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Supercompiler HOSC 1.0: under the hood

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# KELDYSH INSTITUTE OF APPLIED MATHEMATICS Russian Academy of Sciences

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Supercompiler HOSC 1.0: under the hood

Moscow 2009

### Ilya G. Klyuchnikov. Supercompiler HOSC 1.0: under the hood

The paper describes the internal structure of HOSC, an experimental supercompiler dealing with programs written in a higher-order functional language. A detailed and formal account is given of the concepts and algorithms the supercompiler is based upon.

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# Илья Г. Ключников. Суперкомпилятор HOSC 1.0: внутренняя структура

В работе описана внутренняя структура экспериментального суперкомпилятора HOSC. Дано полное описание всех существенных понятий и алгоритмов суперкомпилятора, работающего с функциональным языком высшего порядка.

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## 1 Introduction

Supercompilation is a program transformation technique, suggested by V.F. Turchin in the early 1970-s [21, 22].

This paper describes the specializer HOSC, an experimental supercompiler dealing with programs written in a simple higher-order functional language.

A number of supercompilers for higher-order functional languages have already being described in the literature [8, 10, 11, 12, 15, 14]. However, the published descriptions of concepts and algorithms related to higher-order supercompilation suffer from being incomplete: a number of essential details are either omitted or presented in sketchy and informal style.

The goal of the present work is to provide a *detailed* and *formal* description of the concepts and algorithms the supercompiler HOSC is based upon. The source code of HOSC is publicly available and can be found at http://hosc.googlecode.com.

## 2 Notation

For the sake of conciseness and readability we use overlining to denote constructs that may be repeated, so that  $\overline{A_i}$  means  $A_0A_1 \dots A_n$ , if  $A_0$  is defined, or  $A_1 \dots A_n$  otherwise, where n is determined by the context. Thus, an overlined construct should contain parts with subscripts and, in a sense, is similar to a list comprehension. The doubly overlined text is unfolded "inside out" with respect to overlines and "from left to right" with respect to indices. If the overlined text is found inside the parentheses denoting a set or a list ({}}

```
tDef ::= data \ tCon = \overline{dCon_i};
                                                                     type definition
tCon ::= tn \ \overline{tv_i}
                                                                     type constructor
dCon := c \ \overline{type_i}
                                                                    data constructor
type := tv \mid tCon \mid type \rightarrow type \mid (type)
                                                                     type expression
proq := \overline{tDef_i} \ e \ \mathbf{where} \ \overline{f_i = e_i};
                                                                    program
                                                                    variable
e ::= v
      |c| \overline{e_i}
                                                                    constructor
       \mid f \mid
                                                                    function
       \lambda \ \overline{v_i} \to e
                                                                     \lambda-abstraction
       e_1 e_2
                                                                    application
       | case e_0 of \{\overline{p_i \to e_i};\}
                                                                    case-expression
      | letrec f = \lambda \overline{v_i} \rightarrow e_0 in e_1
                                                                    local function
                                                                    parenthesized expression
p := c \overline{v_i}
                                                                    pattern
```

Figure 1: HLL grammar

or [], respectively), we assume that delimiters (commas) are inserted correctly during unfolding. For example:

```
\begin{array}{l} let \ \overline{v_i=e_i;} \ in \ e \Rightarrow let \ v_1=e_1; \ v_2=e_2; \ \dots; \ v_n=e_n; \ in \ e \\ [\gamma, \ \overline{\gamma_i}] \Rightarrow [\gamma, \ \gamma_1, \ \gamma_2, \ \dots, \ \gamma_n] \\ case \ e_0 \ of \ \{\overline{c_i} \ \overline{v_{ik}} \rightarrow e_i;\} \Rightarrow \\ case \ e_0 \ of \ \{c_1 \ \overline{v_{1k}} \rightarrow e_1; \ \dots; \ c_n \ \overline{v_{nk}} \rightarrow e_n;\} \Rightarrow \\ case \ e_0 \ of \ \{c_1 \ v_{11} \ \dots \ v_{1m_1} \rightarrow e_1; \ \dots; \ c_n \ v_{n1} \ \dots \ v_{nm_n} \rightarrow e_n;\} \end{array}
```

## 3 Input language

HOSC transforms programs written in HLL, a simple higher-order lazy language. HLL is statically typed using the Hindley-Milner polymorphic typing system [4].

A program in HLL consists of a number of data type definitions, an expression to be evaluated and a set of function definitions.

A left-hand side of a data type definition is a type name (more precisely, a type constructor name) followed by a list of type variables. The right-hand side consists of one or more constructor declarations (delimited by a vertical bar,  $\overline{dCon_i} \Rightarrow dCon_1 \mid \ldots \mid dCon_n$ ).

The grammar of HLL is shown in Fig. 1. An expression is either a variable, a constructor, a  $\lambda$ -abstraction, an application, a case-expression, a local func-

```
data List a = Nil | Cons a (List a); 

(compose (map f)(map g)) xs where 

compose = \lambda f g x \rightarrow f (g x); 

map = \lambda f xs \rightarrow case xs of { 

    Nil \rightarrow Nil; 

    Cons x1 xs1 \rightarrow Cons (f x1) (map f xs1); 

};
```

Figure 2: A simple HLL program

tion definition or an expression in parenthesis. A function definition binds a global variable to a  $\lambda$ -abstraction.

The target expression e in a program prog may contain free variables. All local variables in global function definitions should be bound.

An expression  $e_0$  within a case expression is called a selector, and expressions  $\overline{e_i}$  are called branches. We require *all* constructors of a corresponding data type to be enumerated within a case expression, so that patterns in the case expression are *non-overlapping* and *exhaustive*.

We will use two different notations for applications:  $e_1$   $e_2$  and  $e_0$   $\overline{e_i}$ . In the first case, an expression  $e_1$  may be any well-typed expression. In the second case, we require the list of arguments to be non-empty and the expression  $e_0$  not to be an application.

An example of a simple HLL program is shown in Fig. 2.

#### 3.1 Basic definitions

The following definitions will be used throughout the whole paper.

#### 3.1.1 Free and bound variables

A variable is bound if it is either (1) the argument of an enclosing  $\lambda$ -abstraction or (2) is defined in the pattern of an enclosing branch of a case-expression. Variables that are not bound are considered to be free. To avoid problems caused by possible name clashes, we require all variable names to be unique in the following sense: for any variable v and expression e all occurrences of the variable v in the expression e must be either free or bound.

