# Arrows for Parallel Computations

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#### **Abstract**

Arrows are a general interface for computation and therefore form an alternative to monads for API design. We express parallelism using this concept in a novel way: We define an arrows-based language for parallelism and implement it using multiple parallel Haskells. In this manner we are able to bridge across various parallel Haskells.

Additionally, our way of writing parallel programs has the benefit of being portable across flavours of parallel Haskells. Furthermore, as each parallel computation is an arrow, which means that they can be composed and transformed as such. We introduce some syntactic sugar to provide parallelism-aware arrow combinators.

To show that our arrow-based language is on par with the existing parallel languages, we also define several parallel skeletons with our framework. Benchmarks show that our framework does not induce too much overhead performance-wise.

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#### **Contents**

#### 1 Introduction

#### OL: todo, reuse 5.5, and more

blablabla arrows, parallel, haskell.

### Contribution OL: HIT HERE REALLY STRONG

**MB:** different, how? We wrap parallel Haskells inside of our *ArrowParallel* interface, but why do we aim to abstract parallelism this way and what does this approach do better than the other parallel Haskells?

- **Arrow DSL benefits**: With the *ArrowParallel* typeclass we do not lose any benefits of using arrows as *parEvalN* is just yet another arrow combinator. The resulting arrow can be used in the same way a potential serial version could be used. This is a big advantage of this approach, especially compared to the monad solutions as we do not introduce any new types. We can just 'plug' in parallel parts into sequential arrow-based programs without having to change anything.
- **Abstraction**: With the *ArrowParallel* typeclass, we abstract all parallel implementation logic away from the business logic. This leaves us in the beautiful situation of being able to write our code against the interface the typeclass gives us without being bound to any parallel Haskell. So as an example, during development, we can run the program in a simple GHC-compiled variant and afterwards deploy it on a cluster by converting it into an Eden version, by just replacing the actual *ArrowParallel* instance.

**Structure** The remaining text is structures as follows. Section 2 briefly introduces known parallel Haskell flavours and gives an overview of Arrows to the reader (Sec. 2.2). Section 3

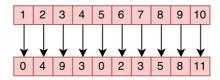


Figure 1: Schematic illustration of parEvalN.

discusses related work. Section 4 defines Parallel Arrows and presents a basic interface. Section 5 defines Futures for Parallel Arrows, this concept enables better communication. Section 6 presents some basic algorithmic skeletons (parallel *map* with and without load balancing, *map* – *reduce*) in our newly defined dialect. More advanced ones are showcased in Section 7 (*pipe*, *ring*, *torus*). Section 8 shows the benchmark results. Section 9 discusses future work and concludes.

#### 2 Background

## 2.1 Short introduction to parallel Haskells

There are already several ways to write parallel programs in Haskell. As we will base our parallel arrows on existing parallel Haskells, we will now give a short introduction to the ones we use as backends in this paper.

In its purest form, parallel computation (on functions) can be looked at as the execution of some functions  $a \to b$  in parallel or  $parEvalN :: [a \to b] \to [a] \to [b]$ , as also Figure 1 symbolically shows. Before we go into detail on how we can use this idea of parallelism for parallel Arrows, as a short introduction to parallelism in Haskell we will now implement parEvalN with several different parallel Haskells.

## 2.1.1 Multicore Haskell

Multicore Haskell (Marlow *et al.*, 2009; Trinder *et al.*, 1998) is way to do parallel processing found in standard GHC.<sup>1</sup> It ships with parallel evaluation strategies for several types which can be applied with *using*::  $a \rightarrow Strategy \ a \rightarrow a$ .

For parEvalN this means that we can just apply the list of functions  $[a \rightarrow b]$  to the list of inputs [a] by zipping them with the application operator \$. We then evaluate this lazy list [b] according to a Strategy[b] with the  $using::a \rightarrow Strategy[a \rightarrow a]$  operator. We construct this strategy with parList::Strategy[a] and rdeepseq::NFData[a] as rdeepseq::NFData[a] where the latter is a strategy which evalutes to normal form. To ensure that programs that use parEvalN have the correct evaluation order, we annotate the computation with  $pseq::a \rightarrow b \rightarrow b$  which forces the compiler to not reorder multiple parEvalN computations. This is particularly necessary in circular communication topologies like in the torus or ring (Chap. 7), where a wrong execution order would result in deadlock scenarios when executed without pseq.

Multicore Haskell on Hackage is available under https://hackage.haskell.org/package/parallel-3.2.1.0, compiler support is integrated in the stock GHC.

The resulting code and a graphical representation can be found in Fig. 2 and Fig. 3, respectively.

```
parEvalN :: (NFData\ b) \Rightarrow [a \rightarrow b] \rightarrow [a] \rightarrow [b]
parEvalN fs \ as = let \ bs = zipWith \ (\$) fs \ as
   in (bs 'using' parList rdeepseq) 'pseq' bs
```

Figure 2: Multicore version of parEvalN.

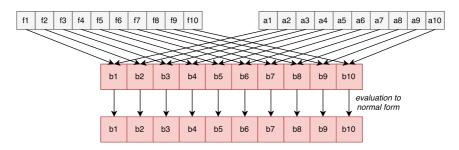


Figure 3: Dataflow of the Multicore Haskell parEvalN version.

# 2.1.2 ParMonad

The Par monad<sup>2</sup> introduced by Marlow et al. (2011a), is a monad designed for composition of parallel programs.

The Par Monad version of our parallel evaluation function parEvalN can be defined by zipping the list of  $[a \to b]$  with the list of inputs [a] with the application operator \$ just like with Multicore Haskell. Then, we map over this not yet evaluated lazy list of results [b] with  $spawnP :: NFData \ a \Rightarrow a \rightarrow Par \ (IVar \ a)$  to transform them to a list of not yet evaluated forked away computations [Par (IVar b)], which we convert to Par [IVar b] with sequenceA. We wait for the computations to finish by mapping over the *IVar b* values inside the *Par* monad with get. This results in Par[b]. We execute this process with runPar to finally get the final [b].

Again, the resulting code and a graphical representation can be found in Fig. 4 and Fig. 5, respectively.

MB: explain problems with laziness here. Problems with torus

## 2.1.3 Eden

Eden (Loogen et al., 2005; Loogen, 2012) is a parallel Haskell for distributed memory and comes with a MPI and a PVM backends.<sup>3</sup> It is targeted towards clusters, but also functions

<sup>&</sup>lt;sup>2</sup> It can be found in the monad-par package on hackage under https://hackage.haskell.org/ package/monad-par-0.3.4.8/.

See also http://www.mathematik.uni-marburg.de/~eden/ and https://hackage. haskell.org/package/edenmodules-1.2.0.0/.

### Arrows for Parallel Computations

```
parEvalN :: (NFData\ b) \Rightarrow [a \rightarrow b] \rightarrow [a] \rightarrow [b]

parEvalN\ fs\ as = runPar\ \$

(sequenceA\ \$map\ (spawnP)\ \$zipWith\ (\$)\ fs\ as) >>= mapM\ get
```

Figure 4: Par Monad version of parEvalN.

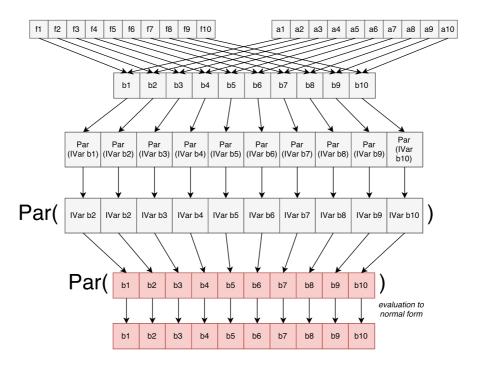


Figure 5: Dataflow of the Par Monad parEvalN version.

well in a shared-memory setting with a further simple backend. However, in contrast to many other parallel Haskells, in Eden each process has its own heap. This seems to be a waste of memory, but with distributed programming paradigm and individual GC per process, Eden yields good performance results also on multicores (Berthold *et al.*, 2009a; Aswad *et al.*, 2009).

