CENG444: Compiling Functional Languages

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This is a short summary of Simon Peyton Jones's 1987 book *The Implementation of Functional Programming Languages*, chapters 2, 12, 13 and 16.

- λ-calculus: Uniform representation and evaluation of functions. Program evaluation = reduction.
- Graph reduction: A simple Virtual Machine for FPLs.
- Supercombinators: Life without global variables.
- SK-machine: Life without variables and intermediate abstractions.

It is a real machine, not abstract or toy! (See Clarke et al. 1980)

 λ -calculus is a uniform way to represent functions, without a need for names.

$$f(x, y) = x + y$$

$$\lambda x.\lambda y. + xy$$

Application of functions to arguments:

$$f(3,4) = 7$$

Lambda calculus

$$\lambda x.\lambda y.(+xy)34 = \lambda y.(+3y)4 = (+34) = 7$$

Functions can be arguments:

$$f(g,x)=g(x)$$

 $\lambda g.\lambda x.gx$

Lambda calculus

Function notation (λ notation)

Function abstraction and application (β conversion)

Function equivalence (α and η conversions)

 β -conversion: $(\lambda x.E)M \leftrightarrow_{\beta} E[M/x]$

E[M/x]: expression E where M is substituted for **free** occurrences of x.

CONS =
$$\lambda a. \lambda b. \lambda f. f a b$$

$$\mathbf{HEAD} = \lambda c.c(\lambda a.\lambda b.a)$$

TAIL =
$$\lambda c.c(\lambda a.\lambda b.b)$$

So that we have

Lambda calculus oo●ooo

CONSx y = list with head x and tail y

$$\mathsf{HEAD}(\mathsf{CONS}x\ y) = x$$

$$TAIL(CONSx y) = y$$

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Lambda calculus

Substitute definitions to obtain expected behaviour:

$$\mathsf{HEAD}(\mathsf{CONS}x\ y) = (\lambda c. c(\lambda a. \lambda b. a))(\mathsf{CONS}x\ y)$$

= **CONS**
$$xy(\lambda a.\lambda b.a) = (\lambda a.\lambda b.\lambda f.f \ a \ b) \ x \ y \ (\lambda a.\lambda b.a)$$

$$= (\lambda b.\lambda f.f \times b)y(\lambda a.\lambda b.a) = (\lambda f.f \times y)(\lambda a.\lambda b.a)$$

$$= (\lambda a.\lambda b.a) \times y = (\lambda b.x) y = x$$

All common PL expressions can be given a λ -calculus interpretation.

Expected behaviour:

Lambda calculus

IFTRUE
$$E_1$$
 $E_2 = E_1$

IFFALSE
$$E_1$$
 $E_2 = E_2$

The following definitions give that behaviour:

$$\mathbf{IF} = \lambda f. \lambda d. \lambda e. fde$$

TRUE =
$$\lambda a. \lambda b. a$$

FALSE =
$$\lambda a. \lambda b. b$$

Ex: **IFTRUE**x y = x

Lambda calculus 000000

IFTRUE $x y = (\lambda f. \lambda d. \lambda e. f)$ **TRUE**x y

 $= (\lambda d.\lambda e.\mathsf{TRUE} d e) x y = (\lambda e.\mathsf{TRUE} x e) y$

 $= (\lambda e.\mathsf{TRUE} x \ e) y = \mathsf{TRUE} x \ y$

 $= (\lambda a.\lambda b.a)xy = (\lambda b.x)y = x$

What if we defined **IF** = $\lambda f.f$?

A combinator is a lambda term without free variables.

We can eliminate renaming of variables, because it's a theorem of λ-calculus.

The following are equivalent functions (modulo variable names). They behave the same.

$$\lambda x$$
. + 1 x λy . + 1 y

Combinators •00000000

 α -conversion: $(\lambda x.M) \leftrightarrow_{\alpha} (\lambda y.M[y/x])$ if y is not free in M

These are equivalent too, because they behave the same as well:

$$\lambda x + 1 x + 1$$

 η -conversion: $\lambda x.F x \leftrightarrow_{\eta} F$ (if x does not occur free in F)

note: $\lambda x. * x x$ is not eta reducible to (*x)

How do we do recursion without names? The fixpoint combinator Y takes care of that:

FAC =
$$\lambda n.$$
IF(= $n \ 0)1(* n($ **FAC**(- $n \ 1)))$

 β conversion in the direction of abstraction gives:

$$\mathbf{H} =_{\beta} \lambda fac.(\lambda n.\mathbf{IF}(= n 0) 1 (* n (fac(- n 1))))$$
 FAC

FAC = **H FAC**, where **H** is
$$\lambda fac.(\cdots)$$

FAC is called the fixpoint of **H**.