#### 3.1.2 Equal expressions

Expressions  $e_1$  and  $e_2$  are considered to be equal, if they differ only in the names of bound variables and there is a one-to-one correspondence between the bound variables of  $e_1$  and  $e_2$ . The equality of expressions  $e_1$  and  $e_2$  is denoted as  $e_1 \equiv e_2$ .

#### 3.1.3 Free variables

fv[e] denotes the set of the free variables of the expression e and is defined inductively as follows:

#### 3.1.4 Bound variables

bv[e] denotes the set of the bound variables of the expression e and is defined inductively as follows:

$$\begin{array}{lll} bv \llbracket v \rrbracket & = \{\} \\ bv \llbracket c \, \overline{e_i} \rrbracket & = \bigcup bv \llbracket e_i \rrbracket \\ bv \llbracket \lambda v \to e \rrbracket & = bv \llbracket e_1 \rrbracket \cup \{v\} \\ bv \llbracket case \, e_0 \, of \, \{ \overline{c_i \, \overline{v_{ik}}} \to e_i; \} \rrbracket & = bv \llbracket e_0 \rrbracket \cup \bigcup bv \llbracket e_i \rrbracket \cup \bigcup v_{ik} \\ bv \llbracket letrec \, f = e_1 \, in \, e_2 \rrbracket & = bv \llbracket e_1 \rrbracket \cup bv \llbracket e_2 \rrbracket \cup \{f\} \end{array}$$

#### 3.1.5 Substitution

A substitution is a finite list of pairs of the form

$$\theta = \{v_1 := e_1, v_2 := e_2, \dots, v_n := e_n\}$$

each pair binding a variable  $v_i$  to its value  $e_i$ .

The domain of a substitution  $\theta$  is denoted by  $domain(\theta)$  and is defined as follows:

$$domain(\{v_1 := e_1, v_2 := e_2, \dots, v_n := e_n\}) = \{v_1, v_2, \dots, v_n\}$$

#### 3.1.6 Application of a substitution

 $e \theta$  denotes the result of applying a substitution  $\theta$  to an expression e, which, informally speaking, is obtained in the following way.

- All free variables of expression e, which are in  $domain(\theta)$ , are replaced by their values from  $\theta$ .
- All other parts of the expression e remain unchanged.

To avoid problems with variable name clashes, we require the domain of  $\theta$  not to contain bound variables of e, i.e.

$$domain(\theta) \cap bv[e] = \emptyset$$

More formally,  $e \theta$  is defined as follows:

$$\begin{array}{lll} v\theta & = & e & \text{if } v := e \in \theta \\ & = & v & \text{otherwise} \\ (c \, \overline{e_i})\theta & = & c \, \overline{(e_i\theta)} \\ (\lambda v \to e)\theta & = & \lambda v \to (e\theta) \\ (e_1 \, e_2)\theta & = & (e_1\theta) \, (e_2\theta) \\ (case \, e_0 \, of \, \{\overline{p_i \to e_i};\})\theta & = & case \, (e_0\theta) \, of \, \{\overline{p_i \to (e_i\theta)};\} \\ (letrec \, f = e_1 \, in \, e_2)\theta & = & letrec \, f = (e_1\theta) \, in \, (e_2\theta) \end{array}$$

#### 3.1.7 Instance of an expression

An expression  $e_2$  is said to be an instance of an expression  $e_1$ , or  $e_1 \leqslant e_2$ , if there is a substitution  $\theta$ , such that  $e_1\theta \equiv e_2$ . As far as the language HLL is concerned, the substitution  $\theta$  is unique and is denoted by  $e_1 \otimes e_2$ .

#### 3.1.8 Renaming

An expression  $e_2$  is a renaming of an expression  $e_1$ ,  $e_1 \simeq e_2$ , if  $e_1 \lessdot e_2$  and  $e_2 \lessdot e_1$ . That is,  $e_1$  and  $e_2$  differ in names of free variables only.

#### 3.1.9 Generalization

A generalization of expressions  $e_1$  and  $e_2$  is a triple  $(e_g, \theta_1, \theta_2)$ , where  $e_g$  is an expression and  $\theta_1$  and  $\theta_2$  are substitutions, such that  $e_g\theta_1 \equiv e_1$  and  $e_g\theta_2 \equiv e_2$ . The set of generalizations of expressions  $e_1$  and  $e_2$  is denoted by  $e_1 \smallfrown e_2$ .

```
\begin{array}{lll} obs & ::= & v \; \overline{e_i} \mid c \; \overline{e_i} \mid (\lambda v \rightarrow e) \\ con & ::= & \langle \rangle \mid con \; e \mid case \; con \; of \; \{\overline{p_i \rightarrow e_i};\} \\ red & ::= & f \mid (\lambda v \rightarrow e_0) \; e_1 \mid case \; v \; \overline{e'_j} \; of \; \{\overline{p_i \rightarrow e_i};\} \\ & \mid & case \; c \; \overline{e'_j} \; of \; \{\overline{p_i \rightarrow e_i};\} \end{array}
```

Figure 3: Observables, contexts, redexes

#### 3.1.10 Most specific generalization

A generalization  $(e_g, \theta_1, \theta_2) \in e_1 \frown e_2$  is a most specific one, if for any generalization  $(e'_g, \theta'_1, \theta'_2) \in e_1 \frown e_2$  holds that  $e'_g \lessdot e_g$ , that is  $e_g$  is an instance of  $e'_g$ . We suppose that there is defined an operation  $\sqcap$ , such that  $e_1 \sqcap e_2$  is a most specific generalization of  $e_1$  and  $e_2$ .

## 3.2 $\lambda$ -lifting

Note that the construction where is no more than syntactic sugar, since the global function definitions may always be transformed into letrec-s and moved to the target expression. Thus, any program written in HLL may always be replaced by an equivalent program consisting of type definitions and a *self-sufficient* target expression. In some cases it is more convenient to deal with such programs. For example, the problem of checking the equivalence of two programs is reduced to checking the equivalence of two expressions.

