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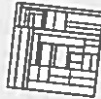
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Lambda-Lifting in Quadratic Time

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Lambda-Lifting in Quadratic Time

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Abstract Lambda-lifting is a program transformation used in compilers and in partial evaluators and that operates in cubic time. In this article, we show how to reduce this complexity to quadratic time.

Lambda-lifting transforms a block-structured program into a set of recursive equations, one for each local function in the source program. Each equation carries extra parameters to account for the free variables of the corresponding local function and of all its callees. It is the search for these extra parameters that yields the cubic factor in the traditional formulation of lambda-lifting, which is due to Johnsson. This search is carried out by a transitive closure.

Instead, we partition the call graph of the source program into strongly connected components, based on the simple observation that all functions in each component need the same extra parameters and thus a transitive closure is not needed. We therefore simplify the search for extra parameters by treating each strongly connected component instead of each function as a unit, thereby reducing the time complexity of lambda-lifting from $\mathcal{O}(n^3 \log n)$ to $\mathcal{O}(n^2 \log n)$, where n is the size of the program. Since a lambda-lifter can output programs of size $\mathcal{O}(n^2)$, we believe that our algorithm is close to optimal.

1 Lambda-lifting

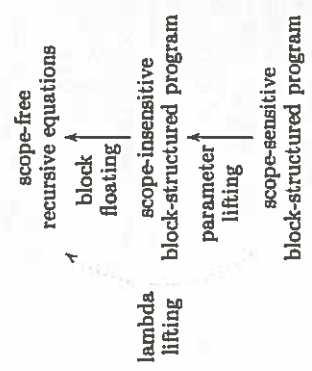
1.1 Setting and background

Lambda-lifting: what. In the mid 1980's, Augustsson, Hughes, Johnsson, and Peyton Jones devised 'lambda-lifting', a meaning-preserving transformation from block-structured programs to recursive equations [6,16,17,24].

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Recursive equations provide a propitious format because they are scope free. Today, a number of systems use lambda-lifting as an intermediate phase. For example, partial evaluators such as Schism, Similix, and Pell-Mell lambda-lift source programs and generate scope-free recursive equations [8,10,21]. Compilers such as Larceny and Moby use local, incremental versions of lambda-lifting in their optimizations [9,25], and so did an experimental version of the Glasgow Haskell Compiler [26]. Program generators such as Bakewell and Runciman's least general common generalization operate on lambda-lifted programs [7].

Lambda-lifting: how. Lambda-lifting operates in two stages: *parameter lifting* and *block floating*.



A block-structured program is scope-sensitive because of free variables in local functions. Parameter lifting makes a program scope-insensitive by passing extra variables to each function. These variables account both for the free variables of each function but also for variables occurring free further in the call path. Block floating erases block structure by making each local function a global recursive equation.

Parameter lifting. Parameter-lifting a program amounts to making all the free variables of a function formal parameters of this function. All callers of the function must thus be passed these variables as arguments as well. A set of solutions is built by traversing the program. A solution pairs each function with a least set of additional parameters. Each block of locally defined functions gives rise to a collection of set equations describing which variables should be passed as arguments to its local functions. The names of functions, however, are not included in the sets, since all functions become globally visible when the lambda-lifting transformation is complete. The solution of each set equation extends the current set of solutions, which is then used to analyze the header and the body of the block.

Block floating. After parameter lifting, a program is scope insensitive. Block floating is thus straightforward: the program is merely traversed, all local functions are collected and all blocks are replaced by their bodies. The collected

function definitions are then appended to the program as global mutually recursive functions, making all functions globally visible.

Lambda-lifting: when. In a compiler, the effectiveness of lambda-lifting hinges on the tension between passing many actuals vs. passing few actuals, and between referring to an actual parameter vs. referring to a free variable.

In practice, though, programmers often stay away both from recursive equations and from maximally nested programs. Instead, they write in a mixed style that both abides by Perlis's epigram "If you have a procedure with ten parameters, you probably missed some." and by Turner's recommendation that good Miranda style means little nesting. In this mixed style, and to paraphrase another of Perlis's epigrams, one man's parameter is another man's free variable.

1.2 Three examples

We first illustrate lambda-lifting with the classical `foldr` functional, and then with two examples involving multiple local functions and mutual recursion. Throughout, we use Standard ML.