While Eden also comes with a monad PA for parallel evaluation, it also ships with a completely functional interface that includes a spawnF::  $(Trans\ a, Trans\ b) \Rightarrow [a \rightarrow b] \rightarrow [a] \rightarrow [b]$  function that allows us to define parEvalN directly:

```
parEvalN :: (Trans\ a, Trans\ b) \Rightarrow [a \rightarrow b] \rightarrow [a] \rightarrow [b]

parEvalN = spawnF
```

A simplistic graphical depiction of this definition can be found in Fig. 6.

**Eden TraceViewer.** To comprehend the efficiency and the lack thereof in a parallel program, an inspection of its execution is extremely helpful. While some large-scale solutions exist (Geimer *et al.*, 2010), the parallel Haskell community mainly utilises the

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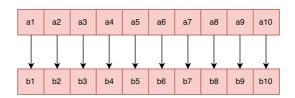


Figure 6: Dataflow of the Eden parEvalN version.

tools Threadscope (Wheeler & Thain, 2009) and Eden TraceViewer<sup>4</sup> (Berthold & Loogen, 2007). In the next sections we will present some *traces*, the post-mortem process diagrams of Eden processes and their activity.

In a trace, the *x* axis shows the time, the *y* axis enumerates the machines and processes. A trace shows a running process in green, a blocked process is red. If the process is 'runnable', i.e. it may run, but does not, it is yellow. The typical reason for then is GC. An inactive machine where no processes are started yet, or all are already terminated, is shows as a blue bar. A comminication from one process to another is represented with a black arrow. A stream of communications, e.g. a transmitted list is shows as a dark shading between sender and receiver processes. An example trace can be found in Fig. 18.

#### 2.2 Arrows

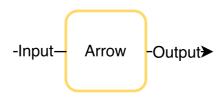


Figure 7: Schematic depiction of an arrow.

Arrows were introduced by Hughes (2000) as a general interface for computation. An arrow  $arr\ a\ b$  represents a computation that converts an input a to an output b. This is defined in the Arrow shown in Fig. 8.

Its arr operation is used to lift an ordinary function to the specified arrow type, similarly to the monadic return. The >>> operator is analogous to the monadic composition >>= and combines two arrows  $arr\ a\ b$  and  $arr\ b\ c$  by "wiring" the outputs of the first to the inputs to the second to get a new arrow  $arr\ a\ c$ . Lastly, the first operator takes the input arrow from b to c and converts it into an arrow on pairs with the second argument untouched. It allows us to to save input across arrows.

The most prominent instances of this interface are regular functions  $(\rightarrow)$ :

<sup>&</sup>lt;sup>4</sup> See http://hackage.haskell.org/package/edentv on Hackage for the last available version of Eden TraceViewer.



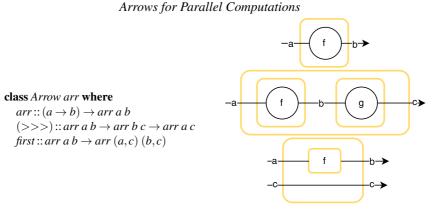


Figure 8: Arrow class definition.

Figure 9: schematic depiction of *Arrow* combinators *arr*, >>> and *first*.

```
instance Arrow (\rightarrow) where arr f = f

f >>> g = g \circ f

first f = \lambda(a,c) \rightarrow (f \ a,c)
```

and the Kleisli type:

```
data Kleisli m a b = Kleisli \{run :: a \rightarrow m b\}

instance Monad m \Rightarrow Arrow (Kleisli m) where

arr f = Kleisli (return \circ f)

f >>> g = Kleisli (\lambda a \rightarrow f a >== g)

first f = Kleisli (\lambda (a, c) \rightarrow f a >== \lambda b \rightarrow return (b, c))
```

With this typeclass in place, Hughes also defined some syntactic sugar (Fig. 10): The

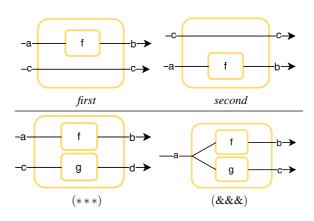


Figure 10: Visual depiction of syntactic sugar for arrows.

combinators second, \*\*\* and &&&. The definition of second, the mirrored version of first, is:

```
second :: Arrow arr \Rightarrow arr ab \rightarrow arr (c,a) (c,b)
second f = arr swap >>> first f >>> arr swap
where swap (x,y) = (y,x)
```

The \*\*\* combinator that combines *first* and *second* to handle two inputs in one arrow, is defined as

```
(***)::Arrow arr \Rightarrow arr a b \rightarrow arr c d \rightarrow arr (a,c) (b,d) f ***g = first <math>f >>> second g
```

, while the &&& combinator, that constructs an arrow which outputs two different values like \*\*\*, but takes only one input, is:

```
(&&&)::Arrow arr \Rightarrow arr \ a \ b \rightarrow arr \ a \ c \rightarrow a \ a \ (b,c)
f \&\&\& g = arr \ (\lambda a \rightarrow (a,a)) >>> (f ***g)
```

A first short example given by Hughes on how to use arrows is addition with arrows:

```
add:: Arrow \ arr \Rightarrow arr \ a \ Int \rightarrow arr \ a \ Int \rightarrow arr \ a \ Int
add \ f \ g = (f \&\&\& \ g) >>> arr \ (\lambda(u,v) \rightarrow u+v)
```

The more restrictive interface of arrows (a monad can be *anything*, an arrow is a process of doing something, a *computation*) allows for more elaborate composition and transformation combinators. One of the major problems in parallel computing is composition of parallel processes.

# 3 Related Work

#### **OL: arrows or Arrows?**

# 3.1 Parallel Haskells

Of course, the three parallel Haskell flavours we have presented above: the GpH (Trinder *et al.*, 1996, 1998) parallel Haskell dialect and its multicore version (Marlow *et al.*, 2009), the *Par* monad (Marlow *et al.*, 2011b; Foltzer *et al.*, 2012), and Eden (Loogen *et al.*, 2005; Loogen, 2012) are related to this work. We use these languages as backends: our DSL can switch from one to other at user's command.

HdpH (Maier *et al.*, 2014; Stewart *et al.*, 2016) is an extension of *Par* monad to heterogeneous clusters. LVish (Kuper *et al.*, 2014) is a communication-centred extension of *Par* monad. Further parallel Haskell approaches include pH (Nikhil & Arvind, 2001), research work done on distributed variants of GpH (Trinder *et al.*, 1996; Aljabri *et al.*, 2014, 2015), and low-level Eden implementation (Berthold, 2008; Berthold *et al.*, 2016). Skeleton composition (Dieterle *et al.*, 2016), communication (Dieterle *et al.*, 2010a), and generation of process networks (Horstmeyer & Loogen, 2013) are recent in-focus research topics in Eden. This also includes the definitions of new skeletons (Hammond *et al.*, 2003; Berthold & Loogen, 2006; Berthold *et al.*, 2009b,c; Dieterle *et al.*, 2010b; de la Encina *et al.*, 2011; Dieterle *et al.*, 2013; Janjic *et al.*, 2013).

More different approaches include data parallelism (Chakravarty *et al.*, 2007; Keller *et al.*, 2010), GPU-based approaches (Mainland & Morrisett, 2010; Svensson, 2011), software

22:28

### 3.2 Algorithmic skeletons

Algorithmic skeletons were introduced by Cole (1989). Early efforts include (Darlington *et al.*, 1993; Botorog & Kuchen, 1996; Danelutto *et al.*, 1997; Gorlatch, 1998; Lengauer *et al.*, 1997). Rabhi & Gorlatch (2003) consolidated early reports on high-level programming approaches. The effort is ongoing, including topological skeletons (Berthold & Loogen, 2006), special-purpose skeletons for computer algebra (Berthold *et al.*, 2009c; Lobachev, 2011, 2012; Janjic *et al.*, 2013), iteration skeletons (Dieterle *et al.*, 2013). The idea of Linton *et al.* (2010) is to use a parallel Haskell to orchestrate further software systems to run in parallel. Dieterle *et al.* (2016) compare the composition of skeletons to stable process networks.