Combinators 00000000

Note that **H** does not refer to **H**, so our recurring name problem is solved.

All we need is a definition **Y** that takes a function and returns its fixpoint:

$$YH = H(YH)$$

It's called the fixpoint combinator:

Combinators 000000000

$$\mathbf{Y} = \lambda h.(\lambda x.h(x x)) (\lambda x.h(x x))$$

Now, def. of **FAC** is non-recursive because

FAC = H FAC and

YH = H(YH)

Therefore.

FAC = Y H

$$FACp = YHp =$$

$$H(YH) p =$$

$$\lambda fac.(\lambda n.IF(= n 0) 1 (* n (fac(- n 1))))(YH) p =$$

$$\lambda n.$$
IF $(= n \ 0) \ 1 \ (* n \ ((YH) \ (- n \ 1))) \ p =$

$$IF(= p 0) 1 (* p ((YH) (- p 1))) =$$

In case $p \neq 0$, rewrite **Y H** as **H** (**Y H**)

So that **H** can get its argument (-p 1)

```
FAC = YH
FAC 1 = Y H 1 =
H (Y H) 1 =
\lambda f \lambda n. IF (= n \ 0) \ 1 \ (\times n \ (f \ (-n \ 1))) \ (\mathbf{Y} \ \mathbf{H}) \ 1 =
\lambda n. IF (= n \cdot 0) 1 (× n \cdot (\mathbf{Y} \cdot \mathbf{H} \cdot (-n \cdot 1))) 1 =
IF (= 1 \ 0) \ 1 \ (\times \ 1 \ (Y \ H \ (-1 \ 1))) =
\times 1 (Y H 0) =
\times 1 (H (Y H) 0) =
\times 1 ((\lambda f \lambda n. IF (= n 0) 1 (\times n (f (- n 1)))) (Y H) 0) =
\times 1 ((\lambda n. IF (= n 0) 1 (\times n (Y H (- n 1)))) 0) =
\times 1 (IF (= 0 0) 1 (\times 0 (Y H (- 0 1)))) =
\times 11 =
```

$$\mathbf{YH} = (\lambda h.(\lambda x.h(x x)) (\lambda x.h(x x))) \mathbf{H} =$$

$$(\lambda x.\mathbf{H}(x x))(\lambda x.\mathbf{H}(x x)) = (\text{apply 1st to 1nd}):$$

$$H((\lambda x.H(x x))(\lambda x.H(x x))) =$$

H(Y H)

Summary:

- all data can be represented as lambda abstractions.
- all functions can be represented as lambda abstractions.

How do we evaluate them?

Methods for β -reduction:

With variables: Graph reduction

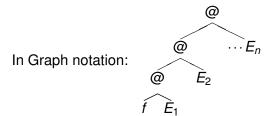
With piece-meal internal abstractions: Supercombinators

Without variables, with only compile-time abstractions: **SK** machine

Graph reduction as program execution

The strategy of choice in LISP.

General template of a program in FPL: $f E_1 E_2 \cdots E_n$



f may be 1) data; 2) built-in function with k < n arguments; 3) lambda abstraction; 4) variable

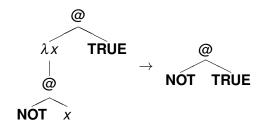
case 1: Program is done.

case 2: Redex is $(f E_1 \cdots E_k)$

case 3: Redex is $(f E_1)$

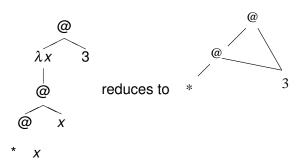
case 4: error (because the variable is free—it's leftmost)

ex: a graph with lambdas: $(\lambda x.NOTx)$ TRUE \rightarrow_{β} NOT TRUE



Why is it called *graph* reduction?

 $(\lambda x. * x x)3 = 9$: Substitute pointers to the argument for formal parameter x



Without pointer substitution, we would evaluate M four times for no reason, because it will always yield the same value.

Y reduction is best described as a graph as well:

$$\begin{array}{ccccc}
& & & & & & & & & & \\
Y & & & & & & & & & \\
Y & f & = f & (Y & f) & = f & (f & (Y & f)) & = & \\
f & (f & (f & (Y & f))) & & & & & \\
\end{array}$$

Supercombinators:

In Graph Reduction, finding a lambda body to reduce during execution is costly; it requires a tree walk at every step.