However, in the process of supercompilation, global function definitions are easier to deal with than letrec-expressions. For this reason, at the first stage of supercompilation, HOSC removes all letrec-expressions using the well-known method of  $\lambda$ -lifting [9].

Thus, in the following, we assume that source HLL programs do not contain letrec-expressions. (Ironically, residual programs constructed by HOSC never contain global function definitions.)

#### 3.3 Observables, contexts, redexes

The syntax of observables, redexes and contexts is shown in Fig. 3.

An observable is either:

- a local variable,
- a local variable application,
- a constructor,
- a  $\lambda$ -abstraction.

```
 \begin{array}{lll} \mathcal{E}[\![c\,\overline{e_i}]\!] & \Rightarrow & c\,\,\overline{e_i} \\ \mathcal{E}[\![\lambda v_0 \to e_0]\!] & \Rightarrow & \lambda v_0 \to e_0 \\ \mathcal{E}[\![con\langle f_0\rangle]\!] & \Rightarrow & \mathcal{E}[\![con\langle unfold(f_0)\rangle]\!] \\ \mathcal{E}[\![con\langle(\lambda v \to e_0)\ e_1\rangle]\!] & \Rightarrow & \mathcal{E}[\![con\langle e_0\{v:=e_1\}\rangle]\!] \\ \mathcal{E}[\![con\langle case\ c_j\ \overline{e_k'}\ of\ \{\overline{c_i\ \overline{v_{ik}}\ \to e_i};\}\rangle]\!] & \Rightarrow & \mathcal{E}[\![con\langle e_j\{v_{jk}:=e_k'\}\rangle]\!] \end{array}
```

Figure 4: Operational semantics of HLL (call by name)

A redex is either:

- a global variable,
- a  $\lambda$ -abstraction application,
- a case-expression with an observable as a selector.

A context is an "expression with a hole". It is either:

- a hole (empty context)  $\langle \rangle$ ,
- a context application,
- a case-expression with a context as a selector.

If con is a context, then  $con\langle e \rangle$  denotes the result of replacing the hole in con with e. Essentially, the hole  $\langle \rangle$  is used as a meta-variable.

It can be shown that any HLL-expression e is either an observable or can be represented as a context con whose hole is replaced with a redex r, so that  $e = con\langle r \rangle$ . This fact is known as the *unique decomposition* property.

## 3.4 Operational semantics

HLL operational semantics is defined as normal-order reduction to a weak-head normal form [3] (Fig. 4). Note that operational semantics is only defined for expressions without free variables.

# 4 Homeomorphic embedding

In the definition of the "classical" homeomorphic embedding [13, 6, 17, 18] all variables are supposed to be free and considered to be embedded into each other. However, bound variables may appear in HLL expressions, for which reason HOSC takes a more subtle approach to the embedding of variables:

- a free variable is embedded into a free variable,
- a bound variable is embedded into a corresponding bound variable,
- a global variable v is embedded into itself only.

The correspondence between bound variables is recorded in a table  $\rho$ :

$$\rho = \{ (v_1', v_1''), \dots, (v_n', v_n'') \}$$

When two  $\lambda$ -abstractions are coupled, the correspondence between their arguments is recorded in  $\rho$ . When two case-expressions are coupled, the correspondence between their pattern variables is also recorded in  $\rho$ .

When checking whether an expression e can dive into the body of a  $\lambda$ -abstraction  $\lambda v \to e_1$ , the bound variable v does not correspond to any variables in e. So the pair  $(\bullet, v)$  is put into  $\rho$ , where  $\bullet$  is a special pseudo-variable (placeholder) for a missing variable.

Checking for an expression e diving into a branch of a case-expression is performed in the same manner - we remember that pattern variables do not correspond to any variables in e.

#### 4.1 Extended homeomorphic embedding

Let  $\rho$  be a table recording the correspondence of bound variables. Then extended homeomorphic embedding relation  $\leq |_{\rho}$  for expressions e' and e'' (constrained by  $\rho$ ) is defined inductively as follows:

$$e' \unlhd e'' \mid_{\rho} \quad \text{ if } e' \unlhd_v e'' \mid_{\rho} \text{ or } e' \unlhd_d e'' \mid_{\rho} \text{ or } e' \unlhd_c e'' \mid_{\rho}$$

Thus we distinguish the following kinds of embedding:

- $e' \leq_v e'' \mid_{\rho}$  embedding of variables
- $e' \leq_d e'' \mid_{\rho}$  diving
- $e' \leq_c e'' \mid_{\rho}$  coupling

We will also use the following abbreviations:

$$\begin{array}{cccc} e' \mathrel{\unlhd} e'' & \stackrel{\scriptscriptstyle \mathrm{def}}{=} & e' \mathrel{\unlhd} e'' \mid_{\{\}} \\ e' \mathrel{\unlhd}_v e'' & \stackrel{\scriptscriptstyle \mathrm{def}}{=} & e' \mathrel{\unlhd}_v e'' \mid_{\{\}} \\ e' \mathrel{\unlhd}_d e'' & \stackrel{\scriptscriptstyle \mathrm{def}}{=} & e' \mathrel{\unlhd}_d e'' \mid_{\{\}} \\ e' \mathrel{\unlhd}_c e'' & \stackrel{\scriptscriptstyle \mathrm{def}}{=} & e' \mathrel{\unlhd}_c e'' \mid_{\{\}} \end{array}$$

## 4.2 Embedding of variables

$$\begin{array}{ll} f \leq_v f\mid_{\rho} \\ v_1 \leq_v v_2\mid_{\rho} & \text{if } (v_1, \ v_2) \in \rho \\ v_1 \leq_v v_2\mid_{\rho} & \text{if } v_1 \notin domain(\rho) \text{ and } v_2 \notin range(\rho) \end{array}$$

#### 4.3 Diving

For an expression e to be able to dive into another expression e' under constraints  $\rho$ , we require that

$$\forall v \in fv(e) : v \notin domain(\rho)$$

The rationale behind this requirement will be given in the next section. The diving relation is defined as follows:

