Compiling functional languages

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Spring school "Semantics of programming languages" Agay, 24-29 March 2002 Goal of this lecture: survey implementation techniques for functional languages, and show some of the (syntactic) theories explaining these techniques.

What is a functional language?

- Examples: Caml, Haskell, Scheme, SML, ...
- "A language where functions are taken seriously".
- A language that supports the manipulation of functions as first-class values, roughly as predicted by the λ -calculus.

Recipe for a functional language:

- Fix an evaluation strategy on the λ -calculus.
 - Weak evaluation (no reductions under λ).
 - Call-by-name or call-by-value.
- Add constants and primitive operations
 - Integers and arithmetic operations.
 - Booleans, strings, characters, input/output, ...
- Add primitive data structures
 - Predefined: tuples, lists, vectors, ...
 - User-defined: records, sums, recursive data types.

- 1. Representing functions as values: closures and environments.
- 2. Abstract machines for functional languages.
- 3. Optimizing functions: control-flow analysis.
- 4. Optimizing data representations.

Part 1

Function closures

let scale = λ n. λ x. n * x

```
let scale_by_2 = scale 2
```

```
let scale_by_10 = scale 10
```

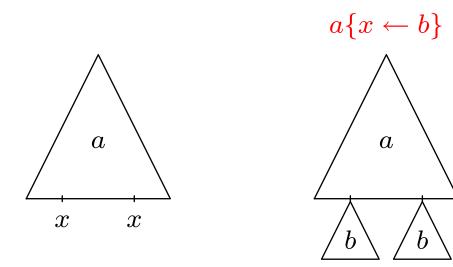
scale 2 should return a function computationally equivalent to $\lambda x \cdot x * 2$, in accordance with the β -reduction rule:

 $(\lambda x.a) \ b \to a\{x \leftarrow b\}$

 $(\lambda x.a) \ b \rightarrow a\{x \leftarrow b\}$

Naive substitution involves a full copy of the term a, replacing occurrences of x by b.

(If b has free variables, a copy of b may also be necessary to avoid variable captures. However, with head reductions, b is guaranteed to be closed.)



The problem gets worse in a compiled setting

In a compiled setting, the standard representation for a function value is a pointer to a piece of compiled code that:

- expects its argument x in a register arg;
- computes the function body *a*;
- leaves the result of a in a register res;
- returns to caller.

If we naively apply this "function as code pointer" model to a functional language, functions that return functions should dynamically generate a piece of compiled code representing the returned function. Example:

```
let scale = \lambdan. \lambdax. n * xscale 2 returnsscale 10 returnsmov res, arg<br/>mul res, 2<br/>returnmov res, arg<br/>mul res, 10<br/>return
```

Problem: this requires run-time code generation, which is complex and expensive (in time and in space for the generated code blocks). Remark: the generated blocks of code share the same "shape": they differ only in the value of the variable n that is free in the returned function λx . x * n.

mov res, arg
mul res, <the value of n passed to scale>
return

Idea: share the common code and put the varying parts (i.e. the values of free variables) in some separate data structure called an environment.

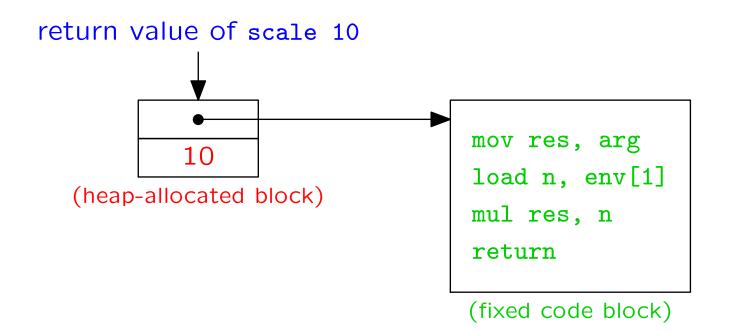
The same idea applies to term-level substitution: to represent $a\{x \leftarrow b\}$, just keep a unchanged and record separately the binding of x to b in an environment.

All functional values are represented by closures.

Closures are dynamically-allocated data structures containing:

- a code pointer, pointing to a fixed piece of code computing the function result;
- an environment: a record providing values for the free variables of the function.

To apply a closure, move the environment part in a conventional register env and call the code pointer.



Closures can also be viewed as terms

$(\lambda x.a)[e]$

where $\lambda x.a$ is a function with free variables y, z, \ldots , and the environment e is an explicit substitution binding terms b, c, \ldots to y, z, \ldots

This is the key to efficient interpreters for functional languages, as we shall now see via the following route:

- Standard semantics: small-step (reduction-based), classical substitutions.
- Big-step semantics with classical substitutions.
- Big-step semantics with environments.
- Small-step semantics with environments / explicit substitutions.

Small-step semantics with classical substitutions

Terms:	$a ::= x \mid \lambda x.a \mid a_1 \mid a_2 \mid cst \mid op(a_1, \dots, a_n)$
Constants:	$c ::= 0 \mid 1 \mid \ldots$
Operators:	$op ::= + pair fst snd \dots$
Values:	$v ::= \lambda x.a \mid cst$

Reduction rules:

$$(\lambda x.a) \ v \ \rightarrow \ a\{x \leftarrow v\} \qquad (\beta_v) op(v_1, \dots, v_n) \ \rightarrow \ v \text{ if } v = \overline{op}(v_1, \dots, v_n) \quad (\delta)$$

Strategy: head reductions, call-by-value, left-to-right

$$\frac{a \to a'}{a \ b \to a' \ b} \qquad \qquad \frac{b \to b'}{v \ b \to v \ b'}$$
$$a \to a'$$

 $op(v_1, \ldots, v_{k-1}, a, b_{k+1}, \ldots, b_n) \to op(v_1, \ldots, v_{k-1}, a', b_{k+1}, \ldots, b_n)$

Big-step semantics with classical substitutions

Rather than chain elementary reductions $a \rightarrow a_1 \rightarrow \ldots \rightarrow v$, define a "big step" evaluation relation $a \Rightarrow v$ that "jumps" from a to its value v. (a must be closed.)

$$(\lambda x.a) \Rightarrow (\lambda x.a) \qquad cst \Rightarrow cst$$

$$\frac{a \Rightarrow (\lambda x.c) \qquad b \Rightarrow v \qquad c\{x \leftarrow a\} \Rightarrow v'}{a \ b \Rightarrow v'}$$

$$a_1 \Rightarrow v_1 \qquad \dots \qquad a_n \Rightarrow v_n \qquad v = \overline{op}(v_1, \dots, v_n)$$

$$op(a_1, \dots, a_n) \Rightarrow v$$

Theorem: if a is closed, $a \xrightarrow{*} v$ if and only if $a \Rightarrow v$.

