parameter passing, and in conditional branches. Each regression test contained assertions of the form “at the end of method m , variable with name x should be (not) constant”.

We also tested the implementation of the IDE algorithm on an IDE problem generated by conversion from an IFDS problem.

To do that, we implemented an IFDS instance for taint analysis.

Recall from Section 3.2 that taint analysis aims to discover variables that are secret at a given program point called a sink.

We used assertions of the form “at program statement n , variable x should be (not) secret” by defining the sink of a secret value through special `isSecret` and `notSecret` methods. Those methods asserted that the parameter passed to them is secret and not secret, respectively. To define a source secret value we created a static `secret()` method that returned a string.

Example 8. Listing 4 illustrates the use of the `isSecret` and `notSecret` assertions.

```
public static void main(String[] args) {
    String n = "not_secret";
    notSecret(n); // assert that n is not secret
    String s1 = f(n);
    isSecret(s1);
    String s2 = f(secret());
}

static String f(String str) {
    isSecret(str);
    return str;
}

public static String secret() { // the secret source
    return "secret";
}
```

Fig. 4: Example usage of `isSecret` and `notSecret` assertions in regression tests

We tested data flow through

- method calls and returns;
- conditional branches and loops, including nested constructions, the ternary operator, and `switch` statements;
- arrays and fields¹¹;

¹¹ In Java, arrays are allocated on the heap, and array elements can be aliases of each other. Hence, if any array element gets assigned a secret value, we considered all

- static and instance class members;
- classes and interfaces that involve inheritance, overriding, and overloading;
- recursion;
- library calls¹²;
- string concatenation and usage of the `StringBuffer` and `StringBuilder` classes¹³;
- generics, type conversions through castings, and exception handling.

Our taint analysis implementation becomes unsound in the presence of static initializers. If a static field is initialized to a secret value, our analysis will not detect it as such.

A static initializer is invoked only once, before the instance creation of a class or the access of a static member of that class. Static initializers are invoked lazily by the Java Virtual Machine [11]. This makes finding out at which program point a static initializer is invoked undecidable [7]. To account for static initializers in the analysis would require modifying WALA’s control-flow supergraph (which does not have edges to static initializers) or using a data-flow analysis for static initialization. Since the primary purpose of the taint-analysis implementation was to test the correlated-call analysis, we did not include a static-initializer analysis in this work.

Correlated-Calls-Analysis Correctness We tested the implementation of the correlated-calls analysis by converting the taint analysis into an IDE problem with an implementation of $\mathcal{T}_{R^{\infty}}$.

Since none of the test cases in the previous section contained correlated calls, we used the same tests with the same assertions to ensure that the correlated-calls analysis produces the same results as an IFDS-equivalent analysis in the absence of correlated calls.

We then added test cases that contained correlated calls. We added a new assertion method, `notSecretCC`. For the IFDS-equivalent analysis, the method asserted that the argument passed to it was secret, and for the correlated-calls analysis, it asserted that the argument was not secret.

Separately, we used unit tests to check the implementation correctness of micro functions. We wrote assertions for the results of the equality, meet, and composition operations on all possible combinations of the identity, top, bottom, and constant functions.

elements of any `String` or `Object` array in the program secret. For the same reason, if a field `f` of an object of class `A` is assigned a secret value, then we considered the field `f` of any object of class `A` secret.

¹² We created a specification for library functions that allowed us to indicate under which conditions a library function returned a secret value. This let us avoid the expensive analysis of library functions.

¹³ Using mutation, objects of these classes can be converted into wrappers around secret strings. This is why we added a special handling for `StringBuffer` and `StringBuilder` objects. For instance, if a field had the `StringBuilder` type, it was considered secret.

We need to decide what exactly we want to include in the evaluation. The numbers of correlated calls are interesting. But I think we want to leave out all mention of the taint analysis. Perhaps we should have no empirical results at all (and only the theory/proofs)? (O)

Benchmark Testing To assess the benefit of the correlated-calls analysis, we counted the frequencies of correlated-call occurrences in the Dacapo benchmarks [2]. We then ran the normal- and correlated-call-taint analysis on the Dacapo benchmarks to see what improvement we would get from the correlated-calls analysis.

Occurrences of Correlated Calls Our goal was to obtain an upper bound on the number of redundant IFDS-result nodes that could be potentially removed by our analysis. We counted the number of correlated calls that occurred in programs of the Dacapo benchmarks, as shown in Table 1.

In the table, the number of all call sites in a program is denoted as C . Polymorphic call sites are denoted as C_P , and correlated call sites as C^\subseteq . The first four columns indicate the overall number of various call sites and correlated-call receivers in a program. The last three columns indicate the ratio of polymorphic to all call sites, the ratio of correlated to polymorphic call sites, and the ratio of correlated call sites to correlated-call receivers.

Table 1: Frequencies of correlated-call occurrences in the Dacapo benchmarks

| Benchmark | $ C $ | $ C_P $ | $ C^\subseteq $ | $ R^\subseteq $ | $\frac{ C_P }{ C }$ | $\frac{ C^\subseteq }{ C_P }$ | $\frac{ C^\subseteq }{ R^\subseteq }$ |
|------------|--------------|------------|-----------------|-----------------|---------------------|-------------------------------|---------------------------------------|
| antlr | 7,610 | 428 | 299 | 70 | 6% | 70% | 4 |
| bloat | 18,157 | 933 | 429 | 119 | 5% | 46% | 4 |
| chart | 18,101 | 466 | 195 | 61 | 3% | 42% | 3 |
| eclipse | 3,222 | 100 | 35 | 10 | 3% | 35% | 4 |
| fop | 4,831 | 129 | 40 | 12 | 3% | 31% | 3 |
| hsqldb | 3,573 | 81 | 35 | 10 | 2% | 43% | 4 |
| python | 12,149 | 487 | 129 | 54 | 4% | 26% | 2 |
| luindex | 7,190 | 188 | 79 | 29 | 3% | 42% | 3 |
| lusearch | 9,043 | 350 | 126 | 47 | 4% | 36% | 3 |
| pmd | 10,972 | 219 | 68 | 23 | 2% | 31% | 3 |
| xalan | 3,889 | 110 | 35 | 10 | 3% | 32% | 4 |
| Geom. mean | 7,572 | 240 | 91 | 29 | 3% | 38% | 3 |

We can see that on average, 3% of all call sites C are polymorphic call sites C_P . Out of those call sites, 38% are correlated call sites C^\subseteq . We also see that for one correlated-call receiver, there are on average three correlated calls.

Experiments We ran the analysis on the Dacapo benchmarks to test if the taint analysis would benefit from the improved, correlated-calls based, analysis. We defined any user input string to be considered a secret source and compared the overall number of results in the original and correlated-call taint analyses. If the number of secret values in the original result were larger than in the correlated-call result, we would see a practical benefit from our analysis.

However, even when we considered each program point as a sink, the “improved” analysis revealed the same number of secret values as the original taint analysis.

A correlated call that could affect a taint-analysis result could most likely occur in the following scenario:

- there is a receiver with at least two polymorphic calls;
- at least one of the calls c_1 returns a string — this would mean that the method potentially returns a secret value;
- at least one of the calls c_2 takes a string parameter — this would mean that a secret value could potentially be propagated to the method as an argument.