## 4.4 Coupling

$$\begin{array}{ll} c \ \overline{e_i'} \ \unlhd_c \ c \ \overline{e_i''} \mid_{\rho} & \text{if} \ \forall i : e_i' \ \unlhd \ e_i'' \mid_{\rho} \\ \lambda v_1 \to e_1 \ \unlhd_c \ \lambda v_2 \to e_2 \mid_{\rho} & \text{if} \ e_1 \unlhd e_1 \mid_{\rho \cup \{(v_1, v_2)\}} \\ e' \ \overline{e_i'} \ \unlhd_c \ e'' \ \overline{e_i''} \mid_{\rho} & \text{if} \ e' \unlhd \ e'' \mid_{\rho} \ \text{and} \ \forall i : e_i' \unlhd \ e_i'' \mid_{\rho} \\ case \ e' \ of \ \{\overline{c_i} \ \overline{v_{ik}'} \to e_i';\} \ \unlhd_c \ case \ e'' \ of \ \{\overline{c_i} \ \overline{v_{ik}''} \to e_i'';\} \mid_{\rho} \\ & \text{if} \ e' \unlhd \ e'' \mid_{\rho} \ \text{and} \ \forall i : e_i' \unlhd \ e_i'' \mid_{\rho \cup \{\overline{(v_{ik}', v_{ik}'')}\}} \end{array}$$

Note that in the third rule (diving of applications) e' and e'' themselves are not applications! This ensures that a global variable can be embedded into itself only.

## 4.5 Examples

Here are some examples of embedding and non-embedding (data types and global functions are taken from the programs presented in Fig. 2 and Fig. 9):

- $x \leq y$
- $map \not \supseteq compose$

- $map f \not \leq compose map f$
- $map \ f \leq_c map \ (compose \ f \ g)$
- $\lambda x \to Nil \leq_c \lambda x \to Cons \ a \ Nil$
- $\lambda x \to Nil \not \supseteq \lambda x \to Cons \ x \ Nil$
- case x of  $\{Z \to x; S b \to f b; \} \leq_c case x of \{Z \to x; S b \to (f g) b; \}$
- case x of  $\{Z \rightarrow x; S b \rightarrow f b; \} \not \leq case x of \{Z \rightarrow x; S b \rightarrow f (g b); \}$

#### 5 Generalization

The specializer HOSC described in the paper never tries to find a generalization for two expressions  $e_1$  and  $e_2$ , unless one of them is embedded into another by coupling. For this reason, we describe an algorithm for finding a most specific generalization (MSG) dealing only with expressions  $e_1$  and  $e_2$ , such that  $e_1 \leq_c e_2$ .

#### 5.1 Most specific generalization of coupled expressions

A most specific generalization of coupled expressions  $e_1$  and  $e_2$  (that is  $e_1 \leq_c e_2$ ) is computed by exhaustively applying the following rewrite rules (common functor rule and common sub-expressions rule) to the initial trivial generalization  $(v, \{v := e_1\}, \{v := e_2\})$ . The requirement from the subsection 4.3 comes from the fact that  $\theta_1$  and  $\theta_2$  should be correct substitutions - that is, substitutions operating on free variables. This requirement guarantees that substitutions will be correct, if  $e_1 \leq_c e_2$ .

#### 5.2 Common functor rule

#### 5.2.1 Variable

$$\begin{pmatrix} e \\ \{v_1 := v_2\} \cup \theta_1 \\ \{v_1 := v_2\} \cup \theta_2 \end{pmatrix} \Rightarrow \begin{pmatrix} e\{v_1 := v_2\} \\ \theta_1 \\ \theta_2 \end{pmatrix}$$

#### 5.2.2 Constructor

$$\left( \begin{array}{c} e \\ \{v := c \ \overline{e_1'}\} \cup \theta_1 \\ \{v := c \ \overline{e_1''}\} \cup \theta_2 \end{array} \right) \Rightarrow \left( \begin{array}{c} e\{v := c \ v_1 \ \dots \ v_n\} \\ \{\overline{v_i := e_1'}\} \cup \theta_1 \\ \{\overline{v_i := e_1''}\} \cup \theta_2 \end{array} \right)$$

#### 5.2.3 $\lambda$ -abstraction

$$\begin{pmatrix} e \\ \{v := \lambda v' \to e'\} \cup \theta_1 \\ \{v := \lambda v'' \to e''\} \cup \theta_2 \end{pmatrix} \Rightarrow \begin{pmatrix} e\{v := \lambda v_0 \to v_1\} \\ \{v_1 := e'\{v' := v_0\}\} \cup \theta_1 \\ \{v_1 := e''\{v'' := v_0\}\} \cup \theta_2 \end{pmatrix}$$

#### 5.2.4 Application

$$\begin{pmatrix} e \\ \{v := e'_0 \, \overline{e'_i}\} \cup \theta_1 \\ \{v := e''_0 \, \overline{e''_i}\} \cup \theta_2 \end{pmatrix} \Rightarrow \begin{pmatrix} e\{v := v_0 \, \overline{v_i}\} \\ \{v_0 := e'_0, \, \overline{v_i} := e'_i\} \cup \theta_1 \\ \{v_0 := e''_0, \, \overline{v_i} := e''_i\} \cup \theta_2 \end{pmatrix}$$

#### 5.2.5 Case-expression

$$\begin{pmatrix} e \\ \{v := case \ e'_0 \ of \ \{\overline{c_i \ \overline{v'_{ik}} \to e'_i};\}\} \cup \theta_1 \\ \{v := case \ e''_0 \ of \ \{\overline{c_i \ \overline{v''_{ik}} \to e''_i};\}\} \cup \theta_2 \end{pmatrix}$$

$$\downarrow \qquad \qquad \downarrow \qquad \qquad \qquad \downarrow$$

$$\begin{pmatrix} e\{v := case \ v_0 \ of \ \{\overline{c_i \ \overline{v_{ik}} \to v_i};\}\} \\ \{v_0 := e'_0, \ \overline{v_i := e'_i \{\overline{v'_{ik} := v_{ik}}\}} \cup \theta_1 \\ \{v_0 := e''_0, \ \overline{v_i := e''_i \{\overline{v'_{ik} := v_{ik}}\}} \cup \theta_2 \end{pmatrix}$$

## 5.3 Common sub-expression rule

$$\begin{pmatrix} e \\ \{v_1 := e', \ v_2 := e'\} \cup \theta_1 \\ \{v_1 := e'', \ v_2 := e''\} \cup \theta_2 \end{pmatrix} \Rightarrow \begin{pmatrix} e\{v_1 := v_2\} \\ \{v_2 := e'\} \cup \theta_1 \\ \{v_2 := e''\} \cup \theta_2 \end{pmatrix}$$

#### 5.4 Let-expression

Let-expressions are used for representing results of generalizing expressions. The syntax of a let-expression has the form:

let 
$$\overline{v_i = e_i}$$
; in  $e_q$ 

where variables  $v_i$  may appear in  $e_g$ , the value of  $v_i$  being obtained by evaluating  $e_i$ . Since HLL is a lazy language, a let-expression is semantically equivalent to  $e_q\{\overline{v_i := e_i}\}$ .

