



MENG INDIVIDUAL PROJECT

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TODO

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Contents

1	Introduction	4
1.1	Motivation	4
1.2	Objectives	4
2	Background	6
2.1	Arrows	6
2.1.1	Definition	6
2.1.2	Example: Calculate the mean	7
2.1.3	Application in parallel computation	8
2.2	Recursion Schemes	8
2.2.1	Definition	8
2.2.2	Example: Merge sort	9
2.3	Multiparty session types	10
2.3.1	Global types and local types	10
2.3.2	Applications in parallel computing	12
2.4	Message passing concurrency	12
2.4.1	Primitives for message-passing concurrency	12
2.4.2	Concurrency Monads	12
2.5	Free monad	14
2.5.1	Definition	14
2.5.2	Example	15
2.5.3	Applications	16
3	Alg: Algebraic Functional Language	17
3.1	Alg	17
3.1.1	Syntax	17
3.1.2	Example: Merge sort	18
3.2	ParAlg: Alg + role annotations	19
3.2.1	Inferring global types	19
3.2.2	Example: Parallel merge sort	19
3.3	Conclusion	19
4	SPar: A session typed free monad EDSL for concurrency	20
4.1	Computation: The Core EDSL	20
4.1.1	Syntax	20
4.1.2	Representation of recursive data structures	21
4.2	Communication: The Proc EDSL	21
4.3	Concurrent computation: A group of Proc	22
4.3.1	Operational semantics	22
4.3.2	Session types and duality checking	23
4.4	Conclusions	24
5	SPar: Implementation	25
5.1	Session types	25
5.1.1	Representations of session types in Haskell	25
5.1.2	Value-level duality check	25
5.1.3	Type-indexed Free Monad	26

5.1.4	Type-level duality check	26
5.2	SPar interpreter	26
5.2.1	Overview	26
5.2.2	Implementation	26
6	ArrowPipe: An arrow interface for writing SPar expressions	28
6.1	Syntax	28
6.1.1	Arrow interface	29
6.1.2	Example: Parallel programming patterns	29
6.2	Implementation of arrow combinators	31
6.3	Strategies for optimized role allocations	32
6.4	Satisfaction of arrow laws	34
6.5	Conclusions	34
7	Type-safe code generation from SPar	35
7.1	Instr: A low-level EDSL for channel communication	35
7.1.1	Syntax and semantic	35
7.1.2	Representation types	36
7.2	Compilation from SPar to Instr	37
7.2.1	Transformation from Proc to Instr	37
7.2.2	Strategies for channel allocations	38
7.2.3	Monad for code generation	38
7.3	Code generation to C: from Instr to C	39
7.3.1	Representations of Core data type in C	39
7.3.2	Compiling from Core to C	40
7.3.3	The structure of generated C code	41
7.4	Conclusion	41
8	Parallel algorithms and evaluation	42
8.1	Parallel algorithms	42
8.1.1	Four steps to write parallel algorithms in SPar	42
8.1.2	Example: Merge sort	43
8.1.3	Example: Word count	43
8.2	Benchmarks	43
8.2.1	Benchmarks against generated Haskell code	43
8.2.2	Benchmarks against C implementation	43
8.3	Evaluation	43
8.3.1	Design choices: Why Haskell?	43
9	Conclusions and future work	44
9.1	Future work	44

Chapter 1

Introduction

1.1 Motivation

Writing parallel software is not a trivial task. Parallel code is hard to write because it is usually written in low level languages with verbose and non-idiomatic decorations, hard to debug because machines, where code is written, are usually different from machines where code is intended to run and hard to maintain and reuse because even though the underlying algorithms are not changed, multiple version of parallel code is needed to tackle various platform and evolution of architectures.

There are many on-going pieces of research aimed at helping programmers write correct parallel programs smoothly. A common approach is to develop a higher level language and compile programmes in this language to required parallel code. There are many high-level frameworks for parallel programming (e.g. algorithmic skeletons[1], domain-specific languages for parallelism[2] or famous MapReduce parallel model[3]). An example is to use arrow terms (Section 2.1) to describe data flow implicitly and hence generate parallel code.

The workflow of writing parallel code has evolved from writing it directly in the target platform to writing software in a high-level language designed for parallel computation and then compiling to the target platform. In this project, we present a method to improve the backend of parallel code generation by introducing a monadic domain-specific language to act as a bridge between high-level and target low-level parallel languages.

This specific language needs to be general enough so that it supports multiple high-level parallel programming frameworks. It can be used to generate different parallel code, e.g. MPI¹, Cuda. Moreover, it can be interpreted with a simulator to aid debugging parallel programs.

With the help of this intermediate languages, the implementation complexity is reduced from $O(M \times N)$, where each of the M high-level languages needs to implement N compilers to generate parallel code in N different platforms, to $O(M + N)$, where each compiler of a high-level language implements a translation rule to the intermediate language which implements one compiler and N backend to generate different target languages.

In addition, it couples with multiparty session type (MPST) [5]. It takes advantages of properties of MPST to enable aggressive optimisation but ensuring code correctness and allow more meaningful static analysis; e.g. cost modelling for parallel programming.

1.2 Objectives

1. **Design:** Design the intermediate languages, argue its generality and build its connection with MPST.
2. **Translation:** Define a translate rule from the language in a high-level parallel framework to our language.
3. **Simulator:** Build a simulator to prove the correctness of code generation and act as a playground for experiments.

¹Message Passing Interface (MPI) is a standardised and portable message-passing standard designed by a group of researchers from academia and industry to function on a wide variety of parallel computing architectures [4].

4. **Code generation:** Generate parallel code in C.
5. **Static Analysis:** Session typing the language to obtain properties guaranteed by session types

More details of the objectives will be explained in the project planning at Section ??

Chapter 2

Background

This section is an overview of techniques that influence the design choices of our monadic language for parallel computation. First of all, we give an overview of techniques applied in the high-level parallel programming framework: arrows (Section 2.1) and recursion schemes (Section 2.2). We then introduce several techniques for message-passing concurrency: multiparty session types (Section 2.3) and monadic languages for concurrency (Section 2.4). In the end, we introduce free monads (Section 2.5), a technique valuable in implementing embedded domain-specific languages (EDSL).

2.1 Arrows

Arrow is a general interface to describe computation. It can ease the process of writing structured code suitable for parallelising. It also demos a common feature of the frameworks: parallelizability is empowered by underlying implicit but precise data-flow. On the other hand, converting to low-level message-passing code, which requires programmers to define communication using message-passing function and primitives, makes the data-flow explicit.

2.1.1 Definition

Listing 1 shows the Arrow definition in Haskell. Intuitively, an arrow type $y \ a \ b$ (that is, the application of the parameterised type y to the two parameter types b and c) can be regarded as a computation with input of type b and output of type b [6]. Visually, arrows are like pipelines (shown in Figure 2.1). In Haskell, an arrow y is a type that implements the following interface (type classes in Haskell are roughly interfaces). `arr` converts an arbitrary function into an arrow. `>>>` sequences two arrows (illustrated in Figure 2.1b). Taking two input, `first` apply the arrow to the first input while keeping the second untouched (Figure 2.1a). Conversely, `second` modifies the second input and keeps the first one unchanged. `***` applies two arrows to two input side by side (Figure 2.1d). `&&&` takes one input and applies two separate arrows to the input and its duplications (Figure 2.1c).

The simplest instance of arrow class is the function type (shown in Listing 2). It is worth noticing that only `arr` and `***` need to be implemented. The reset of function in the arrow type class can be defined in terms of the two functions. For example, `f &&& g = (f *** g) . arr (\b -> (b, b))` and `first = (***) id`

```
1  class Arrow y where
2      arr :: (a -> b) -> y a b
3      first :: y a b -> y (a, c) (b, c)
4      second :: y a b -> y (c, a) (c, b)
5      (***) :: y a c -> y b d -> y (a, b) (c, d)
6      (&&&) :: y a b -> y a c -> y a (b, c)
```

Listing 1: Arrow class in Haskell



Figure 2.1: The visual representations of arrow combinators[7]

```

1  instance Arrow (->) where
2    arr f = f
3    (***) f g ~ (x,y) = (f x, g y)

```

Listing 2: (\rightarrow) instance of Arrow class

2.1.2 Example: Calculate the mean

Consider the a function to calculate the mean from a list of floating number, we will compare the usual, arrows implementations. Implementation using arrows can be regarded as point-free programming. Point-free programming is programming paradigm where function definitions only involve combinators and function composition without mentioning variables[8].

```

1  mean :: [Float] -> Float
2  mean xs = sum xs / (fromIntegral . length) xs
3
4  mean' :: [Float] -> Float
5  mean' = (sum &&& (length >>> fromIntegral)) >>> uncurry (/)

```

The arrows implementation can be visualised in Figure 2.2.

```

1  mean'' :: [Float] -> Float
2  mean'' = liftM2 (/) sum (fromIntegral . length)

```

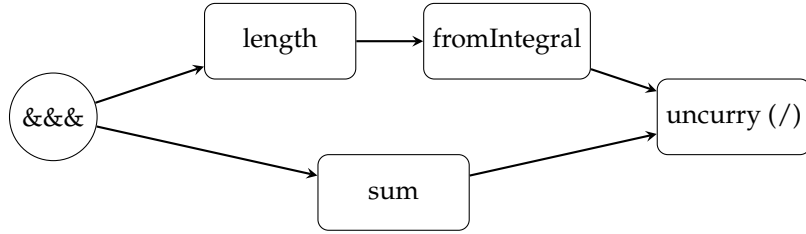


Figure 2.2: Visualization of mean'

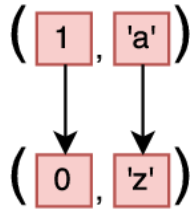
Arrows are not the only way to form point-free programs. The above code snippet is the more traditional approach form of point-free mean function in Haskell. We can argue this form of point-free function is more difficult to understand compared to arrows because it involves knowledge of monads (`liftM2`) and does not map to intuitive data-flow.

The simple example demos that arrows combinators make writing point-free programs easier. Arrows union the implementation of algorithm and data-flow in the algorithm.

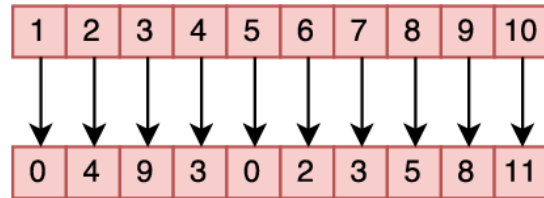
2.1.3 Application in parallel computation

From the previous example, the data flow of programs written regarding arrow combinators can be easily visualised (shown in Figure 2.1). It is intuitive to recognise that the clean separation between the flow of data and actual computation will be useful in generating parallel code. Indeed, arrow describes data flow implicitly, and it is an example of the so-called algebraic pattern. Many works [9, 10, 11] has been done to generate parallel code from algebraic patterns. In particular, details of [11] are introduced in later section.

We will use some figures to explain the idea behind arrows as a framework for parallel computation. For example, as shown in the Figure 2.3a, `f *** g` means computations of `f` and `g` happened in parallel. Figure 2.3b shows that it can be extending to parallel map in terms of arrows, taking an arrow computation `arr ab` and returning a list of computation in parallel (`arr [a] [b]`).



(a) Visualization of parallel `***`[9]



(b) Visualization of `parMap` [9]

2.2 Recursion Schemes

Recursion schemes are patterns for expressing general computation. In particular, they are like high order function abstracting recursion so that programmer can express any kind of recursion by data structures combined with recursion schemes instead of writing explicit recursive functions.

2.2.1 Definition

We will introduce three typical recursion schemes: catamorphisms, anamorphisms and hylomorphisms (seen in Listing 4). As mentioned before, recursion schemes express recursion with the help of data structures, in particular, the fixed point of data structures (seen in Listing 3)

```

1 newtype Fix f = Fix { unfix :: f (Fix f) }
2
3 data TreeF a =
4     Node a a
  
```



```

5 | Leaf int
6 | Empty
7 deriving Functor
8
9 type Tree = Fix TreeF

```

Listing 3: Definition of fix point of data structures

Anamorphisms takes a function from a to $f\ a$ (called the co-algebra) and a value a and return the $\text{Fix}\ f$. Used `Tree` as an example, anamorphisms takes a single value a and applies the co-algebra to the value. It continues to apply itself to the branches of the `TreeF` recursively and finally expands a single value to a complete tree. Intuitively, anamorphism unfolds a single value to a complicated data structure top-down.

Catamorphisms is the reverse of anamorphisms, folding a data structure to a single value bottom-up. It takes a function from $f\ a$ to a (called the algebra) and $\text{Fix}\ f$ to fold and return a single value a . Catamorphisms and anamorphisms describe the process globally (from a to $\text{Fix}\ f$ and from $\text{Fix}\ f$ to a) while co-algebra and algebra capture what happened locally. The elegant part is while co-algebra and algebra do not involve with any recursion data structure (`TreeF` is not recursive), catamorphisms consumes recursive data structure while anamorphism builds them.

Hylomorphisms applies anamorphism followed by catamorphisms. It is the most common pattern to use. We will use an example to illustrate its usefulness. It can be thought of as an abstract divide and conquer algorithm.

```

1 ana :: Functor f => (a -> f a) -> a -> Fix f
2 ana coalg = Fix . fmap (ana coalg) . coalg
3
4 cata :: Functor f => (f a -> a) -> Fix f -> a
5 cata alg = alg . fmap (cata alg) . unfix
6
7 hyllo :: (f b -> b) -> (a -> f a) -> b -> a
8 hyllo g f = f . fmap (hyllo f g) . g

```

Listing 4: Recursion schemes in haskell

2.2.2 Example: Merge sort

We can write merge sort recursively. First of all, we split the list in half and then apply the merge sort recursively to both parts and finally we merge two lists into a single list.