Function values are closures $(\lambda x.a)[e]$.

Terms: $a ::= x \mid \lambda x.a \mid a_1 \mid a_2 \mid cst \mid op(a_1, \dots, a_n)$ Values: $v ::= (\lambda x.a)[e] \mid cst$ Environments: $e ::= \{x_1 \leftarrow v_1, \dots, x_n \leftarrow v_n\}$

The evaluation relation becomes $e \vdash a \Rightarrow v$.

a may now contain free variables, but those must be bound by the environment e.

$$e \vdash x \Rightarrow e(x) \qquad e \vdash (\lambda x.a) \Rightarrow (\lambda x.a)[e] \qquad e \vdash cst \Rightarrow cst$$

$$\frac{e \vdash a \Rightarrow (\lambda x.c)[e'] \qquad e \vdash b \Rightarrow v \qquad e' + \{x \leftarrow v\} \vdash c \Rightarrow v'}{e \vdash a \ b \Rightarrow v'}$$

$$\frac{e \vdash a_1 \Rightarrow v_1 \qquad \dots \qquad e \vdash a_n \Rightarrow v_n \qquad v = \overline{op}(v_1, \dots, v_n)}{e \vdash op(a_1, \dots, a_n) \Rightarrow v}$$

De Bruijn notation: rather than identify variables by names, identify them by position.

Environments become sequences of values $v_1 \dots v_n \varepsilon$, accessed by position: variable number n is bound to v_n . Terms: $a ::= n | \lambda a | a_1 | a_2 | cst | op(a_1, ..., a_n)$ Values: $v ::= (\lambda a)[e] | cst$ Environments: $e ::= \varepsilon | v.e$

$$e \vdash n \Rightarrow e(n) \qquad e \vdash (\lambda a) \Rightarrow (\lambda a)[e] \qquad e \vdash cst \Rightarrow cst$$

$$\frac{e \vdash a \Rightarrow (\lambda c)[e'] \qquad e \vdash b \Rightarrow v \qquad v.e' \vdash c \Rightarrow v'}{e \vdash a \ b \Rightarrow v'}$$

$$\frac{e \vdash a_1 \Rightarrow v_1 \qquad \dots \qquad e \vdash a_n \Rightarrow v_n \qquad v = \overline{op}(v_1, \dots, v_n)}{e \vdash op(a_1, \dots, a_n) \Rightarrow v}$$

All this leads to the canonical efficient interpreter for call-by-value λ -calculus:

```
type term = Var of int | Lambda of term | App of term * term
          | Constant of int | Primitive of primitive * term list
and value = Int of int | Closure of term * environment
and environment = value list
and primitive = value list -> value
let rec eval env term =
  match term with
    Var n -> List.nth env n
  | Lambda a -> Closure(a, env)
  | App(a, b) ->
      let (Closure(c, env')) = eval env a in
      let v = eval env b in
      eval (v :: env') c
  | Constant n \rightarrow
      Int n
  Primitive(p, arguments) ->
      p (List.map (eval env) arguments)
```

Small-step semantics with explicit substitutions

Big-step semantics are good for implementing interpreters, but small-step semantics allow easier reasoning on non-terminating or "stuck" evaluations.

The notion of environment can be internalized in a small-step, reduction-based semantics: λ -calculus with explicit substitutions.

Terms: $a ::= n | \lambda a | a_1 | a_2 | a[e]$ Environments: $e ::= \varepsilon | a.e$

Basic reduction rules:

More complete calculi of explicit substitutions

The previous rules are the minimal rules we need to describe weak call-by-value reduction.

To describe other strategies, including strong reduction, richer calculi of explicit substitutions are needed:

Explicit substitutions, M. Abadi, L. Cardelli, P.L. Curien, J.J. Lévy, Journal of Functional Programming 6(2), 1996.

Confluence properties of weak and strong calculi of explicit substitutions, P.L. Curien, T. Hardin, J.J. Lévy, Journal of the ACM 43(2), 1996.

Closure representation strategies

• When compiling an application, nothing is known about the closure being called (this can be the closure of any function in the program).

 \rightarrow The code pointer must be at a fixed, predictable position in the closure block.

• The environment part of a closure is not accessed during application. Its structure matters only to the code that builds the closure and the code for the function body.

 \rightarrow Considerable flexibility in choosing the layout for the environment part.

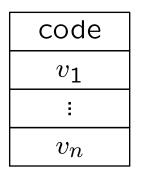
• The environment used for evaluation $(e \text{ in } e \vdash a \Rightarrow v)$ need not have the same structure as the environment put in closures $(e' \text{ in } (\lambda a)[e'])$.

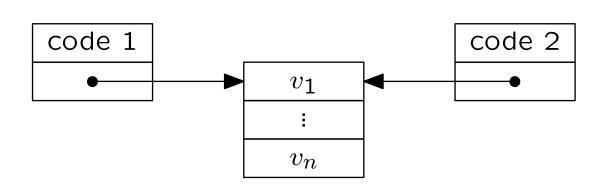
 \rightarrow Even more flexibility in putting part of the evaluation environment e in registers or on the stack rather than in a heap-allocated block.

Examples of closure representations

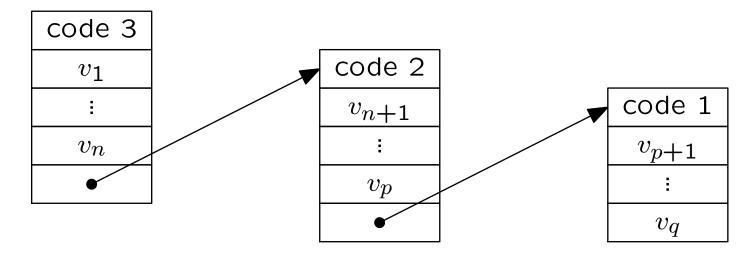
One-block closure

Two-block closures (with environment sharing)





Linked closures



Time trade-off:

- One-block closures: slower to build; faster access to variables.
- Linked closures: faster to build; slower access to variables.

Space trade-off:

- Minimal environments: (bind only the free variables) fewer opportunities for sharing; avoid space leaks.
- Larger environments: (may bind more variables) more opportunities for sharing; may cause severe space leaks.

Modern implementations use one-block closures with minimal environments.