Let-expressions are not allowed in input programs, but may be generated by HOSC when constructing a partial process tree.

#### 5.5 Examples

Here are examples of generalization of expressions embedded by coupling (see examples in 4.5):

- $map \ f \sqcap map \ (compose \ f \ g) = (map \ h, \{h := f\}, \{h := (compose \ f \ g)\})$
- $\lambda x \to Nil \cap \lambda x \to Cons \ a \ Nil = (\lambda x \to y, \{y := Nil\}, \{y := Cons \ a \ Nil\})$
- case x of  $\{Z \to x; S b \to f b; \} \cap case x$  of  $\{Z \to x; S b \to (f g) b; \} = (case x of <math>\{Z \to x; S b \to h b; \}, \{h := f\}, \{h := f g\})$

# 6 Metacomputation

A process tree is a directed (possibly infinite) tree whose nodes are labeled with HLL expressions. A process tree is built by HOSC in the process of metacomputation ([5, 6, 1, 2]), which is performed by symbolically evaluating expressions that, unlike the case of ordinary evaluation, may contain free variables.

The expression a node n is labeled with will be denoted by n.expr.

Given a source program, the construction of its process tree starts with creating a single-node tree whose root is labeled with the program's target expression. Then the process tree is incrementally built by adding children to its leaves according to the *driving* rules presented below.

## 6.1 Driving step

The operator  $\mathcal{D}$  (Fig. 5) takes as argument an expression in a leaf node and returns zero or more expressions  $e_1, \ldots, e_k$ . Then the supercompiler adds k child nodes to the leaf labeling them with the expressions  $e_1, \ldots, e_k$ .

The last rule applies to let-expressions which are not present in source programs but may appear as a result of generalization.

# 7 Partial process tree

In the general case, metacomputation may result in constructing an infinite process tree. The task of supercompilation is to turn this tree into a finite (possibly cyclic) graph, which is referred to as a *partial process tree*. A partial process tree differs from a process tree in the following points:

• Nodes of a partial process tree may contain let-expressions.

$$\mathcal{D}\llbracket v \ \overline{e_i} \rrbracket \qquad \Rightarrow \qquad \llbracket v, \ \overline{e_i} \rrbracket$$

$$\mathcal{D}\llbracket c \ \overline{e_i} \rrbracket \qquad \Rightarrow \qquad \llbracket e_0 \rrbracket$$

$$\mathcal{D}\llbracket con\langle f_0 \rangle \rrbracket \qquad \Rightarrow \qquad \llbracket con\langle unfold(f_0) \rangle \rrbracket$$

$$\mathcal{D}\llbracket con\langle (\lambda v_0 \to e_0) \ e_1 \rangle \rrbracket \qquad \Rightarrow \qquad \llbracket con\langle e_0 \{v_0 := e_1\} \rangle \rrbracket$$

$$\mathcal{D}\llbracket con\langle case \ c_j \ \overline{e_k'} \ of \ \{\overline{c_i \ \overline{v_{ik}} \to e_i};\} \rangle \rrbracket \qquad \Rightarrow \qquad \llbracket con\langle e_i \{v_j := e_k'\} \rangle \rrbracket$$

$$\mathcal{D}\llbracket con\langle case \ v \ \overline{e_j'} \ of \ \{\overline{p_i \to e_i};\} \rangle \rrbracket \qquad \Rightarrow \qquad \llbracket e, \ \overline{con\langle e_i \{v \ \overline{e_j'} := p_i\} \rangle} \rrbracket$$

$$\mathcal{D}\llbracket let \ \overline{v_i = e_i}; \ in \ e \rrbracket \qquad \Rightarrow \qquad \llbracket e, \ \overline{e_i} \rrbracket$$

Figure 5: Driving step

• A partial process tree may have cycles. Let a node  $\alpha$  be an ancestor of a node  $\beta$  and  $\beta.expr$  a renaming of  $\alpha.expr$  ( $\beta.expr \simeq \alpha.expr$ ). Then it is possible to create a "return" edge  $\beta \rightrightarrows \alpha$  from the repeat (or recursive) node  $\beta$  to the function (or base) node  $\alpha$ .

Thus, a partial process tree is a directed tree (whose edges are denoted by  $\rightarrow$ ) supplemented with "return" edges (denoted by  $\rightrightarrows$ ) turning it into a directed graph.

The goal of supercompilation is to construct a partial process tree of a reasonable size.

Fig. 6 shows a number of operations used by HOSC for constructing partial process trees and generating residual programs.

A node  $\beta$  is referred to as processed, if one of the following conditions holds:

- $\beta$  is the source of a "return" edge  $\beta \Rightarrow \alpha$ ,
- $\beta.expr$  is a nullary constructor,
- $\beta.expr$  is a local variable.