To write merge sort in terms of recursion scheme, we need to define the recursive structure to represent the control structure. By the definition of merge sort, this structure must have a case with two branches, a base case representing a singleton list and a base case representing an empty list hence this structure is the `TreeF` we defined above. Splitting a list is like co-algebra while merging is like algebra. We use hylomorphisms to combine them hence getting a sorted list (seen in Listing 5).

```

1 mergeSort :: [Int] -> [Int]
2 mergeSort = hyllo merge split where
3   merge Empty      = []
4   merge (Leaf c)   = [c]
5   merge (Node l r) = usualMerge l r
6
7   split [] = Empty
8   split [x] = Leaf x
9   split xs = Node l r where
10     (l, r) = splitAt (length xs `div` 2) xs

```

Listing 5: Merge sort using hylomorphisms

2.3 Multiparty session types

In complicated distributed systems, participants agree on a protocol, specifying type and direction of data exchanged. Multiparty session types are a branch of behavioural types specifically targeted at describing protocols in distributed systems based on asynchronous communication [5]. They are a type formalism used to model communication-based programming by codifying the structure of communication. The evolution of computing from the era of data processing to the era of communication witnessed the growth and significance of the theory of session types.

The theory of multiparty session type contains three main elements. Global types (seen in Section 2.3.1), local (session) types and processes. Processes are the concrete descriptions of the behaviour of the peers involved in the distributed system [5] using a formal language. Usually, the most used and the original language is π -calculus [12]. However, for the simplicity, we will not introduce π -calculus. The coming sections are an intuitive introduction of session types by examples.

2.3.1 Global types and local types

Global type is at the most abstract level, describing a communication protocol from a neutral viewpoint between two or more participants[5]. The syntax of the global types is shown in Table 2.1 and an example of global types is shown in Table 2.3.

Local types or session types characterise the same communication protocol as the global type, but from the viewpoint of each peer [5]. Each process is typed by local type. The syntax of local types is shown in Table 2.2 and an example of local type is shown in Table 2.4.

The relationship between global types and local types are established by the projection operator (seen in the Section 2.3.1.1), and a type system performs syntactic checks, ensuring that processes are typed by their corresponding local types. Hence, at the compile time, three important properties follow [5].

- **communication safety:** Mismatches between the types of sent and expected messages, despite the same communication channel is used for exchanging messages of different types, do not exist [5].
- **protocol fidelity:** The interactions that occur are accounted for by the global type and therefore are allowed by the protocol [5].
- **progress:** Every message sent is eventually received, and every process waiting for a message eventually receives one [5].

We will learn that these properties are valuable not only in the distributed system but also in the domain of parallel computing in Section 2.3.2.

$G ::=$	Global types
$p \rightarrow q : \langle S \rangle . G$	Value exchange
$p \rightarrow q : \langle T \rangle . G$	Channel exchange
$p \rightarrow q : \{l_i : G_i\}_{i \in I}$	Branching
$\mu t . G \mid t \mid \text{end}$	Recursion/End

Table 2.1: Global types

$S ::=$	Sorts	$T ::=$	Session types/local types
$\text{bool} \mid \text{nat} \mid \text{string}$		$! \langle p, S \rangle . T$	Send value
\dots		$! \langle p, T \rangle . T$	Send channel
		$? \langle p, T \rangle . T$	Channel Receive
		$? \langle p, S \rangle . T$	Sorts Receive
		$\oplus \langle p, \{l_i : T_i\}_{i \in I} \rangle$	Selection
		$\& \langle p, \{l_i : T_i\}_{i \in I} \rangle$	Branching
		$\mu t . T \mid t \mid \text{end}$	Recursion/End

Table 2.2: Session types/local types

1. Customer(0) sends an order number to Agency(1), and Agency sends back a quote to the customer.
2. If Customer is happy with the price then Customer selects accept option and notifies Agency.
3. If Customer thinks the price is too high then Customer terminate the trade by selecting reject.
4. If accept is selected, Agency notify both Customer and Agency2(2).
5. Customer sends an address to Agency2 and Agency2 sends back a delivery date.

$$\begin{aligned}
G = & \\
& 0 \rightarrow 1 : \langle \text{string} \rangle. \\
& 1 \rightarrow 0 : \langle \text{int} \rangle. \\
& 0 \rightarrow 1 : \{ \text{accept} : \\
& \quad 1 \rightarrow \{0, 2\} : \langle \text{string} \rangle. \\
& \quad 0 \rightarrow 2 : \langle \text{string} \rangle. \\
& \quad 2 \rightarrow 0 : \langle \text{int} \rangle. \text{end}, \\
& \quad \text{reject} : \text{end} \}
\end{aligned}$$

Table 2.3: An example of a protocol described by global types G

$$\begin{aligned}
S &\triangleq \mu t. (\& \{ \text{balance} : ![\text{nat}]; t, \\
& \quad \text{deposit} : ?[\text{nat}]; ![\text{nat}]; t, \\
& \quad \text{exit} : \text{end} \}) \\
C &\triangleq \oplus \{ \text{balance} : ?[\text{nat}]; \text{end}, \\
& \quad \text{deposit} : ![\text{nat}]; ?[\text{nat}]; \text{end} \}
\end{aligned}$$

Table 2.4: Session types of client and server end point of a ATM service

2.3.1.1 Projection between global types and local types

Projection is the formalisation of the relationship between global and local types. It is an operation extracting the local type of each peer from the global type [5]. The definition of projection is shown in Table 2.5.

As an example, a projection of global type in Table 2.3 is

$$G \upharpoonright 0 = !\langle 1, \text{string} \rangle; ?\langle 1, \text{string} \rangle; \&(1, \{ \text{accept} : ?\langle 1, \text{string} \rangle; !\langle 2, \text{string} \rangle; ?\langle 2, \text{int} \rangle, \text{reject} : \text{end} \})$$

$$\begin{aligned}
(\mathbf{p} \rightarrow \mathbf{p}' : \langle U \rangle. G') \upharpoonright \mathbf{q} &= \begin{cases} !\langle \mathbf{p}', U \rangle. (G' \upharpoonright \mathbf{q}) & \text{if } \mathbf{q} = \mathbf{p}, \\ ?\langle \mathbf{p}, U \rangle. (G' \upharpoonright \mathbf{q}) & \text{if } \mathbf{q} = \mathbf{p}', \\ G' \upharpoonright \mathbf{q} & \text{otherwise.} \end{cases} \\
(\mathbf{p} \rightarrow \mathbf{p}' : \{ l_i : G_i \}_{i \in I}) \upharpoonright \mathbf{q} &= \begin{cases} \oplus \langle \mathbf{p}', \{ l_i : T_i \}_{i \in I} \rangle & \text{if } \mathbf{q} = \mathbf{p} \\ \&(\mathbf{p}, \{ l_i : G_i \upharpoonright \mathbf{q} \}_{i \in I}) & \text{if } \mathbf{q} = \mathbf{p}' \\ G_{i_0} \upharpoonright \mathbf{q} & \text{where } i_0 \in I \text{ if } \mathbf{q} \neq \mathbf{p}, \mathbf{q} \neq \mathbf{p}' \\ & \text{and } G_i \upharpoonright \mathbf{q} = G_j \upharpoonright \mathbf{q} \text{ for all } i, j \in I. \end{cases} \\
(\mu t. G) \upharpoonright \mathbf{q} &= \begin{cases} \mu t. (G \upharpoonright \mathbf{q}) & \text{if } G \upharpoonright \mathbf{q} \neq t, \\ \text{end} & \text{otherwise.} \end{cases} \quad t \upharpoonright \mathbf{q} = t \quad \text{end} \upharpoonright \mathbf{q} = \text{end}.
\end{aligned}$$

Table 2.5: The definition of projection of a global type G onto a participants \mathbf{q} [5]

2.3.1.2 Duality of session types

In binary session types where all protocols are pairwise, duality formalises the relationship between the types of opposite endpoints. For a type T , its dual or co type, written \bar{T} is defined inductively as in Table 2.6.

$$\begin{array}{llll}
\overline{?[S]; T} = \overline{![S]; T} & \overline{\oplus\{l_i : T_i\}_{i \in I}} = \overline{\&\{l_i : T_i\}_{i \in I}} & \overline{?[T]; T'} = \overline{![T]; T'} & \\
\overline{![S]; T} = \overline{?[S]; T} & \overline{\&\{l_i : T_i\}_{i \in I}} = \overline{\oplus\{l_i : T_i\}_{i \in I}} & \overline{![T]; T'} = \overline{?[T]; T'} & \\
\overline{\text{end}} = \text{end} & \overline{\mu t. T} = \mu t. \overline{T} & \overline{\bar{t}} = \bar{t} &
\end{array}$$

Table 2.6: Inductive definition of duality

Duality is essential for checking type compatibility. Compatible types mean that each common channel k is associated with complementary behaviour: this ensures that the interactions on k run without errors.

In order to apply duality into multiparty session types in which more than two participants are allowed, the partial projection operation (seen in [5]) from multiparty session type to binary session type was introduced to allow reusing the definition of duality after applying the partial projection.

2.3.2 Applications in parallel computing

Multiparty session types not only have rich applications in distributed systems but also value in the domain of parallel computation.

Existing work[13] has shown how to generate MPI¹ programs using session types. Users describe the communication topology as a skeleton using a protocol language which is type checked by session types. After that, an MPI program is generated by merging the skeleton and user-provided kernels for each peer. The parallel code obtained in this way is guaranteed to be deadlock-free and progressing.

2.4 Message passing concurrency

This section introduces some interfaces for message passing concurrency from the primitive case: channel to more advanced one: monad for message passing concurrency.

For simplicity, they are represented in Haskell, but in general, most languages can implement similar interfaces.

2.4.1 Primitives for message-passing concurrency

In Section 2.3, channels are bi-directional and used for communication between two parties. In Haskell, channel primitives are represented in Listing 6. However, just using these primitives cannot guarantee progress or communication safety. For example, a program that has one thread writing channel once combined with another thread reading channel twice is type-correct but will cause deadlock. Many kinds of research to encode MPST using Haskell's type system are presented in [14] so that an (MPST) type-correct Haskell program assures progress, communication safety and session fidelity.

```

1 data Chan a
2 newChan :: IO (Chan a)
3 writeChan :: Chan a -> a -> IO ()
4 readChan :: Chan a -> IO a
5 dupChan :: Chan a -> IO (Chan a)

```

Listing 6: Channel primitives in Haskell

2.4.2 Concurrency Monads

The work done by [15] constructs a monad to express concurrent computation. The definition is in Listing 7. `Action` is the algebraic datatype representing basic concurrency primitives. `Atom`,

¹Message Passing Interface (MPI) is a standardised and portable message-passing standard designed by a group of researchers from academia and industry to function on a wide variety of parallel computing architectures [4].

the atomic unit of computation, is a computation (wrapped in the `IO` monad) followed by an action. `Fork` is two parallel action. `Stop` is the termination of an action. type `C` is a special case of the continuation monad. The continuation monad is an encapsulation of computations in continuation-passing style (CPS)². So `C a` is a CPS computation that produces an intermediate result of type `a` within a CPS computation whose final result type is `Action`. With the help of the monad `C`, sequencing and composing actions can use monadic `bind`.

```

1 data Action =
2     Atom (IO Action)
3   | Fork Action Action
4   | Stop
5
6 newtype C a = C { runC :: (a -> Action) -> Action } ⑥
7
8 instance Monad C where
9     (>=) :: C a -> (a -> C b) -> C b ⑨
10    m >= f = C $ \k -> runC m (\v -> runC (f v) k)
11    return :: a -> C a
12    return x = C $ \k -> k x

```

Listing 7: The definition of concurrency monad

The idea is using continuation to represent the "future" so that computation can pause and resume as well as expressing sequential computation. `Atom` wraps the actual computation and `Fork` is responsible for spawning threads. In addition, in order to write programmes in a monadic way easier, some helper functions are defined (shown in Listing 8). `atom` lifts an `IO` computation to `C`. And `fork` takes a computation in `C` and return a `C` which involves the `Fork` action. Given a `C a`, `action` gives the result of running the CPS computation. We use `_.` `Stop` to represent the final continuation (`Stop` action is the last action).

```

1 atom :: IO a -> C a
2 atom m = C $ \k -> Atom $ do
3     r <- m
4     return $ k r
5
6 fork :: C () -> C ()
7 fork m = C $ \k -> Fork (runC m (const Stop)) (k ())
8
9 action :: C a -> Action
10 action m = m (\_.
```

Listing 8: Helper functions

An example of programme written in the concurrency monad is shown below.

```

1 example :: C ()
2 example = do
3     atom $ putStrLn "Hello"
4     name <- atom getLine
5     fork $ atom $ putStrLn "World"
6     atom $ putStrLn name

```

We can easily define a round-robin scheduler for programmes in this monad. We can regard a list of action as a queue of threads that are running concurrently. `schedule` will pattern match on the head of the list. If it is `Atom` then the scheduler will run the computation (seen a `<- ioa` at ⑦) and pause its remaining computation and put it at the end of the thread queue (seen at ⑧). If it is `Fork` then the scheduler will spawn the thread and put the new thread and the current thread to the bottom of the queue (seen at ⑨). Finally, If it is `Stop` then it means this thread has finished and the scheduler will resume with the rest of threads in the queue. For example, to run the above example, we call `schedule [action example]`.

