Arrows for Parallel Computation

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Abstract

Arrows are a general interface for computation and an alternative to Monads for API design. In contrast to Monad-based parallelism, we explore the use of Arrows for specifying generalised parallelism. Specifically, we define an Arrow-based language and implement it using multiple parallel Haskells.

As each parallel computation is an Arrow, such parallel Arrows (PArrows) can be readily composed and transformed as such. To allow for more sophisticated communication schemes between computation nodes in distributed systems, we utilise the concept of Futures to wrap direct communication.

To show that PArrows have similar expressive power as existing parallel languages, we implement several algorithmic skeletons and four benchmarks. Benchmarks show that our framework does not induce any notable performance overhead. We conclude that Arrows have considerable potential for composing parallel programs and for producing programs that can execute on multiple parallel language implementations.

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1 Introduction

Parallel functional languages have a long history of being used for experimenting with novel parallel programming paradigms. Haskell, which we focus on in this paper, has several mature implementations. We regard here in-depth Glasgow parallel Haskell or short GpH (its Multicore SMP implementation, in particular), the *Par* Monad, and Eden, a distributed memory parallel Haskell. These languages represent orthogonal approaches. Some use a Monad, even if only for the internal representation. Some introduce additional language constructs. Section 3.2 gives a short overview over these languages.

A key novelty in this paper is to use Arrows to represent parallel computations. They seem a natural fit as they can be thought of as a more general function arrow (\rightarrow) and serve as general interface to computations while not being as restrictive as Monads (Hughes, 2000). Section 3.1 gives a short introduction to Arrows.

We provide an Arrows-based type class and implementations for the three above mentioned parallel Haskells. Instead of introducing a new low-level parallel backend to implement our Arrows-based interface, we define a shallow-embedded DSL for Arrows. This DSL is defined as a common interface with varying implementations in the existing parallel Haskells. Thus, we not only define a parallel programming interface in a novel manner – we tame the zoo of parallel Haskells. We provide a common, very low-penalty programming interface that allows to switch the parallel implementations at will. The induced penalty is in the single-digit percent range, with means typically under 2% overhead in measurements over the varying cores configuration (Section 7). Further implementations, based on HdpH or a Frege implementation (on the Java Virtual Machine), are viable, too.

Contributions. We propose an Arrow-based encoding for parallelism based on a new Arrow combinator $parEvalN :: [arr\ a\ b] \to arr\ [a]\ [b]$. A parallel Arrow is still an Arrow, hence the resulting parallel Arrow can still be used in the same way as a potential sequential version. In this paper we evaluate the expressive power of such a formalism in the context of parallel programming.

- We introduce a parallel evaluation formalism using Arrows. One big advantage of this specific approach is that we do not have to introduce any new types, facilitating composability (Section 4).
- We show that PArrow programs can readily exploit multiple parallel language implementations. We demonstrate the use of GpH, a *Par* Monad, and Eden. We do not re-implement all the parallel internals, as this functionality is hosted in the *ArrowParallel* type class, which abstracts all parallel implementation logic. The implementations can easily be swapped, so we are not bound to any specific one.

This has many practical advantages. For example, during development we can run the program in a simple GHC-compiled variant using GpH and afterwards deploy it on a cluster by converting it into an Eden program, by just replacing the *ArrowParallel* instance and compiling with Eden's GHC variant (Section 4).

- We extend the PArrows formalism with *Futures* to enable direct communication of data between nodes in a distributed memory setting similar to Eden's Remote Data (*RD*, Dieterle *et al.*, 2010a). Direct communication is useful in a distributed memory setting because it allows for inter-node communication without blocking the master-node. (Section 5)
- We demonstrate the expressiveness of PArrows by using them to define common algorithmic skeletons (Section 6), and by using these skeletons to implement four benchmarks (Section 7).
- We practically demonstrate that Arrow parallelism has a low performance overhead compared with existing approaches, e.g. the mean over all cores of relative mean overhead was less than 3.5% and less than 0.8% for all benchmarks with GpH and Eden, respectively. As for *Par* Monad, the mean of mean overheads was in favour of PArrows in all benchmarks (Section 7).

PArrows are open source and are available from https://github.com/s4ke/Parrows.