Then, if the correlated call occurred on an invocation $c_2(c_1())$, there might be a possibility of benefiting from the correlated-calls analysis. Given the relatively rare occurrence of correlated calls, this situation is not likely to appear often. This is illustrated in Table 2 which shows how often correlated calls would invoke methods that either take a string as a parameter *or* return a string. The set of receivers on which there are invocations of such methods is denoted as R^{\subseteq}_S . A situation where one correlated call returned a string, *and* another correlated call on the same receiver took a string parameter, appeared in only one case in the **antlr** benchmark. However, the strings invoked were not designated as secret.

This explains why, specifically for a taint analysis as the client analysis, and specifically for the Dacapo benchmarks, the correlated call analysis did not make a difference.

Table 2: Frequency of correlated-call receivers for which at least one of the correlated calls takes a string as a parameter or returns a string

| Benchmark | $ R^{\subseteq}_S $ | $ R^{\subseteq} $ | $\frac{ R^{\subseteq}_S }{ R^{\subseteq} }$ |
|-------------------|---------------------|-------------------|---------------------------------------------|
| antlr | 43 | 70 | 62% |
| bloat | 0 | 119 | 0% |
| chart | 1 | 61 | 2% |
| eclipse | 0 | 10 | 0% |
| fop | 0 | 12 | 0% |
| hsqldb | 0 | 10 | 0% |
| jython | 6 | 54 | 23% |
| luindex | 0 | 29 | 0% |
| lusearch | 2 | 47 | 6% |
| pmd | 1 | 23 | 3% |
| xalan | 0 | 10 | 0% |
| Geom. mean | 3 | 29 | 9 |

5.3 Future Work

We should convert this section into two or three sentences to be added to the conclusion. They should focus especially on the interprocedurally-correlated receivers. (O)

In this section we point out the limitations of the correlated-calls analysis and suggest improvements to the analysis for future work.

One limitation of the analysis is that it only works for IFDS problems like taint analysis, reachable definitions, or available expressions. The correlated-call analysis is not applicable to IDE problems like copy- or linear-constant propagation. Therefore, a possible direction for future work is to create a correlated-calls analysis that transforms an original IDE problem into one that considers correlated calls (with a modified lattice and edge function definition), and then transforms the correlated-calls result into a more precise result of the original IDE problem.

```
class A {
    String string;

    public static void main(String[] args) {
        A a = args == null ? new A() : new B();
        a.setString();
        propagate(a);
    }

    static void propagate(A a) {
        a.printString();
    }

    void setString() {
        string = secret();
    }

    void printString() {
        System.out.println("not_secret");
    }
}

class B extends A {
    void setString() {
        string = "not_secret";
    }

    void printString() {
        System.out.println(a);
    }
}
```

Fig. 5: Inter-procedurally-correlated calls

Another constraint of the algorithm is that it only accounts for intra-procedurally-correlated receivers, or receivers on which correlated calls occur within one method. For example, in Listing 5, `a` is a correlated-call receiver, since there are two polymorphic method invocations on `a`. However, the first one, `a.setString()`, is inside method `main`, and the second one, `a.printString()`, is inside method `propagate`. Therefore, we do not treat `a` as a correlated-call receiver, and the analysis would not improve the original IFDS result for this program.

Finally, correlated calls can occur on multiple receivers and other scenarios discussed in [?] that are not handled in this work.

6 Related Work

IFDS is a version of the functional approach to data-flow analysis developed by M. Sharir and A. Pnueli [21]. Their algorithm is based on computing *summary functions* that return the data-flow value at the end of a procedure, given the data-flow value at the start of the procedure. IFDS problems form a more restricted set of data-flow problems: unlike in the functional approach, IFDS flow functions have to be distributive, and the set of data-flow facts D has to be finite. However, the IFDS algorithm is more general than Sharir’s and Pnueli’s algorithm in that it can handle programs containing local variables and parameters in recursive methods.

IFDS has been used to encode a variety of data-flow problems. More complex examples of applications include typestate analysis (determining which operations can be performed on an object at a given program point) [13] or shape analysis (detecting errors and validating properties of programs at compile time) [10].

IFDS is implemented for two popular static-analysis frameworks, the T.J. Watson Libraries for Analysis (WALA) [6] and Soot [25].

WALA is a framework for static analysis on Java bytecode developed by the IBM T.J. Watson Research Center. In the implementation of our work, we use WALA to build and traverse the supergraph (a special kind of control-flow graph) of a Java program¹⁴.

Soot is a framework for program analysis and optimization on Java bytecode, developed by the Sable Research Group at McGill University. Unlike WALA, Soot also has an implementation of the IDE algorithm. The IFDS and IDE implementations for Soot are part of the Heros project [3].

Whereas one advantage of Soot’s IFDS implementation (and other static analysis tools) is ease of use and extensibility, WALA’s primary focus is efficiency. For example, WALA uses bit-vectors to represent some of the analysis data types, like local variables and parameters. Another difference is that WALA’s intermediate representation of a program uses static single assignment (SSA) form [5]. SSA form is a representation of the program in which each variable has only one definition (assignment). SSA can make dataflow analysis simpler and more efficient [1].

¹⁴ However, we do not use WALA’s IFDS implementation, as explained in Section 5.

Work on improving the IFDS algorithm includes Practical Extensions by N. Naeem and O. Lhoták [14]. Their paper presents four extensions to the IFDS algorithm. Two of the extensions improve the efficiency of the IFDS analysis for certain classes of IFDS problems. Another extension widens the class of problems applicable for the IFDS analysis. However, those extensions do not affect the precision of IFDS problems. Our analysis, in contrast, does not improve the efficiency or generality of IFDS, but it allows us to solve IFDS problems more precisely.

The fourth extension is targeted towards programs that are represented in SSA form. Executing the IFDS analysis on such programs results in loss of precision in the presence of control-flow constructs (e.g. conditionals and loops), compared to programs in non-SSA form. The extension makes the IFDS analysis on programs in SSA form as precise as on programs that are not represented in SSA form. In contrast, the correlated-calls analysis is applicable to programs in both SSA and non-SSA forms. Even if applied to a program in SSA form, our analysis and the extension improve the precision of IFDS in unrelated situations: the first analysis handles correlated calls, and the latter handles control-flow constructs. Thus, an IFDS analysis could benefit from both precision improvements independently.

Another work on improving the efficiency of the IFDS algorithm is E. Bodden et al.’s framework for the analysis of software products lines [4]. Their paper uses transformations from IFDS to IDE problems, a technique we also employ. Finally, J. Rodriguez and O. Lhoták implemented a concurrent version of the IFDS algorithm using actors [19]. However, neither of those works is concerned with improving the precision of IFDS results.

The correlated-calls analysis improves the precision of a data-flow analysis by eliminating a special type of infeasible paths. This is similar to the idea of context-sensitive analysis: just as a context-sensitive analysis eliminates infeasible paths from the end of a procedure to the call sites that do not match the given procedure call, the correlated-calls analysis eliminates infeasible paths caused by correlated method calls.

The idea of using correlated calls to remove infeasible paths in data-flow analyses of object-oriented programs was introduced by F. Tip [?]. The possibility of using IDE to achieve this is mentioned, but not elaborated upon. Our work presents a concrete solution to the problem and an implementation of that solution.

