Trace Analysis for Predicting the Effectiveness of Partial Evaluation*

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1 Introduction

The main goal of partial evaluation [3] is program specialization. Essentially, given a program and part of its input data—the so called static data—a partial evaluator returns a new, residual program which is specialized for the given data. An appropriate residual program for executing the remaining computations—those that depend on the so called dynamic data—is thus the output of the partial evaluator. Despite the fact that the main goal of partial evaluation is improving program efficiency (i.e., producing faster programs), there are very few approaches devoted to formally analyze the effects of partial evaluation, either a priori (prediction) or a posteriori. Recent approaches (e.g., [1,5]) have considered experimental frameworks for estimating the best division (roughly speaking, a classification of program parameters into static or dynamic), so that the optimal choice is followed when specializing the source program.

Here, we introduce an alternative, *symbolic* approach for predicting the potential effects of partial evaluation (which is, in principle, computationally less expensive). Basically, we first generate a finite representation that safely describes all possible *call traces* (i.e., sequences of predicate calls) for a given program. Then, we analyze how this finite representation would change by a particular partial evaluation. By comparing the original and the transformed representations, one may in some cases predict the effects of running the partial evaluator. A more detailed description of our approach can be found in [6].

2 Trace Analysis for Logic Programs

We consider a fixed domain of predicate symbols Π . We assume that Π do not contain occurrences of the same predicate name with different arities. Furthermore, we consider a fixed computation rule for call traces, namely Prolog's leftmost computation rule, which we denote by \mathcal{R}_{left} . We label SLD resolution steps with the predicate symbol of the selected atom, i.e., we write $Q_0 \stackrel{p_0}{\leadsto} Q_1 \stackrel{p_1}{\leadsto} \dots$ with $\operatorname{pred}(\mathcal{R}_{left}(Q_i)) = p_i \in \Pi$, $i \geq 0$, where $\operatorname{pred}(A)$ returns the predicate symbol of atom A.

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Definition 1 (call trace). Let P be a program and Q_0 a query. We say that $\tau = p_0 \ p_1 \dots p_{n-1} \in \Pi^*, \ n \geq 1$, is a call trace for Q_0 with P iff there exists a successful SLD derivation $Q_0 \stackrel{p_0}{\leadsto} Q_1 \stackrel{p_1}{\leadsto} \dots \stackrel{p_{n-1}}{\leadsto} Q_n$.

The first step of our trace analysis consists in producing a context-free grammar (CFG) associated to the considered program. A CFG is a tuple $G = \langle \Sigma, N, R, S \rangle$, where Σ and N are two disjoint sets of terminals and non-terminals, respectively, $S \in N$ is the start symbol, and R is a set of rules. In the following, given a predicate symbol $p \in \Pi$, we denote by $\overline{p} \notin \Pi$ a fresh symbol representing the non-terminal associated to p. Furthermore, we let $\overline{pred}(A) = \overline{p}$ if $A = p(t_1, \ldots, t_n)$. Also, we let $\overline{\Pi}$ denote the set $\{\overline{p} \mid p \in \Pi\}$ of non-terminals associated to predicate symbols. In contrast, we directly use predicate symbols from Π as terminals. We let START be a fresh symbol not in $\Pi \cup \overline{\Pi}$ which we use as a generic start symbol for CFGs.

Definition 2 (trace CFG, CFG_q^P). Let P be a program and $q \in \Pi$ a predicate symbol. The associated trace CFG is $CFG_q^P = \langle \Pi, \overline{\Pi} \cup \{START\}, R, START \rangle$, where

$$R = \{ \underbrace{\text{START}}_{} \to \overline{q} \}$$

$$\cup \{ \underbrace{pred}_{}(A_0) \to pred}_{}(A_0) \underbrace{pred}_{}(B_1) \dots \underbrace{pred}_{}(B_n) \mid A_0 \leftarrow B_1, \dots, B_n \in P, n \geq 0 \}$$

Roughly speaking, the trace CFG associated to a logic program mimics the execution of the original program by replacing queries (sequences of atoms) by sequences of non-terminals and by producing a terminal with the predicate symbol of the selected atom at each SLD-resolution step.