\$S is called a supercombinator if

$$S = \lambda x_1 . \lambda x_2 ... \lambda x_n . E$$

and

- 1) \$S has no free variables
- 2) E is not a lambda abstraction
- 3) any lambda expression in E is a supercombinator
- 4) n > 0

Because of (2), supercombinator's arguments can be supplied all at once.

ex:
$$\lambda x. * x x$$

$$\lambda x.\lambda y. - x y$$

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The following are not supercombinators:

$$\lambda x. - y x$$

$$\lambda f.f(\lambda x.f \times x)$$

Combinators have by convention names, such as \mathbf{B} , \mathbf{S} , \mathbf{K} , \mathbf{I} , \mathbf{C} , \mathbf{T} etc.

Less than a handful is enough to write ANY lambda expression (assuming no free variables) WITHOUT VARIABLES, as Curry and Feys showed in 1958.

Amazingly, two are enough: S and K.

Supercombinators have names made-up during compilation; they are not primitives (hence the \$X notation).

Combinatory Logic of Curry & Feys is equivalent to λ -calculus and Turing Machines.

The (boxing) strategy in Supercombinator Compilation:

- 1) Derive a set of supercombinator definitions (upper box)
- 2) Single Expression to be evaluated (lower box)
- 3) Compile until bottom expression has no lambdas, just evaluations.

Supercombinator definitions

Expresion to be evaluated

Ex: Program $(\lambda x.\lambda y.* x y)$ 3 4 compiles to

\$**XY** *x y* = * *x y* \$**XY**3 4

All arguments must be supplied; expression below is not a supercomb.

$$\$XY \times y = *X y$$

$$\$XY3$$

Ex: Compiling the program: FAC 5

FAC =
$$\lambda n.$$
IF(= $n \cdot 0) \cdot 1 (* n ($ **FAC**(- $n \cdot 1)))$

 β conversion in the direction of abstraction gives:

$$=_{\beta} \lambda fac.(\lambda n.\mathbf{IF}(=n0) 1 (*n(fac(-n1))))$$
 FAC

Let
$$F = \lambda fac.(\lambda n.\mathbf{IF}(= n 0) 1 (* n (fac(-n 1))))$$

Not a supercombinator: inner lambda expr has a free variable (fac)

By β -abstraction, inner lambda body is equivalent to

\$N
$$fac = \lambda w.\lambda n.$$
IF $(= n \ 0) \ 1 \ (* n \ (w \ (-n \ 1)))) fac$

and \$N is a supercombinator

The compiled boxes look like:

FAC = ···
\$Nw
$$n = IF(= n 0) 1 (* n (w (- n 1)))$$

 $\lambda fac.(\lambda n.$N fac n)$ **FAC** 5

Compile until bottom expression has no lambdas, just evaluations.

This is possible because the lambda body $\lambda fac.(\cdots)$ is a supercombinator.

```
FAC = \cdots
Nw n = IF(= n 0) 1 (* n (w (- n 1)))
\mathbf{SNF} fac n = \mathbf{SN} fac n
$NF FAC 5
```

By η -conversion, we can conclude $\mathbf{NF} = \mathbf{N}$, and eliminate CENG444 One of them

Having **FAC** at the bottom might look circular, but We know that FAC = Y H where Y is the fixpoint combinator, and

$$\mathbf{H} = \lambda fac.(\lambda n.\mathbf{IF}(= n \, 0) \, 1 \, (* n \, (fac(-n \, 1))) = \$ \mathbf{N} \, fac \, n$$

We can revise the definitions to reflect that.

$$Y = \cdots$$

\$H fac n = \$N fac n
\$Nw n = IF(= n 0) 1 (* n (w (- n 1)))
\$N (Y H) 5

Therefore $\mathbf{H} = \mathbf{N}$.

Final compilation is:

```
\mathbf{Y} = \cdots
Nw n = IF(= n 0) 1 (* n (w (- n 1)))
$N (Y $N) 5
```

- 1) Derive a set of supercombinator definitions (upper box)
- 2) Single Expression to be evaluated (lower box)
- 3) Compile until bottom expression has no lambdas, just evaluations.

Supercombinator definitions

Expresion to be evaluated

Now fast-forward to Javascript: Wouldn't it be fun to do the upperboxes with arrow functions, and make program body simple use of these arrow functions (lower box).

SK Machine

Miranda and SASL use SK virtual machines.

If we don't want ANY variables, including those in the abstractions, we can use the universal combinators **S** and **K**.

Note that, Supercombinator Method's overhead for keeping environment is minimal but NOT ZERO: there are only local variables, values of which are all supplied at the same time.

SK systems allow any function to be written without variables, in terms of **S**, **K** and built-in functions (like **Y**) and constants.

It is easier to see them at work, starting with I as well.