The idea of eliminating infeasible paths caused by correlated calls is similar to M. Sridharan et al.’s work on improving the precision of pointer analysis for JavaScript programs [22]. For each pointer, a pointer analysis determines the possible set of objects (the *points-to* set) that the pointer can reference at a given program point. In JavaScript, it is challenging to compute the points-to set of fields because in general, field names can be derived from arbitrary expressions and bound at runtime. As a result, an imprecise data-flow analysis will include infeasible paths between values of the form $o[p]$ (access of a property p of object o), where at compile time, p can be bound to different values. The idea of the paper is to track all dynamic property accesses (reads and writes) on an object

o with property name p . The code snippets containing the references $o[p]$ are then extracted into a separate function f . The analysis is then run so that for each possible value of p , f is analyzed separately; therefore, for a given property name, all correlated objects with that name are analyzed together.

The differences between this method of tracking correlated calls and our analysis are the following.

- *Type of target data-flow analysis* whose precision is to be improved. Our analysis improves the precision of IFDS data-flow analyses, whereas the JavaScript analysis improves the precision of pointer analysis.
- *Target language*. Our analysis is for object-oriented languages where polymorphic methods, and not property names (which are known at compile time), cause infeasible paths.
- *Different handling of correlated calls*. Extracting code that contains correlated calls into separate methods would not prevent infeasible paths. Instead, our analysis uses IDE flow functions to detect and eliminate infeasible paths caused by correlated calls.

7 Conclusions

We presented a technique to improve the precision of solutions to IFDS problems in the presence of correlated calls. Correlated calls occur when there are multiple polymorphic method invocations on the same receiver. Such method calls cause a data-flow analysis to consider infeasible paths, which makes the data-flow analysis less precise.

Our method of eliminating infeasible paths caused by correlated calls works by transforming an existing IFDS problem into a specialized IDE problem. In this way, we are able to track the classes to which method invocations get dispatched. After solving the specialized IDE problem, we convert its result into an IFDS result that is potentially more precise than the solution to the original IFDS problem. The increase in precision can occur for programs that contain correlated calls. Specifically, if, on a certain data-flow path, there are two polymorphic method invocations on the same receiver that dispatch to incompatible classes, the IDE analysis will consider the path infeasible.

We proved that the correlated-calls analysis is sound and that it improves the precision of IFDS results.

Our Scala implementation of the correlated-calls analysis includes

- an implementation of the IDE analysis, which is based on the WALA static program analysis framework;
- a taint-analysis implementation as an IFDS problem instance;
- a transformer of IFDS problems to equivalent IDE problems, and a second transformer that accounts for correlated calls.

We tested the correlated-calls analysis on our taint analysis implementation by comparing the number of secret values that were leaked when using an IFDS

taint analysis and a taint analysis that accounts for correlated calls. We used the Dacapo benchmarks as input programs. Although the benchmarks contained a number of correlated calls, we were not able to improve the precision of the taint analysis, because the correlated calls did not occur on paths of secret information leaks.

We are hopeful that other analyses can benefit from the extra information provided by the correlated-calls analysis, and plan to test this hypothesis in the future.

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Appendix

In this appendix we present the proofs to the Lemmas introduced in Section ??.

Soundness and Precision

We start by proving the Lemmas of Soundness and Precision of the correlated-calls analysis.

Proof of Lemma 4. The transformation \mathcal{U}^\subseteq is the same as \mathcal{U}^\equiv , except that it can remove data-flow facts from the result:

$$\begin{aligned} \mathcal{U}^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P)))(n) &= \{(n', D'_n(\mathcal{R}(\mathcal{T}_R^\subseteq(P)))) \mid n \in N^*\}(n) \\ &= D_n^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P))) \\ &\subseteq \text{MVP}_F(n) \\ &= \mathcal{R}_{\text{IFDS}}(P)(n). \end{aligned} \quad \square$$

To prove the Soundness Lemma, we first introduce Lemmas 7 and 8. We will denote the top element in the environment lattice as

$$\Omega = \lambda d. \top_\subseteq. \quad (28)$$

For the purpose of the proofs, we will rewrite Equation (23) that defines an edge function as follows:

$$\text{EdgeFn}_S^\subseteq = \lambda e. \begin{cases} \text{id} & \text{if } d_1 = d_2 = A, \\ \lambda m. \varepsilon(e)(\delta(m)) & \text{otherwise,} \end{cases} \quad (29)$$

where $S \subseteq R$, d_1 and d_2 are the source and target facts, and for a map $m \in L_U^\subseteq$, $\delta(m)$ is either m or \perp_\subseteq :

$$\delta(m) = \begin{cases} \perp_\subseteq & \text{if } d_1 = A \\ m & \text{otherwise.} \end{cases} \quad (30)$$

Additionally, for a path $p = [\text{start}_{\text{main}}, \dots]$ and a fact $d \in D$, we will denote the lattice element that is mapped to d according to the flow functions of path p as follows:

$$\xi(p, d) = M_{\text{Env}}(p)(\Omega)(d). \quad (31)$$

The following Lemma shows that the lattice elements (receiver-to-types maps) of a correlated-calls IDE analysis correctly overapproximate the possible types of a receiver in a program execution.

Lemma 7. *Let $p = [\text{start}_{\text{main}}, \dots, n]$ be some concrete execution trace of the program, and let $r \in R$ be a receiver. If after the execution trace p , at node n , r points to an object of runtime type t , and $d \in D$ is a fact such that $d \in M_F(p)(\emptyset)$, then*

$$t \in \xi(p, d)(r). \quad (32)$$

Proof. By induction on the length of the trace.

Basis: $p = [\text{start}_{\text{main}}]$. Then there is no instruction at which a receiver r could be instantiated, and the Lemma is trivially true.

Induction hypothesis: Let $p = [\text{start}_{\text{main}}, \dots, n_{k-1}]$, and let τ be the set of types to which $\xi(p, d_{k-1})$ maps r :

$$\tau = \xi(p, d_{k-1})(r). \quad (33)$$

Assume that for a concrete execution path $p = [\text{start}_{\text{main}}, \dots, n_{k-1}]$, at node (n_{k-1}, d_{k-1}) , the Lemma holds, i.e. $t \in \tau$.

Induction step: Let $p' = [\text{start}_{\text{main}}, \dots, n_{k-1}, n_k]$ and $t' \in T$ be the type to which r is mapped at n_k .