²In continuation-passing style function result is not returned, but instead is passed to another function, received as a parameter (continuation)[16]

```

1 schedule :: [Action] -> IO ()
2 schedule [] = return ()
3 schedule (a:as) = sched as a
4
5 sched :: [Action] -> Action -> IO ()
6 sched as (Atom ioa) = do
7     a <- ioa ⑦
8     schedule $ as ++ [a] ⑧
9 sched as (Fork a1 a2) = schedule $ as ++ [a2, a1] ⑨
10 sched as Stop = schedule as

```

The concurrency monad can be extended to support many features. For example, work done by [17] modifies the definition of Action as well as implements a work-stealing parallel scheduler (seen in Listing 9) to build a monad for parallel computation.

Besides, extending the concurrency monad to monad for message-passing concurrency can be done by adding channel primitives like newChan, writeChan and readChan into the Action. Since channel primitives are possible to represent in this monad, we naturally think of its prospect in connecting with MPST (will be discussed in the later section).

```

1 newtype IVar a = IVar (IORef (IVarContents a)) ①
2 data IVarContents a = Full a | Blocked [a -> Action.]
3
4 data Action . =
5     Fork Action Action
6     | Stop
7     | forall a . Get (IVar a) (a -> Action) ⑦
8     | forall a . Put (IVar a) a Action ⑧
9     | forall a . New (IVar a -> Action)

```

Listing 9: Par Monad

- ① Parent threads and child threads communicate data via IVar
- ⑦ Get operation blocks when the underlying IVarContents is Blocked
- ⑧ Put operation updates the underlying IVarContents to Full with the result a and resume the list of blocking threads by applying a to the continuation.

In summary, many techniques and ideas like continuation presented in the implementation of this monad afford us inspirations in designing our intermediate language.

2.5 Free monad

Free monad[18] is a concept from category theory. Intuitively, a free monad as a programming abstraction is a technique for implementing EDSLs, where a functor represents basic actions of the EDSL and the free monad of this Functor provides a way to sequence and compose actions. Speaking of the advantages, we are particularly interested in its benefits in flexible interpretations which will be illustrated by an example (Section 2.5.2) and discussed further (Section 2.5.3).

2.5.1 Definition

In practice, a free monad in Haskell can be defined as an algebraic data type(ADT) (shown in Listing 10). `Free f` is the monad produced given a functor `f`. `Free` has two type constructors: `Pure` and `Free`. `Monad (Free f)` is the Haskell implementation of the `Monad` interface for `Free f`. Many useful helper functions are derived from the simple definition of the free monad (shown in Listing 11). `liftF` lift the functor to its free monad representations. `freeM` maps a natural transformation of functor (`f a -> g a`) to the natural transformation of their free monad versions. Given `m` is a monad, `freeM` is a special case of interpreting `Free m a`: to the `m` monad itself. Finally, `interpret` shows the power of free monad. We can interpret the free monad version of a functor `f` to any monad `m` given a natural transformation from `f` to `m`.

```

1 data Free f a
2   = Pure a
3   | Free f (Free f a)
4
5 instance Functor f => Monad (Free f) where
6   return = pure
7   (Pure x) >>= fab = fab x
8   (Free fx) >>= fab = Free $ fmap (>>= fab) fx

```

Listing 10: Free monad in Haskell

```

1 liftF :: Functor f => f a -> Free f a
2 liftF = Free . fmap Pure
3
4 freeM :: (Functor f, Functor g) => (f a -> g a) -> (Free f a) -> (Free g a)
5 freeM phi (Pure x) = Pure x
6 freeM phi (Free fa) = Free $ phi (fmap (freeM phi) fa)
7
8 monad :: Monad m => Free m a -> m a
9 monad (Pure x) = pure x
10 monad (Free mfx) = mfx >>= monad
11
12 interpret :: (Functor f, Monad m) => (f a -> m a) -> (Free f a -> m a)
13 interpret phi = monad . freeM phi

```

Listing 11: Helper functions based on free monad

2.5.2 Example

Free monad is useful in interpreting an abstract syntax tree (AST). In order to apply free monad to a given AST, we can follow a routine [18].

1. Create an AST, usually represented as an ADT
2. Implement functor for the ADT
3. Create helper constructors to Free ADT for each type constructor in ADT by liftF
4. Write a monadic program using helper constructors. It is essentially a program written in DSL operations.
5. Build interpreters for Free ADT by interpreting
6. Interpret the program by the interpreter.

We will demo the above procedure by a made-up example. We would like to build a simple EDSL for getting customers' name and greeting customers. First of all, we build a functor `GreetingF` to represent the basic operations: getting the name and greeting. Then we wrap the functor with `Free` constructor so that a program written in our EDSL can be regarded as a Haskell expression with type `Free GreetingF a`.

```

1 data GreetingF next
2   = Getname (String -> next)
3   | Greet String next
4   deriving Functor
5
6 type Greeting = Free GreetingF

```

Then we create helper functions of `Greeting` using `liftF`.

```

1 getName = liftF $ Getname id
2 greet str = liftF $ Greet str ()

```

Then we can write a simple program using operations provided by Greeting.

```
1 exampleProgram :: Greeting ()
2 exampleProgram = do
3     a <- getName
4     greet a
5     b <- getName
6     greet b
```

Then we can easily implement an interpreter for the example program

```
1 goodMorningInterpreter :: Greeting a -> IO a
2 goodMorningInterpreter = interpret helper
3     where
4         helper (Getname next) = fmap next getLine
5         helper (Greet str next) = putStrLn ("Good morning " ++ str) >> return next
```

Finally, execute the program.

```
ghci:> goodMorningInterpreter examplePrograe
Tom
Good morning Tom
Mary
Good morning Mary
```

2.5.3 Applications

As illustrated by the example (Section 2.5.2), free monad decouple the abstract syntax tree of domain specific language (DSL) and the interpreter. Interpreters with different purposes can be implemented without changing the syntax.

In the project, we apply free monad to the intermediate language so not only we make the languages monadic for free but also benefits from decoupling the interpreter and the syntax to implement different interpreters, e.g. Simulator, code generators to different platforms easily.

Chapter 3

Alg: Algebraic Functional Language

Algebraic Functional Language (Alg) is an example of the high-level languages to generate parallel code, proposed in the paper [11]. The work done by [11] also proposes a method to do the code generation. Part of this project is about implementing an alternative code generation backend for the PAL. In the evaluation section, we will compare the speed of generated code of our method against the original method. We will give an overview of the language in this section.

3.1 Alg

3.1.1 Syntax

$F_1, F_2 ::=$		$t_1, t_2 ::=$	Type
I	Identity functor	$() \mid \text{int} \mid \dots$	Primitive types
Kt	Constant functor	$a \rightarrow b$	Function types
$F_1 + F_2$	Sum functor	$a + b$	Sum types
$F_1 \times F_2$	Product functor	$a \times b$	Product types
		$F t_1$	Functor types
		$\mu.F$	Recursive types

$e_1, e_2 ::=$		Expression
$f \mid v \mid \text{const } e \mid e_1 \circ e_2 \mid \pi_i \mid e_1 \Delta e_2 \mid e_1 \nabla e_2 \mid l_i \mid F e \mid \text{in}_F \mid \text{out}_F \mid \text{rec}_F e_1 e_2$		

Table 3.1: Syntax of Alg language

```
newtype L = K () + K Int * I
type List = Rec L
```

Listing 12: Type of integer list in PAL

The syntax of Alg is shown in Table 3.1. In terms of the syntax of expressions, f represents atomic functions which are functions of which we only know their types [11]. The presence of atomic function is important in the code generation which we will discuss in the later chapter. v is the primitive value like integer 1. F represents functor and a, b are types. Besides primitives type, function types and recursive type constructor μ , Alg use four functors to form more types hence representing complicated data structures by composing them. For example, a list of integer is expressed in Listing 12. One thing we should learn from the syntax is that the Alg and its combinators avoid mentioning variables explicitly and hence Alg programs are known as point-free programs.

The typing rule for alg is expressed in Table 3.2 and the semantic for alg is shown in Table 3.3.

An important feature of Alg is the lack of usual control flow like if branch or while loop to build algorithms, instead, it uses the flow of transformation of data structures to replace conventional control flow. For example, ∇ combinator is the case operation whose types is $a + b \rightarrow c$.

$\frac{f : a \rightarrow b \in \Gamma}{\vdash f : a \rightarrow b}$
$\frac{\vdash e : a}{\vdash \text{cons } e : b \rightarrow a}$
$\frac{}{\vdash \text{id} : a \rightarrow a}$
$\frac{}{\vdash \text{in}_F : F \mu F \rightarrow \mu F}$
$\frac{}{\vdash \text{out}_F : \mu F \rightarrow F \mu F}$
$\frac{\vdash e_1 : b \rightarrow c, \quad \vdash e_2 : a \rightarrow b}{e_1 \circ e_2 : a \rightarrow b}$
$\frac{i \in [1, 2]}{\pi_i : a_1 \times a_2 \rightarrow a_i}$
$\frac{i \in [1, 2]}{l_i : a_i \rightarrow a_1 + a_2}$
$\frac{\vdash e_1 : a \rightarrow b, \quad \vdash e_2 : a \rightarrow b}{e_1 \Delta e_2 : a \rightarrow b \times c}$
$\frac{\vdash e_1 : a \rightarrow c, \quad \vdash e_2 : b \rightarrow c}{e_1 \nabla e_2 : a + b \rightarrow c}$
$\frac{\vdash e : a \rightarrow b}{F e : F a \rightarrow F b}$
$\frac{\vdash e_1 : F b \rightarrow b, \quad \vdash e_2 : a \rightarrow F a}{\text{rec}_F e_1 e_2 : a \rightarrow b}$

Table 3.2: Typing rules for Alg

It can be seen as analogy of if branch in normal programming languages. Δ combinator represents split operation. More importantly, the combinator rec_F uses the idea of recursion schemes (explained in Section 2.2) to build divided-and-conquer algorithms.

To summary, algorithms in Alg are represented as a series of transformations of data, making it easy to transform the Alg programs to programs in arrows mechanically since arrows also express flow of data and their transformations naturally. This property allows us to generate parallel code without burdens.

3.1.2 Example: Merge sort

$$\begin{aligned}
\text{ms} &= \text{rec}_T \text{mrg spl} = \text{mrg} \circ T (\text{rec}_T \text{mrg spl}) \circ \text{spl} \\
&= \text{mrg} \circ (\text{id} + \text{id} + (\text{rec}_T \text{mrg spl}) \times (\text{rec}_T \text{mrg spl})) \circ \text{spl} \\
&= \text{mrg} \circ (\text{id} + \text{id} + \text{ms} \times \text{ms}) \circ \text{spl}
\end{aligned}$$

Listing 13: Merge sort in Alg

By making the hylomorphism as a built-in combinator rec_F , Alg can express merge sort similarly as the example shown in Section 2.2.2. We use the functor $T = K() + Ka + I \times I$ to be substitute the functor F in rec_F and the type Ls to represent a list of elements whose type is a . We also have two atomic function $\text{spl} : Ls \rightarrow T Ls$ and $\text{mrg} : T Ls \rightarrow Ls$. Finally we express merge sort as $\text{ms} = \text{rec}_T \text{mrg spl}$. It is shown in Listing 13. From the example, we observe that ms can expanded infinitely. Later, we will exploit this property to generate parallel code.

Constant, Identity and Composition

Products

$\pi_i = \lambda(x_1, x_2). x_i \text{ if } i \in [1, 2]$

Coproducts

$\iota_i = \lambda x. \text{inj}_i x$

Functors

$\mathsf{I} a = a$
 $\mathsf{I} e = e$

Recursion

$\text{const } e = \lambda x. e \quad \text{id} = \lambda x. x \quad e_1 \circ e_2 = \lambda x. e_1 (e_2 x)$

$e_1 \triangle e_2 = \lambda x. (e_1 x, e_2 x) \quad e_1 \times e_2 = (e_1 \circ \pi_1) \triangle (e_2 \circ \pi_2)$

$e_1 \nabla e_2 = \lambda x. e_i y \text{ if } x = \text{inj}_i y \quad e_1 + e_2 = (\iota_1 \circ e_1) \nabla (\iota_2 \circ e_2)$

$(F_1 \times F_2) a = F_1 a \times F_2 a \quad (F_1 + F_2) a = F_1 a + F_2 a$
 $(F_1 \times F_2) e = F_1 e \times F_2 e \quad (F_1 + F_2) e = F_1 e + F_2 e$

$\text{in}_F = \lambda x. \text{in}_F x \quad \text{out}_F = \lambda x. y \text{ where } \text{in}_F y = x \quad \text{rec}_F e_1 e_2 = f \text{ where } f = e_1 \circ F f \circ e_2$

Table 3.3: Semantics of Alg expression

3.2 ParAlg: Alg + role annotations

3.2.1 Inferring global types

3.2.2 Example: Parallel merge sort

3.3 Conclusion

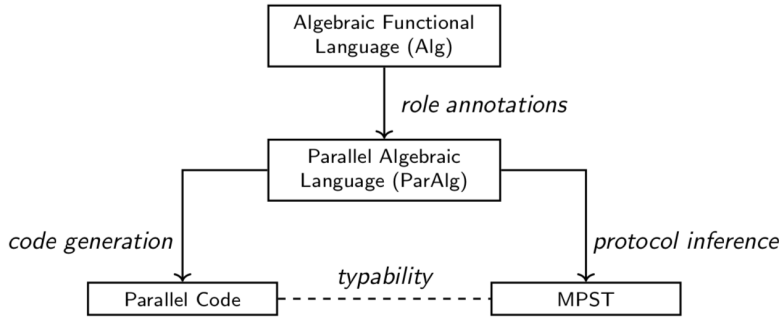


Figure 3.1: Overview of code generating pipeline from Alg[11]

A visualization of the pipeline can be seen in the Figure 3.1. The content of the section is a skim of Alg and ParAlg. Curious readers should read about [11] for more details.