#### 3.3 Arrows

Arrows were introduced by Hughes (2000), basically they are a generalised function arrow  $\rightarrow$ . Hughes (2005a) is a tutorial on arrows. Some theoretical details on arrows (Jacobs *et al.*, 2009; Lindley *et al.*, 2011; Atkey, 2011) are viable. Paterson (2001) introduced a new notation for arrows. Arrows have applications in information flow research (Li & Zdancewic, 2006, 2010; Russo *et al.*, 2008), invertible programming (Alimarine *et al.*, 2005), and quantum computer simulation (Vizzotto *et al.*, 2006). But perhaps most prominent application of arrows is arrow-based functional reactive programming, AFRP (Hudak *et al.*, 2003).**OL: cite more!** 

Liu *et al.* (2009) formally define a more special kind of arrows that capsule the computation more than regular arrows do and thus enable optimizations. Their approach would allow parallel composition, as their special arrows would not interfere with each other in concurrent execution. In contrast, we capture a whole parallel computation as a single entity: our main instantiation function *parEvalN* makes a single (parallel) arrow out of list of arrows. **OL: ugh, take care!** Huang *et al.* (2007) utilise arrows for parallelism, but strikingly different from our approach. They basically use arrows to orchestrate several tasks in robotics. We, however, propose a general interface for parallel programming, remaining completely in Haskell.

## 3.4 Other languages

Although this work is centred on Haskell implementation of arrows, it is applicable to any functional programming language where parallel evaluation and arrows can be defined. Experiments with our approach in Frege language <sup>5</sup> (which is basically Haskell on the JVM) were quite successful, we were able to use typical Java libraries for parallelism MB:

<sup>&</sup>lt;sup>5</sup> GitHub project page at https://github.com/Frege/frege

not really tested completely. basic parEvalN worked, we didn't test for different Java **ExecutorServices, though...** However, it is beyond the scope of this work.

Achten et al. (2004, 2007) use an arrow implementation in Clean for better handling of typical GUI tasks. Dagand et al. (2009) used arrows in OCaml in the implementation of a distributed system.

#### 4 Parallel Arrows

We have seen what Arrows are and how they can be used as a general interface to computation. In the following section we will discuss how Arrows constitute a general interface not only to computation, but to parallel computation as well. We start by introducing the interface and explaining the reasonings behind it. Then, we discuss some implementations using exisiting parallel Haskells. Finally, we explain why using Arrows for expressing parallelism is beneficial.

## 4.1 The ArrowParallel typeclass

As we have seen earlier, in its purest form, parallel computation (on functions) can be seen as the execution of some functions  $a \to b$  in parallel, parEvalN (Chap. 2.1). Translating this into arrow terms gives us a new operator parEvalN that lifts a list of arrows [arr a b] to a parallel arrow arr[a][b]. This combinator is similar to our utility function *listApp* from Appendix A, but does parallel instead of serial evaluation.

```
parEvalN :: (Arrow \ arr) \Rightarrow [arr \ a \ b] \rightarrow arr \ [a] \ [b]
```

With this definition of parEvalN, parallel execution is yet another arrow combinator. But as the implementation may differ depending on the actual type of the arrow arr and we want this to be an interface for different backends, we introduce a new typeclass ArrowParallel arr a b:

```
class Arrow arr \Rightarrow Arrow Parallel arr a b where
   parEvalN :: [arr \ a \ b] \rightarrow arr [a] [b]
```

Sometimes parallel Haskells require or allow for additional configuration parameters, e.g. an information about the execution environment or the level of evaluation (weak head normal form vs. normal form). For this reason we also introduce an additional conf parameter to the function. We also do not want *conf* to be a fixed type, as the configuration parameters can differ for different instances of ArrowParallel. So we add it to the type signature of the typeclass as well: **OL**: ArrowParallel arr a b conf **or** ArrowParallel conf arr a b? **MB**: does it really matter?

```
class Arrow arr \Rightarrow Arrow Parallel arr a b conf where
   parEvalN :: conf \rightarrow [arr \ a \ b] \rightarrow arr \ [a] \ [b]
```

Note that we don't require the *conf* parameter in every implementation. If it is not needed, we usually just default the *conf* type parameter to () and even blank it out in the parameter list of the implemented parEvalN, as we will see in the implementation of the Multicore Haskell and the Par Monad backends.

```
instance (NFData b,ArrowApply arr,ArrowChoice arr) \Rightarrow ArrowParallel arr a b () where parEvalN _fs = listApp fs >>> arr (withStrategy (parList rdeepseq)) && arr id>>> arr (uncurry pseq)
```

Figure 11: Fully evaluating ArrowParallel instance for the Multicore Haskell backend.

## 4.2 ArrowParallel instances

#### 4.2.1 Multicore Haskell

The Multicore Haskell implementation of this class is implemented in a straightforward manner by using listApp (Appendix A) combined with the withStrategy::  $Strategy \ a \to a \to a$  and pseq::  $a \to b \to b$  combinators from Multicore Haskell, where withStrategy is the same as using::  $a \to Strategy \ a \to a$  but with flipped parameters. For most cases a fully evaluating version like in Fig. 11 would probably suffice, but as the Multicore Haskell interface allows the user to specify the level of evaluation to be done via the Strategy interface, our DSL should allow for this. We therefore introduce the Conf a data-type that simply wraps a Strategy a:

```
data Conf \ a = Conf \ (Strategy \ a)
```

We can't directly use the *Strategy a* type here as GHC (at least currently) does not allow type synonyms in type class instances. To get our configurable *ArrowParallel* instance, we simply unwrap the strategy and pass it to *parList* like in the fully evaluating version (Fig. 12).

```
\begin{array}{l} \textbf{instance} \ (NFData \ b, ArrowApply \ arr, ArrowChoice \ arr) \Rightarrow \\ ArrowParallel \ arr \ a \ b \ (Conf \ b) \ \textbf{where} \\ parEvalN \ (Conf \ strat) \ fs = \\ listApp \ fs >>> \\ arr \ (withStrategy \ (parList \ strat)) \&\&\& \ arr \ id >>> \\ arr \ (uncurry \ pseq) \end{array}
```

Figure 12: Configurable ArrowParallel instance for the Multicore Haskell backend.

# 4.2.2 Par Monad

OL: introduce a newcommand for par-monad, "arrows", "parrows" and replace all mentions to them to ensure uniform typesetting The Par monad implementation (Fig. 13) makes use of Haskells laziness and Par monad's  $spawnP::NFData\ a\Rightarrow a\rightarrow Par\ (IVar\ a)$  function. The latter forks away the computation of a value and returns an IVar containing the result in the Par monad.

We therefore apply each function to its corresponding input value with I and then fork the computation away with *arr spawnP* inside a *zipWithArr* (Fig. A 3) call. This yields a

list [Par (IVar b)], which we then convert into Par [IVar b] with arr sequenceA. In order to wait for the computation to finish, we map over the IVars inside the Par monad with arr (>= mapM get). The result of this operation is a Par [b] from which we can finally remove the monad again by running arr runPar to get our output of [b].

```
instance (NFData b, ArrowApply arr, ArrowChoice arr) \Rightarrow
  ArrowParallel arr a b conf where
     parEvalN \_fs =
       (arr \$ \lambda as \rightarrow (fs, as)) >>>
       zipWithArr (app>>> arr spawnP)>>>
       arr sequenceA>>>
       arr (> = mapM get) >> >
       arr runPar
```

Figure 13: ArrowParallel instance for the Par Monad backend.

```
4.2.3 Eden
```

For both the Multicore Haskell and Par Monad implementations we could use general instances of ArrowParallel that just require the ArrowApply and ArrowChoice typeclasses. With Eden this is not the case as we can only spawn a list of functions and we cannot extract simple functions out of arrows. While we could still manage to have only one class in the module by introducing a typeclass:

```
class (Arrow arr) \Rightarrow Arrow Unwrap arr where
   arr\ a\ b \rightarrow (a \rightarrow b)
```

We don't do it here for aesthetic resons, though. For now, we just implement ArrowParallel for normal functions:

```
instance (Trans a, Trans b) \Rightarrow ArrowParallel (\rightarrow) a b conf where
   parEvalN \_fs \ as = spawnF \ fs \ as
and the Kleisli type:
   instance (Monad m, Trans a, Trans b, Trans (m b)) \Rightarrow
      ArrowParallel (Kleisli m) a b conf where
   parEvalN conf fs =
      (arr parEvalN conf (map (\lambda(Kleislif) \rightarrow f) fs)) >>>
      (Kleisli $ sequence)
```

# 4.3 Extending the Interface

With the ArrowParallel typeclass in place and implemented, we can now implement some further basic parallel interface functions. These are algorithmic skeletons that, however, mostly serve as a foundation to further, more specific algorithmic skeletons.