For each i , let e_i be the edge $((n_{i-1}, d_{i-1}), (n_i, d_i))$. Note that

$$e_1 = ((\text{start}_{\text{main}}, \Lambda), (n_1, d_1)).$$

Observe that

$$\begin{aligned} \xi(p', d) &= M_{\text{Env}}(p')(\Omega)(d) \\ &= (M_{\text{Env}}(e_k) \circ M_{\text{Env}}(e_{k-1}) \circ \dots \circ M_{\text{Env}}(e_1))(\Omega)(d) \\ &= M_{\text{Env}}(e_k)(M_{\text{Env}}(p)(\Omega))(d). \end{aligned}$$

According to (18),

$$\begin{aligned} &M_{\text{Env}}(e_k)(M_{\text{Env}}(p)(\Omega))(d)(r) \\ &= \left(\text{EdgeFn}_R^{\subseteq}((n_{k-1}, \Lambda), (n_k, d))(\top_{\subseteq}) \right) \sqcap \\ &\quad \left(\bigcap_{d' \in D} \text{EdgeFn}_R^{\subseteq}((n_{k-1}, d'), (n_k, d))(M_{\text{Env}}(p)(\Omega)(d')) \right)(r) \\ &\supseteq \bigcap_{d' \in D} \text{EdgeFn}_R^{\subseteq}((n_{k-1}, d'), (n_k, d))(M_{\text{Env}}(p)(\Omega)(d'))(r) \\ &\supseteq \text{EdgeFn}_R^{\subseteq}((n_{k-1}, d_{k-1}), (n_k, d))(\xi(p, d_{k-1}))(r). \end{aligned}$$

Therefore,

$$\text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) \subseteq \xi(p', d)(r). \quad (34)$$

We will now show that

$$t' \in \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r),$$

which, due to (34), means that the Lemma holds.

According to (29), there are two cases in which $\text{EdgeFn}_R^{\subseteq}(e_k)$ could fail.

If $d_{k-1} = d_k = \Lambda$, then $d_k \notin M_F(p)(\emptyset)$, since it does not belong to the set D , and the Lemma trivially holds.

Otherwise,

$$\text{EdgeFn}_R^{\subseteq}(e_k) = \lambda m. \varepsilon(e_k)(\delta(m)).$$

It follows that

$$\begin{aligned} \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) &= (\lambda m. \varepsilon(e_k)(\delta(m)))(\xi(p, d_{k-1}))(r) \\ &= \varepsilon(e_k)(\delta(\xi(p, d_{k-1}))(r)). \end{aligned} \quad (35)$$

Let us denote the lattice element $\delta(\xi(p, d_{k-1}))$ with Δ :

$$\Delta = \delta(\xi(p, d_{k-1})).$$

Note that since Δ , according to (30), can be either \perp_{\subseteq} or $\xi(p, d_{k-1})$, it always maps r to a set containing t :

$$t \in \Delta(r). \quad (36)$$

Note also that unless the instruction at n_{k-1} contains an assignment for r , r is mapped to the same object of type t as at node n_{k-1} , and $t = t'$. Therefore, for the non-assignment instructions, it is sufficient to prove that $t \in \Delta(r)$.

Depending on the instructions at the nodes n_{k-1} and n_k , there are four cases:

1. The instruction at n_{k-1} is an assignment for a receiver $r' \in R$. Since $\varepsilon_R(e_k) = \lambda m. m[r' \rightarrow \perp_T]$,

$$\begin{aligned} \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) &= (\lambda m. m[r' \rightarrow \perp_T])(\Delta)(r) \\ &= \Delta[r' \rightarrow \perp_T](r). \end{aligned}$$

In the resulting map, r' is mapped to \perp_T . Then

- (a) if $r = r'$, then $\text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) = \perp_T$, which contains t' .
 - (b) If $r \neq r'$, then r has not been reassigned a value, and still maps to the same object of type t . The receiver r is mapped to $\Delta(r)$, which, according to (36), contains t . Since $t = t'$, $\Delta(r)$ contains t' .
2. e_k is a call-start edge with signature $s_{\mathcal{F}}$, and $f \in \mathcal{F}$ is the called procedure. Then

$$\begin{aligned} \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) &= (\lambda m. m[r' \rightarrow m(r') \cap \tau(s_{\mathcal{F}}, f)])(\Delta)(r) \\ &= \Delta[r' \rightarrow \Delta(r') \cap \tau(s_{\mathcal{F}}, f)], \end{aligned}$$

where r' is the receiver of the call.

- If $r' = r$, then $\Delta(r') = \Delta(r)$ which contains t . Since $t \in \tau(s_{\mathcal{F}}, f)$, it follows that $t \in \Delta(r) \cap \tau(s_{\mathcal{F}}, f)$, and $t \in \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r)$.
 - If $r' \neq r$, see (1b).
3. e_k is an end-return edge, $r_1, \dots, r_k \in R$ are the local variables in the callee method, r' is the receiver of the call site corresponding to the return node n_k , and $f \in \mathcal{F}$ is the called method with signature $s_{\mathcal{F}}$. Then

$$\varepsilon_R(e_k) = \lambda m. m[r' \rightarrow m(r') \cap \tau(s_{\mathcal{F}}, f)][r_1 \rightarrow \perp_T] \dots [r_k \rightarrow \perp_T].$$

If $r \in \{r_1, \dots, r_k\}$, see Case 1. Otherwise, the case is analogous to Case 2.

4. The node contains any other instruction. Then

$$\text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) = \text{id}(\Delta)(r) = \Delta(r),$$

which contains t according to (36). \square

We will now show that on a node of a concrete execution path, the correlated-calls analysis does not map receivers to \top_T . In other words, the analysis never considers nodes of a concrete execution path unreachable.

Lemma 8. *Let $p = [\text{start}_{\text{main}}, \dots, n]$ be a concrete execution path, $r \in R$ a receiver, and $d \in D$ a data-flow fact. Then if $d \in M_F(p)(\emptyset)$,*

$$\xi(p, d)(r) \neq \top_T. \quad (37)$$

Proof. By induction on the length of the execution trace.

Basis: Let $p = [\text{start}_{\text{main}}]$. Since the only realizable path corresponding to p is $[(\text{start}_{\text{main}}, \Delta)]$, there is no fact $d \in D$ such that $d \in M_F(p)(\emptyset)$, and the claim follows immediately.

Induction hypothesis: Let $p = [\text{start}_{\text{main}}, \dots, n_{k-1}]$. Let τ be the set of types to which r is mapped by $\xi(p, d_{k-1})$:

$$\tau = \xi(p, d_{k-1})(r). \quad (38)$$

Assume the Lemma holds for that for a concrete execution path

$$p = [\text{start}_{\text{main}}, n_1, \dots, n_{k-1}],$$

i.e. $\tau \neq \top_T$ for an arbitrary $r \in R$ and $d_{k-1} \in D$.

Induction step: Let $p' = [\text{start}_{\text{main}}, n_1, \dots, n_{k-1}, n_k]$ be a concrete execution path.

Let $e_k = ((n_{k-1}, d_{k-1}), (n_k, d))$. As shown in (34),

$$\xi(p', d)(r) \supseteq \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r).$$

From Definition 5, we can see that unless e_k is a call-start edge or an end-return edge, the result follows from the induction hypothesis. More formally, if e_k is not a call-start or end-return edge, then for all $m \in L_R^{\subseteq}$,

$$\text{EdgeFn}_R^{\subseteq}(e_k)(m) \subseteq m.$$

The edge function corresponding to the call-start and end-return edges is the only place in which the set of types that a receiver maps to can be reduced.

Assume that e_k is an end-return edge with a call on the receiver $r' \in R$ with a signature $s_{\mathcal{F}}$ to a function $f \in \mathcal{F}$.

$$\begin{aligned} \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) &= (\lambda m. m[r' \rightarrow m(r) \cap \tau(s_{\mathcal{F}}, f)][r_1 \rightarrow \perp_T] \dots [r_l \rightarrow \perp_T])(\xi(p, d_{k-1}))(r) \\ &= (\xi(p, d_{k-1})[r' \rightarrow \tau \cap \tau(s_{\mathcal{F}}, f)][r_1 \rightarrow \perp_T] \dots [r_l \rightarrow \perp_T])(r), \end{aligned}$$

where $r_1, \dots, r_l \in R$ are the local variables in the called method.