Example 1: We consider the classical fold function for lists, defined with a local function.

```
fun foldr f b xs
  = let fun walk nil
        = b
          | walk (x :: xs)
            = f (x, walk xs)
        in walk xs
    end
```

Lambda-lifting this block-structured program yields two recursive equations: the original entry point, which now serves as a wrapper to invoke the other function, and the other function, which has been extended with two extra parameters.

```
fun foldr f b xs
  = foldr_walk f b xs
  and foldr_walk f b []
  = b
    | foldr_walk f b (x :: xs)
  = f (x, foldr_walk f b xs)
```

Example 2: The following token program adds its two parameters.

```
fun main x y
  = let fun add p
        = add_to_x p
          and add_to_x q
            = q + x
        in add y
    end
```

Lambda-lifting this block-structured program yields three recursive equations:

```

fun main x y
  = main_add x y
and main_add x p
  = main_add_to_x x p
and main_add_to_x x q
  = q + x

```

As a local function, `add_to_x` has a free variable, `x`, and thus it needs to be passed the value of `x`. Since `add_to_x` calls `add_to_x`, it needs to pass the value of `x` to `add_to_x` and thus to be passed this value, even though `x` is not free in the definition of `add`. During parameter lifting, each function thus needs to be passed not only the value of its free variables, but also the values of the free variables of all its callees.

Example 3: The following token program multiplies its two parameters with successive additions, using mutual recursion.

```

fun mul x y
  = let fun loop z
        = if z=0 then 0 else add_to_x z
        and add_to_x z
          = x + loop (z-1)
        in loop y
    end

```

Again, lambda-lifting this block-structured program yields three recursive equations:

```

fun mul x y
  = mul_loop x y
and mul_loop x z
  = if z=0 then 0 else mul_add_to_x x z
and mul_add_to_x x z
  = x + mul_loop x (z-1)

```

As before, the free variable `x` of `add_to_x` has to be passed as a formal parameter, through its caller `loop`. When `add_to_x` calls `loop` recursively, it must pass the value of `x` to `loop`, so that `loop` can pass it back in the recursive call.

This third example illustrates our key insight: during parameter lifting, mutually recursive functions must be passed the same set of variables as parameters.

1.3 Overview

Lambda-lifting, as specified by Johnson, takes cubic time (Section 2). In this article, we show how to reduce this complexity to quadratic time (Section 3).

Throughout the main part of the article, we consider Johnson's algorithm [17,18]. Other styles of lambda-lifting, however, exist: we describe them as well, together with addressing related work (Section 4).

2 Lambda-lifting in cubic time

2.1 Johnson's parameter-lifting algorithm

Johnson's algorithm descends recursively through the program structure, calculating the set of variables that are needed by each function. This is done by solving set equations describing the dependencies between functions. These dependencies may be arbitrarily complex, since a function can depend on any variable or function that is lexically visible to it. In particular, mutually recursive functions depend upon each other, and so they give rise to mutually recursive set equations.

The mutually recursive set equations are solved using fixed-point iterations. A program containing m function declarations gives rise to m set equations. In a block-structured program the functions are distributed across the program, so we solve the set equations in groups, as we process each block of local functions. Each set equation unifies $\mathcal{O}(m)$ sets of size $\mathcal{O}(n)$, where n is the size of the program. However, the total size of all equations is bounded by the size of the program n , so globally each iteration takes time $\mathcal{O}(n \log n)$. The number of set union operations needed is $\mathcal{O}(n^2)$, so the time needed to solve all the set equations is $\mathcal{O}(n^3 \log n)$, which is the overall running time of lambda-lifting.

2.2 An alternative specification based on graphs

Rather than using set equations, one can describe an equivalent algorithm using graphs. We use a graph to describe the dependencies between functions. Peyton Jones names this representation a *dependency graph* [24], but he uses it for a different purpose (see Section 4.1). Each node in the graph corresponds to a function in the program, and is associated with the free variables of this function. An edge in the graph from a node f to a node g indicates that the function f depends on g , because it refers to g . Mutually recursive dependencies give rise to cycles in this graph. Rather than solving the mutually recursive equations using fixed-point iteration, we propagate the variables associated with each node backwards through the graph, from callee to caller, merging the variable sets, until a fixed point is reached.

2.3 Example

Figure 1 shows a small program, defined using three mutually recursive functions, each of which has a different free variable.