Combinators and the syntax of Alg makes every Alg program to be point-free. Without the use of explicit variables, point-free programs express the underlying data-flow of the computation clearly. Adding role annotations transforms Alg programs to ParAlg programs converting the implicit data-flow to explicit communication. Communication will aid us generate parallel code using message-passing concurrency. We will introduce the our method from next chapter.

Chapter 4

SPar: A session typed free monad EDSL for concurrency

In order to generate parallel code from ParAlg, we first introduce the syntax of our intermediate language, the session-typed free monad EDSL for concurrency hosted in Haskell (SPar). SPar are compromised of two components: Core and Proc. Core is the language expressing sequential computation while Proc is a monadic language with message-passing primitives, communicating Core expression between different roles. We use a group of Proc interacting with each other to represent parallel computations. In addition, session typing the group of Proc ensure the computation is deadlock-free.

4.1 Computation: The Core EDSL

Core is the elemental computation. The syntax of Core is mostly inspired by Alg [11] and FunC, a demo DSL defined in the work done by [19]. For this project, we choose to implement Core syntax as small as possible without sacrificing expressibility.

4.1.1 Syntax

```
data Core a where
  Lit :: a -> Core a
  Var :: Int -> Core a
  Prim :: String -- XXX: name of the "primitive function"
        -> a -- Semantics in Haskell of the primitive function
        -> Core a

  Ap :: Core (a -> b) -> Core a -> Core b
  Id :: Core (a -> a)
  Const :: Core a -> Core (b -> a)

  Fst :: Core ((a, b) -> a)
  Snd :: Core ((a, b) -> b)
  Pair :: Core a -> Core b -> Core (a, b)

  Inl :: Core (a -> Either a b)
  Inr :: Core (b -> Either a b)
```

Listing 14: The syntax of Core

The syntax of Core is shown in Listing 14. Inl and Inr are for the construction of sum type. Pair is responsible for constructing value of product type while Fst and Snd extract value from product type. Supporting sum type, product type and inductive type (see next section) is enough

to express any data structure in any computation. In addition to these actions manipulating basic data structure, we have `Lit` which is used in the Haskell interpreter (see Section 5.2) and benchmarking (see Section 8) and `Var`, a constructor which is useful when what we do not actually evaluation the Core expression but inspect its static structure. It is used in the code generation (see Section 7.3.2) and session typing (see Section 5.1). `Id` is similar to identity function and `Const` is similar to the `const` function in Haskell. `Prim` representing user defined functions takes two field name and the haskell implementation. The first field will be useful in the code generation (see Section 7.3.2) when applying user-defined function calls. The second field will be used in the interpreter implementation.

4.1.2 Representation of recursive data structures

Core has primitives to operate on sum and product types. Representing recursive type like $\mu \text{list}(). + \text{Int} \times \text{list}$ will be covered in this section. The method are taken from the implementation of Alg language in [11]. First of all, we extend the core with the following two operation. `In` represents the fold operation on iso-recursive types and `Out` represents the unfold operation on iso-recursive types. Consider a recursive type $\mu \alpha. \tau$, the type parameter τ is equivalent to α , $f : @ : \tau$ is equivalent to τ and the typeclass `Data f t` associate two types which is equivalent to $\mu \alpha. \tau$. $:@ :$ is a type family which is a function acting on types instead of values. $:@ :$ converts f to sum and product type in Haskell which are `(,)` and `Either`.

```
data Core a where
  -- omit other constructors
  In  :: Data f t => Core (f :@: t -> t)
  Out :: Data f t => Core (t -> f :@: t)

type family (@:) (a :: Poly Type) (b :: Type) :: Type where
  'PId :@: x = x
  'PK y :@: _ = y
  'PProd f g :@: x = (f :@: x, g :@: x)
  'PSum f g :@: x = Either (f :@: x) (g :@: x)

class Data (f :: Poly Type) t | t -> f where
  roll  :: f :@: t -> t
  unroll :: t -> f :@: t
```

A concrete example is a list shown below. We know a list has recursive type: $\mu \alpha. () + a \times \alpha$. So the `f` is `('PSum ('PK ()) ('PProd ('PK a) 'PId))` and we use haskell list type `[a]` to present α (equivalent to `t` in `f :@: t`). `f :@: t` is evaluated to the type `Either () (a, [a])`.

```
instance Data ('PSum ('PK ()) ('PProd ('PK a) 'PId)) [a] where
  roll (Left _) = []
  roll (Right (a, b)) = a : b

  unroll [] = Left ()
  unroll (x:xs) = Right (x, xs)
```

However, even though this way of representing recursive types is very smart but later we found out they are seldomly used in the code generation of parallel algorithms. Because functions processing complicated recursive data structure is usually represented by the `Prim` constructor hiding the implementation details so `In` and `Out` are too low-level operations and hence user are difficult to write complicated function by them. So in the conclusion, Core provides basic functionality, i.e `Fst` or `Inl` to process data and more importantly, `Prim`, a black box constructor hiding the implementation details of the sequential computation. `Prim` turn out to be a very useful abstraction in terms of code generation. More details will be discussed in the Section 7.3.2.

4.2 Communication: The Proc EDSL

Proc will be a free monad EDSL for message passing. As introduced in the free monad section in the background, the first thing to do it to define the algebra of message-passing concurrency:

```

data ProcF next where
  Send :: Nat -> Core a -> next -> ProcF next

  Recv :: Nat -> (Core a -> next) -> ProcF next

  Select :: Nat
    -> Core (Either a b)
    -> (Core a -> Proc c)
    -> (Core b -> Proc c)
    -> next
    -> ProcF next

  Branch :: Nat -> Proc c -> Proc c -> (Core c -> next) -> ProcF next

  Broadcast :: [Nat]
    -> Core (Either a b)
    -> (Core a -> Proc c)
    -> (Core b -> Proc c)
    -> next
    -> ProcF next

type Proc a = Free ProcF (Core a)

```

Listing 15: The algebra for message-passing

`ProcF` and `Proc` is defined using free monad constructor and `ProcF`. The definition is shown in the Listing 15. Careful reader might notice that `Proc` and `ProcF` are defined mutually with each others in `Branch`, `Select` and `Broadcast`.

The semantics will be defined in the next subsection in terms of a group of `Proc` programs. A single `Proc` program is either sequential or deadlock. Operational semantics is only worth discussing when given a group of `Proc` programs interacting with each others.

4.3 Concurrent computation: A group of Proc

We have introduced syntax for computation and communication. We also know that a single `Proc` expression is meaningless since there does not exist another party to interact with this `Proc` hence the computation has no progress. Naturally, we use a group of `Proc` to represent concurrent computations. To be more precise, a collection of `Proc` with their own role identifiers can be treated as a system of roles executing their own programs concurrently. In most of the cases, in order to make the group of `Proc` meaningful, we will allocate a start role in the system acting as the original data provider and a end role whose `Proc` program will receive data from others, process and output the final computation which is wrapped by the `Pure` constructor at the end of the `Proc` program.

Readers might find it easy to visualize the group of `Proc` as a computation graph. The start role is the source node and the end role is the sink node. A pair of nodes are connected if they communicate data with each others.

4.3.1 Operational semantics

Due to the similarities between `Proc` and multiparty session calculus introduced in [5]. We borrow some syntax and operational semantics rule from multiparty session calculus to define the operational semantics of `Proc`. P, Q denote `Proc` programs. A message queue is h which contains messages (q, p, v) meaning that the sender q sends the receiver p with value v . $h \cdot m$ is a message queue whose bottom element is message m . h are runtime syntax to model the asynchronous message communication where the order of the messages are retained [5]. $e \Downarrow v$ means the evaluation of the `Core` expression to the value v .

$(P_1, r_1) \mid (P_2, r_2) \mid \dots \mid (P_n, r_n) \rightarrow$	(Init)
$(P_1, r_1) \mid (P_2, r_2) \mid \dots \mid (P_n, r_n) \mid \emptyset$	
$(\text{Free } (\text{Send } r_j \ e \ \text{next}), r_i) \mid \dots \mid h \rightarrow$	(Send)
$(\text{next}, r_i) \mid \dots \mid h \cdot (r_i, r_j, v) \quad (e \downarrow v)$	
$(\text{Free } (\text{Recv } r_i \ \text{cont}), r_j) \mid \dots \mid (r_i, r_j, v) \cdot h \rightarrow$	(Recv)
$(p, r_j) \mid \dots \mid h \quad (\text{cont } v \downarrow p)$	
$(\text{Free } (\text{Select } r_j \ e \ \text{cont1} \ \text{cont2} \ \text{next}), r_i) \mid \dots \mid h \rightarrow$	(Sel-Left)
$(\text{cont1 } v \gg \text{next}, r_i) \mid \dots \mid h \cdot (r_i, r_j, L) \quad (e \downarrow v, \text{label}(v) \downarrow L)$	
$(\text{Free } (\text{Select } r_j \ e \ \text{cont1} \ \text{cont2} \ \text{next}), r_i) \mid \dots \mid h \rightarrow$	(Sel-Right)
$(\text{cont2 } v \gg \text{next}, r_i) \mid \dots \mid h \cdot (r_i, r_j, R) \quad (e \downarrow v, \text{label}(v) \downarrow R)$	
$(\text{Free } (\text{Branch } r_i \ \text{next1} \ \text{next2} \ \text{cont}), r_j) \mid \dots \mid (r_i, r_j, L) \cdot h \rightarrow$	(Branch-Left)
$(\text{next1} \gg \text{cont}, r_j) \mid \dots \mid h$	
$(\text{Free } (\text{Branch } r_i \ \text{next1} \ \text{next2} \ \text{cont}), r_j) \mid \dots \mid (r_i, r_j, R) \cdot h \rightarrow$	(Branch-Right)
$(\text{next2} \gg \text{cont}, r_j) \mid \dots \mid h$	
$(\text{Free } (\text{Broadcast } [r_{k_1}, \dots, r_{k_n}] \ e \ \text{cont1} \ \text{cont2} \ \text{next}), r_i) \mid \dots \mid h \rightarrow$	(Broadcast-1)
$(\text{Free } (\text{Select } r_{k_1} \ v \ c \ (\text{Pure } ())), r_i) \mid \dots \mid h$	
where $(e \downarrow v, c = \text{Free } (\text{Broadcast } [r_{k_2}, \dots, r_{k_n}] \ \text{cont1} \ \text{cont2} \ \text{next}))$	
$(\text{Free } (\text{Broadcast } [r_{k_1}] \ e \ \text{cont1} \ \text{cont2} \ \text{next}), r_i) \mid \dots \mid h \rightarrow$	(Broadcast-1)
$(\text{Free } (\text{Select } r_{k_1} \ v \ \text{cont1} \ \text{cont2} \ \text{next}), r_i) \mid \dots \mid h \quad (e \downarrow v)$	
$(\text{Pure } v, r_i) \mid (P_i, r_i) \mid \dots \mid (P_j, r_j) \mid h \rightarrow (P_i, r_i) \mid \dots \mid (P_j, r_j) \mid h$	(Pure)

Table 4.1: Small step semantics for Proc

Table 4.1 shows the small step semantics for Proc. Rule (Init) describes the initialization of a group of Proc programs with an empty message queue at the beginning. Rule (Send) append the value to the message queue. Its complementary rule: (Recv) will recv the value at the top of the message queue. Rule (Branch) is also the complementary rule of the rule (Sel). You might notice that the Rule (Broadcast) is kind of special. Its rule states that Broadcast is expressed in terms of a series of Select operations. This operation broadcasts the label to a group of receivers. Even though its semantics can be expressed in terms of Select, the reason why we still treat broadcast an independent operations in SPar is because 1) this operation is very common in communication, including Broadcast as a primitive operation is beneficial for user to write code 2) In the code generation stages, we can generate more efficient code for Broadcast operation instead of generating a series of Select operations. Broadcast operation can be treated as a syntax sugar in the Proc language.

4.3.2 Session types and duality checking

Immediately, we notice that a Proc program can be typed by session types. Send operation in ProcF corresponds to the type $\langle p, S \rangle.T$, Select corresponds to $\oplus \langle p, \{l_i : T_i\}_{i \in I} \rangle$ and so on. One exception is the broadcast operation which does not mapped to any type of the session types. We will discuss how to handle Broadcast in Section 5.1.

The Table 4.2 shows the session typing rule for a Proc expression. We borrowed some notation from the work done [11]. $\Gamma \vdash e : \text{Proc } L \ a$ is the type of a Proc expression that follows protocol L and returns a value of type a . The types are parameterized by a variable l representing the continuation of a local type L . The typing rules contains all the operations except Broadcast. This is because Broadcast can be expressed in terms of a series of Select as explained in the previous subsection. So its session type $\oplus \{r_j\}_{j \in [1, n]} L : L_1, R : L_2$ can be expanded to $r_1 \oplus \{l : r_2 \oplus \{\dots \oplus \{r_n \dots\}\}, \dots\}_{l \in [L, R]}$.