If $r \in \{r_1, \dots, r_l\}$, then $\text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) = \perp_T \ni t^{15}$.

Otherwise, if $r = r'$, then $\text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) = \tau \cap \tau(s_{\mathcal{F}}, f)$.

According to Lemma 7 and by the induction hypothesis, the runtime type t of r must be contained in $\xi(p, d_{k-1})(r) = \tau$. At the same time, by definition, t is part of $\tau(s_{\mathcal{F}}, f)$. Therefore, $t \in \tau \cap \tau(s_{\mathcal{F}}, f) \subseteq \text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r)$, which means that $\text{EdgeFn}_R^{\subseteq}(e_k)(\xi(p, d_{k-1}))(r) \neq \top_T$.

The same reasoning applies to the case where e_k is a call-start edge. \square

Finally, we can prove the Soundness Lemma.

Proof of Lemma 5. Let $\rho = \mathcal{R}(\mathcal{T}_R^{\subseteq}(P))$. Then

$$\begin{aligned} \mathcal{U}^{\subseteq}(\rho)(n) &= D_n^{\subseteq}(\rho) \\ &= \{d' \mid d' \in \text{MVP}_F(n) \wedge \forall r \in R: \rho(n, d')(r) \neq \top_T\}. \end{aligned}$$

Since $\text{MVP}_F(n) = \bigcap_{q \in \text{VP}(n)} M_F(q)(\emptyset)$, and $p \in \text{VP}(n)$, it follows that

$$\begin{aligned} d &\in M_F(p)(\emptyset) \\ &\subseteq \text{MVP}_F(n). \end{aligned}$$

At the same time, for all receivers $r \in R$,

$$\begin{aligned} \rho(n, d)(r) &= \left(\bigcap_{q \in \text{VP}(n)} \xi(q, d) \right)(r) \\ &= \bigcap_{q \in \text{VP}(n)} \xi(q, d)(r). \end{aligned}$$

According to Lemma 8, $\xi(p, d)(r) \neq \top_T$. Since $p \in \text{VP}(n)$,

$$\xi(p, d)(r) \subseteq \bigcap_{q \in \text{VP}(n)} \xi(q, d)(r).$$

From $\bigcap_{q \in \text{VP}(n)} \xi(q, d)(r) = \rho(n, d)(r)$ it follows that $\xi(p, d)(r) \subseteq \rho(n, d)(r)$. Therefore, $\rho(n, d)(r) \neq \top_T$, and $d \in D_n^{\subseteq}(\rho) = \mathcal{U}^{\subseteq}(\rho)(n)$. \square

Correlated Call Receivers

This appendix contains the proof for Lemma ?? which shows that in a correlated-calls analysis, it is enough to consider only correlated-call receivers R^{\subseteq} .

The following Lemma shows that the types to which a given receiver is mapped in the result of the algorithm is not affected by other receivers and the types to which they are mapped.

¹⁵ In the case of a recursive call, it is possible that both $r \in \{r_1, \dots, r_l\}$ and $r = r'$. In that case, the set to which r will be mapped would be still “overwritten” by \perp_T .

Lemma 9. *Let P be an IFDS problem. Let N^* be the supergraph for P , D the set of data-flow facts, $n \in N^*$ a node, and $p = [\text{start}_{\text{main}}, \dots, n]$ a path in the supergraph. Let $d \in D \cup \{\Lambda\}$. Then for any realizable path $p' \in \text{RP}(p, d)$, set $S \subseteq R$, and receiver $r \in S$,*

$$\text{EdgeFn}_S^{\subseteq}(p')(\top_{\subseteq})(r) = \text{EdgeFn}_{\{r\}}^{\subseteq}(p')(\top_{\subseteq})(r). \quad (39)$$

Proof. By induction on the length of p .

Basis: $p' = [(\text{start}_{\text{main}}, \Lambda)]$. Then $\text{EdgeFn}_S^{\subseteq}(p') = \text{id} = \text{EdgeFn}_{\{r\}}^{\subseteq}(p')$, and the Lemma follows directly.

Induction hypothesis: Suppose that for a path $q = [(\text{start}_{\text{main}}, \Lambda), \dots, (n_{k-1}, d_{k-1})]$, where $q \in \text{RP}(n, d)$, the Lemma holds, i.e. both edge functions map r to the same set of types τ :

$$\begin{aligned} \tau &= \text{EdgeFn}_S^{\subseteq}(q)(\top_{\subseteq})(r) \\ &= \text{EdgeFn}_{\{r\}}^{\subseteq}(q)(\top_{\subseteq})(r). \end{aligned}$$

Induction step: Let $q' = [(\text{start}_{\text{main}}, \Lambda), \dots, (n_{k-1}, d_{k-1}), (n_k, d_k)]$ and the edge $e_k = ((n_{k-1}, d_{k-1}), (n_k, d_k))$.

Observe that for any set $U \subseteq R$ such that $r \in U$,

$$\text{EdgeFn}_U^{\subseteq}(q')(\top_{\subseteq})(r) = \text{EdgeFn}_U^{\subseteq}(e_k)(\text{EdgeFn}_U^{\subseteq}(q)(\top_{\subseteq}))(r). \quad (40)$$

We can see from (29) that there are two cases.

If $d_{k-1} = d_k = \Lambda$, $\text{EdgeFn}_S^{\subseteq}(e_k) = \text{id} = \text{EdgeFn}_{\{r\}}^{\subseteq}(e_k)$, and, due to (40),

$$\begin{aligned} \text{EdgeFn}_S^{\subseteq}(q')(\top_{\subseteq})(r) &= \tau \\ &= \text{EdgeFn}_{\{r\}}^{\subseteq}(q')(\top_{\subseteq})(r). \end{aligned}$$

Otherwise, there are four sub-cases.

1. e_k is a call-start edge, $r'.c()$ is the call site at n_{k-1} with signature $s_{\mathcal{F}}$, $f \in \mathcal{F}$ is the called procedure, and $r' \in U$. Then

$$\text{EdgeFn}_U^{\subseteq}(e_k) = \lambda m. \delta(m)[r' \rightarrow \delta(m)(r) \cap \tau(s_{\mathcal{F}}, f)].$$

There are two sub-cases.

- (a) If $r = r'$, then, according to (40), the resulting set of types

$$\text{EdgeFn}_U^{\subseteq}(q')(\top_{\subseteq})(r) = \delta(\text{EdgeFn}_U^{\subseteq}(q)(\top_{\subseteq}))(r) \cap \tau(s_{\mathcal{F}}, f).$$

If $d_{k-1} = \Lambda$, then $\delta(\text{EdgeFn}_U^{\subseteq}(q)(\top_{\subseteq}))(r) = \perp_{\subseteq}(r) = \perp_T$. If $d_{k-1} \neq \Lambda$, then $\delta(\text{EdgeFn}_U^{\subseteq}(q)(\top_{\subseteq}))(r) = \text{EdgeFn}_U^{\subseteq}(q)(\top_{\subseteq})(r) = \tau$. The set $\tau(s_{\mathcal{F}}, f)$ is the same for either case.