A node  $\beta$  is non-trivial, if  $\beta$ .expr can be decomposed as either

$$con\langle f\rangle \qquad \text{or} \qquad con\langle case\ v\ \overline{e_j}\ of\ \{\overline{p_i\to e_i;}\}\rangle$$

Otherwise, the node  $\beta$  is referred to as *trivial*.

processed (node)	Returns <i>true</i> or <i>false</i> depending on whether the node is processed or not.
trivial (node)	Returns <i>true</i> or <i>false</i> depending on whether the node is trivial or not.
$children(t, \ \alpha)$	Returns an ordered list of child nodes for the node $\alpha$ of a tree $t$ .
$addChildren(t, \beta, es)$	Adds child nodes to the node $\beta$ of the tree $t$ and puts expressions $es$ into them.
$replace(t, \alpha, expr)$	Replaces a subtree with root $\alpha$ by a single node $\beta$ such that $\beta.expr = expr$ .
$ancestor(t, \beta, \simeq)$	Returns a node $\alpha$ - the closest ancestor of $\beta$ , such that $\alpha.expr \simeq \beta.expr$ , or $\bullet$ if $\beta$ doesn't have such ancestor.
$ancestor(t, \ \beta, \ \lessdot)$	Returns a node $\alpha$ - the closest ancestor of $\beta$ , such that $\alpha.expr < \beta.expr$ , or $\bullet$ if $\beta$ doesn't have such ancestor.
$ancestor(t, \ \beta, \ \leq_c)$	Returns a node $\alpha$ - the closest ancestor of $\beta$ , such that $\alpha.expr \leq_c \beta.expr$ , or $\bullet$ if $\beta$ doesn't have such ancestor.
$fold(t, \alpha, \beta)$	Adds a "return" edge $\beta \Rightarrow \alpha$ to form a cycle.
$abstract(t, \ \alpha, \ \beta)$	= $replace(t, \alpha, let \overline{v_i = e_i}; in e_g)$ , where $(e_g, \theta_1, \theta_2) = \alpha.expr \cap \beta.expr$ , $e_g\theta_1 = e_g\{\overline{v_i := e_i}\} = let \overline{v_i = e_i}; in e_g$ .
$\lceil \alpha \downarrow_{s} t \rceil$	Returns all repeat nodes of a node $\alpha$ , $\lceil \alpha \bowtie t \rceil = [\overline{\beta_i}]$ : $\beta_i \rightrightarrows \alpha$ , or $\bullet$ if $\alpha$ is not a function node.
$\lceil \alpha \uparrow \uparrow t \rceil$	Returns a function node for a node $\alpha$ , $\lceil \alpha \uparrow \uparrow t \rceil = \beta : \alpha \Rightarrow \beta$ , or $\bullet$ if $\alpha$ is not a repeat node.
$drive(t, \alpha)$ unprocessedLeaves(t)	= $addChildren(t, \alpha, \mathcal{D}[\![\alpha.expr]\!])$ - executes driving step. Returns a list of all unprocessed leaves of a tree $t$ or $\bullet$ if all leaves are processed.

Figure 6: Operations on partial process tree

# 8 Partial process tree construction

Now we describe the core of the supercompiler HOSC: the algorithm of constructing a partial process tree (see Fig. 7).

Given a source program, the algorithm starts with the creation of a singlenode tree whose root is labeled with the program's target expression. Then, while there is at least one unprocessed leaf  $\beta$ :

- 1. If  $\beta$  is trivial,  $\beta.expr$  is "metaevaluated" by performing a driving step.
- 2. If  $\beta$  has an ancestor  $\alpha$ , such that  $\alpha.expr \simeq \beta.expr$ , then  $\alpha$  becomes a function node for  $\beta$ , i.e. a "return" edge  $\beta \rightrightarrows \alpha$  is added to the tree.

```
while unprocessedLeaves(t) \neq \bullet do
    \beta = unprocessedLeaves(t).head
    if trivial(\beta) then
         t = drive(t, \beta)
    else if ancestor(t, \beta, \simeq) \neq \bullet then
         \alpha = ancestor(t, \beta, \simeq)
         fold(t, \alpha, \beta)
    else if ancestor(t, \beta, \lessdot) \neq \bullet then
         \alpha = ancestor(t, \beta, \lessdot)
         t = abstract(t, \beta, \alpha)
    else if ancestor(t, \beta, \leq_c) \neq \bullet then
         \alpha = ancestor(t, \beta, \leq_c)
         t = abstract(t, \alpha, \beta)
    else
         t = drive(t, \beta)
end
```

Figure 7: Algorithm of partial process tree construction

- 3. If  $\beta$  has an ancestor  $\alpha$ , such that  $\alpha.expr < \beta.expr$ , then  $\beta.expr$  is generalized.
- 4. If  $\beta$  has an ancestor  $\alpha$ , such that  $\alpha.expr \leq_c \beta.expr$ , then  $\alpha.expr$  is generalized.
- 5. Otherwise,  $\beta.expr$  is "metaevaluated" by performing a driving step.

The fourth case is the central point of the algorithm: removing it would cause the supercompiler not to terminate for a lot of source programs!

However, the test for the homeomorphic embedding works as a "whistle" signalling that two expressions have "too much in common", but a "return" edge still cannot be added to the tree. So the supercompiler generalizes the expression labeling the "function" (base) node, thereby preventing the the infinite growth of the process tree.

# 9 Residual program construction

Here we describe an algorithm for transforming a partial process tree t into a program in the language HLL.

The algorithm (presented in Fig. 8) is defined via two mutually recursive operators (functions):  $\mathcal{C}$  and  $\mathcal{C}'$ . A residual program can be extracted from a

$$\begin{array}{lll} & \mathcal{C} & \| \alpha \|_{t,\Sigma} = \mathcal{C}' \| \alpha . expr p^* \|_{t,\alpha,\Sigma} \Rightarrow \mathcal{C} \| \gamma' \| \{\overline{v_i} = \mathcal{C} \| \gamma_i \|_{t,\Sigma} \} \\ & \mathcal{C}' \| tet \ \overline{v_i} = e_i; \ in \ e \|_{t,\alpha,\Sigma} \Rightarrow \mathcal{C} \| \gamma' \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{e_i} \|_{t,\alpha,\Sigma} \Rightarrow v \ \overline{\mathcal{C}} \| \gamma_i \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{e_i} \|_{t,\alpha,\Sigma} \Rightarrow c \ \overline{\mathcal{C}} \| \gamma_i \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{e_i} \|_{t,\alpha,\Sigma} \Rightarrow c \ \overline{\mathcal{C}} \| \gamma_i \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{e_i} \|_{t,\alpha,\Sigma} \Rightarrow c \ \overline{\mathcal{C}} \| \gamma_i \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{e_i} \|_{t,\alpha,\Sigma} \Rightarrow c \ \overline{\mathcal{C}} \| \gamma_i \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{e_i} \|_{t,\alpha,\Sigma} \Rightarrow c \ \overline{\mathcal{C}} \| \gamma' \|_{t,\Sigma} \\ & \mathcal{C}' \| v \ \overline{\mathcal{C}} \|_{t,\alpha,\Sigma} \\ & \mathcal{C} \| v \ \overline{\mathcal{C}} \|_{t,\alpha,\Sigma} \\ & \mathcal{C}' \| v \ \overline{\mathcal{C}} \|_{t,\alpha,\omega} \\ & \mathcal{$$

In order to make the rules less cumbersome the following abbreviations are used. If  $\gamma_i$  appears in the right-hand side, we assume that  $[\overline{\gamma_i}] = children(t, \alpha)$ , and all child nodes are processed in the same way. If  $\gamma'$  and  $\gamma'_i$  appear in the right-hand side, we assume that  $[\gamma', \overline{\gamma'_i}] = children(t, \alpha)$ , and either there is exactly one child node or the first child node requires special treatment.