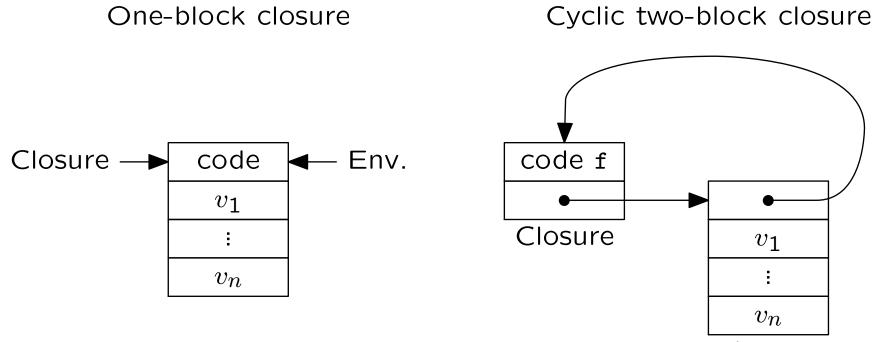
Recursive functions need access to their own closure:

let rec f x = \dots List.map f l \dots

(The body of f needs to pass the closure of f as first argument to the higher-order function List.map.)

This can be done in several ways:

- Reconstruct the closure of f from the current environment.
- Treat f as a free variable of the function body: put a pointer to the closure of f in the environment of that closure (cyclic closures).
- In the one-block approach: the environment passed to f is its own closure, just reuse it.



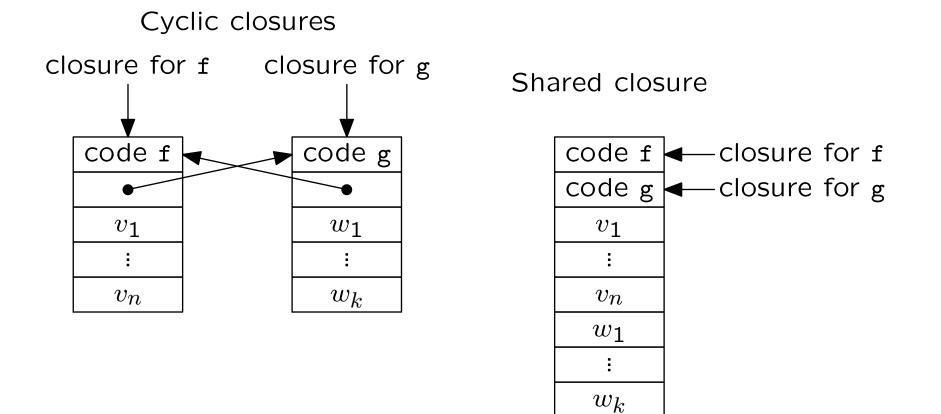
Environment

Mutually recursive functions need access to the closures of all functions in the mutual recursive definition:

let rec f x = ... List.map f l1 ... List.map g l2 ... and g y = ... List.map f l3 ...

This can be done in two ways:

- The closure for f contains a pointer to that of g and conversely (cyclic closures).
- Share a closure between f and g using infix pointers (*Compiling with Continuations*, A. Appel, Cambridge U. Press, 1992.)



Part 2

Abstract machines

• Interpretation:

control (the sequencing of computations) is represented by a source-level, tree-shaped term. The interpreter walks this tree at run-time.

• Native compilation:

control is compiled down to a sequence of machine instructions. These instructions are those of a real processor, and are executed by hardware.

• Compilation to abstract machine code: control is compiled down to a sequence of abstract instructions. These instructions are those of an abstract machine: they do not match existing hardware, but are chosen to match closely the operations of the source language. Arithmetic expressions: $a ::= cst | op(a_1, ..., a_n)$

The machine uses a stack to hold intermediate results.

Compilation is translation to "reverse Polish notation".

$$\mathcal{C}(cst) = \text{CONST}(cst)$$

$$\mathcal{C}(op(a_1, \dots, a_n)) = \mathcal{C}(a_1); \mathcal{C}(a_2); \dots; \mathcal{C}(a_n); I(op)$$

where I(+) = ADD, I(-) = SUB, etc.

The machine has two components:

- a code pointer c giving the next instruction to execute;
- a stack *s* holding intermediate results.

Notation for stacks: top of stack is on the left

push v on $s: s \longrightarrow v.s$ pop v c

pop v off s: $v.s \longrightarrow s$

Transitions of the machine:

State be	fore	State after		
Code	Stack	Code	Stack	
CONST(cst); c	S	С	cst.s	
ADD; c	$n_2.n_1.s$	С	$(n_1 + n_2).s$	
SUB; c	$n_2.n_1.s$	С	$(n_1 - n_2).s$	

Halting state: $code = \varepsilon$ and stack = v.s.Result of computation is v. (Similar to Landin's SECD and Cardelli's FAM.)

Three components in this machine:

- a code pointer c giving the next instruction to execute;
- an environment e giving values to free variables;
- a stack s holding intermediate results and return frames.

Compilation scheme:

$$C(n) = ACCESS(n)$$

$$C(\lambda a) = CLOSURE(C(a); RETURN)$$

$$C(a b) = C(a); C(b); APPLY$$

(Constants and arithmetic: as before.)

Transitions

State	State after				
Code	Env	Stack	Code	Env	Stack
ACCESS(n); c	e	8	c	e	e(n).s
CLOSURE(c'); c	e	8	c	e	[c', e].s
APPLY; c	e	v.[c',e'].s	<i>c</i> ′	v.e'	c.e.s
RETURN; c	e	v.c'.e'.s	c'	e'	v.s

- ACCESS(n): push value of variable number n
- CLOSURE(c'): push closure [c, e] of c with current env.
- APPLY: pop argument, pop closure, push return frame, jump to closure code.
- RETURN: restore saved code and environment from stack.

Code for a stack-based abstract machine can be executed either

• By expansion of abstract machine instructions to real machine instructions, e.g.

CONST(i)	>	pushl i
ADD	>	popl %eax addl 0(%esp), %eax

• By efficient interpretation.