2 Related Work

Parallel Haskells. The non-strict semantics of Haskell, and the fact that reduction encapsulates computations as closures, makes it relatively easy to define alternate parallelisations. A range of approaches have been explored, including data parallelism (Chakravarty *et al.*, 2007; Keller *et al.*, 2010), GPU-based approaches (Mainland & Morrisett, 2010; Svensson, 2011), software transactional memory (Harris *et al.*, 2005; Perfumo *et al.*, 2008). The Haskell–GPU bridge Accelerate (Chakravarty *et al.*, 2011; Clifton-Everest *et al.*, 2014; McDonell *et al.*, 2015) is completely orthogonal to our approach. A good survey of parallel Haskells can be found in Marlow (2013).

Our PArrow implementation uses three task parallel languages as backends: the GpH (Trinder *et al.*, 1996, 1998) parallel Haskell dialect and its multicore version (Marlow *et al.*, 2009), the *Par* Monad (Marlow *et al.*, 2011; Foltzer *et al.*, 2012), and Eden (Loogen *et al.*, 2005; Loogen, 2012). These languages are under active development, for example a combined shared and distributed memory implementation of GpH is available (Aljabri *et al.*, 2014, 2015). Research on Eden includes low-level 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). The definitions of new Eden skeletons is a specific focus (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).

Other task parallel Haskells related to Eden, GpH, and the *Par* Monad include the following. 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.

Algorithmic skeletons. Algorithmic skeletons were introduced by Cole (1989). Early publications on this topic include (Danelutto *et al.*, 1992; Darlington *et al.*, 1993; Botorog & Kuchen, 1996; Lengauer *et al.*, 1997; Gorlatch, 1998). Rabhi & Gorlatch (2003) consolidated early reports on high-level programming approaches. Types of algorithmic skeletons include *map-*, *fold-*, and *scan-*based parallel programming patterns, special applications such as divide-and-conquer or topological skeletons.

The *farm* skeleton (Hey, 1990; Peña & Rubio, 2001; Poldner & Kuchen, 2005) is a statically task-balanced parallel *map*. When tasks' durations cannot be foreseen, a dynamic load balancing (*workpool*) brings a lot of improvement (Rudolph *et al.*, 1991; Hammond *et al.*, 2003; Hippold & Rünger, 2006; Berthold *et al.*, 2008; Marlow *et al.*, 2009). For special tasks *workpool* skeletons can be extended with dynamic task creation (Priebe, 2006; Dinan *et al.*, 2009; Brown & Hammond, 2010). Efficient load-balancing schemes for *workpools* are subject of research (Blumofe & Leiserson, 1999; Acar *et al.*, 2000; van Nieuwpoort *et al.*, 2001; Chase & Lev, 2005; Olivier & Prins, 2008; Michael *et al.*, 2009). The *fold* (or *reduce*) skeleton was implemented in various skeleton libraries (Kuchen, 2002; Karasawa & Iwasaki, 2009; Buono *et al.*, 2010; Dastgeer *et al.*, 2011), as also its inverse, *scan* (Bischof & Gorlatch, 2002; Harris *et al.*, 2007). Google *map-reduce* (Dean & Ghemawat, 2008, 2010) is more special than just a composition of the two skeletons (Lämmel, 2008; Berthold *et al.*, 2009b).

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.

Arrows. Arrows were introduced by Hughes (2000) as a less restrictive alternative to Monads, in essence they are a generalised function arrow \rightarrow . Hughes (2005) presents a tutorial on Arrows. Jacobs et al. (2009); Lindley et al. (2011); Atkey (2011) develop theoretical background of Arrows. 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 probably most prominent application of Arrows is Arrow-based functional reactive programming, AFRP (Nilsson et al., 2002; Hudak et al., 2003; Czaplicki & Chong, 2013). Liu et al. (2009) formally define a more special kind of Arrows that capsule the computation more than regular Arrows do and thus enable optimisations. 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. Huang et al. (2007) utilise Arrows for parallelism, but strikingly different from our approach. They use Arrows to orchestrate several tasks in robotics. We, however, propose a general interface for parallel programming, while remaining completely in Haskell.

Arrows in 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. Basic definitions of PArrows are possible in the Frege language¹ (which is basically Haskell on the JVM). However, they are beyond the scope of this work, as are similar experiments with the Eta language², a new approach to Haskell on the JVM.