Figure 8: Rules for constructing residual program

partial process tree by evaluating:

$$\mathcal{C} [t.root]_{t,\{\}}$$

In order to construct a residual program we traverse a partial process tree top-down starting from the root node. Traversing a function node produces the definition of a recursive function. The correspondence between a function node and the signature of a new function in recorded in  $\Sigma$ . Later  $\Sigma$  is used for constructing a recursive function call.

Let us consider the most important details of the algorithm.

- The rule  $C_0$  defines  $\mathcal{C}$  in terms of  $\mathcal{C}'$ .
- The rules  $C_1$ ,  $C_2$  and  $C_3$  define how to combine supercompiled parts of a let-expression, a call to an unknown function or a constructor expression in order to produce the corresponding residual expression.
- The rule  $C_4$  defines how to construct a residual  $\lambda$ -abstraction from its supercompiled body.
- The rules  $C_5$  and  $C_6$  correspond to the cases  $con\langle(\lambda v_0 \to e_0) \ e_1\rangle$  and  $con\langle case \ c \ \overline{e_j} \ of \ \{\overline{p_i \to e_i};\}\rangle$ , in which a reduction step was completely performed by driving. Thus no actions need to be inserted into the residual program.
- The rules  $C_7$  and  $C_8$  are used when traversing non-trivial nodes:
  - A recursive function definition is generated when traversing a function node (the rules  $C_7^1$  and  $C_8^1$ ). Note that only those free variables which were not preserved in repeat nodes become the arguments of λ-abstractions. This trick allows the arity of residual functions to be reduced. The correspondence between a function node and the signature of a new function is recorded in Σ and passed down the tree.
  - A call to a recursive function defined in a base node is generated when traversing a repeat node (the rules  $C_7^2$  and  $C_8^2$ ).
  - The rules  $C_7^3$  and  $C_8^3$  are used when traversing a node which is neither a "function" nor "repeat" one.

## 10 Related work

Originally supercompilation was formulated for the programming language Refal [20]. Currently, the most advanced Refal supercompiler is SCP4 [16].

In 1990-s supercompilation was reformulated for a number of simpler languages. The main goal of that work was to understand which parts of supercompilation were due to Refal and which ones were of a fundamental character. The first *complete* and *formal* description of the concepts and algorithms essential for any supercompiler (driving, generalization, residual program construction) can be found in Sørensen's Master's Thesis [17] and in the later papers by Sørensen and Glück [6, 19], presenting a supercompiler for a simple first-order functional language (with the call-by-name semantics).

A full description of a supercompiler for the language TSG (a simple first-order functional call-by-name language) is given by Abramov [2].

The supercompilers by Abramov, Sørensen and Glück construct a partial process tree which is then transformed into a residual program.

In recent years there has been a growing interest in supercompilation for higher-order functional languages [8, 10, 11, 12, 15, 14].

The techniques used by Jonsson [10, 11] and Mitchell [15, 14] are similar to that of deforestation [23, 7], since they do not generate a partial process tree at a separate stage, so that the driving (symbolic evaluation) and the residual program generation are closely interleaved.

On the contrary, Hamilton [8] and Kabir [12] prefer to deal with partial process trees.

While the papers by Hamilton, Jonsson, Kabir and Mitchell pay much attention to driving and residual program generation, unfortunately, the problems related to homeomorphic embedding and generalization are not given an account they deserve.

Homeomorphic embedding relation (used as a whistle for generalization) is described in detail in [6, 17, 18] for a first-order functional language without case-expressions. Case-expressions are discussed in [19], however a formal description of homeomorphic embedding of case-expressions is missing. The same situation is with the higher-order languages with case-expressions dealt with in [8, 10, 11, 12, 15, 14].

An algorithm for finding a most specific generalization for a first-order language without case-expressions (that is, without bound variables) is described in already mentioned papers [6, 17, 18]. Bound variables are considered in [19, 10, 8, 15], and the algorithm for finding an MSG is said to be a modification of algorithms from [6, 17, 18], but it is not described in detail.

#### 11 Conclusions

We have described the internal structure of an experimental supercompiler HOSC dealing with programs written in a higher-order functional language.

A complete and formal account has been given of all essential concepts and algorithms the supercompiler HOSC is based upon. Of particular interest are the questions related to homeomorphic embedding and generalization for expressions containing bound variables.

# 12 Acknowledgements

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# A Examples

This Appendix contains two examples of supercompilation performed by HOSC supercompiler. Both examples deal with higher-order functions.

#### A.1 Iterate

```
data List a = Nil | Cons a (List a);
data Nat = Z | S Nat;
iterate (\lambdan \to S n) Z where
iterate = \lambdaf x \to Cons x (iterate f (f x));
```

Figure 9: Iterate: input

The language HLL allows infinite data structures to be processed. Given a function f and an initial value x, the function iterate produces an infinite list (stream) of repeated applications of f to x:

```
iterate f x = Cons x (Cons (f x) (Cons (f (f x)) (Cons ...)))
```

The target expression in the program in Fig. 9 defines the infinite list (stream) of natural numbers by means of the function **iterate**.