We can describe the dependencies between the local block of functions using set equations, as shown in Figure 2. To solve these set equations, we need to perform three fixed-point iterations, since there is a cyclic dependency of size three. Similarly, we can describe these dependencies using a graph, also shown in Figure 2. The calculation of the needed variables can be done using this representation, by propagating variable sets backwards through the graph. A single propagation step is done by performing a set union over the variables

```

fun main x y z n
  = let fun f1 i
      = if i=0 then 0 else x + f2 (i-1)
      and f2 j
      = let fun g2 b = b * j
          in if j=0 then 0 else g2 y + f3 (j-1)
          end
      and f3 k
      = let fun g3 c = c * k
          in if k=0 then 0 else g3 z + f1 (k-1)
          end
      in f1 n
      end

```

Figure 1. Three mutually recursive functions

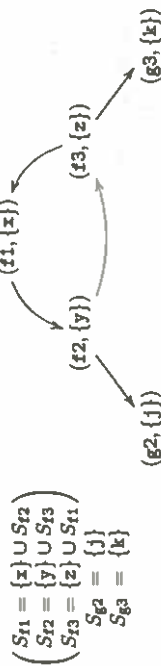


Figure 2. Dependencies between the local functions of the program of Figure 1

```

fun main x y z n
  = f1 x y z n
  and f1 x y z i
  = if i=0 then 0 else x + f2 x y z (i-1)
  and f2 x y z j
  = if j=0 then 0 else g2 y j + f3 x y z (j-1)
  and g2 b j
  = b * j
  and f3 x y z k
  = if k=0 then 0 else g3 z k + f1 x y z (k-1)
  and g3 c k
  = c * k

```

Figure 3. Lambda-lifted counterpart of Figure 1

associated with a node and the variables associated with its successors. Similarly to the case of the set equations, each node must be visited three times before a fixed point is reached.

When the set of needed variables has been determined for each function, solutions describing how each function must be expanded with these variables are added to the set of solutions. The result is shown in Figure 3.

3 Lambda-lifting in quadratic time

We consider the variant of the parameter-lifting algorithm that operates on a dependency graph. It propagates needed variables backwards through the graph since the caller needs the variables of each callee.

It is our observation that functions that belong to the same strongly connected component of the call graph must be parameter-lifted with the same set of variables (as was illustrated in Section 1.2). We can thus treat these functions in a uniform fashion, by coalescing the strongly connected components of the dependency graph. Each function must define at least its free variables together with the free variables of the other functions of the strongly connected component. Coalescing the strongly connected components of the dependency graph produces a DAG with sets of function names for nodes. A breadth-first backwards propagation of variables can then be done in linear time, which eliminates the need for a fixed-point computation.

3.1 Complexity analysis

The parameter-lifting algorithm must first construct the dependency graph, which takes time $\mathcal{O}(n \log n)$, where n is the size of the program. The strongly connected components of the graph can then be computed in time $\mathcal{O}(n)$. The ensuing propagation requires a linear number of steps since we are now operating on a DAG. Each propagation step consists of a number of set-union operations, each of which takes $\mathcal{O}(n \log n)$ time, i.e., the time to unify two sets of variables of size $\mathcal{O}(n)$. Globally, a number of set-union operations linear in the size of the program needs to be performed, yielding a time complexity of $\mathcal{O}(n^2 \log n)$. The overall running time is thus $\mathcal{O}(n^2 \log n)$, where n is the size of the program.

3.2 Lower bound

Consider a function with m formal parameters $\{v_1, \dots, v_m\}$ that declares m mutually recursive local functions, each of which has a different variable from $\{v_1, \dots, v_m\}$ as a free variable. The size of the program n is $\mathcal{O}(m)$. The output program contains the m functions, each of which needs to be expanded with the m formal parameters of the enclosing function. The output program is therefore of size $\mathcal{O}(m^2)$, which is also $\mathcal{O}(n^2)$. One thus cannot perform lambda-lifting faster than $\mathcal{O}(n^2)$. Since one needs $\mathcal{O}(n \log n)$ time to compute the sets of free variables of the program, our complexity of $\mathcal{O}(n^2 \log n)$ must be close to optimal.

3.3 Contribution

Our contribution is

- to characterize the fixed-point operations on the set equations as propagation through the dependency graph, and
- to recognize that functions in the same strongly connected component require the same set of variables.

We can therefore first determine which variables need to be known by each function in a strongly connected component, and then add them as formal parameters to these functions. In each function, those variables not already passed as parameters to the function should be added as formal parameters.