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.

3 Background

This section gives a short overview of Arrows (Section 3.1) and of GpH, the *Par* Monad, and Eden, the three parallel Haskells which we base our DSL on (Section 3.2).

3.1 Arrows

Arrows were introduced by Hughes (2000) as a general interface for computation and a less restrictive generalisation of Monads. Hughes motivates the broader interface of Arrows with the example of a parser with added static meta-information that can not satisfy the monadic bind operator (\gg):: $m \, a \to (a \to m \, b) \to m \, b$ (with $m \to m \, b$).

An Arrow $arr\ a\ b$ represents a computation that converts an input a to an output b. This is defined in the Arrow type class shown in Fig. 1. To lift an ordinary function to an Arrow, arr is used, analogous to the monadic return. Similarly, the composition 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 $arr\ a\ b$ and converts it into an Arrow on pairs $arr\ (a,c)\ (b,c)$ that leaves the second argument untouched. It allows us to to save input across Arrows. Figure 2 shows a graphical representation of these basic Arrow combinators. The most prominent instances of this interface are regular functions (\rightarrow) and the Kleisli type (Fig. 1), which wraps monadic functions, e.g. $a \rightarrow m\ b$.

Hughes also defined some syntactic sugar (Fig. 3): *second*, *** and &&&. *second* is the mirrored version of *first* (Appendix A). The *** function combines *first* and *second* to handle two inputs in one arrow, and is defined as follows:

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

The &&& combinator, which constructs an Arrow that outputs two different values like ***, but takes only one input, is:

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

¹ GitHub project page at https://github.com/Frege/frege

² Eta project page at http://eta-lang.org

In the example a parser of the type *Parser s a* with static meta information *s* and result *a* is shown to not be able to use the static information *s* without applying the monadic function $a \to m b$. With Arrows this is possible.

```
class Arrow\ arr\ where arr::(a \rightarrow b) \rightarrow arr\ a\ b (>>>)::arr\ a\ b \rightarrow arr\ b\ c \rightarrow arr\ a\ c first::arr\ a\ b \rightarrow arr\ (a,c)\ (b,c) instance Arrow\ (\rightarrow) where arr\ f=f f>>>> g=g\circ f first\ f=\lambda(a,c) \rightarrow (f\ a,c) 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))
```

Figure 1: The *Arrow* type class and its two most typical instances.

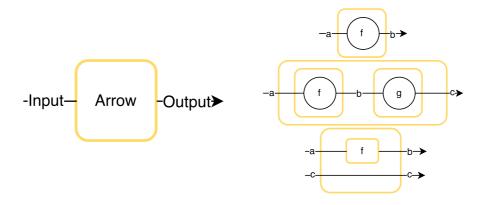


Figure 2: Schematic depiction of an Arrow (left) and its basic combinators *arr*, >>> and *first* (right).