Therefore, the value of $\text{EdgeFn}_U^{\subseteq}(q')(\top_{\subseteq})(r)$ has the same result regardless of U , which means that $\text{EdgeFn}_S^{\subseteq}(q')(\top_{\subseteq})(r) = \text{EdgeFn}_{\{r\}}^{\subseteq}(q')(\top_{\subseteq})(r)$, and the Lemma holds.

(b) If $r \neq r'$, then

$$\text{EdgeFn}_U^{\subseteq}(q')(\top_{\subseteq})(r) = \delta(\text{EdgeFn}_U^{\subseteq}(q)(\top_{\subseteq}))(r), \quad (41)$$

which, as we have seen in Case (1a), does not depend on U , and the Lemma holds.

2. e_k is an end-return edge, $r_1, \dots, r_l \in U$ are the local variables in the callee method, $r'.c()$ is the call corresponding to the return node at n_k , $f \in \mathcal{F}$ is the called method with signature $s_{\mathcal{F}}$, and $r' \in U$. Then

$$\text{EdgeFn}_U^{\subseteq}(e_k) = \lambda m. \delta(m)[r' \rightarrow \delta(m)(r) \cap \tau(s_{\mathcal{F}}, f)][r_1 \rightarrow \perp_T] \dots [r_l \rightarrow \perp_T].$$

There are three sub-cases.

- (a) If $r \in \{r_1, \dots, r_l\}$, then regardless of the value of U ,

$$\text{EdgeFn}_U^{\subseteq}(q')(\top_{\subseteq})(r) = \perp_T,$$

and the Lemma holds.

- (b) Otherwise, if $r = r'$, the case is analogous to Case (1a).
(c) If $r \notin \{r', r_1, \dots, r_l\}$, then see Case (1b).

3. n_{k-1} contains an assignment for $r' \in U$. Then

$$\text{EdgeFn}_U^{\subseteq}(e_k) = \lambda m. \delta(m)[r' \rightarrow \perp_T].$$

If $r = r'$, see Case (2a). If $r \neq r'$, see Case (1b).

4. Otherwise,

$$\text{EdgeFn}_U^{\subseteq}(e_k) = \lambda m. \delta(m),$$

and the case is analogous to Case (1b). \square

The following Lemma shows that the correlated-calls analysis computes the results for each receiver independently, or separately. To compute the set of types to which a receiver r is mapped at each exploded-graph node, we can exclude all other receivers in the program from the analysis (recall from (23) that the set of receivers that are considered in the analysis is specified by the set S in a correlated-calls transformation $\mathcal{T}_S^{\subseteq}$). Therefore, for a given receiver r , the results of a $\mathcal{T}_S^{\subseteq}$ - and a $\mathcal{T}_{\{r\}}^{\subseteq}$ -analysis are the same.

Lemma 10. *Let P be an IFDS problem. Let N^* be the supergraph for P , D the set of data-flow facts, and $S \subseteq R$ a set of receivers. Then for any $n \in N^*$, $d \in D$, and receiver $r \in S$,*

$$\mathcal{R}(\mathcal{T}_S^{\subseteq}(P))(n, d)(r) = \mathcal{R}(\mathcal{T}_{\{r\}}^{\subseteq}(P))(n, d)(r). \quad (42)$$

Proof. According to (??), (9), and (19),

$$\begin{aligned}
\mathcal{R}(\mathcal{T}_S^{\subseteq}(P))(n, d)(r) &= \text{MVP}_{\text{Env}}(n, d)(r) \\
&= \left(\prod_{q \in \text{VP}(n)} M_{\text{Env}}(q)(\Omega)(d) \right)(r) \\
&= \left(\prod_{q \in \text{VP}(n)} \prod_{q' \in \text{RP}(q, d)} \text{EdgeFn}_S^{\subseteq}(q')(\top_{\subseteq}) \right)(r) \\
&= \bigcup_{q \in \text{VP}(n)} \bigcup_{q' \in \text{RP}(q, d)} \text{EdgeFn}_S^{\subseteq}(q')(\top_{\subseteq})(r). \tag{43}
\end{aligned}$$

Then from Lemma 9,

$$\begin{aligned}
\mathcal{R}(\mathcal{T}_S^{\subseteq}(P))(n, d)(r) &= \bigcup_{q \in \text{VP}(n)} \bigcup_{q' \in \text{RP}(q, d)} \text{EdgeFn}_{\{r\}}^{\subseteq}(q')(\top_{\subseteq})(r) \\
&= \mathcal{R}(\mathcal{T}_{\{r\}}^{\subseteq}(P))(n, d)(r). \quad \square
\end{aligned}$$

The next lemma shows that the set of types to which a receiver is mapped in a correlated-calls lattice element can be represented as an intersection of static-type function applications $\tau(s_{\mathcal{F}_i}, f_i)$.

Lemma 11. *For an IFDS problem P , a node $n \in N^*$, and fact $d \in D$, let $p \in \text{RP}(n, d)$ be a realizable path and $r \in R$ a receiver. Then there exists a non-negative number γ of calls on the receiver r with signatures $s_{\mathcal{F}_\gamma}$ to the functions $f_\gamma \in \mathcal{F}_\gamma$, for which*

$$\text{EdgeFn}_{\{r\}}^{\subseteq}(p)(\top_{\subseteq})(r) = \bigcap_{\gamma \geq 0} \tau(s_{\mathcal{F}_\gamma}, f_\gamma).$$

Proof. Let p have the following form¹⁶:

$$p = [(\text{start}_{\text{main}}, \Lambda), (n_1, \Lambda), \dots, (n_k, \Lambda), (n_{k+1}, d_{k+1}), \dots, (n_{k+l}, d_{k+l})],$$

where $l \geq 1$ and the facts for all nodes up to n_k are equal to Λ and $d_{k+i} \in D$ for $0 < i \leq l$.

As previously, for all i , we will denote the edge (n_i, n_{i+1}) by e_i .

From (23) we can infer that

$$\text{EdgeFn}_{\{r\}}^{\subseteq}(p) = \text{EdgeFn}_{\{r\}}^{\subseteq}(e_{k+l}) \circ \dots \circ \text{EdgeFn}_{\{r\}}^{\subseteq}(e_{k+2}) \circ (\lambda m. \beta) \circ \text{id} \circ \dots \circ \text{id},$$

¹⁶ It can be shown from the definition of a pointwise representation in Sagiv et al. [20] that in a realizable path, there is never an edge from a fact of the set D to a Λ fact. Therefore, we can represent p as a sequence of nodes that has a prefix of Λ -fact nodes, after which all nodes are non- Λ facts.

where

$$\beta = \begin{cases} \perp_{\in} [r \rightarrow \tau(s_{\mathcal{F}}, f)] & \text{if } (n_k, n_{k+1}) \text{ is a call-start or end-return edge, and} \\ & \text{the call site } r.c() \text{ with signature } s_{\mathcal{F}} \text{ to the function} \\ & f \in \mathcal{F} \text{ corresponds to the call-start or end-return edge,} \\ \perp_{\in} & \text{otherwise}^{17}. \end{cases}$$

Therefore,

$$\begin{aligned} \text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in}) &= \left(\text{EdgeFn}_{\{r\}}^{\in}(e_{k+l}) \circ \dots \circ \text{EdgeFn}_{\{r\}}^{\in}(e_{k+2}) \right) ((\lambda m. \beta)(\top_{\in})) \\ &= \left(\text{EdgeFn}_{\{r\}}^{\in}(e_{k+l}) \circ \dots \circ \text{EdgeFn}_{\{r\}}^{\in}(e_{k+2}) \circ \text{id} \right) (\beta). \end{aligned} \quad (44)$$

We can now prove the lemma by induction on l .