Example 1. Consider the following program P which defines a procedure for multiplying all elements of a list by a given value:¹

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(c_1) mlist([], I, []).
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- (c_2) $mlist([X|R], I, L) \leftarrow ml(X, R, I, L).$
- (c_3) $ml(X, R, I, [XI|RI]) \leftarrow mult(X, I, XI), mlist(R, I, RI).$
- (c_4) mult(0, Y, 0). (c_5) $mult(s(X), Y, Z) \leftarrow mult(X, Y, Z1)$, add(Z1, Y, Z).
- (c_6) add(X,0,X). (c_7) $add(X,s(Y),s(Z)) \leftarrow add(X,Y,Z)$.

where natural numbers are built from 0 and $s(\cdot)$. The associated trace CFG is $CFG^P_{mlist} = \langle \{mlist, ml, mult, add\}, \{START, MLIST, ML, MULT, ADD\}, R, START\rangle$, where the set of rules R is as follows:

In [6], we prove that CFG_q^P is indeed a correct approximation of the call traces for P w.r.t. the leftmost computation rule \mathcal{R}_{left} .

¹ In the examples, we write non-terminals associated to predicates using capital letters.

Unfortunately, trace CFGs do not always allow us to produce a simple and compact representation of the call traces of a program (e.g., when the associated language is not regular). To overcome this drawback, we use the transformation from [4] to approximate a trace CFG with a *strongly regular* grammar (SRG). The relevance of SRGs is that they can be mapped to equivalent finite-state automata using an efficient algorithm. Moreover, the transformation of [4] guarantees that the result remains readable and mainly preserves the structure of the original CFG, which is particularly useful in our context.

A grammar is *left-linear* if every rule has either the form $(A \to t)$ or $(A \to t)$, where t is a finite sequence of terminals and A, B are non-terminals.

Definition 3 (trace SRG, SRG_q^P). Let P be a program and $q \in \Pi$ a predicate symbol. The associated trace SRG, SRG_q^P , is obtained from CFG_q^P as follows. First, we compute the sets of mutually recursive non-terminals of CFG_q^P . Then, for each set M of mutually recursive non-terminals such that their rules are not all left-linear w.r.t. the non-terminals of M (i.e., considering non-terminals from $(\overline{\Pi} \setminus M)$ as terminals), we apply a grammar transformation as follows:

- 1. For each non-terminal $A \in M$, we introduce a fresh non-terminal A' and add the rule $A' \to \epsilon$ to the grammar (we denote by ϵ the empty sequence).
- 2. For each non-terminal $A \in M$ and each rule $A \to t_0 B_1 t_1 B_2 t_2 \dots B_m t_m$ of CFG_q^P with $m \ge 0, B_1, \dots, B_m \in M, t_0, \dots, t_m \in (\Pi \cup (\overline{\Pi} \setminus M))^*$, we replace this rule by the following set of rules:

$$A \to t_0 B_1$$
 $B'_1 \to t_1 B_1$... $B'_{m-1} \to t_{m-1} B_m$ $B'_m \to t_m A'$

(Note that this set reduces to $A \rightarrow t_0 A'$ when m = 0.)

We let $SRG_q^P = \langle \Pi, \overline{\Pi} \cup N \cup START, R', START \rangle$, where R' are the rules obtained as described above and N are the fresh non-terminals added during this process.

Example 2. Consider the CFG_q^P of Example 1. The sets of mutually recursive nonterminals are $\{\{\mathrm{MLIST},\mathrm{ML}\},\{\mathrm{MDLT}\},\{\mathrm{ADD}\}\}$. Here, the rules for both MLIST and ML are left-linear w.r.t. $\{\mathrm{MLIST},\mathrm{ML}\}$. The rules for ADD are clearly left-linear too. However, the second rule of MULT is not left-linear because, even if ADD is treated as a terminal, it appears to the right of the non-terminal MULT. Therefore, in SRG_{mlist}^P we replace the original rules for MULT by the following ones:

$$\{ \text{MULT}' \to \epsilon, \text{ MULT} \to \text{mult MULT}', \text{ MULT} \to \text{mult MULT}, \text{ MULT}' \to \text{ADD MULT}' \}$$

Once we have an SRG that safely approximates the call traces of a program, there are several possibilities for representing the language generated by this SRG in a compact and intuitive way. Here, we consider the generation of a *finite-state automaton* (FA) that accepts the language generated by the SRG; an alternative approach that produces *regular expressions* can be found in [6].