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)
```

As we can rewrite the monadic bind operation (\gg) with only the Kleisli type into $m \ a \to Kleisli \ m \ a \ b \to m \ b$, but not with a general Arrow $arr \ a \ b$, we can intuitively get an idea of why Arrows must be a generalisation of Monads. While this also means that a general Arrow can not express everything a Monad can, Hughes (2000) shows in his parser example that this trade-off is worth it in some cases.

In this paper we will show that parallel computations can be expressed with this more general interface of Arrows without requiring Monads. We also do not restrict the compatible Arrows to ones which have *ArrowApply* instances but instead only require instances for *ArrowChoice* (for if-then-else constructs) and *ArrowLoop* (for looping). Because of this, we have a truly more general interface as compared to a monadic one.

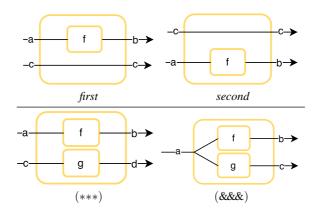


Figure 3: Visual depiction of syntactic sugar for Arrows.

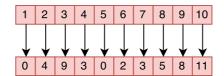


Figure 4: Schematic illustration of *parEvalN*. A list of inputs is transformed by different functions in parallel.

While we could have based our DSL on Profunctors as well, we chose Arrows for this paper since they they allow for a more direct way of thinking about parallelism than general Profunctors because of their composability. However, they are a promising candidate for future improvements of our DSL. Some Profunctors, especially ones supporting a composition operation, choice, and looping, can already be adapted to our interface as shown in Appendix B.

3.2 Short introduction to parallel Haskells

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 4 symbolically shows.

In this section, we will implement this non-Arrow version which will later be adapted for usage in our Arrow-based parallel Haskell.

There exist several parallel Haskells already. Among the most important are probably GpH (based on *par* and *pseq* 'hints', Trinder *et al.*, 1996, 1998), the *Par* Monad (a monad for deterministic parallelism, Marlow *et al.*, 2011; Foltzer *et al.*, 2012), Eden (a parallel Haskell for distributed memory, Loogen *et al.*, 2005; Loogen, 2012), HdpH (a Template Haskell-based parallel Haskell for distributed memory, Maier *et al.*, 2014; Stewart *et al.*, 2016) and LVish (a *Par* extension with focus on communication, Kuper *et al.*, 2014).

As the goal of this paper is not to re-implement yet another parallel runtime, but to represent parallelism with Arrows, we base our efforts on existing work which we wrap as

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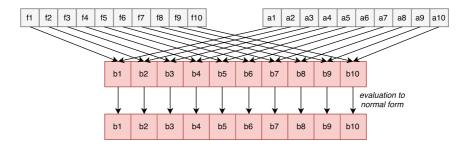


Figure 5: parEvalN (GpH).

backends behind a common interface. For this paper we chose GpH for its simplicity, the *Par* Monad to represent a monadic DSL, and Eden as a distributed parallel Haskell.

LVish and HdpH were not chosen as the former does not differ from the original *Par* Monad with regard to how we would have used it in this paper, while the latter (at least in its current form) does not comply with our representation of parallelism due to its heavy reliance on Template Haskell.

We will now go into some detail on GpH, the *Par* Monad and Eden, and also give their respective implementations of the non-Arrow version of *parEvalN*.

3.2.1 Glasgow parallel Haskell - GpH

GpH (Marlow *et al.*, 2009; Trinder *et al.*, 1998) is one of the simplest ways to do parallel processing found in standard GHC.⁴ Besides some basic primitives (*par* and *pseq*), it ships with parallel evaluation strategies for several types which can be applied with $using :: a \rightarrow Strategy \ a \rightarrow a$, which is exactly what is required for an implementation of parEvalN.

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

parEvalN\ fs\ as = \mathbf{let}\ bs = zipWith\ (\$)\ fs\ as

\mathbf{in}\ bs\ `using\ `parList\ rdeepseq
```

In the above definition of parEvalN we 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 $parList::Strategy[a] \rightarrow Strategy[a]$ and $rdeepseq::NFData[a] \rightarrow Strategy[a]$ where the latter is a strategy which evaluates to normal form. Other strategies like e.g. evaluation to weak head normal form are available as well. It also allows for custom Strategy implementations to be used. Fig. 5 shows a visual representation of this code.

⁴ The Multicore implementation of GpH is available on Hackage under https://hackage.haskell.org/package/parallel-3.2.1.0, compiler support is integrated in the stock GHC.

Arrows for Parallel Computation

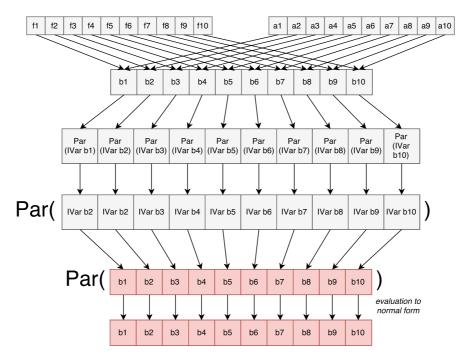


Figure 6: parEvalN (Par Monad).