This approach can be applied locally to work like Johnson's algorithm, processing each block independently. It can also be applied globally to the overall dependency graph. The global algorithm, however, must explicitly limit the propagation of free variables, so that they are not propagated beyond their point of definition.

3.4 The new algorithm

We operate on programs conforming to the simple syntax of Figure 4.

$$\begin{aligned}
 p \in \text{Program} &::= \{d_1, \dots, d_m\} \\
 d \in \text{Def} &::= f \equiv \lambda v_1, \dots, \lambda v_n. e \\
 e \in \text{Exp} &::= e_0 \dots e_n \\
 &| \text{LetRec} \{d_1, \dots, d_k\} e_0 \\
 &| \text{if } e_0 \text{ e}_1 \text{ e}_2 \\
 &| f \\
 &| v \\
 &| \text{literal} \\
 v \in \text{Variable} & \\
 f \in \text{FunctionName} \cup \text{PdefinedFunction} &
 \end{aligned}$$

Figure 4. Simplified syntax of source programs

The set $\text{FV}(f)$ denotes the set of free variables in the function f , and the set $\text{FF}(f)$ denotes the set of free functions in f (note that $\text{FV}(f) \cap \text{FF}(f) = \emptyset$). In our algorithm, we assume variable hygiene, i.e., that no name clashes can occur. Figure 5 shows our (locally applied) $\mathcal{O}(n^2 \log n)$ parameter-lifting algorithm. It makes use of several standard graph and list operations that are described in the appendix. Figure 6 shows the standard linear-time (globally applied) block-floating algorithm. Johnson's original lambda-lifting algorithm includes steps to explicitly name anonymous lambda expressions and replace non-recursive let blocks by applications. These steps are trivial and omitted from the figures.

```

parameterLiftProgram :: Program -> Program
parameterLiftProgram p = map (parameterLiftDef \() p
parameterLiftDef :: Set(FunName, Set(Variable)) -> Def -> Def
parameterLiftDef S (f \equiv \lambda v_1, \dots, \lambda v_n. e) =
  applySolutionToDef S (f \equiv \lambda v_1, \dots, \lambda v_n. (parameterLiftExp S e))

parameterLiftExp :: Set(FunName, Set(Variable)) -> Exp -> Exp
parameterLiftExp S (e_0 \dots e_n) =
  in (e'_0 \dots e'_n)
  where
    parameterLiftExp S (LetRec {d_1, \dots, d_k} e_0) =
      let G = ref (\(), \())
          V_fi = ref (FV(f_i)), for each (d_k \equiv (f_i \equiv \lambda v_1, \dots, \lambda v_n. e)) \in \{d_1, \dots, d_k\}
          P_fi = \{v_1, \dots, v_n\}, for each (d_k \equiv (f_i \equiv \lambda v_1, \dots, \lambda v_n. e)) \in \{d_1, \dots, d_k\}
          in for each f_i \in \{f_1, \dots, f_k\} do
              Graph.add-edge G f_i g
          let (G' as (V', E')) = Graph.coalesceSCC G
              succ_p = \{q \in V' | (p, q) \in E'\}, for each p \in V'
              F_p = \bigcup_{q \in succ_p} q, for each p \in V'
              propagate :: List(Set(FunName)) -> ()
              propagate [] = ()
              propagate (p :: \tau) =
                let V = (\bigcup_{f \in P} V_f) \cup (\bigcup_{p \in F_p} V_p)
                    in for each f \in p do
                        V_f := V \setminus P_f;
                        (propagate r)
                in (propagate (List.reverse (Graph.breadthFirstOrdering G')))
          let S' = S \cup \{(f_1, V_{f_1}), \dots, (f_k, V_{f_k})\}
              f_s = map (parameterLiftDef S') \{d_1, \dots, d_k\}
              e'_0 = parameterLiftExp S' e_0
          in (LetRec f_s e'_0)

parameterLiftExp S (if e_0 e_1 e_2) =
  let e'_i = parameterLiftExp S e_i, for each e_i \in \{e_0, e_1, e_2\}
      in (if e'_0 e'_1 e'_2)

parameterLiftExp S f = applySolutionToExp S f
parameterLiftExp S v = v
parameterLiftExp S (x as literal) = x
applySolutionToDef :: Set(FunName, Set(Variable)) -> Def -> Def
applySolutionToDef (S as \{(f_1, \{v_1, \dots, v_n\}), \dots\}) (f \equiv \lambda v_1, \dots, \lambda v_n. e) =
  (f \equiv \lambda v_1, \dots, \lambda v_n. \lambda v'_1, \dots, \lambda v'_n. e)

applySolutionToDef S d = d
applySolutionToExp :: Set(FunName, Set(Variable)) -> Exp -> Exp
applySolutionToExp (S as \{(f_1, \{v_1, \dots, v_n\}), \dots\}) f = (f v_1 \dots v_n)
applySolutionToExp S e = e