We have argued that a Proc program can be typed by multi-party session types. In order to

Inst	$\frac{\Gamma \vdash e : \forall l. \text{Proc } L \ a}{\Gamma \vdash e : \text{Proc } ([\text{end}/l]L) \ a}$
Gen	$\frac{\Gamma \vdash e : \forall l. \text{Proc } L \ a, \quad \text{fresh } l}{\Gamma \vdash e : \text{Proc } ([l/\text{end}]L) \ a}$
Ret	$\frac{\Gamma \vdash v : a}{\Gamma \vdash \text{Pure } v : \forall l. \text{Proc } l \ a}$
Abs	$\frac{\Gamma, x : a \vdash e : \forall l. \text{Proc } L \ b}{\Gamma \vdash \lambda x. e : a \rightarrow \forall l. \text{Proc } L \ b}$
Bind	$\frac{\Gamma \vdash m : \forall l_1. \text{Proc } L_1 \ a, \quad \Gamma \vdash f : a \rightarrow \forall l_2. \text{Proc } L_2 \ b}{\Gamma \vdash m \gg f : \forall l_2. \text{Proc } [L_2/L_1]L_1 \ b}$
Send	$\frac{\Gamma \vdash v : a}{\Gamma \vdash \text{Free } (\text{Send } r \ v \ (\text{Pure } ())) : \forall l. \text{Proc } (r!(a).l) \ ()}$
Recv	$\frac{}{\Gamma \vdash \text{Free } (\text{Recv } r \ (\backslash v \rightarrow \text{Pure } v)) : \forall l. \text{Proc } (r?(a).l) \ a}$
Select	$\frac{\Gamma \vdash v : a + b, \quad \Gamma \vdash f_1 : a \rightarrow \forall l_1. \text{Proc } L_1 \ c, \quad \Gamma \vdash f_2 : b \rightarrow \forall l_2. \text{Proc } L_2 \ c}{\Gamma \vdash \text{Free } (\text{Select } r \ v \ f_1 \ f_2 \ (\text{Pure } ())) : \forall l. \text{Proc } r \oplus \{L : [l/l_1]L_1, R : [l/l_2]L_2\} \ ()}$
Branch	$\frac{\Gamma \vdash \text{next1} : \forall l_1. \text{Proc } L_1 \ c, \quad \Gamma \vdash \text{next2} : \forall l_2. \text{Proc } L_2 \ c}{\Gamma \vdash \text{Free } (\text{Branch } r \ \text{next1} \ \text{next2} \ (\backslash x \rightarrow \text{Pure } x)) : \forall l. \text{Proc } r \ \& \{L : [l/l_1]L_1, R : [l/l_2]L_2\} \ ()}$

Table 4.2: Typing rules for Proc expressions

utilize this property, we should check the duality of the each pair of Proc in the group. In short, duality check examines whether any pair of Proc in the system are complement with each other. If the duality properties is satisfied, the computation is guaranteed to be deadlock-free. This safety guarantee is useful and powerful in the application of parallel code generation. In this domain, SPar is considered to be the intermediate languages. Passing duality check for the intermediate representation means that as long as we preserve types and communication pattern carefully in later stages of code generation pipeline, the generated code obtained will share the same non-trivial properties: communication safety, protocol fidelity and deadlock-free.

The work done by [5] constructed the theoretic foundation of algorithms for checking dualities and we will give an overview of the implementation in the Section 5.1.

4.4 Conclusions

In this section, we have introduced our intermediate language. It is friendly to use thanks to the monadic interface. In addition, communication and computation are independent in SPar. We can make Proc to parameterized by the type represent sequential computation so users can simply use their construction for sequential computation if they found Core is limited. More importantly, our strategy for parallelism is clear now. In a nutshell, we achieve parallelism by message-passing concurrency: spawning a group of threads on a multi-core CPU where each thread executes its corresponding Proc program.

Before jumping into the code generation, we will use next chapter to give an overview of some implementation challenges related to SPar first.

Chapter 5

SPar: Implementation

5.1 Session types

Haskell does not have session type natively. Other encodings of session types in Haskell is an overkill to this project since Proc does not support all actions that can be typed by session types i.e. channel delegation. We decide to create our representation of session types in Haskell corresponding to set of actions supported by Proc. We will introduce two method for session typing the group of Proc programs followed by duality checking. One is at the value level and another is at the type level.

5.1.1 Representations of session types in Haskell

Session types belong to the family of behavior types. So the type looks very similar to the value that will be typed. We exploit this similarity to defined our session types in terms of free monad, the same method we have used in defining Proc.

```
data STypeF a next where
  S :: Nat -> a -> next -> STypeF a next
  R :: Nat -> a -> next -> STypeF a next
  B :: Nat -> SType a c -> SType a c -> next -> STypeF a next
  Se :: Nat -> SType a c -> SType a c -> next -> STypeF a next

type SType a next = Free (STypeF a) next
```

Listing 16: Session types in Haskell

The Listing 16 shows the definition in Haskell. `SType` is the session type in Haskell. It is parameterized by a type variable `a` so that the value-level session types and the type-level session types can share the same basic definition of session types. `S` is mapped to $!\langle p, S \rangle.T$. `R` is mapped to $?(p, S).T$. `B` is mapped to Branch type and `Se` is mapped to Select type.

5.1.2 Value-level duality check

For the value-level session types, the type variable `a` is instantiated with type `TypeRep`. `TypeRep` reifies types to some extent by associating type representations to types [20]. Due to session types are represented as value expressions in Haskell, session typing a Proc program is the same as writing an interpreter which can be easily done since a Proc program is based on free monad.

We traverses Proc programs converting each operation to its corresponding type in `STypeF` and convert the value to its `TypeRep`. For output actions, we recursively call the substructure to build the rest of session types. For input actions like `Recv`, we will apply the continuation with the Core value constructed by `Var` and recursively call the function on the result. The trick to apply `Var` to the continuation make it possible to inspect the static structure of every Proc programs because `b -> Proc a` is not inspectable i.e, we cannot pattern match on `i`, while `Proc a` can

be inspected. We will also use this trick in the code generation which will be introduced in later chapter.

5.1.3 Type-indexed Free Monad

5.1.4 Type-level duality check

5.2 SPar interpreter

5.2.1 Overview

SPar interpreter is an interpreter that interprets a group of Proc programs in Haskell. It can be considered as the simplest backend for evaluating SPar expressions. It records traces of the executions and the final output values of each Proc programs in the system.

It focused on providing a reference implementation explaining the semantics of SPar expressions not on performance. In addition, SPar interpreter served as a very useful tool in the development of this project. We use it as prototype to quickly verify whether the computation produces the expected results. This feature is useful especially in the early stage of an implementation or during debugging.

5.2.2 Implementation

The implementation of SPar interpreter is standard. It is similar to the implementation of the scheduler we explained in the free monad section of the background chapter. In essence, it is a round robin scheduler for a group of Proc programs. A partial implementation can be seen in

```
data InterpState = InterpState
{
    outputValues :: [(Nat, String)],
    trace :: [String],
    messageQueue :: [(Nat, Nat, String)]
}

interpret :: [(Proc (), Nat)] -> State InterpState ()
interpret [] = return ()
interpret (x : xs) = interp x xs

interp :: (Proc (), Nat) -> [(Proc (), Nat)] -> State InterpState ()
interp (Pure value, role) xs = do
    updateOutputValue role value
    interpret xs
interp (Free (Send receiver value next), role) xs = do
    updateMessageQueue (role, receiver, show $ interpCore value)
    updateTrace "send"
    interpret $ xs ++ [(next, role)]
interp p@(Free (Recv sender cont), role) xs = do
    (x, y, v) <- getTopMessage
    if x == sender && y == role
    then do
        popMessageQueue
        updateTrace "recv"
        let value = Lit (read v)
        interpret $ xs ++ [(cont value, role)]
    else
        interpret $ xs ++ [p]
```

Listing 17: Partial implementation of the SPar interpreter

Listing 17. It takes a list of Proc as a parameter and it maintain a state which is the combination of

message queue, the trace and the list of output values. For the base case, the list is empty which means all processes has exited, it returns the trace and the list of output values. For the recursive case, it pattern matches on the first process at the beginning of the list. If the operation is `Pure`, it will update the list of output values, then abandon this process, and call itself recursively on the tail of the list. If the operation is an output action i.e, `Send`, `Select` or `Broadcast`, it will update the message queue with the corresponding message containing the sender, the receiver and the value, then pop the head proc and append its next step to the end of list of Procs, and call the list recursively. If the operation is an input action i.e, `Recv` or `Branch`, it will first examine whether sender and receiver pairs match the pair from the message at the top of the message queue. If so, it applies the continuation with the value in the message, removes the message from the queue, removes the head of list, moves the result of applying continuation to the end of the list and calls the list recursively. If not, it simply moves the head of list to the end of list without changing the value of head process and call the list recursively. Because we have checked the duality of the processes in the system, it guarantee that for any input action, the required message will eventually appear at the head of the message queue in finite steps.

Also, the interpreter has a helper function that interpret Core expression into haskell value. This helper function is easy to write because operation on product or sum type has its own mapping function in Haskell and for the `Prim` and `Lit`, we can get their underlying Haskell implementation directly. As for `Var`, this constructor is not intended to externally so we will not encounter.

Chapter 6

ArrowPipe: An arrow interface for writing SPar expressions

When trying to express more complicated and interesting parallel patterns, e.g map or reduce, We realize SPar is too low-level so that it is difficult to express simple computation because of overheads of expressing communication patterns by hand. In addition, compilation from Par-Alg to SPar is hard since they are very different domain specific languages.

To solve both issues, we draw inspirations from the Arrow interface (in particular, work done by [9] where they use arrow interface to express parallel computation) and introduce ArrowPipe.

ArrowPipe is an arrow interface for writing SPar expressions. Withe the help from ArrowPipe, Users can use canonical arrow combinators to write algorithms in Arrow without writing any explicit communication and gain parallelized algorithms for free. Similarly, ArrowPipe makes hassle-free compilation from Par-Alg to SPar possible because Par-Alg Proto is also an arrow expression and simply interpreting arrow combinators by the ArrowPipe implementations fills the gap between Par-Alg and SPar.

6.1 Syntax

```
data Pipe a b = Pipe
  { start :: Nat
  , cont  :: a -> Proc
  , env   :: Map Nat Proc
  , end   :: Nat
  }

type ArrowPipe a b = Nat -> Pipe a b

instance Arrow ArrowPipe where
  -- implementation omit
instance ArrowChoice ArrowPipe where
  -- implementation omit
```

Listing 18: Definition of ArrowPipe

The simplified syntax of ArrowPipe can be found in Listing 18. ArrowPipe is a type synonym of `Nat -> Pipe a b`. It consumes `Nat` which means the identifier of a process and output `Pipe a b`. The reason why we use `Nat` as the only parameter is to ensure no duplication of processes name since in most of the time, duplication is bad for parallelization. It will be explained more thoroughly in Section 6.3.

`Pipe a b` data structure is the essential component of ArrowPipe. It regards computation as a pipe where data with type `a` goes into the pipe and data with type `b` get out of the pipe. Internally, it's a record type of four fields. `start` field identifies the process where the input data

is received. `cont` field has the type `a -> Proc` which is a continuation waiting for the input data produced by the last pipe. `env` represents a group of Procs interacting inside the pipe to produce the output data, in other words, it is the parallel computation. `end` indicates the process that produces the output data in the end. We can retrieve the corresponding process by a look up on the `env` with the key `end`. The returned Proc returns a data with type `b`.

6.1.1 Arrow interface

`ArrowPipe` is an instance of `Arrow` typeclass as well as `ArrowChoice` type class. For example, the type signature of the combinators `>>>`, `|||`, `&&&` and `arr` are shown below. The main difference between their type signatures and the usual arrow interface is that in the `arr`, the function is wrapped with `Core`. In general, it captures the same meaning as the usual arrow interfaces. Implementation details of these combinators will be explained in Section 6.2.

```
(>>>) :: (ArrowPipe a b) -> (ArrowPipe b c) -> (ArrowPipe a c)
-- (>>>) :: a b c -> a c d -> a b d
arr :: (Core (a -> b)) -> ArrowPipe a b
-- arr :: (b -> c) -> a b c
(|||) :: (ArrowPipe a c) -> (ArrowPipe b c) -> ArrowPipe (Either a b) c
-- (|||) :: a b d -> a c d -> a (Either b c) d
(&&&) :: (ArrowPipe b c) -> (ArrowPipe b c') -> ArrowPipe b (c, c')
-- (&&&) :: a b c -> a b c' -> a b (c, c')
(***) :: (ArrowPipe b c) -> (ArrowPipe b' c') -> ArrowPipe (b, b') (c, c')
-- (***) :: a b c -> a b' c' -> a (b, b') (c, c')
```

6.1.2 Example: Parallel programming patterns

As an example, we will illustrate some typical computation patterns used in parallel computing.