Basis: If $l = 1$, then $\text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in}) = \text{id}(\beta) = \beta$. There are two cases.

If $\beta = \perp_{\in}$, then

$$\begin{aligned} \text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in})(r) &= \beta(r) \\ &= \perp_T, \end{aligned}$$

and $\gamma = 0$.

If $\beta = \perp_{\in} [r \rightarrow \tau(s_{\mathcal{F}}, f)]$, then

$$\text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in})(r) = \tau(s_{\mathcal{F}}, f),$$

and $\gamma = 1$.

Induction hypothesis: Assume that for a path $p = [(\text{start}_{\text{main}}, \Lambda), \dots, (n_{k+l}, d_{k+l})]$, the Lemma holds for $\gamma = N$, where $N \geq 0$.

Induction step: Let $p' = [(\text{start}_{\text{main}}, \Lambda), \dots, (n_{k+l}, d_{k+l}), (n_{k+l+1}, d_{k+l+1})]$.

Recall that

$$\text{EdgeFn}_{\{r\}}^{\in}(p')(\top_{\in})(r) = \text{EdgeFn}_{\{r\}}^{\in}(e_{k+l+1}) \left(\text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in}) \right) (r).$$

From (22) we can see that unless e_{k+l+1} is a call-start or end-return edge corresponding to a call on the receiver r , then $\text{EdgeFn}_{\{r\}}^{\in}(e_{k+l+1})(r)$ must be equal to either \perp_T or $m(r)$, where $m = \text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in})$.

If $\text{EdgeFn}_{\{r\}}^{\in}(e_{k+l+1})(r) = \perp_T$, then the Lemma holds for $\gamma = 0$.

Otherwise,

$$\begin{aligned} \text{EdgeFn}_{\{r\}}^{\in}(e_{k+l+1})(\top_{\in})(r) &= \text{EdgeFn}_{\{r\}}^{\in}(p)(\top_{\in})(r) \\ &= \bigcap_N \tau(s_{\mathcal{F}_N}, f_N), \end{aligned}$$

¹⁷ Since $d_k = \Lambda$ and $d_{k+1} \neq \Lambda$, the micro function for the edge e_{k+1} is equal to $\lambda m. \varepsilon_{\{r\}}(e_{k+1})(\perp_{\in})$. From the definition of ε_S (22) we can see that the only case where $\varepsilon_{\{r\}}(e_{k+1})(m)$ would not be equal to \perp_{\in} is when e_{k+1} is call-start or end-return edge.

and therefore $\gamma = N$.

Suppose that e_{k+l+1} is a call-start edge with a call on the receiver r with signature s_G to a function $g \in \mathcal{G}$. Then, according to (22),

$$\text{EdgeFn}_{\{r\}}^{\subseteq}(e_{k+l+1}) = \lambda m . m[r \rightarrow m(r) \cap \tau(s_G, g)].$$

Therefore,

$$\begin{aligned} \text{EdgeFn}_{\{r\}}^{\subseteq}(p')(\top_{\subseteq})(r) &= \lambda m . m[r \rightarrow m(r) \cap \tau(s_G, g)] \left(\text{EdgeFn}_{\{r\}}^{\subseteq}(p)(\top_{\subseteq}) \right)(r) \\ &= \text{EdgeFn}_{\{r\}}^{\subseteq}(p)(\top_{\subseteq})(r) \cap \tau(s_G, g) \\ &= \left(\bigcap_N \tau(s_{\mathcal{F}_N}, f_N) \right) \cap \tau(s_G, g), \end{aligned}$$

and the Lemma holds for $\gamma = N + 1$.

The case where e_{k+l+1} is an end-return edge is analogous to the previous case. \square

We now show that a receiver will be only mapped to \top_{\subseteq} if it is the receiver of a correlated call.

Lemma 12. *For an IFDS problem P , let $n \in N^*$ be a node, and $d \in D$ a dataflow fact such that there exists a realizable path $p \in RP(n, d)$. Let T be the set of all types in the program. If there exists a receiver $r \in R$ such that*

$$\text{EdgeFn}_{\{r\}}^{\subseteq}(p)(\top_{\subseteq})(r) = \top_T,$$

then $r \in R^{\subseteq}$.

Proof. According to Lemma 11,

$$\text{EdgeFn}_{\{r\}}^{\subseteq}(p)(\top_{\subseteq})(r) = \bigcap_{\gamma \geq 0} \tau(s_{\mathcal{F}_\gamma}, f_\gamma).$$

Let $\tau_i = \tau(s_{\mathcal{F}_i}, f_i)$. For a given k , let $r.m_k()$ be the call site corresponding to τ_k , and T' the set of types compatible with the static type of r . Recall from Section 3.3 that

- $\tau_k \neq \top_T$;
- if $\tau_k = T'$ then the corresponding call site is monomorphic;
- if $\tau_k \subset T'$ then the call site is polymorphic.

From the conditions of the Lemma,

$$\bigcap_{\gamma \geq 0} \tau_\gamma = \top_T. \tag{45}$$

If all $\tau_k = T'$, then $\bigcap_{\gamma \geq 0} \tau_\gamma$ is also equal to T' . Since $T' \neq \top_T$, this is a contradiction.

If exactly one $\tau_k \subset T'$ and the rest are equal to T' , then $\bigcap_{\gamma \geq 0} \tau_\gamma$ is equal to τ_k , which cannot be \top_T either.

Therefore, there are at least two sets, τ_i and τ_j , which are strict subsets of T' . Since both τ_i and τ_j are non-empty and their intersection equals \top_T , τ_i and τ_j must be disjoint. If τ_i and τ_j are disjoint, they must correspond to different call sites.

In other words, there are at least two calls on the same receiver for which the static-type function is a strict subset of the set of types compatible with a given receiver r . It follows that both calls have to be polymorphic. Therefore, $r \in R^\subseteq$. \square

We will now show that if a receiver ever gets mapped to top, then it is a correlated-calls receiver.

Lemma 13. *For an IFDS problem P , let $n \in N^*$ be a node, and $d \in D$ a dataflow fact such there exists a realizable path $p \in RP(n, d)$. Then, if there exists a receiver $r \in R$, such that*

$$\mathcal{R}(\mathcal{T}_{\{r\}}^\subseteq(P))(n, d)(r) = \top_T,$$

then $r \in R^\subseteq$.

Proof. As shown in (43),

$$\mathcal{R}(\mathcal{T}_{\{r\}}^\subseteq(P))(n, d)(r) = \bigcup_{q \in VP(n)} \bigcup_{q' \in RP(q, d)} \text{EdgeFn}_{\{r\}}^\subseteq(q')(\top_\subseteq)(r).$$

Since the latter is equal to \top_T , it follows that for each realizable path p' to node n , $\text{EdgeFn}_{\{r\}}^\subseteq(p')(\top)(r) = \top_T$. According to Lemma 13, this is only possible if $r \in R^\subseteq$. \square

Finally, we present the proof for Lemma ?? which states that a correlated-calls analysis that considers all receivers computes the same result as an analysis that considers only correlated-call receivers.