```

Figure 5. Parameter lifting: free variables are made parameters

4 Related work

We review alternative approaches to handling free variables in higher-order, block-structured programming languages, namely supercombinator conversion, closure conversion, lambda-dropping, and incremental versions of lambda-lifting and closure conversion. Finally, we address the issue of formal correctness.

4.1 Supercombinator conversion

Peyton Jones's textbook describes the compilation of functional programs towards the G-machine [24]. Functional programs are compiled into supercombinators, which are then processed at run time by graph reduction. Supercombinators are closed lambda-expressions. Supercombinator conversion [16,23] produces a series of closed terms, and thus differs from lambda-lifting that produces a series of mutually recursive equations where the names of the equations are globally visible.

Peyton Jones also uses strongly connected components for supercombinator conversion. First, dependencies are analyzed in a set of recursive equations. The resulting strongly connected components are then topologically sorted and the recursive equations are rewritten into nested letrec blocks. There are two reasons for this design:

1. it makes type-checking faster and more precise; and
2. it reduces the number of parameters in the ensuing supercombinators.

Supercombinator conversion is then used to process each letrec block, starting outermost and moving inwards. Each function is expanded with its own free variables, and made global under a fresh name. Afterwards, the definition of each function is replaced by an application of the new global function to its free variables, including the new names of any functions used in the body. This application is mutually recursive in the case of mutually recursive functions, relying on the laziness of the source language; it effectively creates a closure for the functions.

Peyton Jones's algorithm thus amounts to first applying dependency analysis to a set of mutually recursive functions and then to perform supercombinator conversion. As for dependency analysis, it is only used to optimize type checking and to minimize the size of closures.

In comparison, applying our algorithm locally to a letrec block would first partition the functions into strongly connected components, like dependency analysis. We use the graph structure, however, to propagate information, not to obtain an ordering of the nodes for creating nested blocks. We also follow Johnson's algorithm, where the names of the global recursive equations are free in each recursive equations, independently of the evaluation order. Instead, Johnson's algorithm passes all the free variables that are needed by a function and its callees, rather than just the free variables of the function.

To sum up, Peyton Jones's algorithm and our revision of Johnson's algorithm both coalesce strongly connected components in the dependency graph,

```

blockFloatProgram :: Program -> Program
blockFloatProgram p = foldr makeUnion (map blockFloatDef p) ()
blockFloatDef :: Def -> (Set(Def), Def)
blockFloatDef (f ≡ λv1. ... λvn. e) = let (Fnew, e') = blockFloatExp e
    in (Fnew, f ≡ λv1. ... λvn. e')
blockFloatExp :: Exp -> (Set(Def), Exp)
blockFloatExp (e0 ... en) =
  let (F1, e'1) = blockFloatExp e1, for each ei ∈ {e0, ..., en}
      Fnew = foldr (∪) {F1, ..., Fn} ()
      in (Fnew, e'0 ... e'n)
blockFloatExp (LetRec {d, ...} e0) =
  let (Fnew, e) = blockFloatExp (LetRec {...} e0)
      in ({d} ∪ Fnew, e)
blockFloatExp (LetRec θ e0) = blockFloatExp e0
blockFloatExp (If e0 e1 e2) =
  let (F1, e'1) = blockFloatExp e1, for each ei ∈ {e0, e1, e2}
      in (F0 ∪ F1 ∪ F2, (If e'0 e'1 e'2))
blockFloatExp f = (θ, f)
blockFloatExp v = (θ, v)
blockFloatExp (x as literal) = (θ, x)
makeUnion :: ((Set(Def), Def), Set(Def)) -> Set(Def)
makeUnion ((Fnew, d), S) = Fnew ∪ {d} ∪ S

```

Figure 6. Block floating: block structure is flattened

When parameter-lifting a set of mutually recursive functions $\{f_1, \dots, f_n\}$, and some function f_i defines a variable x that is free in one of its callees f_j , a naive algorithm expands the parameter list of the function with x . The sets P_{f_i} used in our parameter-lifting algorithm serve to avoid this problem.