The interpreter is typically written in C as shown on the next slide, and is about one order of magnitude faster than term-level interpretation.

```
value interpret(int * start_code)
{
 register int * c = start_code;
 register value * s = bottom_of_stack;
 register environment e;
 while(1) {
    switch (*c++) {
      case CONST: *s++ = *c++; break;
      case ADD: s[-2] = s[-2] + s[-1]; s--; break;
      case ACCESS: *s++ = Lookup(e, *c++); break;
      case CLOSURE: *s++ = MakeClosure(*c++, e); break;
      case APPLY: arg = *--sp; clos = *--sp;
                    *sp++ = (value) c; *sp++ = (value) e;
                    c = Code(clos);
                    e = AddEnv(arg, Environment(clos));
                    break;
      case RETURN:
                   res = *--sp; e = (environment) *--sp; c = (int *) *--sp;
                    *sp++ = res; break;
      case STOP:
                   return *--sp;
    }
 }
}
```

An abstract machine for call-by-name: Krivine's machine

As before, three components in this machine: code c, environment e, stack s.

However, the stack does not contain values, but thunks: closures [c, e] representing expressions whose evaluation is delayed until needed.

Compilation scheme:

$$C(n) = ACCESS(n)$$

$$C(\lambda a) = GRAB; C(a)$$

$$C(a b) = PUSH(C(b)); C(a)$$

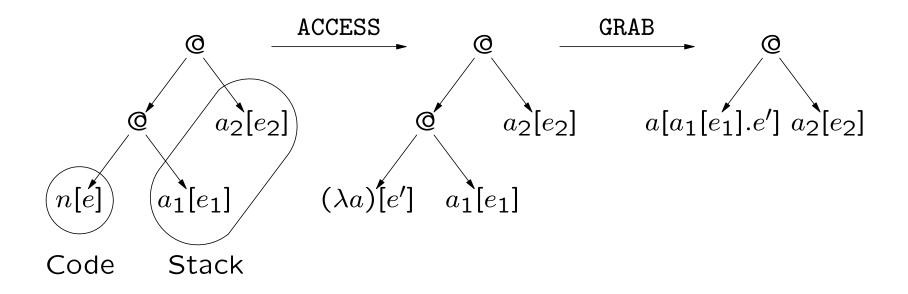
Transitions of Krivine's machine

State before			State after		
Code	Env	Stack	Code	Env	Stack
ACCESS(n); c	e	8	<i>c</i> ′	e'	s if $e(n) = [c', e']$
GRAB; c	e	[c', e'].s	c	[c', e'].e	S
PUSH(c'); c	e	8	c	e	[c',e].s

- ACCESS(n): fetch thunk bound to variable number n, and proceed evaluating it
- GRAB: pop the next argument provided off the stack, and add it to the environment (β -reduction step)
- PUSH(c): build thunk for c and push it

The stack maintains the current spine of applications.

The code is the leftmost outermost part of the spine.



Making call-by-name practical

Realistic abstract machines for call-by-name functional languages are more complex than Krivine's machine in two aspects:

• Constants and strict primitive operations:

Operators such as integer addition are strict in their arguments. Extra machinery is required to reduce sub-expressions to head normal form in a strict way.

• Sharing of evaluations (lazy evaluation):

Call-by-name reduces an expression every time its head normal form is needed. Lazy evaluation performs the reduction the first time it is needed, then caches the result for further reference.

See e.g. *Implementing lazy functional languages on stock hardware: the Spineless Tagless G-machine*, S.L. Peyton Jones, Journal of Functional Programming 2(2), Apr 1992.

Proving the correctness of an abstract machine

At this point, we have two notions of evaluation for terms:

- 1. Source-level evaluation with environments: $a[e] \xrightarrow{*} v$ or $e \vdash a \Rightarrow v$.
- 2. Compilation, then execution by the abstract machine:

$$\begin{pmatrix} c = \mathcal{C}(a) \\ e = \varepsilon \\ s = \varepsilon \end{pmatrix} \xrightarrow{*} \begin{pmatrix} c = \varepsilon \\ e = \dots \\ s = v \dots \end{pmatrix}$$

Do these two notions agree? Does the abstract machine compute the right results?

The compilation scheme is compositional: each sub-term is compiled to code that evaluates it and leaves its value on the top of the stack.

This parallels exactly a derivation of $e \vdash a \Rightarrow v$ in the big-step semantics, which contains sub-derivations $e' \vdash a' \Rightarrow v'$ for each sub-expression a'.

Theorem: if $e \vdash a \Rightarrow v$, then

$$\begin{pmatrix} \mathcal{C}(a); k \\ \mathcal{C}(e) \\ s \end{pmatrix} \xrightarrow{*} \begin{pmatrix} k \\ \mathcal{C}(e) \\ \mathcal{C}(v).s \end{pmatrix}$$

The compilation scheme $C(\cdot)$ is extended to values and environments as follows:

$$C(cst) = cst$$

$$C((\lambda a)[e]) = [(C(a); \text{RETURN}), C(e)]$$

$$C(v_1 \dots v_n . \varepsilon) = C(v_1) \dots C(v_n) . \varepsilon$$

The theorem is proved by induction on the derivation of $e \vdash a \Rightarrow v$. We show the most interesting case: function application.

$$\frac{e \vdash a \Rightarrow (\lambda c)[e'] \qquad e \vdash b \Rightarrow v' \qquad v'.e' \vdash c \Rightarrow v}{e \vdash a \ b \Rightarrow v}$$

 $(\mathcal{C}(a); \mathcal{C}(b); \text{APPLY}; k \mid \mathcal{C}(e) \mid s)$ $\downarrow *$ ind. hyp. on first premise $(\mathcal{C}(b); \text{APPLY}; k \mid \mathcal{C}(e) \mid [(\mathcal{C}(c); \text{RETURN}), \mathcal{C}(e')].s)$ $\downarrow *$ ind. hyp. on second premise (APPLY; $k \mid \mathcal{C}(e) \mid \mathcal{C}(v').[(\mathcal{C}(c); \text{RETURN}), \mathcal{C}(e')].s$) \downarrow APPLY transition $(\mathcal{C}(c); \text{RETURN} \mid \mathcal{C}(v'.e') \mid k.\mathcal{C}(e).s)$ $\downarrow *$ ind. hyp. on third premise (RETURN | $\mathcal{C}(v'.e')$ | $\mathcal{C}(v).k.\mathcal{C}(e).s$) \downarrow **RETURN** transition $(k \mid \mathcal{C}(e) \mid \mathcal{C}(v).s)$