## Arrows for Parallel Computations

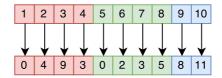


Figure 14: Schematic depiction of parEvalNLazy.

```
parEvalNLazy :: (ArrowParallel\ arr\ a\ b\ conf\ ,ArrowChoice\ arr\ ,ArrowApply\ arr) \Rightarrow conf\ \rightarrow ChunkSize\ \rightarrow [arr\ a\ b]\ \rightarrow (arr\ [a]\ [b]) parEvalNLazy\ conf\ chunkSize\ fs = arr\ (chunksOf\ chunkSize) >>> \\ listApp\ fchunks >>> \\ arr\ concat \mathbf{where}\ fchunks = map\ (parEvalN\ conf\ )\ \$\ chunksOf\ chunkSize\ fs
```

Figure 15: Definition of parEvalNLazy.

### 4.3.1 Lazy parEvalN

The function parEvalN is 100% strict, which means that it fully evaluates all passed arrows. Sometimes this might not be feasible, as it will not work on infinite lists of functions like e.g.  $map\ (arr\circ(+))\ [1..]$  or just because we need the arrows evaluated in chunks. parEvalNLazy (Fig. 14, 15) fixes this. It works by first chunking the input from [a] to [[a]] with the given ChunkSize in  $arr\ (chunksOf\ chunkSize)$ . These chunks are then fed into a list  $[arr\ [a]\ [b]]$  of parallel arrows created by feeding chunks of the passed ChunkSize into the regular parEvalN by using listApp. The resulting [[b]] is lastly converted into [b] with  $arr\ concat$ .

# 4.3.2 Heterogenous tasks



Figure 16: Schematic depiction of parEval2.

We have only talked about the paralellization arrows of the same type until now. But sometimes we want to paralellize heterogenous types as well. However, we can implement such a parEval2 combinator (Fig. 16, B 4) which combines two arrows  $arr\ a\ b$  and  $arr\ c\ d$  into a new parallel arrow  $arr\ (a,c)\ (b,d)$  quite easily with the help of the ArrowChoice typeclass. The idea is to use the +++ combinator which combines two arrows  $arr\ a\ b$  and  $arr\ c\ d$  and transforms them into  $arr\ (Either\ a\ c)\ (Either\ b\ d)$  to get a common arrow type that we can then feed into parEvalN.

## 5 Futures

Consider the parallel arrow combinator in Fig. 17 In a distributed environment, a resulting

```
someCombinator :: (Arrow\ arr) \Rightarrow [arr\ a\ b] \rightarrow [arr\ b\ c] \rightarrow arr\ [a]\ [c]
someCombinator\ fs1\ fs2 = parEvalN\ ()\ fs1 >>> rightRotate >>> parEvalN\ ()\ fs2
```

Figure 17: An example parallel Arrow combinator without Futures.

arrow of this combinator first evaluates all  $[arr\ a\ b]$  in parallel, sends the results back to the master node, rotates the input once and then evaluates the  $[arr\ b\ c]$  in parallel to then gather the input once again on the master node. Such situations arise, e.g. in scientific computations when the data distributed across the nodes needs to be transposed. A concrete example is 2D FFT computation (Gorlatch & Bischof, 1998; Berthold *et al.*, 2009c).

While the example in Fig. 17 could be rewritten into only one *parEvalN* call by directly wiring the arrows properly together, this example illustrates an important problem: When using a *ArrowParallel* backend that resides on multiple computers, all communication between the nodes is done via the master node, as shown in the Eden trace in Figure 18. This can become a serious bottleneck for larger amount of data and number of processes (showcases Berthold *et al.*, 2009c, as, e.g.).

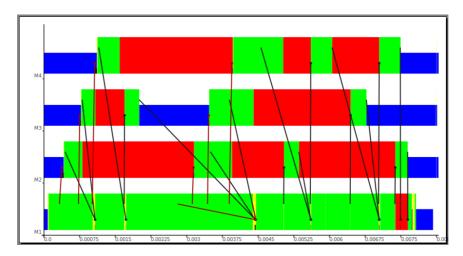


Figure 18: Communication between 4 threads without Futures. All communication goes through the master node. Each bar represents one process. Black lines between processes represent communication. Colors: blue  $\hat{=}$  idle, green  $\hat{=}$  running, red  $\hat{=}$  blocked, yellow  $\hat{=}$  suspended.

OL: more practical and heavy-weight example! fft (I have the code)?

MB: Depends... Are the communications easy to read in such an example?

MB: Keep the description for the different colours, or link to the EdenTV description in 2.1.3

This motivates for an approach that allows the nodes to communicate directly with each other. Thankfully, Eden, the distributed parallel Haskell we have used in this paper so far,

But as we want code written against our API to be implementation agnostic, we have to wrap this context. We do this with the *Future* typeclass (Fig. 19). Since *RD* is only a

```
class Future fut a \mid a \rightarrow fut where

put :: (Arrow arr) \Rightarrow arr \ a \ (fut \ a)

get :: (Arrow arr) \Rightarrow arr \ (fut \ a) \ a
```

Figure 19: Definition of the Future typeclass.

type synonym for communication type that Eden uses internally, we have to use some wrapper classes to fit that definition, though, as seen in Appendix in Fig. B 1. This is due to the same reason we had to introduce a wrapper for *Strategy a* in the Multicore Haskell implementation of *ArrowParallel* in Section 4.2.1.

For our Par Monad and Multicore Haskell backends, we can simply use *MVars* (Jones *et al.*, 1996) (Fig. 20), because we have shared memory in a single node and don't require Eden's sophisticated communication channels. **MB: explain MVars** 

```
{-# NOINLINE putUnsafe #-}

putUnsafe :: a → MVar a

putUnsafe a = unsafePerformIO $ do

mVar ← newEmptyMVar

putMVar mVar a

return mVar

instance (NFData a) ⇒ Future MVar a where

put = arr putUnsafe

get = arr takeMVar >>> arr unsafePerformIO
```

Figure 20: MVar instance of the Future typeclass for the Par Monad and Multicore Haskell backends.

Furthermore, in order for these *Future* types to fit with the *ArrowParallel* instances we gave earlier, we have to give the necessary *NFData* and *Trans* instances, the latter are only needed in Eden. Because *MVar* already ships with a *NFData* instance, we only have to supply two simple instances for our *RemoteData* type:

```
instance NFData (RemoteData a) where rnf = rnf \circ rd instance Trans (RemoteData a)
```

The *Trans* instance does not have any functions declared as the default implementation suffices here.

Going back to our communication example we can use this *Future* concept in order to enable direct communications between the nodes in the following way: In a distributed environment, this gives us a communication scheme with messages going through the master

```
someCombinator :: (Arrow \ arr) \Rightarrow [arr \ a \ b] \rightarrow [arr \ b \ c] \rightarrow arr \ [a] \ [c] someCombinator \ fs1 \ fs2 = parEvalN \ () \ (map \ (>>>put) \ fs1) >>> rightRotate >>> parEvalN \ () \ (map \ (get>>>) \ fs2)
```

Figure 21: The combinator from Fig. 17 in parallel.

node only if it is needed - similar to what is shown in the trace in Fig. 22.OL: Fig. 3 is not really clear. Do Figs 2-3 with a lot of load?