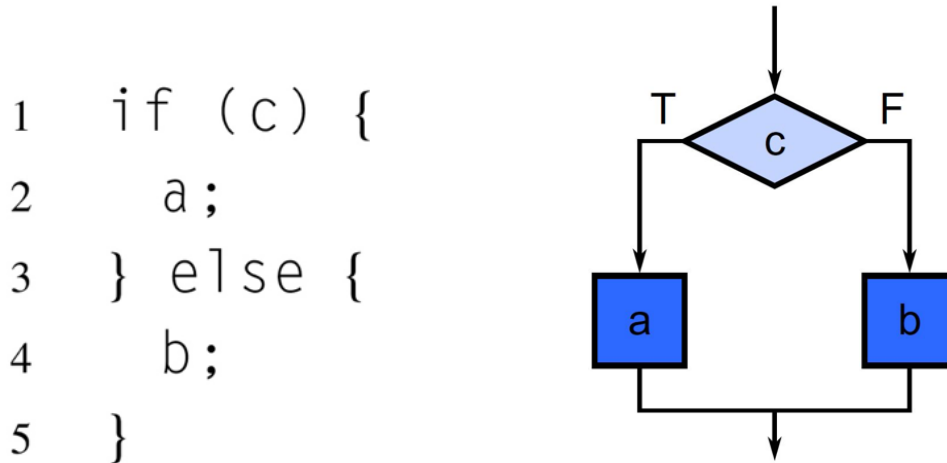
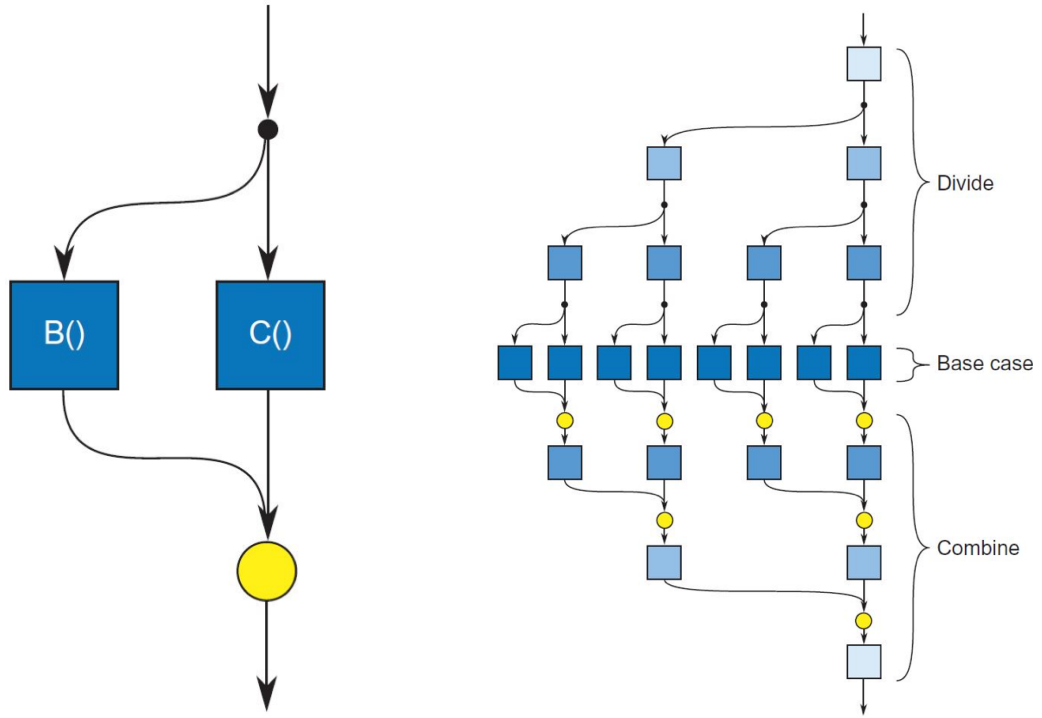


Figure 6.1: Visualization of the select pattern [21]

First of all, the select pattern illustrated by Figure 6.1 is equivalent to an expression formed by `|||` combinators, where the data constructor `Left` means True and the data constructor `Right` means False for the sum type `Either`. Secondly, the fundamental building block of parallel pattern, the fork-join pattern illustrated by Figure 6.2a can be expressed by `&&&` combinator. The arrowPipe produced by `&&&` has the two-ary tuple as the output type collecting the computation result of the main thread and the forked thread and also acts as a synchronization point.

Thirdly, the familiar parallel map pattern illustrated in Figure 6.3 is also a candidate to be expressed in `ArrowPipe`. The code sample is in Listing 19. `pmap` splits the input `a` into 4 chunks using the splitting function `s`, applied the elemental function `f` and the arrow combinator `***` in parallel and finally use the collecting function `c` to collect the results. Usually, the input `a` is a list and `s` splits the list into 4 equal chunks. The number of function `f` applied decides the number of ways of parallelism. We will describe a possible solution to make `pmap` more generic so that we do not need write a new functions for every number of ways of parallelism. (TODO).



(a) Visualization of the fork-join pattern [21]

(b) Fork-Join Pattern for Divide-Conquer [21]

Figure 6.2: Fork-join pattern and divide-and-conquer algorithms

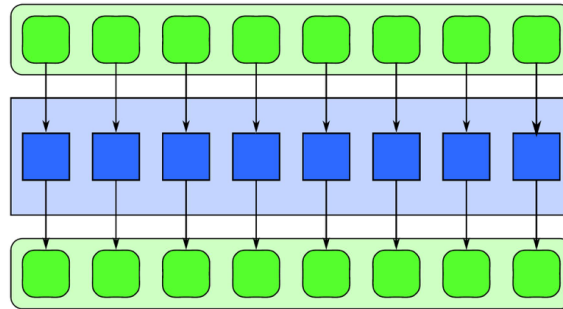


Figure 6.3: Visualization of parallel map [21]

Fourthly, we can apply the similar logic to express parallel reduce pattern shown in Figure 6.4. The code sample is in Listing 20. The result of parallel reduce has similar type signature as the collecting function in `pmap` so it is often used with the `pmap` function. We use nested tuple `(a, (a, (a, a)))` to represent a sized array of data. The `helper` function transforms the array representation of data into a form so that we apply the reduce function `r` to the elements pair-wise and parallel.

Finally, more complicated pattern can be expressed compositely from simpler pattern expressed in ArrowPipe. We can use a typical divide-and-conquer algorithms implemented with fork-join as an example. Figure 6.2b shows a divide-and-conquer algorithms with 2-ways and 3-levels of fork-join. The algorithm can be expressed in arrowPipe shown in Listing 21. The divide-and-conquer pattern can be built recursively in Haskell. For the base case, we simply apply the basic computation. Otherwise, we first call `split` and then call the function recursively with the level decremented by one and, in the end, call the `merge` to combine the results. Every expressions in the function definition are connected using arrow combinators. A 3-level divided-and-conquer algorithm is constructed by passing 3 to the function resulting a algorithm with $2^3 = 8$ -way parallelism.

The implementation demos the power of implementing ArrowPipe as a domain specific language embedded in Haskell. We make full use of Haskell features, i.e high order functions and

```

pmap :: ArrowPipe a b
-> ArrowPipe a (a, (a, (a, a)))
-> ArrowPipe (b, (b, (b, b))) b
-> ArrowPipe a b
pmap f s c = s >>> (f *** (f *** (f *** f))) >>> c

```

Listing 19: Parallel map in ArrowPipe

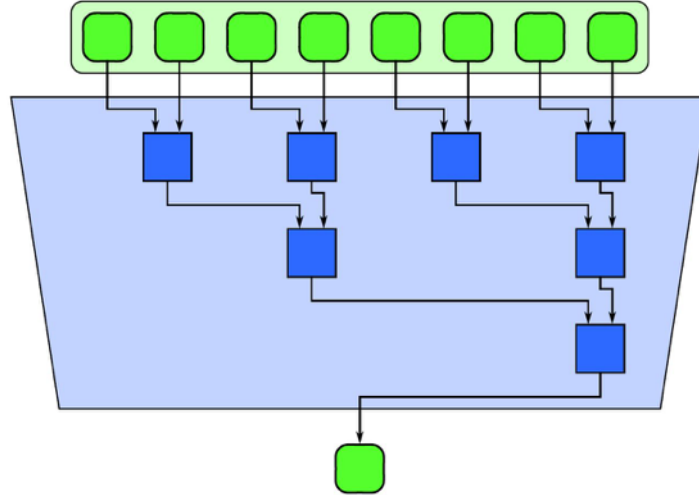


Figure 6.4: Visualization of parallel reduce in ArrowPipe [21]

polymorphic functions to construct expressive, composable and generic computation patterns.

More examples of algorithms formed by ArrowPipe, e.g. dot product or merge sort are shown in the Section 8.

6.2 Implementation of arrow combinators

In this chapter, we will present naive implementation and the optimized solution is introduced in the next section.

The intuition why ArrowPipe is an instance of Arrow comes from the Kleisli arrow of a monad is an instance of Arrow class (shown in Listing 22). The `cont` field in the Pipe has similar type signatures as the `runKleisli` field in the Kleisli arrow. From the previous section, we shown that Proc is a monad so Pipe is just an extended version of Kleisli arrow where computations in Pipe usually finish in one of the process stored in `env` instead of finishing at `cont` like Kleisli arrow. Intuitively, ArrowPipe, a function from the role to the Pipe, should be an instance of Arrow since Pipe looks like an arrow instance and functions are composite.

The essential issue when implementing arrow combinators is how to connect one Pipe by another Pipe. The first problem we need to address is how to deal with the `cont` in the tail Pipe. We know that only one `cont` field exists in the resulting Pipe and it must be that from the head Pipe. Hence the only option is to convert it to a Proc expression and store the converted expression in the updated `env` in the resulting Pipe. A right way to do it bind the `cont` with the action: receive from end in the first Pipe. Also, extend the proc related to the end in the head Porc with action: send to start in the second Pipe. Besides addition of the converted `cont` expression, the new `env` is formed by merging the `env` from the head Pipe and the `env` from the tail Pipe. Merging two `envs` is trivial. When there are duplication, we simply use monadic bind to combine them so that the actions belonging to head Pipe followed by the actions belonging to the tail Pipe. The `start` field in the resulting Pipe is the same as that from the head Pipe and the `end` field will be set the same as that from the tail Pipe. We can apply the Pipe composing function to implement arrow combinators for ArrowPipe. The implementation is just apply the first ArrowPipe with the input role (don't forget ArrowPipe is a function) to get the Pipe and apply the second ArrowPipe with a new role to get the second Pipe (usually in order to avoid duplication of roles, the new role is

```

preduc :: ArrowPipe (a, a) a -> ArrowPipe (a, (a, (a, a))) a
preduc r = helper >>> (r *** r) >>> r
where
  helper = (arr Id *** arr Fst) &&& (arr Snd >>> arr Snd)

```

Listing 20: Parallel reduce in ArrowPipe

```

divConquer
  :: Int
  -> ArrowPipe a b
  -> ArrowPipe a (a, a)
  -> ArrowPipe (b, b) b
  -> ArrowPipe a b
divConquer 0 baseFunc _split _merge = baseFunc
divConquer level baseFunc split merge =
  split
    >>> (    divConquer (level - 1) baseFunc split merge
          *** divConquer (level - 1) baseFunc split merge
        )
    >>> merge

twoWayThreeLevelDq = divConquer 3

```

Listing 21: 2-ways and 3-levels divided-and-conquer algorithm in arrowPipe

set to be maximum role in the first Pipe + 1) and finally apply the Pipe composing functions to both Pipe. A simplified code explanation can be seen in Listing 23. The rest of combinators can be implemented in a similar fashion.

6.3 Strategies for optimized role allocations

From the last section, we know the number of roles in the system is directly related to the number of processes in the final generated code. Hence, role allocating is an essential part in generating efficient parallel programs.

In this section, we propose strategies for optimizing role allocations. We have two goals in mind when optimizing; The first one is we would like to reduce the number of roles (processes) in the computation since the overhead of thread creation and data transmission has negative impact on performance. The second one is we do not want roles duplication when we try to compose ArrowPipes since role duplications means the different computation must be merged in the same role and computations in the same thread is sequential hence role duplications has negative impact on degree of parallelization.

If we only put the first goal in mind, an easy solution will be setup an upper bound of the number roles, and then we cycle through a fixed bound when allocating new roles. Processes corresponding to duplicated roles can be simply merged using binds since Proc is a monadic DSL and duality check ensures binding will not cause deadlocks. However, this strategies is not ideal since duplications of roles will decrease the degree of parallelization in the system.

If we only consider the second goal, naive strategies used in the previous section will satisfy the goal. However, the number of channels required and the number of roles in the system will grow exponentially. In a divided-and-conquer algorithm, the number of channels increases from 10 to 120 and the number of roles increases from 6 to 36 when the level is increased from 1 to 3.

For the purpose of illustration, we use inference rules to explain our proposed strategies for optimized role allocations when composing ArrowPipes. Please see Listing 24 as an example. $x \Rightarrow x$ means the computation start with role x and end with role x .