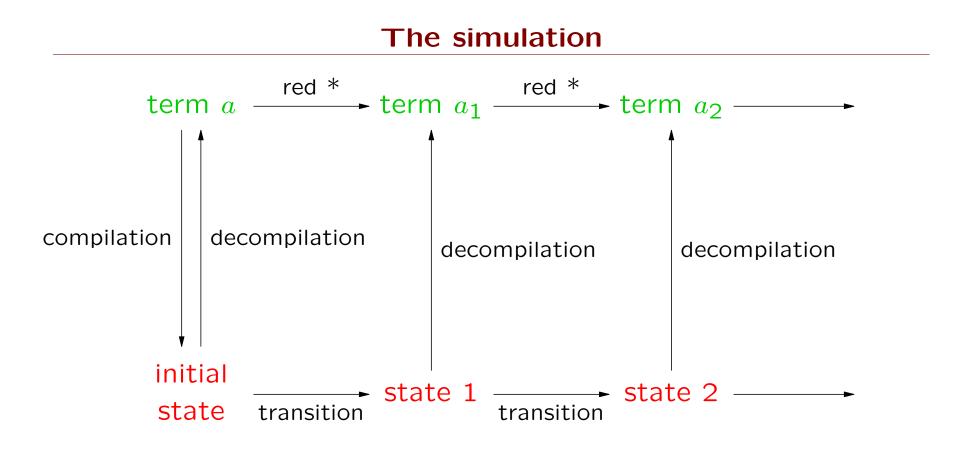
The previous theorem shows the correctness of the abstract machine for terminating terms. However, if the term a does not terminate, $e \vdash a \Rightarrow v$ does not hold, and we do not know anything about the execution of the compiled code.

(It might loop, but it might just as well stop and reply "42".)

To show correctness for all terms (terminating or not), we need to establish a simulation between machine transitions and source-level reductions:

Each transition of the machine corresponds to zero, one, or several source-level reductions.

See Functional Runtimes within the Lambda-Sigma Calculus, T. Hardin, L. Maranget, B. Pagano, J. Func. Prog 8(2), 1998.



Problem: not all intermediate states of the machine correspond to the compilation of a source term.

Solution: define a decompilation function \mathcal{D} : States \rightarrow Terms that is defined on all intermediate states and is left inverse of the compilation function.

Idea: decompilation is a symbolic variant of the abstract machine: it reconstructs source terms rather than performing the computations.

Decompilation of values:

 $\mathcal{D}(cst) = cst$ $\mathcal{D}([c,e]) = (\lambda a)[\mathcal{D}(e)]$ if $c = \mathcal{C}(a)$; RETURN

Decompilation of environments and stacks:

$$\mathcal{D}(v_1 \dots v_n.\varepsilon) = \mathcal{D}(v_1) \dots \mathcal{D}(v_n).\varepsilon$$
$$\mathcal{D}(\dots v \dots c.e \dots) = \dots \mathcal{D}(v) \dots c.\mathcal{D}(e) \dots$$

Decompilation of concrete states:

$$\mathcal{D}(c \mid e \mid s) = \mathcal{D}(c \mid \mathcal{D}(e) \mid \mathcal{D}(s))$$

Decompilation of abstract states (E, S already decompiled):

 $\mathcal{D}(\varepsilon \mid E \mid a.S) = a$ $\mathcal{D}(\text{CONST}(cst); c \mid E \mid S) = \mathcal{D}(c \mid E \mid cst.S)$ $\mathcal{D}(\text{ACCESS}(n); c \mid E \mid S) = \mathcal{D}(c \mid E \mid E(n).S)$ $\mathcal{D}(\text{CLOSURE}(c'); c \mid E \mid S) = \mathcal{D}(c \mid E \mid (\lambda a)[E].S)$ if $c' = \mathcal{C}(a)$; RETURN $\mathcal{D}(\text{RETURN}; c \mid E \mid a.c'.E'.S) = \mathcal{D}(c' \mid E' \mid a.S)$ $\mathcal{D}(\text{APPLY}; c \mid E \mid b.a.S) = \mathcal{D}(c \mid E \mid (a \ b).S)$ $\mathcal{D}(I(op); c \mid E \mid a_n \dots a_1.S) = \mathcal{D}(c \mid E \mid (op(a_1, \dots, a_n)).S)$ Simulation: if $\mathcal{D}(S)$ is defined, and the machine performs a transition from S to S', then $\mathcal{D}(S')$ is defined and $\mathcal{D}(S) \xrightarrow{*} \mathcal{D}(S')$.

Progress: if S is not a final state, and $\mathcal{D}(S)$ is defined and reduces, then the machine can perform a transition from S.

No stuttering: there exists a non-negative measure |S| on machine states such that if the machine does a silent transition from S to S' (i.e. $\mathcal{D}(S) = \mathcal{D}(S')$), then |S| > |S'|.

Initial states: $\mathcal{D}(\mathcal{C}(a) | \varepsilon | \varepsilon) = a$ if a is closed.

Final states: $\mathcal{D}(\varepsilon \mid e \mid v.s) = \mathcal{D}(v)$.

Theorem: Let a be a closed term and $S = (\mathcal{C}(a) | \varepsilon | \varepsilon)$.

- If a ^{*}→ v, then the abstract machine started in state S terminates and returns the value C(v).
- If *a* reduces infinitely, the machine started in state *S* performs an infinite number of transitions.

Part 3

Optimized compilation of functions: flow analysis

Step 1: Replace functions by closures; make explicit the construction, passing, and accessing of the environments.

This brings us to a conventional intermediate language that manipulates code pointers (i.e. closed functions).

(Think of the intermediate language as a subset of C with support for dynamic allocation and garbage collection.)

Step 2: Optimize and generate machine code from the intermediate language.

- Write your own code generator using conventional compiler technology (OCaml, SML/NJ)
- Or use a C compiler (GHC, Bigloo).

```
\mathbf{x} = \mathbf{x}
                                        [cst] = cst
[op(a_1, ..., a_n)] = op([a_1], ..., [a_n])
[[a b]] = let clos = [[a]] in clos[0] (clos, [[b]])
[\lambda x. a] =
    let code_fn (clos, x) =
       let v_1 = clos[1] and ... and v_n = clos[n] in ||a||
    in
       makeblock(code_fn, v_1, ..., v_n)
```

where v_1, \ldots, v_n are the variables free in $\lambda x.a$.

Note: the function code_fn above has no free variables. The occurrence of code_fn in makeblock(code_fn, \ldots) denotes a pointer to its (fixed) code.

From intermediate language to machine code

Standard compiler technology applies largely unchanged. optimizations Register transfer language Intermediate order evaluations with virtual registers language instruction selection register allocation Actual Register transfer language final code assembly with real registers generation code and stack frames optimizations

See any good compiler textbook, e.g. *Modern Compiler Implementation in ML*, Andrew W. Appel, Cambridge Univ. Press, 1998.