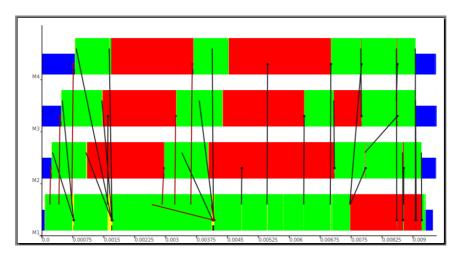


Figure 22: Communication between 4 threads with Futures. Other than in Fig. 18, threads communicate directly (black lines between the bars) instead of always going through the master node (bottom bar).

# 6 Map-based Skeletons

Now we have developed Parallel Arrows far enough to define some algorithmic skeletons useful to an application programmer.

# 6.1 Parallel map

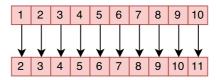


Figure 23: Schematic depiction of parMap.

The *parMap* skeleton (Fig. 23, 24) is probably the most common skeleton for parallel programs. We can implement it with ArrowParallel by repeating an arrow  $arr\ a\ b$  and then passing it into parEvalN to get an arrow  $arr\ [a]\ [b]$ . Just like parEvalN, parMap is 100% strict.

```
parMap :: (ArrowParallel \ arr \ a \ b \ conf) \Rightarrow conf \rightarrow (arr \ a \ b) \rightarrow (arr \ [a] \ [b])
parMap \ conf \ f = parEvalN \ conf \ (repeat \ f)
```

Figure 24: Definition of parMap.

# 6.2 Lazy parallel map

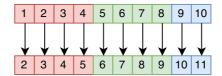


Figure 25: Schematic depiction of parMapStream.

As *parMap* (Fig. 23, 24) is 100% strict it has the same restrictions as *parEvalN* compared to *parEvalNLazy*. So it makes sense to also have a *parMapStream* (Fig. 25, 26) which behaves like *parMap*, but uses *parEvalNLazy* instead of *parEvalN*.

```
parMapStream :: (ArrowParallel\ arr\ a\ b\ conf\ , ArrowChoice\ arr\ , ArrowApply\ arr) \Rightarrow conf \to ChunkSize \to arr\ a\ b \to arr\ [a]\ [b] 
parMapStream\ conf\ chunkSize\ f = parEvalNLazy\ conf\ chunkSize\ (repeat\ f)
```

Figure 26: Definition of parMapStream.

### 6.3 Statically load-balancing parallel map

A parMap (Fig. 23, 24) spawns every single computation in a new thread (at least for the instances of ArrowParallel we gave in this paper). This can be quite wasteful and a farm (Fig. 27, 28) that equally distributes the workload over numCores workers (if numCores is greater than the actual processor count, the fastest processor(s) to finish will get more tasks) seems useful. The definitions of helper functions unshuffle, takeEach, shuffle (shown in Appendix) originate from an Eden skeleton<sup>6</sup>.

Figure 27: Schematic depiction of a farm, a statically load-balanced parMap.

```
farm :: (ArrowParallel\ arr\ a\ b\ conf\ , \\ ArrowParallel\ arr\ [a]\ [b]\ conf\ , ArrowChoice\ arr) \Rightarrow \\ conf\ \to NumCores\ \to arr\ a\ b\ \to arr\ [a]\ [b] \\ farm\ conf\ numCores\ f = \\ unshuffle\ numCores\ >> \\ parEvalN\ conf\ (repeat\ (mapArr\ f))>>> \\ shuffle \\ unshuffle\ :: (Arrow\ arr) \Rightarrow Int\ \to arr\ [a]\ [[a]] \\ unshuffle\ n = arr\ (\lambda xs\ \to [takeEach\ n\ (drop\ i\ xs)\ @\ [a\ i\leftarrow [0\ ..\ n-1]]) \\ takeEach\ :: Int\ \to\ [a]\ \to\ [a] \\ takeEach\ n\ (x:xs) = x: takeEach\ n\ (drop\ (n-1)\ xs) \\ shuffle\ :: (Arrow\ arr) \Rightarrow arr\ [[a]]\ [a] \\ shuffle\ = arr\ (concat\ o\ transpose)
```

Figure 28: The definition of farm.

# 6.4 The farmChunk Skeleton

Since a *farm* (Fig. 27, 28) is basically just *parMap* with a different work distribution, it is, again, 100% strict. So we can define *farmChunk* (Fig. 29, B 2) which uses *parEvalNLazy* instead of *parEvalN*. It is basically the same definition as for *farm*, with *parEvalN* replaced with *parEvalNLazy*, as Appendix shows.

## 6.5 Map and reduce

A simple *map–reduce* can be written like in Figure 30. Notice that the performance of the >>> combinator is essential for the performance of the skeleton. A definitive version would use Futures.

OL: it appears STRANGE. are the data really left alone and noded after map and taken from there by reduce? It makes sense only is no communication through master takes place. ELSE: CUT!

MB: this requires some work. Either change this to use futures or cut, yes. OL: now rewritten as motivation for futures. maybe still cut?

<sup>6</sup> Available on Hackage under https://hackage.haskell.org/package/edenskel-2.1.0.0/docs/src/Control-Parallel-Eden-Map.html.



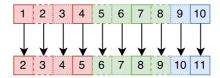


Figure 29: Schematic depiction of farmChunk.

```
parMapReduceDirect:: (ArrowParallel \ arr \ [a] \ b \ conf, ArrowApply \ arr, ArrowChoice \ arr) \Rightarrow conf \rightarrow ChunkSize \rightarrow arr \ a \ b \rightarrow arr \ (b,b) \ b \rightarrow b \rightarrow arr \ [a] \ b parMapReduceDirect \ conf \ chunkSize \ mapfin \ foldfin \ neutral = arr \ (chunkSof \ chunkSize) >>> parMap \ conf \ (mapArr \ mapfin >>> foldlArr \ foldfin \ neutral) >>> foldlArr \ foldfin \ neutral
```

Figure 30: Definition of parMapReduceDirect.

# 7 Topological Skeletons

Even though many algorithms can be expressed by parallel maps, some problems require more sophisticated skeletons. The Eden library leverages this problem and already comes<sup>7</sup> with more predefined skeletons, among them a *pipe*, a *ring*, and a *torus* implementations (Loogen, 2012). These seem like reasonable candidates to be ported to our Arrow-based parallel Haskell. We aim to showcase that we can express more sophisticated skeletons with Parallel Arrows as well.

## 7.1 Parallel pipe

The parallel *pipe* skeleton is semantically equivalent to folding over a list  $[arr\ a\ a]$  of arrows with >>>, but does this in parallel, meaning that the arrows do not have to reside on the same thread/machine. We implement this skeleton using the ArrowLoop typeclass which gives us the  $loop::arr\ (a,b)\ (c,b)\to arr\ a\ c$  combinator which allows us to express recursive fix-point computations in which output values are fed back as input. For example MB: das kann man hier so lassen, oder?OL: sicherlich!

```
loop (arr (\lambda(a,b) \rightarrow (b,a:b))) which is the same as loop (arr snd \&\&\& arr (uncurry (:)))
```

defines an arrow that takes its input a and converts it into an infinite stream [a] of it. Using this to our advantage gives us a first draft of a pipe implementation (Fig. 31) by plugging in the parallel evaluation call parEvalN conf fs inside the second argument of &&& and then only picking the first element of the resulting list with arr last.

Available on Hackage: https://hackage.haskell.org/package/edenskel-2.0.0.2/docs/Control-Parallel-Eden-Topology.html.

```
\begin{aligned} &pipeSimple :: (ArrowLoop\ arr, ArrowParallel\ arr\ a\ a\ conf) \Rightarrow \\ &conf \rightarrow [arr\ a\ a] \rightarrow arr\ a\ a \\ &pipeSimple\ conf\ fs = \\ &loop\ (arr\ snd\ \&\&\& \\ &(arr\ (uncurry\ (:) >>> lazy) >>> parEvalN\ conf\ fs)) >>> \\ &arr\ last \\ &lazy :: (Arrow\ arr) \Rightarrow arr\ [a]\ [a] \\ &lazy = arr\ (\lambda \sim (x:xs) \rightarrow x: lazy\ xs) \end{aligned}
```

Figure 31: A first implementation of the *pipe* skeleton expressed with Parallel Arrows. Note that the use of *lazy* is essential as without it programs using this definition would never halt. We need to enforce that the evaluation of the input [a] terminates before passing it into *parEvalN*.