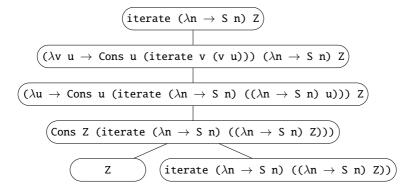


Figure 10: Iterate: metacomputation before generalization

Driving this target expression, HOSC produces the process tree shown in Fig. 10. At this step, it finds out that the expression in the root node is embedded into the expression in the rightmost leaf:

iterate (
$$\lambda n \to S$$
 n) Z  $\leq_c$  iterate ( $\lambda n \to S$  n) (( $\lambda n \to S$  n) Z)

However, the second expression is not an instance of the first one! For this reason, HOSC generalizes the expression in the root node, replacing

iterate (
$$\lambda n \rightarrow S n$$
) Z

with the let-expression

**let** 
$$z = Z$$
 **in** iterate  $(\lambda n \rightarrow S n) z$ 

and deleting the whole subtree. Further driving (after this generalization and till the next whistle) is shown in Fig. 11:

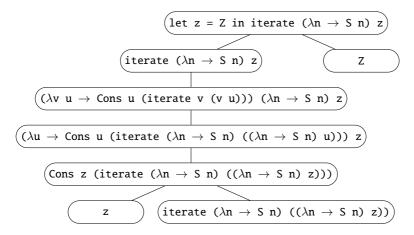


Figure 11: Iterate: metacomputation after the first generalization

At this step an embedding is detected:

iterate 
$$(\lambda n \to S n)$$
 z  $\leq_c$  iterate  $(\lambda n \to S n)$   $((\lambda n \to S n) z)$ 

This time the second expression is an instance of the first one:

iterate 
$$(\lambda n \to S n) z \in \text{iterate } (\lambda n \to S n) ((\lambda n \to S n) z)$$

Thus the second expression is generalized by replacing the expression

iterate (
$$\lambda n \rightarrow S n$$
) (( $\lambda n \rightarrow S n$ ) z)

with the let-expression

**let** 
$$y = (\lambda n \rightarrow S n) z in iterate (\lambda n \rightarrow S n) ((\lambda n \rightarrow S n) z)$$

Further metacomputation results in a complete partial process tree shown in Fig. 12. The residual program constructed from this tree is presented in Fig. 13. It is worth to note that the function

$$(\lambda n \rightarrow S n)$$

was inlined into the body of the recursive function f.

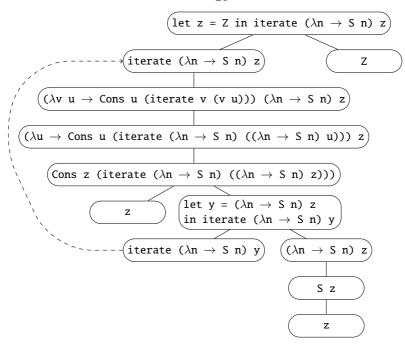


Figure 12: Iterate: metacomputation after the second generalization

Figure 13: Iterate: output

#### A.2 Church numerals

Church numerals are the representations of natural numbers under Church encoding. The higher-order function that represents a natural number n is the function that maps any other function f to its n-fold composition. In simpler terms, the "value" of the numeral is equivalent to the number of times the function encapsulates its argument.

$$f^n = f \circ f \circ f \circ \dots \circ f$$

Church numerals 0, 1, 2, ..., n are defined as follows in the lambda calculus:

```
data Nat = Z | S Nat;
data Boolean = False | True;
eq (add x y) (unchurch(churchAdd (church x) (church y))) where
eq = \lambdam y \rightarrow case m of {
  Z \rightarrow case \ n \ of \ \{Z \rightarrow True; \ S \ n1 \rightarrow False; \} ;
  S m1 \rightarrow case n of {Z \rightarrow False; S n1 \rightarrow eq m1 n1;};
};
church = \lambda n \rightarrow case n of {
  Z \rightarrow \lambda f t \rightarrow t;
  S n1 \rightarrow \lambda f t \rightarrow f (church n1 f t);
unchurch = \lambda n \rightarrow n \ (\lambda t \rightarrow S \ t) \ Z;
churchAdd = \lambdam n \rightarrow (\lambdaf t \rightarrow m f (n f t));
add = \lambda m \ n \rightarrow case \ m \ of  {
  Z \rightarrow n:
  S m1 \rightarrow S (add m1 n);
};
```

Figure 14: Church addition: input

```
0 = \lambda f \ x \to x

1 = \lambda f \ x \to f \ x

2 = \lambda f \ x \to f \ (f \ x)

...

n = \lambda f \ x \to f^n \ x
```

The addition function churchAdd m n = m + n is based on the identity  $f^{(m+n)}$   $x = f^m$   $(f^n x)$ .

The program shown in Fig. 14 contains definitions of Boolean and Nat datatypes. Nat represents numbers in Peano encoding by means of constructors Z (zero) and S (successor). The function eq tests Peano numbers for equality. The church and unchurch functions convert between Peano numbers and Church numerals. The churchAdd and add functions perform addition of Church numerals and Peano numbers correspondingly.

The target expression of the program tests whether the sum of two Peano numbers  $\boldsymbol{x}$  and  $\boldsymbol{y}$  is equal to the decoded sum of corresponding Church numerals **church**  $\boldsymbol{x}$  and **church**  $\boldsymbol{x}$ . Note that the target expression may return False, since this constructor appears in the definition of the function eq.

The result of supercompiling this program is shown in Fig. 15. Now it is evident that the residual program cannot return False since this constructor never appears in the body of the program.

The structure of the original program has been simplified: it is almost

```
data Nat = Z | S Nat;
data Boolean = False | True ;
case x of {
  Z \rightarrow case y of {
      Z \rightarrow True;
      S y1 \rightarrow letrec f= \lambda z \rightarrow
                     case z of { Z \rightarrow True; S z1 \rightarrow f z1;}
                  in f y1;
      }:
  S x1 \rightarrow
     letrec g = \lambda v \rightarrow
          case v of {
             Z \rightarrow case y of {
                  Z \rightarrow True;
                  S y2 \rightarrow letrec h= \lambdaw \rightarrow
                               case w of { Z \rightarrow True; S w1 \rightarrow h w1; }
                          in h y2;
             };
             S v1 \rightarrow g v1;
     in g x1;
}
```

Figure 15: Church addition: output

evident that the residual program just destructs  $\boldsymbol{x}$  and  $\boldsymbol{y}$  and finally returns True.

It can also be shown that this program returns True, if x and y are "good" values (finite Peano numbers), since the result of program execution depends on successful destruction of x and y only.