The rule for the rest of combinators are shown in Listing 25. Notice that for compose, id, arr and ArrowChoice we do not introduce any new roles, in other words, there is no parallelization for these combinators. Reader may find it strange that we do not intent to parallelize arr combi-


```

newtype Kleisli m a b = Kleisli { runKleisli :: a -> m b }

instance Monad m => Arrow (Kleisli m) where
  id = Kleisli return
  (Kleisli f) . (Kleisli g) = Kleisli (\b -> g b >=> f)
  arr f = Kleisli (return . f)
  first (Kleisli f) = Kleisli (\ ~(b,d) -> f b >=> \c -> return (c,d))
  second (Kleisli f) = Kleisli (\ ~(d,b) -> f b >=> \c -> return (d,c))

```

Listing 22: The implementation of arrow instance for Kleisli arrow of a monad

```

(>>>) :: (ArrowPipe a b) -> (ArrowPipe b c) -> (ArrowPipe a c)
(>>>) leftArrow rightArrow start = compose firstPipe secondPipe
where
  firstPipe = leftArrow start
  secondPipe = rightArrow (end leftP + 1)

```

Listing 23: The simplified implementation of >>>

nator which lifts a sequential computation represented by Core $(a \rightarrow b)$ into ArrowPipe. It makes sense to introduce a new role to execute the computation and hence parallelize computational heavy tasks. We use this strategy in the first place but later we found a more suitable strategy exists which will be introduced in the later paragraph. Also, another reason not to introduce new role when encountering arr combinators is that we gained function fusion for free. Simple function i.e. fst, inject left or snd are automatically fused into more complicated user defined functions. Introducing new roles for these simple functions will damage performance.

For the class of combinators belonging to arrow choice, we do not introduce any new role. The expressions at the lhs and at the rhs starts with the same role x because when only one code path will be executed as the name choice suggested so we should not use separate roles for two expressions that will never be executed simultaneously. In the end, we decided the computation end in the role $\max(y, z)$. Max guarantees that there will not be role duplications when we compose expressions formed by ArrowChoice combinators with other combinators. For the implementation, all process in both left and right ArrowPipe expression are wrapped inside a branch operation separately. Assume $\max(y, z) = y$, the process at the role y will be extended with actions that receive data from $\min(y, z) = z$ role at its right branch. Finally, applying inject left and inject right at left and right branches gives us Either type as the output.

Finally, we discovered the right place to allocate new roles is &&& combinator. As shown in type signature, product types mean computation at both branches will both be executed and they are independent. In order to make sure both computation are executed simultaneously, we constraint that the right ArrowPipe expression must start with a role greater than the end role of the left ArrowPipe expression. This ensures no role duplications hence maximize parallelism. The combined expression ends in the end role of the right ArrowPipe expression instead of introducing a unnecessary new role. For the implementations, the process corresponded to end role z are extended with actions that receive data from the end role y of the left process and store the computation of ArrowPipe expression at the right side and finally output a pair.

In conclusions, even though from the implementation point of view, arrowPipe composition with the optimized role allocations is ad-hoc and less elegant to implement because we need to consider composition by send-and-receive and composition by local monadic bind and more edge cases to be dealt with compared to the naive solution in the last section where composition is done solely by send-and-receive and role allocations is mindless. We believe the effort is worthy because for a n-level divided-and-conquer algorithms, the optimized role allocation strategies allocate 2^n roles in total which is the same as the number of way of parallelization in theory. All the roles are used to maximize parallelism instead of wasting the valuable resources to create roles that merely transmit data.

$$\text{id} \frac{x : \text{Role}, \quad a : \text{Type}}{id : \text{ArrowPipe } a \ a, x \Rightarrow x}$$

Listing 24: Role allocation for id

$$\text{compose} \frac{e1 : \text{ArrowPipe } a \ b, x \Rightarrow y, \quad e2 : \text{ArrowPipe } b \ c, y \Rightarrow z}{e1 >>> e2 : \text{ArrowPipe } a \ c, x \Rightarrow z}$$

$$\text{arr} \frac{f : \text{Core } (a \rightarrow b), \quad x : \text{Role}}{\text{arr } f : \text{ArrowPipe } a \ b, x \Rightarrow x}$$

$$\text{arrow choice: } ||| \frac{e1 : \text{ArrowPipe } a \ c, x \Rightarrow y, \quad e2 : \text{ArrowPipe } b \ c, x \Rightarrow z}{e1 ||| e2 : \text{ArrowPipe } (\text{Either } a \ b) \ c, x \Rightarrow \max(y, z)}$$

$$\text{arrow choice: } +++ \frac{e1 : \text{ArrowPipe } a \ c, x \Rightarrow y, \quad e2 : \text{ArrowPipe } b \ d, x \Rightarrow z}{e1 +++ e2 : \text{ArrowPipe } (\text{Either } a \ b) \ (\text{Either } c \ d), x \Rightarrow \max(y, z)}$$

$$\text{arrow: } \&\&\& \frac{e1 : \text{ArrowPipe } a \ b, x \Rightarrow y, \quad e2 : \text{ArrowPipe } a \ c, (y + 1) \Rightarrow z}{e1 \&\&\& e2 : \text{ArrowPipe } a \ (b, c), x \Rightarrow z}$$

$$\text{arrow: } *** \frac{e1 : \text{ArrowPipe } a \ b, x \Rightarrow y, \quad e2 : \text{ArrowPipe } a' \ c, (y + 1) \Rightarrow z}{e1 *** e2 : \text{ArrowPipe } (a, a') \ (b, c), x \Rightarrow z}$$

Listing 25: Rules fo role allocations of different combinators

6.4 Satisfaction of arrow laws

6.5 Conclusions

Arrow interface is the perfect interface to express general computation for this project because not only is it intuitive to understand and visualize but also its combinators `***` and `&&&` has built-in parallel natural.

So far, we've used ArrowPipe to help us compile Par-Alg to SPar. The remained challenge is the code generation from SPar to a target platform. In the next chapter, we will introduce our methods for code generation and specifically code generation to C. Once we achieved this, every computation in ArrowPipe can be transformed into parallel C code in one step.

Chapter 7

Type-safe code generation from SPar

SPar has two components: Core representing the unit of computation and Proc as a skeleton of the computation, describing the communication patterns. Naturally, the process of code generation from SPar should be divided into two parts correspondingly. We choose to make two parts independent of each other so that it's possible to swap the code generation strategy of one component without modifying another one.

The procedure of code generation is standard: transformation. We start our programs in a high level DSL and run a series of transformation to low-level DSL. SPar expressions are converted to a low-level EDSL which is then transformed to an abstract syntax tree (AST) of C (TODO cite the package). The generated code is obtained by pretty printing the AST.

7.1 Instr: A low-level EDSL for channel communication

In Proc, we have high-level actions like select, broadcast and branch abstracting implementation details, i.e variable declaration, variable assignment, channel initialization, channel communication and channel deletion. Hence, we need to define a EDSL containing instructions related to these low-level operations. We name it Instr. A SPar programs will be translated to a sequence of Instr.

When we design Instr, we keep the simplicity in mind so Instr is not coupled with any fancy language feature related to some specific target languages. Any reasonable target language with a channel communication library can be easily converted from Instr.

7.1.1 Syntax and semantic

The definition of Instr is seen in Listing 26. Channel is our abstract representation of Channel in Instr. It is indexed by a type `a` from the reified type `ReprType a`. More details of this reified type will be introduced in Section 7.1.2. This type parameter preserves type in channel initialization hence make sure the value to be sent or received in this channel has the same type as this channel. This is necessary because for some target languages, the channel are typed. Similarly, type parameters in `Exp` have the same functionality. `Exp` is just a wrapper of the expression in Core. In later stages, we will take care of code generation of `Exp` along with `Instr`. `Instr` defines the set of statements that will be generated and `Exp` represents the sequential computation, which is a value that will be generated.

The semantic of Instr operations are similar to what their names suggest. `CInitChan` represents operations that initialize a channel according to the given type and `cid`. `CDeleteChan` will destroy a channel. `CSend` operation sends the value `Exp a` through the Channel. `CRecv` action means the value received in the channel will be assigned to variable whose postfix name is the int field. `CEnd` means the instruction exits with the value `Exp a`. `CDecla` and `CAssgn` are instructions for variable declaration and assignment. The type of the variable is determined by `ReprType a` and the value is `Exp a`. `CBranch` and `CSelect` are used to express conditional control flow of the Instr languages. SPar action like broadcast are built on top of these operations. For `CBranch`, the first field represents the value of Either type to be received via the channel and two `Seq Instrs` represents the sequence of Instrs in the left branch and the right branch. For `CSelect`, the first field represents the variable containing Either value and the second field field

```

data Channel a where
  Channel :: CID -> ReprType a -> Channel a

data Exp a where
  Exp :: Core a -> ReprType a -> Exp a

data Instr where
  CInitChan :: Channel a -> Instr
  CDeleteChan :: Channel a -> Instr
  CSend :: Channel a -> Exp a -> Instr
  CRecv :: Channel a -> Int -> Instr
  CEnd :: Exp a -> Instr
  CDecla :: Int -> ReprType a -> Instr
  CAssgn :: Int -> Exp a -> Instr
  CBranch :: Int -> Seq Instr -> Seq Instr -> Instr
  CSelect :: Int -> Int -> Seq Instr -> Seq Instr -> Instr

```

Listing 26: The syntax of Instr in Haskell with accompanying low-level data types

```

data ReprType a where
  NumReprType :: NumType a -> ReprType a
  LabelReprType :: ReprType Label
  SumReprType :: ReprType a -> ReprType b -> ReprType (Either a b)
  UnitReprType :: ReprType ()
  ProductReprType :: ReprType a -> ReprType b -> ReprType (a, b)
  ListReprType :: ReprType a -> ReprType [a]

```

Listing 27: The definition of representation types

represents the variable whose value is assigned by the end results of instructions from either the left branch or the right branch. The third and fourth fields represent the instructions in the left branch and the right branch.

7.1.2 Representation types

SPar programs cannot be fully parametric since the target languages of code generation from SPar are usually less expressive, i.e, they do not treat function type $a \rightarrow b$ as a value, and are less efficient when processing with some specific form of data, i.e, languages targeting GPUs are usually more productive in dealing with array of floating point number while slow in working with aggregate structures [22]. Hence, we need to restrict the set of types allowed in SPar. We achieve this using the type class `Repr` and corresponding reified type `ReprType` (shown in Listing 27). `Repr` determines the set of type allow in SPar. Reified type `ReprType` will be used

```

constToCExpr :: ReprType a -> a -> CExpr
constToCExpr (NumReprType numType) v = numTypeToCExpr numType v
constToCExpr LabelReprType          v = case v of
  Le -> cVar "LEFT"
  Ri -> cVar "RIGHT"
constToCExpr s@(ProductReprType a b) v = defCompoundLit
  (show s)
  [ ([], initExp $ constToCExpr a (fst v))
  , ([], initExp $ constToCExpr b (snd v))
  ]
-- omit other cases

```

Listing 28: An example usage of reified type in the code generation

to alter the behaviors of code generation based on the type. This can be simply done by pattern matching because reified types are values in Haskell [23]. To be more concrete, Listing 28 gives a demo. `constToCExpr` is function that handle code generation from constant value to expression in C programming languages. By pattern matching, we vary the behaviors of code generation so that constants with different types has their own way to be represented in C.

In conclusion, we allow the following type: numerical type like `Float` and `Int`, the unit type `()`, the label type which is used in the code generation of select and branch and the aggregate type: list, product and sum that are built recursively, to be expressed in `SPar`.

7.2 Compilation from `SPar` to `Instr`

7.2.1 Transformation from `Proc` to `Instr`

As described in the previous section, `Instr` contains a data type called `Exp` which is a wrapper of Core expression. Compiling Core to `Instr` is hence not difficult. The challenge of compiling Core is mainly how to compile it to a specific target language. This will be discussed in the next subsection.

In this section, we will explain how we transform operations in `Proc` to `Instr`. Generally speaking, each `Proc` operation is mapped to a sequence of actions in `Instr`. The transformation algorithm from a `Proc` expression to a sequence of `Instr` can be implemented easily by traversing `Proc` expression, applying the mapping and collecting the results by concatenation. This is an advantage of using free monad technique to build the AST because `Proc` expression can be treated as a data structures and traversing recursive data structures can be easily done in Haskell. In addition, operations like `Recv` which involves continuation whose type is `Core a -> next` in their constructors are treated differently than those operations whose constructors only have a value type `next`. The latter is easy to implement, we can simply call the traversing function recursively. For the former, we have to pass an expression whose type is `Core a` to the continuation so that we can recursively call the traverse function recursively on the results of applying value to the continuation. The answer to what Core expression should be used is `Var` denoting variable. Passing a unique variable to the continuation gives us `next` inexpensively and we will define where does values of variables comes from for each operations in `Proc`.

We have introduced the general principle to the readers. Now let us dig into details of translation rules for each operations.

- Pure:** It is the base case in free monad. Hence it is mapped to the `CEnd` instruction.
- Send:** It is mapped to a sequence of three instructions. First of all, We declared a temporary variable using `CDecla` and then assign the value that will be sent to this variable using `CAssign` and send the content of the variable via the specific channel. The problem of how make sure the same channel is used in a send-and-receive pair will be discussed in the next sub chapter.
- Recv:** It is the reverse of send operation. Firstly, it declares a new variable with `CDecla` and use `CRecv` to assign the value received from the sender to this variable. Notice that the `recv` has a continuation, we will pass the variable declared in the first `instr` to the continuation to traverse the `Proc` expression recursively as discussed above.
- Select:** It is a more complicated operations. Its constructor contains two continuation: one for left branch and one for right branch. Hence for this instruction, we need to declared two variable to passed into continuations. The value of both variables is assigned by the Core expression whose value is a `Either` type. Besides, we need to send label indicating whether the execution of the left branch or the right branch of the receiver is selected. Its value is assigned by the either value as well and the sending operation is done by `CSend`. Finally, we call the transforming function recursively on the left branch and right branch and combine the results using `CSelect`.
- Broadcast:** The mapping from Broadcast is similar to that of Select. The only difference is that the former sends to a list of receivers while the latter sends to a single receiver. So in this operation, we will have multiple `CSend` corresponding to each receiver.