A finite-state automaton (FA) is specified by a tuple $\langle Q, \Sigma, \delta, s_0, F \rangle$, where Q is a set of states, Σ is an input alphabet, $\delta \subseteq Q \times \Sigma \times Q$ is a set of transitions, $s_0 \in Q$ is the start state and $F \subseteq Q$ is a set of final states. For constructing a

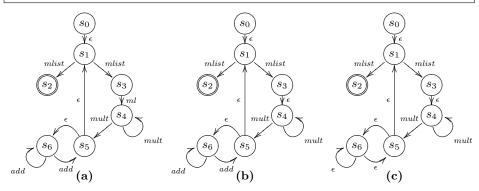


Fig. 1. Transformation of trace FAs

finite automaton FA(G) from an SRG G, called $trace\ FA$, we follow the classical approach: there is a start state associated to the start symbol of the SRG; for each reduction $w \to w'$ with a rule $A \to t B$ of the SRG, we have a transition (s, α, s') in the FA, where states s, s' are associated with the sequence of non-terminals in w, w' and character α is set to the sequence t in the applied rule.

Example 3. Consider the SRG SRG_{mlist}^P of Example 2. The associated FA is shown in Fig. 1 (a), where the final state s_2 is denoted with a double circle.

3 Towards Predicting the Speedup of Partial Evaluation

The trace analysis gives us the *context* where every predicate call appears. Now, we informally describe two transformations (a more formal definition can be found in [6]) that modify the computed traces to account for the potential effects of a partial evaluation. By analyzing the traces before/after partial evaluation, one can extract useful conclusions on its effectiveness.

The first transformation is used to eliminate *intermediate* predicates. Basically, for every state with exactly one input transition and one output transition, we replace the label of the output transition by ϵ (i.e., we delete calls to predicates which are called from a single program point). Consider the trace FA of the program of Ex. 1 which is shown in Fig. 1 (a). After the elimination of intermediate states (the case of s_3), we get the trace FA shown in Fig. 1 (b).

Our second transformation is parameterized by the output of a binding-time analysis (BTA), which annotates each predicate with either unfold or memo, where unfold means that the predicate can be safely unfolded (i.e., without entering an infinite loop), and memo means that it should be specialized (i.e., a residual predicate is produced). Basically, our second transformation replaces the labels of unfoldable predicates with ϵ . Consider, e.g., that the output of a BTA annotates mlist, ml, and mult as memo and add as unfold—this is the case

when the second argument of the initial call to *mlist* is static. We then get the trace FA of Fig. 1 (c). Here, we achieve a significant improvement since, in every iteration for *mlist*, we save the (recursive) evaluation of the calls to *add*.

Clearly, we could eliminate those states whose transitions are all labeled with ϵ . However, we think that keeping the structure of the original trace FA may help the user—and automated analysis tools—to formally compare the original and transformed trace FAs.

4 Discussion

The closest approach to our trace analysis is that of [2], though we offer a different trade-off between analysis cost and accuracy. Basically, they generate trace *terms* abstracting computation trees independently of a computation rule, while we generate sequences of predicate calls for a specific computation rule; also, they do not include a technique for enumerating the (possibly infinite) set of trace terms of a program, while this is a key ingredient of our approach (though one could also apply the transformation from [4] to the CFG associated to the trace terms of [2] to obtain a finite representation). A deeper comparison with the approach of [2] is an interesting topic for further research.

A proof-of-concept implementation of our technique, called PEPE, is publicly available from http://german.dsic.upv.es/pepe.html. Our approach can be seen as a first step for the development of automated techniques and tools for predicting the potential speedup of partial evaluation, thus it opens a number of interesting lines for further research.

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