However, using this definition directly will make the master node a potential bottleneck in distributed environments as described in Section 5. Therefore, we introduce a more sophisticated version that internally uses Futures and get the final definition of *pipe*:

```
pipe :: (ArrowLoop arr, ArrowParallel arr (fut a) (fut a) conf, Future fut a) \Rightarrow conf \rightarrow [arr a a] \rightarrow arr a a pipe conf fs = unliftFut (pipeSimple conf (map liftFut fs))
```

Sometimes, this *pipe* definition can be a bit inconvenient, especially if we want to pipe arrows of mixed types together, i.e. *arr* a b and *arr* b c. By wrapping these two arrows inside a common type we obtain *pipe2* (Fig. 32).

#### **OL:** I swapped the type classes here:

```
\begin{array}{l} \textit{pipe2} :: (ArrowLoop\ arr, ArrowChoice\ arr, Future\ fut\ (([a],[b]),[c]),\\ ArrowParallel\ arr\ (fut\ (([a],[b]),[c]))\ (fut\ (([a],[b]),[c]))\ conf) \Rightarrow\\ conf \to arr\ a\ b \to arr\ b\ c \to arr\ a\ c\\ \textit{pipe2}\ conf\ f\ g =\\ (arr\ return\ \&\&\ arr\ (const\ []))\ \&\&\&\ arr\ (const\ [])>>>\\ \textit{pipe}\ conf\ (replicate\ 2\ (unify\ f\ g))>>>\\ \textit{arr\ snd}>>> \textit{arr\ head\ where}\\ \textit{unify}:: (ArrowChoice\ arr) \Rightarrow \textit{arr\ a}\ b \to \textit{arr\ b}\ c \to \textit{arr\ }(([a],[b]),[c])\ (([a],[b]),[c])\\ \textit{unify}\ f\ g =\\ (\textit{mapArr\ f}\ ***\textit{mapArr\ g}) ***\textit{arr\ }(\setminus \to [])>>>\\ \textit{arr\ }(\lambda((a,b),c) \to ((c,a),b)) \end{array}
```

Figure 32: Definition of pipe2.

Note that extensive use of this combinator over *pipe* with a hand-written combination data type will probably result in worse performance because of more communication overhead from the many calls to parEvalN. Nonetheless, we can define a parallel piping operator *parcomp* (Fig. 33, which is semantically equivalent to >>> similarly to other parallel syntactic sugar from Section C.

#### Arrows for Parallel Computations

## OL: swapped type classes

```
(|>>>|) :: (ArrowLoop\ arr, ArrowChoice\ arr, Future\ fut\ (([a],[b]),[c]), ArrowParallel\ arr\ (fut\ (([a],[b]),[c]))\ (fut\ (([a],[b]),[c]))\ ()) \Rightarrow arr\ a\ b \to arr\ b\ c \to arr\ a\ c \\ (|>>>|) = pipe2\ ()
```

Figure 33: Definition of parcomp.

## 7.2 Ring skeleton

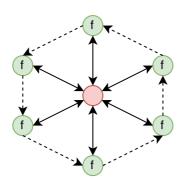


Figure 34: Schematic depiction of the ring skeleton.

Eden comes with a ring skeleton<sup>8</sup> (Fig. 34) implementation that allows the computation of a function  $[i] \rightarrow [o]$  with a ring of nodes that communicate in a ring topology with each other. Its input is a node function  $i \rightarrow r \rightarrow (o, r)$  in which r serves as the intermediary output that gets send to the neighbour of each node. This data is sent over direct communication channels, so called 'remote data'. We depict it in Appendix, Fig. B 3.

We can rewrite this functionality easily with the use of *loop* as the definition of the node function, arr(i,r)(o,r), after being transformed into an arrow, already fits quite neatly into the *loop*'s  $arr(a,b)(c,b) \rightarrow arr\ a\ c$ . In each iteration we start by rotating the intermediary input from the nodes  $[fut\ r]$  with  $second\ (rightRotate >>> lazy)$ . Similarly to the pipe from Chapter 7.1 (Fig. 31), we have to feed the intermediary input into our lazy arrow here, or the evaluation would hang. **OL:** meh, wording The reasoning is explained by Loogen (2012):

Note that the list of ring inputs ringIns is the same as the list of ring outputs ringOuts rotated by one element to the right using the auxiliary function rightRotate. Thus, the program would get stuck without the lazy pattern, because the ring input will only be produced after process creation and process creation will not occur without the first input. Next, we zip the resulting  $([i],[fut\,r])$  to  $[(i,fut\,r)]$  with arr (uncurry zip) so we can feed that into a our input arrow arr (i,r) (o,r), which we transform into arr  $(i,fut\,r)$   $(o,fut\,r)$  before lifting it to arr  $[(i,fut\,r)]$   $[(o,fut\,r)]$  to get a list  $[(o,fut\,r)]$ . Finally we unzip this list into  $([o],[fut\,r])$ . Plugging this arrow arr  $([i],[fut\,r])$   $([o],fut\,r)$  into the definition of loop

<sup>8</sup> Available on Hackage: https://hackage.haskell.org/package/edenskel-2.0.0.2/docs/Control-Parallel-Eden-Topology.html

from earlier gives us arr[i][o], our ring arrow (Fig. 35). This combinator can, for example, be used to calculate the shortest paths in a graph using Warshall's algorithm. **OL: let's do it?** 

```
ring:: (ArrowLoop arr, Future fut r, ArrowParallel arr (i, fut r) (o, fut r) conf) \Rightarrow conf \rightarrow arr (i,r) (o,r) \rightarrow arr [i] [o] ring conf f = loop (second (rightRotate >>> lazy) >>> arr (uncurry zip) >>> parMap conf (second get >>> f >>> second put) >>> arr unzip) rightRotate:: (Arrow arr) \Rightarrow arr [a] [a] rightRotate = arr \$ \lambda list \rightarrow case list of [] \rightarrow [] xs \rightarrow last xs: init xs lazy:: (Arrow arr) \Rightarrow arr [a] [a] lazy = arr (\lambda~(x:xs) \Rightarrow x: lazy xs)
```

Figure 35: Final definition of the *ring* skeleton.

#### 7.3 Torus skeleton

If we take the concept of a ring from 7.2 one dimension further, we get a torus (Fig. 36, 37). Every node sends ands receives data from horizontal and vertical neighbours in each communication round. With our Parallel Arrows we re-implement the *torus* combinator<sup>9</sup> from Eden—yet again with the help of the *ArrowLoop* typeclass.

Similar to the *ring*, we once again start by rotating the input, but this time not only in one direction, but in two. This means that the intermediary input from the neighbour nodes has to be stored in a tuple ([ $[fut\ a]]$ , [ $[fut\ b]$ ]) in the second argument (loop only allows for two arguments) of our looped arrow arr ([[c]], ([ $[fut\ a]]$ , [ $[fut\ b]$ ])) ([[d]], ([ $[fut\ a]$ ], [ $[fut\ b]$ ])) and our rotation arrow becomes second ( $(mapArr\ rightRotate >>> lazy)***(<math>arr\ rightRotate >>> lazy)****(<math>arr\ rightRotate >>> lazy)***(<math>arr\ rightRotate >>> lazy)**(<math>arr\ rightRotate >>> lazy)***(<math>arr\ rightRotate >>> lazy)**(<math>arr\ rightRotate >>> lazy)**(<math>arr\ rightRotate >>> lazy)**(<math>arr\ rightRotate >>> lazy)**(<math>arr\ rightRotate >> lazy)**(<math>arr\ rightRo$ 

<sup>9</sup> Available on Hackage: https://hackage.haskell.org/package/edenskel-2.0.0.2/docs/Control-Parallel-Eden-Topology.html.

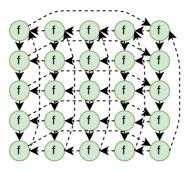


Figure 36: Schematic depiction of the torus skeleton.