Branch: It is the reverse of the Select operation. So it will use `CRecv` to receive a label from the sender and call recursively on two branches and finally use `CBranch` to collect results.

7.2.2 Strategies for channel allocations

Channel allocations is important because correction allocation is essential in making sure the correctness and deadlock-free of generated code. Besides correctness concerns, we are also want to reduce the amount of channels creations hence increase performances.

In the first iteration, inspired by the linearity of channels in the π calculus, we choose to allocate a one-time channel for each send-and-receive pair. All channels' buffer size is one because of the linearity. Send action will initialize a channel and Receive action will destroy this channel once it receive the value. We use the ensure the same channel is used for the pair. During the transformation, we use a map of queue whose key is a pair of sender id and receiver id. When visiting the send action, it will push the channel into the queue and the corresponding received operation will pop the channel from the queue. Because we've ensured the duality of all processes in the system, we can claim the channel is right for each send-and-receiver pair. However, we realize this strategy is complicate to implement and not resource efficient since too many channels are initialized.

In the second iteration, we come up with a more efficient and simple strategy. The buffer size of all channels is still one due to the same reason about linearity. However, we decide to only allocate one channel for a pair of sender and receiver. We will not destroy the channel after the value has received and will reuse the channel for the next communication. When all processes have returned, we will destroy all channels at once. For this strategy, we have simplified the state from a map of queues of channels to a map of channels.

7.2.3 Monad for code generation

```
data CodeGenState = CodeGenState
{
  chanTable :: Map ChanKey CID,
  varNext   :: Int,
  chanNext  :: CID,
  dataStructCollect :: Set AReprType
}

newtype CodeGen m a =
  CodeGen { runCodeGen :: StateT CodeGenState m a }
  deriving (Functor,
            Applicative,
            Monad,
            MonadState CodeGenState,
            MonadIO)
```

Listing 29: States required during the traversal

From the last two subsections, we need to maintain a number of states during the compilation process. Hence we define a state monad to be used during the traversal. The `CodeGenState` is the collection of states with different purposes and it is shown in Listing 29. `chanTable` is the map we required during the channel allocation. `varNext` represents the next variable id to used. It will increment by one every time we declare a new variable. It helps us make sure the variable names are unique. `chanNext` has the similar functionality ensuring the uniqueness of channel names. `dataStructCollect` is the set of compound type we encounter during the traversal. With this states, we know what kind of data structures whose definitions will be generated before the generation of executing code.

7.3 Code generation to C: from Instr to C

The last piece of jigsaw is compilation from a sequence of Instr to C. This process is rather simple but trivial. This is simply done by transforming the sequence of Instr to a C AST. We used an open source library: language-c [24] to represent C AST in Haskell and pretty printing the C AST gives us the generated code. This method can be generalized to any target language. As for the implementation of channel communication in C, we used another open source library: chan [25] whose internal is based on shared memory. In the following subsections, we will present some design choices during this final step.

7.3.1 Representations of Core data type in C

```
typedef enum Label {
    LEFT, RIGHT
} Label;

typedef struct Prod_int_int {
    int fst; int snd;
} Prod_int_int;

typedef struct Sum_unit_Prod_int_int {
    Label label;
    union {
        int left; Prod_int_int right;
    } value;
} Sum_unit_Prod_int_int;
```

Listing 30: Compound data type in C

The first challenge we face is how to represent data structure in C. For primitive data type like Int or Float, a simple one-to-one mapping is sufficed. It is hard to deal with compound data type in C. First of all, C does not support polymorphic type. Hence, we choose to generate specific data type for every different compound data type even though they have the same structure. We have a way to name the generated data type to avoid name duplication. The naming simply reflects the structure of the data types with its elemental type. For example, $(\text{Int}, (\text{Int}, ()))$ is converted to `Prod_Int_Prod_Int_Unit` and `Either (Either () (Int, Int)) (Int, Int)` is converted to `Sum_Sum_unit_Prod_int_int_Prod_int_int`. In this project, all compound types are formed by sum type and product type. The product type will be converted to a struct with two field in C. The sum type is represented by the tagged union type. Tagged union is a struct with two field. The value of the first field indicates whether it is a left value or right value and the value of second field is a union type containing either the left value or right value. We also implement a sorting algorithm based on the depth of compound type so that all necessary data types have been defined before the definition of the compound types. An example of what `Either Int (Int, Int)` will be converted to is shown in Listing 30.

Another challenge is representation of the recursive type. From the type theory, we learn that a list of int can be expressed as $\mu a.() + \text{Int} \times a$. We might reuse the idea from the last paragraph to generated recursive type in terms of sum and product type. Hence a list of Int will look like the code below.

```
typedef struct Sum_unit_Prod_int_a {
    Label label;
    union {
        int left;
        Sum_unit_Prod_int_a *right;
    } value;
} Sum_unit_Prod_int_a
```

However, we believe expressing typical recursive data structures like list of int in this way is bad for performances. Obviously, C has a more efficient way to represent list of int using an array. So

we decided to have two ways to represent recursive data structures. For a set of specific recursive data structures, users can write their own representation to exploit the advantages of the target language. For example, a list of int are encoded in C using a wrapper of pointer type (shown in Listing 31). This way is not very generic but friendly for performances. For other types of recursive data structures where user do not specify their optimized versions in C, we simply apply the method in the last paragraph to encode them in C. This way is generic but not efficient.

```
typedef struct List_int {
    size_t size; int * value;
} List_int;
```

Listing 31: Optimized represent of List in C

7.3.2 Compiling from Core to C

$$\begin{array}{c}
\text{Var} \frac{}{\llbracket \text{Var } n \rrbracket = \text{vn}} \qquad \text{Lit} \frac{}{\llbracket \text{Lit val} \rrbracket = \text{toC(val)}} \\
\text{Fst} \frac{\llbracket a \rrbracket = c}{\llbracket \text{Fst 'ap' } a \rrbracket = c.\text{fst}} \qquad \text{Snd} \frac{\llbracket a \rrbracket = c}{\llbracket \text{Snd 'ap' } a \rrbracket = c.\text{snd}} \\
\text{Inl} \frac{\llbracket a \rrbracket = c}{\llbracket \text{Inl 'ap' } a \rrbracket = \{\text{LEFT}, c\}} \qquad \text{Inr} \frac{\llbracket a \rrbracket = c}{\llbracket \text{Inr 'ap' } a \rrbracket = \{\text{RIGHT}, c\}} \\
\text{Pair} \frac{\llbracket a \rrbracket = c_1, \quad \llbracket b \rrbracket = c_2}{\llbracket \text{Pair } a \text{ } b \rrbracket = \{c_1, c_2\}} \\
\text{Id} \frac{\llbracket a \rrbracket = c}{\llbracket \text{Id 'ap' } a \rrbracket = c} \qquad \text{Const} \frac{\llbracket a \rrbracket = c, \quad \llbracket v \rrbracket = b}{\llbracket (\text{Const } v) \text{ 'ap' } a \rrbracket = b} \\
\text{Prim} \frac{\llbracket a \rrbracket = c}{\llbracket (\text{Prim fname fimpl}) \text{ 'ap' } a \rrbracket = \text{fname}(c)}
\end{array}$$

Listing 32: Rules for compilation from Core to C

Core has a concise syntax so it does not require too much work to write a function that generate C expressions from Core expressions. Not surprisingly, the compilation is a traversal of the Core expression. Pattern matching on the structure of the core expression alters the behaviors of compilation. `ap` (apply) constructor is used with an expression whose type is `Core (a -> b)` and another expression whose type is `Core b`. The code generation for `ap` depends on the what the function expression is. The code generation rule is explained by the inference rules shown in Listing 32. $\llbracket a \rrbracket$ means the C code generated by Core expression `a`. `toC` is function that convert constant value in Haskell to constant value in C.

Var, Lit: Code generation of `Lit` simply convert Haskell value to its corresponding C value. As for `Var`, it will be converted to a string literal composed of the variable identity prefixed by `'v'`.

Fst, Inl, Pair... For `Inl` and `Inr` and `Pair`, we used C99 style to initialize struct and union. The rule for generating corresponding struct are explained in the previous subsection. For `Fst` and `Snd`, we simply access the specific value using the designator.

Id, Const v: Code generation of `Id` constructor is same as the code generation of the argument of the `Id`. Code generation of `Const` ignores the argument and use code generation of `v` expression instead.

Prim: `Prim` constructor represents user defined functions. The code generation for function call of `Prim` is converted to function call by the name of the primitive function. Users can implement those functions in C and include them in the main generated file.


```

void proclRt ()
{
    List_int v0;
    chan_recv_buf(c1, &v0, sizeof(List_int));
    List_int v1;
    v1 = sort(v0);
    chan_send_buf(c2, &v1, sizeof(List_int));
}
void * procl ()
{
    proclRt ();
    return NULL;
}
List_int proc0(List_int v0)
{
    c1 = chan_init(1);
    c2 = chan_init(1);
    pthread_t th1;
    pthread_create(&th1, NULL, procl, NULL);
    chan_send_buf(c1, &v0, sizeof(List_int));
    List_int v1;
    chan_recv_buf(c2, &v1, sizeof(List_int));
    pthread_join(th1, NULL);
    chan_dispose(c1);
    chan_dispose(c2);
    return v1;
}

```

Listing 33: An example of generated code

7.3.3 The structure of generated C code

We have covered the code generation algorithm for each process. We will now tackle how the generated code is structured from a group of interacting processes representing a parallel computation. For each process, we will generate its own C functions that take no argument. We decide to generate an addition C function since function is portable and acts like a black box hiding all the execution details so users can interact with generated code by calling the function. This function will take a parameter as input data and return the computation result. Inside, the function, it will spawn the same number of threads as the number of roles in the system. Each thread will execute the code which is generated from their corresponding Proc expression. In addition to that, the function will send the input parameter to the starting role in the group of Proc to kick off the computation using channel and waiting for the the ending role to send back the computation results. After it receive the result and all the threads have returned, it will return the result. We called this function `proc0`.

An example of the generated code is shown in Listing 33. The code contains one process which received a list of int from `proc0`, sort the input by an user defined function and then send the result to `proc0`. We omit global channels declarations and data type declaration for simplicity.

7.4 Conclusion

With the completion of code generation, we deliver the results we promised in the introduction section. We have implemented a end-to-end process that will generate low-level deadlock-free parallel code from an expressive high-level languages embedded with a flexible backend that can target multiple languages with ease. Now, it is time to evaluate the performances of our achievement with quantitative measurements.

Chapter 8

Parallel algorithms and evaluation

What we have implemented so far not only can be used as a backend for compiling ParAlg but also a stand-alone tool to express parallel algorithms. In this section, we will give an overview of how to use SPar for parallel algorithms, benchmarking the performance of the generated code for various computation and analyzing the design choices of the project.

8.1 Parallel algorithms

The biggest advantage of writing parallel programs in SPar is that user can express the computation similar to the sequential one without worrying about any low-level primitives for parallel computations. We will give a recommended recipe of expressing computation in SPar to follow and two concrete examples for the explanation.

8.1.1 Four steps to write parallel algorithms in SPar

Usually, divide-and-conquer algorithms are the best candidates for SPar to parallelize. The recipe is below:

1. Understand the algorithm: We recommend programmers to express the algorithms in the recursion schemes (see Section 2.2) Recursion schemes is recommended because it separates the split function, the merge function and the structure of divide-and-conquer from each others so it helps you familiar with the building blocks of the divide-and-conquer algorithms, i.e, the data structures involved, the type signature of the split function, the type signature of merge functions and their implementations.
2. Systematic parallelization: In the first iteration, we can express the split, merge functions and other necessary helper functions in terms of the ArrowPipe by combing `arr` constructor and `Prim` constructor. Every computation wrapped by `arr` with `Prim` is consider to be sequential. We will substitute them into high-level parallel patterns provided by SPar. For example, divide-and-conquer algorithms can be parallelized by `divConquer` (see Listing 21) helper functions. Notice that the parameter for number of ways of parallelized divide-and-conquer algorithms should be set accordingly by the number of cores of the execution machine.
3. Specific parallelization: The above step is generic and can be applied to any divide-and-conquer algorithm. In this stage, what will be done is determined by the specific implementations. Programmer should inspect the implementation of these function wrapped by `arr` with `Prim` to see whether there are any parallelism to exploit. If so, programmer should rewrite these functions by the parallel patterns provided by SPar or arrow combinators. For example, the split function of the Quickhull [26] to solve convex hull problems will use a for-loop to find the point whose distance to the line is at the maximum. This step can be expressed by the parallel map-and-reduce pattern (see Listing 20). Programmers can apply this step iteratively until all possible parallelisms are exploited or programmers think it is enough. All the sequential computation are left in the form: `arr with Prim`.

4. Wrap up: Before the code generation, programmer need to implement all the `Prim` functions in terms of the target language. In the scope of this project, programmer will write them in C and create a header file. The generated code will include the header file and from then on, programmers obtain the parallelized version of the algorithm that is guaranteed to be deadlock-free.

8.1.2 Example: Merge sort

8.1.3 Example: Word count

8.2 Benchmarks

8.2.1 Benchmarks against generated Haskell code

8.2.2 Benchmarks against C implementation

8.3 Evaluation

8.3.1 Design choices: Why Haskell?

Chapter 9

Conclusions and future work

9.1 Future work

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