3.5 Revisiting the example of Section 2.3

Applying the algorithm of Figure 5 to the program of Figure 1 processes the `main` function by processing its body. The letrec block of the body is processed by first constructing a dependency graph similar to that shown in Figure 2 (except that we simplify the description of the algorithm to not include the sets of free variables in the nodes). Coalescing the strongly connected components of this graph yields a single node containing the three functions. Since there is only a single node, the propagation step only serves to associate each function in the node with the union of the free variables of each of the functions in the component. These variable sets directly give rise to a new set of solutions.

Each of the functions defined in the letrec block and the body of the letrec block are traversed and expanded with the variables indicated by the set of solutions. Block floating according to the algorithm of Figure 6 yields the program of Figure 3.

but for different purposes, our purpose being to reduce the time complexity of lambda-lifting from cubic to quadratic.

4.2 Closure conversion

The notion of closure originates in Landin's seminal work on functional programming [19]. A closure is a functional value and consists of a pair: a code pointer and an environment holding the denotation of the variables that occur free in this code. Efficient representations of closures are still a research topic today [30].

Closure conversion is a key step in Standard ML of New Jersey [4,5], and yields scope-insensitive programs. It is akin to supercombinator conversion, though rather than creating a closure through a mutually recursive application, the closure is explicitly created as a vector holding the values of the free variables of the possibly mutually recursive functions.

In his textbook [24], Peyton Jones concluded his discussion between lambda-lifting and supercombinator/closure conversion by pointing out a tension between

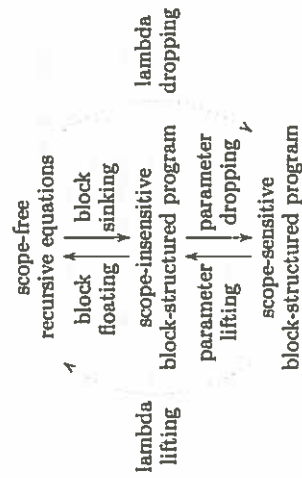
- passing all the [denotations of the] free variables of all the callees but not the values of the mutually recursive functions (in lambda-lifting), and
- passing all the values of the mutually recursive functions but not the free variables of the callees (in closure conversion).

He left this tension unresolved, stating that future would tell which algorithm (lambda-lifting or closure conversion) would prevail.

Today we observe that in the compiler world (Haskell, ML, Scheme), closure conversion has prevailed, with only one exception in Scheme [9]. Conversely, in the program-transformation world [8,10,21], lambda-lifting has prevailed. We also observe that only for lambda-lifting has an inverse transformation been developed: lambda-dropping.

4.3 Lambda-dropping

Lambda-dropping is the inverse of lambda-lifting [13]:



Block floating is reversed by block sinking, which creates block structure by making functions used in only one function local to this function. Parameter lifting is reversed by parameter dropping, which removes redundant formal parameters that are originally defined in an outer scope and that always take on the same value.

Lambda-lifting simplifies the structure of a program. However, a program transformation that employs lambda-lifting as a preprocessing phase tends to output a lambda-lifted program rather than a block-structured one. For one point, the resulting programs are less readable. For another point, compilers are often geared for source programs with few parameters.¹ Therefore, increased numbers of formal parameters often form a major overhead in procedure invocation at run time. Against these odds, lambda-dropping can be applied to re-create block structure and reduce the number of formal parameters.

A few years ago, Appel has pointed out a correspondence between imperative programs in SSA form and functional programs using block structure and lexical scope [2]. Specifically, he has shown how to transform an SSA program into its functional representation.² We were struck by the fact that this transformation corresponds to performing block sinking on the recursive equations defining the program. As for the transformation into optimal SSA form (which diminishes the number of ϕ -nodes), it is equivalent to parameter dropping. This made us conclude that lambda-dropping can be used to transform programs in SSA form into optimal SSA form [13].

This conclusion prompted us to improve the complexity of the lambda-dropping algorithm to $\mathcal{O}(n \log n)$, where n is the size of the program, by using the dominance graph of the dependency graph. We then re-stated lambda-lifting in a similar framework using graph algorithms, which led us to the result presented in the present article.