[lazy)) instead of the singular rotation in the ring as we rotate  $[[fut\ a]]$  horizontally and  $[[fut\ b]]$  vertically. Then, we once again zip the inputs for the input arrow with  $arr\ (uncurry3\ zipWith3\ lazyzip3)$  from  $([[c]],([[fut\ a]],[[fut\ b]]))$  to  $[[(c,fut\ a,fut\ b)]]$ , which we then feed into our parallel execution.

This, however, is more complicated than in the ring case as we have one more dimension of inputs to be transformed. We first have to *shuffle* all the inputs to then pass it into *parMap conf* (*ptorus f*) which yields us  $[(d, fut\ a, fut\ b)]$ . We can then unpack this shuffled list back to its original ordering by feeding it into the specific unshuffle arrow we created one step earlier with  $arr\ length >>> arr\ unshuffle$  with the use of  $app::arr\ (arr\ a\ b,a)\ c$  from the ArrowApply typeclass. Finally, we unpack this matrix  $[[[(d, fut\ a, fut\ b)]]$  with  $arr\ (map\ unzip3) >>> arr\ unzip3 >>> threetotwo\ to\ get\ ([[d]], ([[fut\ a]], [[fut\ b]])).$ 

#### OL: swapped type classes

```
torus:: (ArrowLoop arr, ArrowChoice arr, ArrowApply arr, Future fut a, Future fut b,
         ArrowParallel arr (c, \text{fut } a, \text{fut } b) (d, \text{fut } a, \text{fut } b) \text{ conf}) \Rightarrow
         conf \rightarrow arr\left(c,a,b\right)\left(d,a,b\right) \rightarrow arr\left[\left[c\right]\right]\left[\left[d\right]\right]
torus conf f =
         loop\ (second\ ((mapArr\ rightRotate >>> lazy) *** (arr\ rightRotate >>> lazy)) >>> lazy) *** (arr\ rightRotate >>> lazy)) >>> lazy) *** (arr\ rightRotate >>> lazy)) >>> lazy) >>> lazy) >>> lazy) *** (arr\ rightRotate >>> lazy)) >>> lazy) >>> lazy) >>> lazy) *** (arr\ rightRotate >>> lazy)) >>> lazy) >> lazy) >>> lazy) >> lazy) >>> lazy) >> lazy) >> lazy) >> lazy) >> lazy) >>> lazy) >>> lazy) >> lazy) >> lazy) >>> lazy) >>> lazy) >> lazy) >>> lazy) >> lazy) >>> lazy) >> lazy) >>
         arr (uncurry3 (zipWith3 lazyzip3))>>>
         (arr length >>> arr unshuffle) &&& (shuffle >>> parMap conf (ptorus f)) >>> app >>>
         arr\left(map\ unzip3\right)>>> arr\ unzip3>>> threetotwo)
ptorus :: (Arrow \ arr, Future \ fut \ a, Future \ fut \ b) \Rightarrow
         arr(c,a,b)(d,a,b) \rightarrow arr(c,fut\ a,fut\ b)(d,fut\ a,fut\ b)
ptorus f = arr (\lambda \sim (c, a, b) \rightarrow (c, get \ a, get \ b)) >>> f >>> arr (\lambda \sim (d, a, b) \rightarrow (d, put \ a, put \ b))
uncurry3::(a \rightarrow b \rightarrow c \rightarrow d) \rightarrow (a,(b,c)) \rightarrow d
uncurry3f(a,(b,c)) = f a b c
lazyzip3::[a] \rightarrow [b] \rightarrow [c] \rightarrow [(a,b,c)]
lazyzip3 as bs cs = zip3 as (lazy bs) (lazy cs)
threetotwo::(Arrow\ arr) \Rightarrow arr\ (a,b,c)\ (a,(b,c))
threetotwo = arr \$ \lambda \sim (a, b, c) \rightarrow (a, (b, c))
```

Figure 37: Definition of the torus skeleton.

As an example of using this skeleton (Loogen, 2012) showed the matrix multiplication using the Gentleman algorithm (Gentleman, 1978). Their instantiation of the skeleton *nodefunction* can be adapted as shown in Fig. 38. If we compare the trace from a call using

```
nodefunction::Int \rightarrow ((Matrix,Matrix),[Matrix],[Matrix]) \rightarrow ([Matrix],[Matrix],[Matrix]) \\ nodefunction \ n \ ((bA,bB),rows,cols) = ([bSum],bA:nextAs,bB:nextBs) \\ \textbf{where} \ bSum = foldl' \ matAdd \ (matMult \ bA \ bB) \ (zipWith \ matMult \ nextAs \ nextBs) \\ nextAs = take \ (n-1) \ rows \\ nextBs = take \ (n-1) \ cols
```

Figure 38: Adapted *nodefunction* for matrix multiplication with the *torus* from Fig. 37.

our arrow definition of the torus (Fig. 39) with the Eden version (Fig. 40) we can see that the behaviour of the arrow version is comparable. OL: much more details on this!

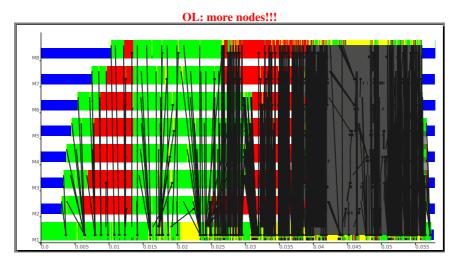


Figure 39: Matrix Multiplication with a torus (PArrows).

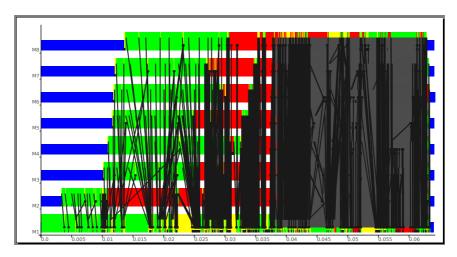
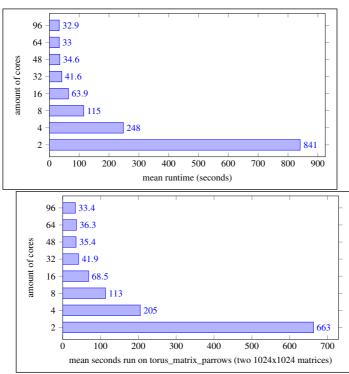


Figure 40: Matrix Multiplication with a torus (Eden).

# Arrows for Parallel Computations

## 8 Benchmarks



## 9 Conclusion

Arrows are a generic concept that allows for powerful composition combinators. To our knowledge we are the first ones to represent parallel computation with arrows. OL: that strange arrows-based robot interaction paper from 1993 or so! clearly discuss in related work done!

Arrows turn out to be a useful tool for composing in parallel programs. We do not have to introduce new monadic types that wrap the computation. Instead use arrows just like regular sequential pure functions. This work features multiple parallel backends: the already available parallel Haskell flavours. Parallel Arrows feature an implementation of the *ArrowParallel* interface for Multicore Haskell, *Par* Monad, and Eden. With our approach parallel programs can be ported across these flavours with no effort. Performancewise, Parallel Arrows are on par with existing parallel Haskells, as they do not introduce any notable overhead. OL: PROOFS. Many proofs in benchmarks!

MB: ArrowApply (or equivalent) are needed because we basically want to be able to produce intermediary results, this is by definition of the parallel evaluation combinators

OL: Remove websites from citations, put them into footnotes!

OL: Parrows + accelerate = love? Metion port to Frege. Mention the Par monad troubles.

# 9.1 Future Work

Our PArrows interface can be expanded to futher parallel Haskells. More specifically we target HdpH (Maier *et al.*, 2014) a modern distributed Haskell that would benefit from our Arrows notation. Future-aware special versions of Arrow combinated can be extended and further improved. We would look into more transparency of the API, it should basically infuse as little overhead as possible.

Of course, definitions of further skeletons are viable and needed. We are looking into more experiences with seamless porting of parallel PArrow-based programs across the backends.

Accelerate (Chakravarty *et al.*, 2011) is not related to our approach. It would be interesting to see a hybrid of both APIs.

OL: replace API with DSL globally?

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### **A Utility Functions**

To be able to go into detail on parallel arrows, we introduce some utility combinators first, that will help us later: map, foldl and zipWith on arrows.