[[a b]] = let clos = [[a]] in clos[0] (clos, [[b]])

Our translation scheme transforms every function application into

- one load clos[0]
- one call to a computed address clos[0](...)

Calls to a computed address are expensive on modern processors:

- the destination address is usually not predicted in advance;
- this stalls the pipeline (cannot fetch and start executing instructions from the called function while earlier instructions complete).

Typically, a factor of 10 more expensive than a call to a statically known address.

The overhead of calls to computed addresses can be avoided in many practical situations:

let succ = λx . x + 1 in succ (succ 2)

The two applications of succ "obviously" call the code for $\lambda x.x + 1$, and no other code.

let rec f = λx f arg ...

The application of f always calls the code for the current function.

```
let sort_list = \lambda ordering. \lambda list. ... in
... sort_list (\lambda x.\lambda y. compare(x,y)) some_list ...
```

If there are no other calls to Sort.list in the program, all applications of ordering in the body of sort_list call the code for $\lambda x.\lambda y.compare(x, y)$.

In all of these cases:

- application of a function in the static scope of its definition
- recursive calls
- higher-order functions applied only once

we could (and should):

- Generate calls to statically-known code addresses.
- Or if the called function is small, perform inline expansion (compile-time β -reduction) of its body, e.g.

succ (succ 2) \Rightarrow (2 + 1) + 1

A program analysis is needed to discover those opportunities for closure optimization.

Control-flow analyses (Shivers, PLDI 1988) approximate at each application point the set of functions that can be called here (in other terms, the set of function values that can flow to this application point).

If that set is a singleton $\{\lambda x.a\}$, we can generate a direct call to the code for a, or inline it if a is small enough.

In all cases, we also get an approximation of the call graph for the program (who calls who?), required for later interprocedural optimizations (e.g. global register allocation).

Since functions are first-class values, CFA is actually a data-flow analysis that keeps track of the flow of functional values and determines control-flow along the way.

CFA sets up a system of constraints of the form

$V(\ell_1) \subseteq V(\ell_2)$

meaning that all values at program point ℓ_1 can flow to point ℓ_2 .

Solve that system into a flow graph:

producer point \longrightarrow consumer point

(constant, $\lambda x.a$ (function parameter, result of operator, function part of appl., data constructor) argument to operator) For (if a^m then b^n else c^p)^{ℓ}: add the constraints

 $V(n) \subseteq V(\ell)$ (the then branch flows to the result) $V(p) \subseteq V(\ell)$ (the else branch flows to the result)

For (let $x = a^m \text{ in } b^n)^{\ell}$: add

$$V(m) \subseteq V(x)$$
 $V(n) \subseteq V(\ell)$

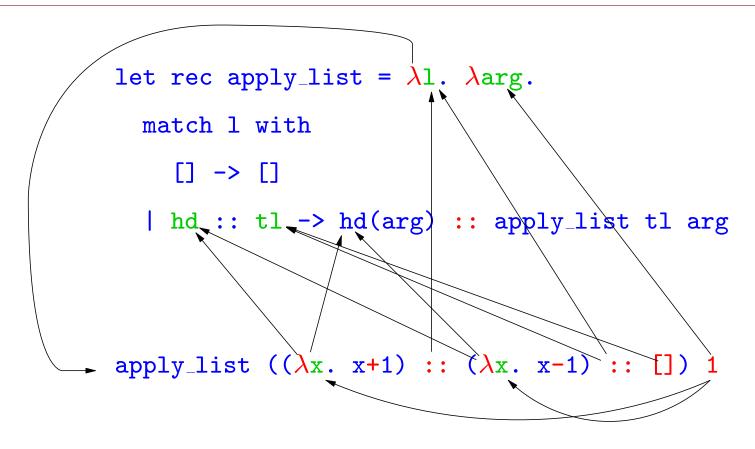
For $(a^m(b^n))^{\ell}$:

for each function λx . c^q in V(m), add the constraints

 $V(n) \subseteq V(x)$ (the argument flows to the parameter) $V(q) \subseteq V(\ell)$ (the function result flows to the application result)

Note: need to interlace constraint building and constraint solving, and iterate till fixpoint is reached.

An example of CFA



(Not all edges are shown.)

We have determined the two functions called from apply_list, and also that they are always applied to 1.

Basic algorithm (0-CFA) is $O(n^3)$ (*n* is the size of the program). Main applications:

- Optimize function calls in functional languages. (As seen before.)
- Optimize method dispatch in object-oriented languages.
 (Similar problems to that of closures. Roughly, an object is a closure with multiple entry points.)
- Eliminate run-time type tests in dynamically-typed languages such as Scheme.

(E.g. if all values flowing to + are integers, + does not need to check the type of its arguments.)

• More applications later...

More precise analyses:

- Polyvariant analyses (*n*-CFA, polymorphic splitting, ...): distinguish between different call sites of the same function.
- Finer approximation of values (Heintze's set-based analysis): capture the shapes of data structures using grammars.

Less precise (faster) analyses:

- Coarser representations of sets of values:
 Ø or {v} (singletons) or ⊤ (all values).
- Do not iterate till fixpoint: (Ashley, ICFP 1997) start with ⊤ on all variables and do 1 or 2 iterations.
- Use equality constraints (unification) in addition to inclusion constraints.

Digression on inlining: why not compile-time β **-reductions?**

An analysis like CFA might seem overkill for function inlining: the effect of inlining can also be achieved by β -reductions at compile-time.

> let succ = λx . x + 1 in succ (succ 2) $\xrightarrow{\beta}$ (λx . x + 1) ((λx . x + 1) 2) $\xrightarrow{\beta}$ (λx . x + 1) (2 + 1) $\xrightarrow{\beta}$ (2 + 1) + 1

Unrestricted compile-time reductions may cause program size explosion (or even execution of the whole program at compile-time!); various size-based heuristics control when to perform them. Problem 1: these heuristics are difficult to calibrate.

let p =
$$(\lambda x. small)$$
, $(\lambda y. HUGE)$ in fst(p)(1)
 $\xrightarrow{\beta}$ fst($(\lambda x. small)$, $(\lambda y. HUGE)$)(1)
 $\xrightarrow{\beta}$ $(\lambda x. small)$ (1) $\xrightarrow{\beta}$ $small{x \leftarrow 1}$

An intermediate reduct can be huge, then collapse to a much smaller term; shall the compiler lose the opportunity for inlining, or risk explosion?