Even with the improvement presented in this article, we are still left in an asymmetric situation where lambda-lifting and lambda-dropping do not have the same time complexity. With some thought, though, this asymmetry is not so surprising, since lambda-dropping is applied to the output of lambda-lifting, and the complexity is measured in terms of the size of the output program. Measuring the complexity of lambda-dropping in terms of the size of the program before lambda-lifting yields a relative time complexity of lambda-dropping of $\mathcal{O}((n^2) \log(n^2))$, which is $\mathcal{O}(n^2 \log n)$, a fitting match for the $\mathcal{O}(n^2 \log n)$ time complexity of lambda-lifting.

4.4 Mixed style

In order to preserve code locality, compilers such as Twobit [9] or Moby [25] often choose to lift parameters only partially. The result is in the mixed style described at the end of Section 1.1.

¹ For example, the magic numbers of parameters, in OCaml, are 0 to 7.

² The point is made comprehensively in his SIGPLAN Notices note, which is also available in his home page [3].

In more detail, rather than lifting all the free variables of the program to become formal parameters, parameter lifting is used incrementally to transform programs by lifting only a subset of the free variables of each function. If a function is to be moved to a different scope, however, it needs to be passed the free variables of its callees as parameters. As was the case for global lambda-liftings, propagating the additional parameters through the dependency graph requires cubic time. To improve the time complexity, our quadratic-time parameter-lifting algorithm can be applied to the subsets of the free variables instead. The improvement in time complexity for incremental lambda-lifting is the same as what we observed for the global algorithm.

We note that a partial version of closure conversion also exists, namely Stecker and Wand's [31], that leaves some variables free in a closure because this closure is always applied in the scope of these variables. We also note that combinator-based compilers [33] could be seen as using a partial supercombinator conversion.

4.5 Correctness issues

Only idealized versions of lambda-lifting and lambda-dropping have been formally proven correct. Danvy has related lambda-lifted and lambda-dropped functions and their fixed point [12]. Fischbach and Hannan have capitalized on the symmetry of lambda-lifting and lambda-dropping to formalize them in a logical framework, for a simply typed and recursion-free source language [14].

Overall, though, and while there is little doubt about Johnson's original algorithm, its semantic correctness still remains to be established.

5 Conclusion and future work

We have shown that a transitive closure is not needed for lambda-lifting. In this article, we have reformulated lambda-lifting as a graph algorithm and improved its time complexity from $\mathcal{O}(n^3 \log n)$ to $\mathcal{O}(n^2 \log n)$, where n is the size of the program. Based on a simple example where lambda-lifting generates a program of size $\mathcal{O}(n^2)$, we have also demonstrated that our improved complexity is close to optimal.

The quadratic-time algorithm can replace the cubic-time instances of lambda-lifting in any partial evaluator or compiler, be it for global or for incremental lambda-lifting.

As for future work, we are investigating lambda-lifting in the context of object-oriented languages. Although block structure is instrumental in object-oriented languages such as Java, Beta and Simula [11,15,20], existing work on partial evaluation for object-oriented languages has not addressed the issue of block structure [27,28,29]. Problems similar to those found in partial evaluation for functional languages appear to be unavoidable: residual methods generated in a local context may need to be invoked outside of the scope of their class. Side effects, however, complicate matters.

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A Graph algorithms

The description of the algorithm for parameter lifting (Figure 5) makes use of a number of graph and list algorithms. We give a short description of each of these algorithms in Figure 7.

```

Graph.add-edge :: Graph( $\alpha$ )  $\rightarrow$  ( $\alpha$ ,  $\alpha$ )  $\rightarrow$  ( $\alpha$ ,  $\alpha$ )
Graph.add-edge  $G$  ( $n_1$ ,  $n_2$ ) : Updates  $G$  to contain the nodes  $n_1$  and  $n_2$  as well as an
edge between the two.
Graph.coalesceSCC :: Graph( $\alpha$ )  $\rightarrow$  Graph(Set( $\alpha$ ))
Graph.coalesceSCC  $G$  : Returns  $G$  with its strongly connected components coalesced
into sets [1].
Graph.breadthFirstOrdering :: Graph( $\alpha$ )  $\rightarrow$  List( $\alpha$ )
Graph.breadthFirstOrdering  $G$  : Returns a list containing the nodes of  $G$ , in a
breadth-first ordering.
List.reverse :: List( $\alpha$ )  $\rightarrow$  List( $\alpha$ )
List.reverse  $L$  : Returns  $L$  with its elements reversed.

```

Figure 7. Graph and list functions.

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