The mapArr combinator (Fig. A 1) lifts any arrow arr a b to an arrow arr [a][b] (Hughes, 2005b). Similarly, we can also define foldlArr (Fig. A 2) that lifts any arrow arr (b, a) b with a neutral element b to arr [a] b.

```
mapArr :: ArrowChoice \ arr \Rightarrow arr \ a \ b \rightarrow arr \ [a] \ [b]
mapArrf =
  arr listcase >>>
  arr (const []) ||| (f * * * mapArr f >>> arr (uncurry (:)))
listcase [] = Left ()
listcase(x:xs) = Right(x,xs)
```

Figure A 1: The definition of *map* over Arrows and the *listcase* helper function.

Finally, with the help of mapArr (Fig. A 1), we can define zipWithArr (Fig. A 3) that lifts any arrow arr(a,b) c to an arrow arr([a],[b]) [c].

These combinators make use of the ArrowChoice type class which provides the combinator. It takes two arrows arr a c and arr b c and combines them into a new arrow

```
foldlArr :: (ArrowChoice\ arr, ArrowApply\ arr) \Rightarrow arr\ (b,a)\ b \rightarrow b \rightarrow arr\ [a]\ b
foldlArrfb =
    arr listcase>>>
    arr (const b) |||
        (\textit{first} \; (\textit{arr} \; (\lambda a \rightarrow (b, a)) >>> f >>> \textit{arr} \; (\textit{foldlArr} \; f)) >>> \textit{app})
```

Figure A 2: The definition of *foldl* over Arrows.

```
zipWithArr::ArrowChoice\ arr \Rightarrow arr\ (a,b)\ c \rightarrow arr\ ([a],[b])\ [c]
zipWithArrf = (arr \$ \lambda(as,bs) \rightarrow zipWith(,) as bs) >>> mapArrf
```

Figure A 3: *zipWith* over arrows.

arr (Either a b) c which pipes all Left a's to the first arrow and all Right b's to the second arrow.

```
(|||) :: Arrow Choice arr a c \rightarrow arr b c \rightarrow arr (Either a b) c
```

With the zipWithArr combinator we can also write a combinator *listApp*, that lifts a list of arrows  $[arr \ a \ b]$  to an arrow  $arr \ [a] \ [b]$ .

```
listApp :: (ArrowChoice\ arr, ArrowApply\ arr) \Rightarrow [arr\ a\ b] \rightarrow arr\ [a]\ [b]
listApp\ fs = (arr \$ \lambda as \rightarrow (fs, as)) >>> zipWithArr\ app
```

Note that this additionally makes use of the ArrowApply typeclass that allows us to evaluate arrows with  $app :: arr (arr \ a \ b, a) \ c$ .

## **B** Omitted Funtion Definitions

We have omitted some function definitions in the main text for brevity, and redeem this here. We warp Eden's build-in Futures in PArrows as in Figure B 1. The full definition of farmChunk is in Figure B 2. Eden definition of ring skeleton following (Loogen, 2012) is in Figure B 3.

```
data RemoteData \ a = RD \{ rd :: RD \ a \}
instance (Trans\ a) \Rightarrow Future\ RemoteData\ a\ where
  put = arr (\lambda a \rightarrow RD \{ rd = release a \})
  get = arr rd >>> arr fetch
```

Figure B 1: RD-based RemoteData version of Future for the Eden backend.

Furthermore, parEval2 (Fig. B 4) is achieved as follows: We start by transforming the (a,c) input into a two-element list [Either a c] by first tagging the two inputs with Left and Right and wrapping the right element in a singleton list with return so that we can combine them with arr (uncurry (:)). Next, we feed this list into a parallel arrow running on two instances of f + + + g as described above. After the calculation is finished, we convert the

```
\begin{split} & \textit{farmChunk} :: (\textit{ArrowParallel arr a b conf}, \textit{ArrowParallel arr } [a] \ [b] \ \textit{conf}, \\ & \textit{ArrowChoice arr}, \textit{ArrowApply arr}) \Rightarrow \\ & \textit{conf} \rightarrow \textit{ChunkSize} \rightarrow \textit{NumCores} \rightarrow \textit{arr a b} \rightarrow \textit{arr } [a] \ [b] \\ & \textit{farmChunk conf chunkSize numCores} f = \\ & \textit{unshuffle numCores} >>> \\ & \textit{parEvalNLazy conf chunkSize (repeat (mapArrf))} >>> \\ & \textit{shuffle} \end{split}
```

Figure B 2: Definition of farmChunk.

```
\begin{array}{l} \textit{ringSimple} :: (\textit{Trans } i, \textit{Trans } o, \textit{Trans } r) \Rightarrow (i \rightarrow r \rightarrow (o, r)) \rightarrow [i] \rightarrow [o] \\ \textit{ringSimple } f \textit{ is} = os \\ \textbf{where} \textit{ } (os, \textit{ringOuts}) = \textit{unzip } \textit{ } (\textit{parMap } \textit{ } (\textit{toRD} \$ \textit{uncurry } f) \textit{ } (\textit{zip } \textit{is} \$ \textit{lazy ringIns})) \\ \textit{ringIns} = \textit{rightRotate ringOuts} \\ \textit{toRD} :: (\textit{Trans } i, \textit{Trans } o, \textit{Trans } r) \Rightarrow ((i, r) \rightarrow (o, r)) \rightarrow ((i, RD \ r) \rightarrow (o, RD \ r)) \\ \textit{toRD} \textit{ } f \textit{ } (i, \textit{ringIn}) = (o, \textit{release ringOut}) \\ \textbf{where} \textit{ } (o, \textit{ringOut}) = f \textit{ } (i, \textit{fetch ringIn}) \\ \textit{rightRotate} :: [a] \rightarrow [a] \\ \textit{rightRotate } xs = \textit{last } xs : \textit{init } xs \\ \textit{lazy} :: [a] \rightarrow [a] \\ \textit{lazy} \sim (x : xs) = x : \textit{lazy } xs \\ \end{array}
```

Figure B 3: Eden's definition of the *ring* skeleton.

resulting [Either b d] into ([b],[d]) with arr partitionEithers. The two lists in this tuple contain only one element each by construction, so we can finally just convert the tuple to (b,d) in the last step.

```
parEval2 :: (ArrowChoice\ arr, \\ ArrowParallel\ arr\ (Either\ a\ c)\ (Either\ b\ d)\ conf) \Rightarrow \\ conf \to arr\ a\ b \to arr\ c\ d \to arr\ (a,c)\ (b,d) \\ parEval2\ conf\ f\ g = \\ arr\ Left *** (arr\ Right >>> arr\ return) >>> \\ arr\ (uncurry\ (:)) >>> \\ parEvalN\ conf\ (replicate\ 2\ (f+++g)) >>> \\ arr\ partitionEithers >>> \\ arr\ head *** arr\ head
```

Figure B 4: Definition of parEval2.

# C Syntactic Sugar

For basic arrows, we have the \*\*\* combinator (Fig. 10) which allows us to combine two arrows  $arr\ a\ b$  and  $arr\ c\ d$  into an arrow  $arr\ (a,c)\ (b,d)$  which does both computations at

once. This can easily be translated into a parallel version \*\*\* with the use of *parEval2*, but for this we require a backend which has an implementation that does not require any configuration (hence the () as the *conf* parameter):

```
(|***|):: (ArrowChoice\ arr, ArrowParallel\ arr\ (Either\ a\ c)\ (Either\ b\ d)\ ())) \Rightarrow arr\ a\ b \to arr\ c\ d \to arr\ (a,c)\ (b,d) (|***|) = parEval2\ ()
```

We define the parallel &&& in a similar manner to its sequential pendant &&& (Fig. 10):

```
(|\&\&\&|) :: (ArrowChoice\ arr, ArrowParallel\ arr\ (Either\ a\ a)\ (Either\ b\ c)\ ()) \Rightarrow \\ arr\ a\ b \to arr\ a\ c \to arr\ a\ (b,c) \\ (|\&\&\&|)\ f\ g = (arr\,\$\,\lambda a \to (a,a)) >>> f\ |\ *** |\ g
```