Problem 2: β -reduction is not selective enough

let p =
$$(\lambda x.small)$$
, $(\lambda y.HUGE)$ in
fst(p)(1), snd(p)(2), snd(p)(3)
 $\xrightarrow{\beta*} small\{x \leftarrow 1\}$, $HUGE\{y \leftarrow 2\}$, $HUGE\{y \leftarrow 3\}$

There is no way to inline *small* without duplicating *HUGE*.

CFA avoids both problems by simulating what β -reduction would do without actually performing it.

Connections between CFA and type systems

CFA can be used as a type system if enriched with safety checks (e.g. fail if an integer flows to an application site).

Conversely, many type systems (and type inference algorithms) can be viewed as checking / approximating the flow of data in a program.

Palsberg and O'Keefe (TOPLAS 1995) show equivalence between:

- 0-CFA with safety checks;
- the Amadio-Cardelli type system (subtyping + recursive types).

Provides an efficient type inference algorithm for that system.

Also: type inference algorithms for type systems with subtyping are based on inclusion constraints similar to those used by CFA. (Aiken and Wimmers, FPCA 1993; Smith et al, MFPS 1995; Pottier, ICFP 1998.)

Part 4

Optimizing the representation of data

Representations for high-level data structures

High-level data structures (such as ML's datatypes) leave considerable flexibility to the compiler in deciding a data representation.

 \rightarrow clever representation tricks are feasible.

(Would be hard to do by hand in C, at least portably.)

Examples:

- For dynamically-typed languages (Scheme): clever tagging scheme (to embed the type of an object in its bit pattern).
- For ML's datatypes: clever encodings of the constructor.

```
type expr = Const of int | Var
    | Sum of expr * expr | Prod of expr * expr
```

Constant constructors are represented by odd integers 1, 3, ... (Bit pattern: $\dots xxx1$)

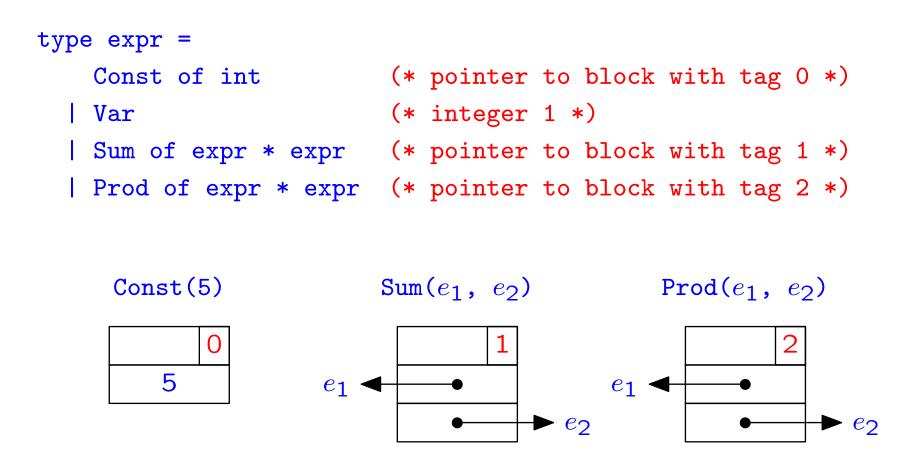
Constructors with arguments are represented by word-aligned pointers to heap blocks.

(Bit pattern: $\ldots xx00$)

The heap block contains one byte (the "tag" byte) representing the number of the constructor.

(This byte is stored at no extra cost in the header word required by the garbage collector.)

Example of datatype representation



Without static typing (Scheme):

- Need tagging to implement run-time type tests.
- All data types must fit a common format (usually one word).
 - \rightarrow floats are boxed (heap-allocated);
 - \rightarrow records are boxed;
 - \rightarrow arrays are arrays of pointers to boxed elements.
- All functions must use the same calling conventions: e.g. argument in R0; result in R0.

With monomorphic static typing (Pascal, C)

- No need to support run-time type tests.
- Different data types can have different sizes.
 - \rightarrow unboxed floats
 - \rightarrow unboxed records (if small enough)
 - \rightarrow flat arrays

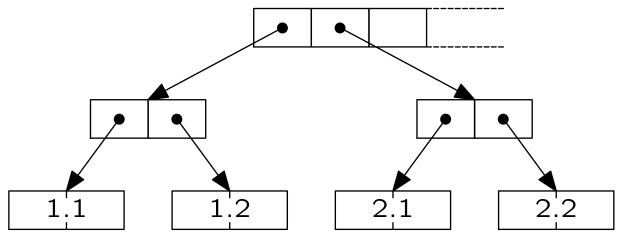
The compiler determines the size from the static type:

|int| = 1 word |float| = 2 words $|\tau \times \sigma| = |\tau| + |\sigma|$

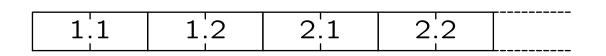
 Functions of different types can use different calling conventions. E.g. use floating-point registers for float arguments and results.

float \rightarrow floatargument in FP0, result in FP0int \times int \rightarrow intargument in R0 and R1, result in R0

In Scheme:



In C:



The type system guarantees type safety, but does not assign a unique type to every value at compile-time:

Polymorphism:	Type abstraction:
fun x -> x : $\forall \alpha. \ \alpha \rightarrow \alpha$	type t
	val x : t
	val f : t -> t
Actual type of x: any	Actual type of x: unknown
Size of x: variable	Size of x: unknown
Calling conventions: variable	Calling conventions: unknown

• Restrict polymorphism and type abstraction.

Modula: abstract types must be pointer types. Java: cannot coerce integers and floats to/from type Object. Problem: unnatural.

• Code replication.

Ada, C++: compile a specialized version of a generic function for each type it is used with.

Problem: code size explosion; link-time code generation.

Revert to Scheme-style representations.

Problem: inefficient; lots of boxing and unboxing.

- Use run-time type inspection: pass type information at run-time to polymorphic code; use this information to determine sizes and layouts at run-time.
- Mix C-style representations for monomorphic code and Scheme-style representations for polymorphic code.
- Combine Scheme-style representations with local unboxing, partial inlining, and special treatment of arrays.

In order to reconstruct exact types of data structures at run-time, polymorphic function must receive as extra arguments the types to which they are specialized.

let f x = x let f α x = x let g x = f (x, x) let g β x = f $\langle \beta \times \beta \rangle$ (x, x) g 5 g $\langle \text{int} \rangle$ 5

In this example, this allows f to determine at run-time that its x parameter has actual type $int \times int$.