Sf
$$g x \rightarrow f x (g x)$$

$$\mathbf{K} \times \mathbf{y} \rightarrow \mathbf{x}$$

$$\mathbf{I} x \to x$$

S allows us to see that the following functions are equivalent:

$$F_1 = \lambda x. (e_1 e_2)$$

$$F_2 = \mathbf{S} (\lambda x.e_1) (\lambda x.e_2)$$

$$F_1 a = F_2 a$$

If e_1 and e_2 are lambda expressions, **S** allows us to push down the variable one level more, until it is no longer possible:

$$\lambda x. (\lambda y.e_3 e_4) (\lambda y.e_5 e_6)$$

= $\lambda x. (\mathbf{S} (\lambda y.e_3) (\lambda y.e_4))(\mathbf{S} (\lambda y.e_5) (\lambda y.e_6))$

$$\lambda x.x = 1$$

$$\lambda x.c = \mathbf{K} c$$

We can use these equivalences as *program transformations*.

ex:
$$(\lambda x. * x x)$$
 5. This is same as $(\lambda x. ((*x) x))$ 5

$$= (\mathbf{S} (\lambda x. * x)(\lambda x. x)) 5 = (\mathbf{S} (\mathbf{S} (\lambda x. *)(\lambda x. x))(\lambda x. x)) 5$$

= (S(S(K*)I)I)5. By associativity this is same as

= **S** (**S** (**K** *) **I**) **I** 5. Now reduction, outermost combinator first:

= (**S** (**K** *) **I** 5) (**I** 5). By associativity same as

= **S** (**K** *) **I** 5 (**I** 5). By reduction (**K** * 5) (**I** 5) (**I** 5)

= * (15) (15) = *55

Try I a and S K Ka

In case of multiple lambdas, you must push in the innermost one first: take $\lambda x \lambda y + xy$

```
\lambda x \lambda y + xy = \lambda x (\lambda y ((+x)y))
                                                            by associativity
\lambda x(S(\lambda y. + x)(\lambda y. y)) =
\lambda x(S(S(\lambda y.+)(\lambda y.x))(SKK)) =
                                                        now push x inward
S(\lambda x.S(S(K+)(Kx)))(\lambda x.SKK) =
S(\lambda x.S(S(K+)(Kx)))(S(\lambda x.SK)(\lambda x.K)) =
S(\lambda x.S(S(K+)(Kx)))(S(S(KS)(KK))(KK)) =
S(S(\lambda x.S)(\lambda x.S(K+)(Kx)))(S(S(KS)(KK))(KK)) =
S(S(KS)(S(\lambda x.S(K+))(\lambda x.Kx)))(S(S(KS)(KK))(KK)) =
                            you can use eta-reduction too; try on \lambda x.\mathbf{K}x
S(S(KS)(S(\lambda x.S(K+))K))(S(S(KS)(KK))(KK)) =
S(S(KS)(S(S(\lambda x.S)(\lambda x.K+))K))(S(S(KS)(KK))(KK)) =
S(S(KS)(S(KS)(S(\lambda x.K)(\lambda x.+)))K))(S(S(KS)(KK))(KK)) =
Phew!!
Now, try (\lambda x.\lambda y. + xy)34 = +34 and above to see their
```

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equivalence.

Generally, I am more lenient in the exams, so I might ask something like:

$$(\lambda x. * x x)$$
 5. This is same as $(\lambda x. ((*x) x))$ 5

$$= (\mathbf{S} (\lambda x.* x)(\lambda x.x)) 5 = (\mathbf{S} (\mathbf{S} (\lambda x.*)(\lambda x.x))(\lambda x.x)) 5$$

= (S(S(K*)I)I)5. By associativity this is same as

= **S** (**S** (**K** *) **I**) **I** 5. Now reduction, outermost combinator first:

= (**S** (**K** *) **I** 5) (**I** 5). By associativity same as

= **S** (**K** *) **I** 5 (**I** 5). By reduction (**K** * 5) (**I** 5) (**I** 5)

= * (15) (15) = *55

Y = SSK(S(K(SS(S(SSK))))K)

It is not pretty, but it does the job.

Note that although the programmer can use variables in the source language for convenience,

the compiler completely eliminates them via program transformations.

There is no environment to keep during run-time.

For an entertaining story of Starling and Kestrel, try Ray Smullyan's 1985 book.

The Turkish equivalents are, imo, Saka and Kerkenez.

References

Clarke, T. J., P. J. Gladstone, C. D. MacLean, and A. C. Norman (1980). SKIM - the S, K, I reduction machine. In *Proceedings* of the 1980 ACM conference on LISP and functional programming, LFP '80, New York, NY, USA, pp. 128–135. ACM.

Smullyan, R. (1985). *To Mock a Mockingbird*. New York: Knopf.