Proof of Lemma 6. From (??) we know that

$$\mathcal{U}^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P))) = \{(n, D_n^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P)))) \mid n \in N^*\}.$$

According to (??) and Lemma 10, for a given $n \in N^*$,

$$\begin{aligned} D_n^\subseteq(\mathcal{R}(\mathcal{T}_R^\subseteq(P))) &= \left\{ d \mid d \in MVP_F(n) \wedge \forall r \in R : \left\{ (r, \mathcal{R}(\mathcal{T}_{\{r\}}^\subseteq(P))(n, d)(r)) \mid r \in R \right\} (r) \neq \top_T \right\} \\ &= \left\{ d \mid d \in MVP_F(n) \wedge \forall \mathbf{r} \in \mathbf{R} : \mathcal{R}(\mathcal{T}_{\{\mathbf{r}\}}^\subseteq(P))(n, d)(\mathbf{r}) \neq \top_T \right\}. \end{aligned}$$

Since, according to Lemma 13, $\mathcal{R}(\mathcal{T}_{\{r\}}^{\mathbb{E}}(P))(n, d)(r)$ can only be equal to \top_T when $r \in R^{\mathbb{E}}$, we can conclude that

$$\begin{aligned} D_n^{\mathbb{E}}(\mathcal{R}(\mathcal{T}_R^{\mathbb{E}}(P))) \\ &= \left\{ d \mid d \in \text{MVP}_F(n) \wedge \forall r \in R^{\mathbb{E}} : \mathcal{R}(\mathcal{T}_{\{r\}}^{\mathbb{E}}(P))(n, d)(r) \neq \top_T \right\} \\ &= D_n^{\mathbb{E}}(\mathcal{R}(\mathcal{T}_{R^{\mathbb{E}}}^{\mathbb{E}}(P))). \end{aligned}$$

Therefore,

$$\begin{aligned} \mathcal{U}^{\mathbb{E}}(\mathcal{R}(\mathcal{T}_R^{\mathbb{E}}(P))) &= \{ (n, D_n^{\mathbb{E}}(\mathcal{R}(\mathcal{T}_{R^{\mathbb{E}}}^{\mathbb{E}}(P)))) \mid n \in N^* \} \\ &= \mathcal{U}^{\mathbb{E}}(\mathcal{R}(\mathcal{T}_{R^{\mathbb{E}}}^{\mathbb{E}}(P))). \end{aligned} \quad \square$$

Representation of Micro Functions

In this part of the appendix we present the proofs to the Lemmas related to the representation of micro functions with update maps.

Proof of Lemma ??. Let $\text{update}_{f,r}^* = \langle I, U \rangle$. For any $\tau \in T$,

$$\begin{aligned} \llbracket \mathcal{N}(\text{update}_{f,r}^*) \rrbracket (\tau) &= \llbracket \mathcal{N}(\langle I, U \rangle) \rrbracket (\tau) \\ &= \llbracket \langle I \cup U, U \rangle \rrbracket (\tau) \\ &= \tau \cap (I \cup U) \cup U \\ &= (\tau \cap I) \cup (\tau \cap U) \cup U \\ &= \tau \cap I \cup U \\ &= \llbracket \langle I, U \rangle \rrbracket (\tau) \\ &= \llbracket \text{update}_{f,r}^* \rrbracket (\tau). \end{aligned}$$

Thus, $\llbracket \mathcal{N}(\text{update}_{f,r}^*) \rrbracket = \llbracket \text{update}_{f,r}^* \rrbracket$. \square

Proof of Lemma ??. Let us show that there always exists a set $\tau \subseteq T$ such that $\llbracket \langle I, U \rangle \rrbracket (\tau) \neq \llbracket \langle I', U' \rangle \rrbracket (\tau)$. There are two cases:

1. $U \neq U'$. Then for the empty set $\tau = \emptyset$,

$$\llbracket \langle I, U \rangle \rrbracket (\tau) = \llbracket \langle I, U \rangle \rrbracket (\emptyset) = (\emptyset \cap I) \cup U = U,$$

whereas

$$\llbracket \langle I', U' \rangle \rrbracket (\tau) = \llbracket \langle I', U' \rangle \rrbracket (\emptyset) = (\emptyset \cap I') \cup U' = U'.$$

Hence, $\llbracket \langle I, U \rangle \rrbracket \neq \llbracket \langle I', U' \rangle \rrbracket$.

2. $I \neq I'$. Then for the set of all types $\tau = T$,

$$\llbracket \langle I, U \rangle \rrbracket (\tau) = \llbracket \langle I, U \rangle \rrbracket (T) = (T \cap I) \cup U = I \cup U.$$

Since $\langle I, U \rangle$ is normalized, $U \subseteq I$, and

$$I \cup U = I.$$

At the same time,

$$\llbracket \langle I', U' \rangle \rrbracket (\tau) = \llbracket \langle I', U' \rangle \rrbracket (T) = (T \cap I') \cup U' = I' \cup U' = I'.$$

Since $I \neq I'$, it follows that $\llbracket \langle I, U \rangle \rrbracket \neq \llbracket \langle I', U' \rangle \rrbracket$. \square

Proof of Lemma ??. 1. The identity function is represented as

$$\llbracket \text{id} \rrbracket = \{(r, \langle \perp_T, \top_T \rangle) \mid r \in R^\subseteq\};$$

the top function is represented as

$$\llbracket \lambda m. \top_\subseteq \rrbracket = \{(r, \langle \top_T, \top_T \rangle) \mid r \in R^\subseteq\}.$$

2. Equations (??) and (??) show that the representation of micro functions is closed under composition and meet.
3. To show that our representation for micro functions forms a lattice with finite height, let us first show that $L_{R^\subseteq}^\subseteq : R^\subseteq \rightarrow 2^T$ forms a lattice. Since T is a finite set, $(2^T, \subseteq)$ is a finite-height lattice. R^\subseteq is a finite set. Hence, the mapping

$$R^\subseteq \mapsto 2^T = \{(r, t) \mid r \in R^\subseteq, t \in 2^T\} = L_{R^\subseteq}^\subseteq$$

also forms a finite-height lattice [15].

Furthermore, $L_{R^\subseteq}^\subseteq$ is a finite set. Every element of $L_{R^\subseteq}^\subseteq$ can be applied to $|R^\subseteq|$ receivers, where each receiver is mapped to a set of types. There are $|R^\subseteq| \cdot 2^{|T|}$ different possibilities to form those mappings, so

$$|L_{R^\subseteq}^\subseteq| = |R^\subseteq| \cdot 2^{|T|}.$$

Therefore, $L_{R^\subseteq}^\subseteq \mapsto L_{R^\subseteq}^\subseteq$ also forms a finite-height lattice.

4. All operations can be computed in $O(R^\subseteq \times T)$ time. Note that the R^\subseteq and T sets are an input to the correlated-calls analysis, and the time it takes to compute the meet or composition of micro functions is independent of the representation of the specific operand micro functions.
5. The space bound is $O(R^\subseteq \times T)$. \square