The TIL approach (Harper, Morrisett, et al, PLDI 1996):

- Use C-style, "flat", multi-word representations of data structures (just like in a monomorphic type system).
- In polymorphic code, compute size information, data layout, and calling conventions from the run-time type information.
- (In monomorphic code, this information is computed at compile-time.)

Source code:

```
let assign_array a b i = b.(i) <- a.(i)
Generated code, Scheme style:
assign_array(a, b, i) {
    load one word from a + i * 4;
    store this word at b + i * 4;
}
Generated code, TIL style:
assign_array(a, a, b, i) {
    s = size_of_type(a);
    copy s bytes
    from a + i * s
    to b + i * s;
    }
</pre>
```

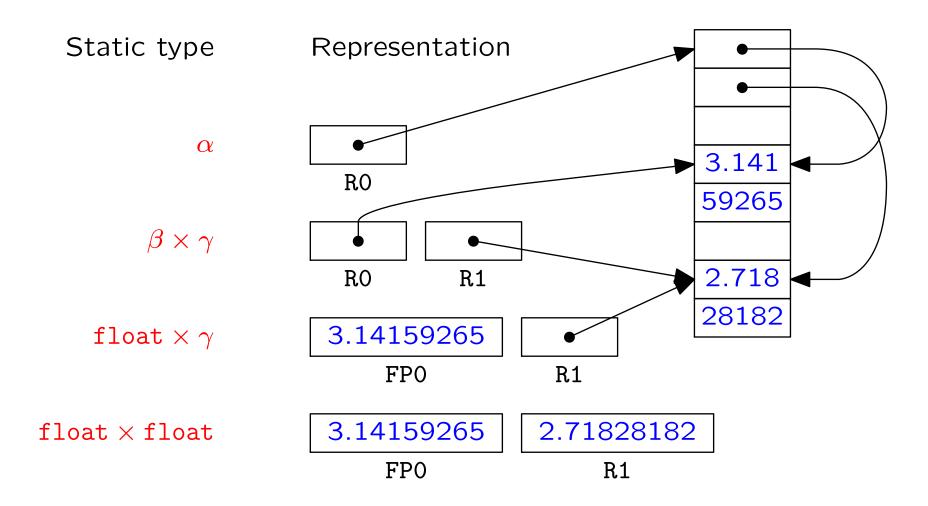
Variant (Ohori, Lisp.Symb.Comp.1993): pass only the size of types at run-time, not representations of whole type expressions.

(Leroy, POPL 1992; Shao and Appel, PLDI 1995; the SML/NJ compiler)

Use C-style representations for data whose exact type is known at compile-time (i.e. inside monomorphic code).

Revert to Scheme-style representations for manipulating data whose type is not completely known at compile-time (i.e. inside polymorphic code).

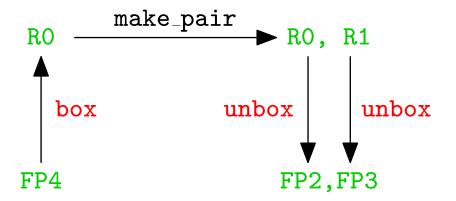
Insert coercions between the two representations at interface points.



Source code:

let make_pair x = (x, x) in ... make_pair 3.41519

Coercion diagram:



Generated code:

let make_pair x = (x, x) in ...
let (fst, snd) = make_pair(box_float(3.14159)) in
 (unbox_float(fst), unbox_float(snd))

When using a value v of type $\forall \alpha. \tau$ with type $\tau' = \tau \{ \alpha \leftarrow \sigma \}$, insert the coercion $[\tau \Rightarrow \tau'](v)$.

When implementing an abstract type $\exists \alpha. \tau$ by a value v of type $\tau' = \tau \{ \alpha \leftarrow \sigma \}$, insert the coercion $[\tau' \Rightarrow \tau](v)$.

(Objective Caml; Glasgow Haskell; Bigloo Scheme.)

Instead of basing the data representations on the types, use Scheme-style representations by default, plus:

• Perform intra-function unboxing by standard dataflow analysis:

let x = box(f) in \Rightarrow let x = f in ... unbox(x) ... unbox(x) ... \therefore x ... x ... x ...

- Extend it to inter-function unboxing using control-flow analysis or partial inlining.
- Use simple tagging schemes and tag testing to support important special cases of generic data structures (e.g. float arrays).

Partial inlining (a.k.a. the worker-wrapper technique)

(Peyton-Jones and Lauchbury, FPCA 1991; Goubault, SAS 1994.)

Split a function into:

- a worker function taking and returning unboxed data;
- a wrapper function performing the boxing and unboxing around the worker.

At call sites, try to inline the wrapper function (typically small) and hope its boxing and unboxing cancel out with those of the call context.

Example

```
let worker_f a b =
  (* a and b are unboxed floats *)
  (* compute result *)
  (* return unboxed float result *)
let f a b = box(worker_f (unbox a) (unbox b))
... unbox(f (box 3.14) (box 2.71)) ...
```

After inlining of f and simplifications:

... worker_f 3.14 2.71 ...

(Crucially depends on the availability of a good inlining pass in the compiler.)

Conclusions and perspectives

Other relevant topics not addressed here

Memory management and garbage collection. *Uniprocessor Garbage Collection Techniques*, P. Wilson, ftp://ftp.cs.utexas.edu/pub/garbage/bigsurv.ps.

Garbage collection support in intermediate code and code generation.

The C-- intermediate language, N. Ramsey and S. Peyton Jones.

Optimized compilation of pattern-matching. *Two techniques for Compiling Lazy Pattern Matching*, L. Maranget, INRIA research report 2385.

Relevance and adaptation of classic compiler optimizations.

Automatic parallelization.

Conclusions from an engineering standpoint

Compilation technology for functional languages is relatively mature:

- On comparable programs, achieve at least 50% of the performance of optimizing C compilers.
- Match or exceed the performance of C on allocation-intensive programs.

Still more work to do:

- Getting rid of the last factor of 2 is difficult.
- Truly efficient functional programs still require programmers to be conscious of performance issues while writing their code.

Functional languages promote software reliability:

clean semantics \Rightarrow formal methods \Rightarrow reliable programs

But: proving the correctness of source code is useless if the compiler is incorrect.

 $\rightarrow \text{Certified compilers} \\ Proof \vdash \forall Prog. \quad Comp(Prog) \equiv Prog$

→ Certifying compilers (proof-carrying code) $\forall Prog. \ Certif(Prog) \vdash Comp(Prog) \equiv Prog$

Today: we are able to certify realistic bytecode compilers and abstract machines.

Tomorrow: certification of optimizing native-code compilers?