3.2.2 Par Monad

The *Par* Monad⁵ introduced by Marlow *et al.* (2011), is a Monad designed for composition of parallel programs. Let:

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

parEvalN\ fs\ as = runPar\ \$

(sequenceA\ (map\ (return \circ spawn)\ (zipWith\ (\$)\ fs\ as))) > mapM\ get
```

The Par Monad version of our parallel evaluation function parEvalN is defined by zipping the list of $[a \rightarrow b]$ with the list of inputs [a] with the application operator \$ just like with GpH. Then, we map over this not yet evaluated lazy list of results [b] with $spawn:NFData\ a\Rightarrow Par\ a\rightarrow Par\ (IVar\ a)$ 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 [b]. While we used spawn in the definition above, a head-strict variant can easily be defined by replacing spawn with $spawn_::Par\ a\rightarrow Par\ (IVar\ a)$. Fig. 6 shows a graphical representation.

The Par monad can be found in the monad-par package on Hackage under https://hackage.haskell.org/package/monad-par-0.3.4.8/.

3.2.3 Eden

Eden (Loogen *et al.*, 2005; Loogen, 2012) is a parallel Haskell for distributed memory and comes with MPI and PVM as distributed backends.⁶ It is targeted towards clusters, but also functions 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 on multicores, as well (Berthold *et al.*, 2009a; Aswad *et al.*, 2009).

While Eden 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
```

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 tools Threadscope (Wheeler & Thain, 2009) and Eden TraceViewer⁷ (Berthold & Loogen, 2007). In the next sections we will present some *trace visualisations*, the post-mortem process diagrams of Eden processes and their activity.

The trace visualisations are colour-coded. In such a visualisation (Fig. 14), the *x* axis shows the time, the *y* axis enumerates the machines and processes. The visualisation 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 this is GC. An inactive machine, where no processes are started yet, or all are already terminated, shows as a blue bar. A communication 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.

4 Parallel Arrows

While Arrows are a general interface to computation, we introduce here specialised Arrows as a general interface to *parallel computations*. We present the *ArrowParallel* type class and explain the reasoning behind it before discussing some parallel Haskell implementations and basic extensions.

The projects homepage can be found at http://www.mathematik.uni-marburg.de/~eden/. The Hackage page is at https://hackage.haskell.org/package/edenmodules-1.2.0.0/.

⁷ See http://hackage.haskell.org/package/edentv on Hackage for the last available version of Eden TraceViewer.

4.1 The ArrowParallel type class

A parallel computation (on functions) can be seen as execution of some functions $a \to b$ in parallel, as our parEvalN prototype shows (Section 3.2). 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 evalN 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 - or even the input a and output b - and we want this to be an interface for different backends, we introduce a new type class 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 introduce an additional *conf* parameter as we do not want *conf* to be a fixed type, as the configuration parameters can differ for different instances of *ArrowParallel*.

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

By restricting the implementations of our backends to a specific *conf* type, we also get interoperability between backends for free. We can parallelize one part of a program using one backend, and parallelize the next with another one.

4.2 ArrowParallel instances

With the type class defined, we will now give implementations of it with GpH, the *Par* Monad and Eden.

4.2.1 Glasgow parallel Haskell

The GpH implementation of *ArrowParallel* is implemented in a straightforward manner in Fig. 7, but a bit different compared to the variant from Section 3.2.1. We use *evalN*:: $[arr\ a\ b] \to arr\ [a]\ [b]$ (definition in Appendix A, think *zipWith* (\$) on Arrows) combined with *withStrategy*:: *Strategy* $a \to a \to a$ from GpH, where *withStrategy* is the same as *using*:: $a \to Strategy\ a \to a$, but with flipped parameters. Our *Conf* a datatype simply wraps a *Strategy* a, but could be extended in future versions of our DSL.

4.2.2 Par Monad

As for GpH we can easily lift the definition of *parEvalN* for the *Par* Monad to Arrows in Fig. 8. To start off, we define the *Strategy a* and *Conf a* type so we can have a configurable instance of ArrowParallel:

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

instance (ArrowChoice\ arr) \Rightarrow

ArrowParallel\ arr\ a\ b\ (Conf\ b) where

parEvalN\ (Conf\ strat)\ fs =

evalN\ fs >>>

arr\ (withStrategy\ (parList\ strat))
```

Figure 7: GpH ArrowParallel instance.

```
type Strategy a = a \rightarrow Par (IVar a) data Conf a = Conf (Strategy a)
```

Now we can once again define our ArrowParallel instance as follows: First, we convert our Arrows $[arr\ a\ b]$ with $evalN\ (map\ (>>>> arr\ strat)\ fs)$ into an Arrow $arr\ [a]\ [(Par\ (IVar\ b))]$ that yields composable computations in the Par monad. By combining the result of this Arrow with $arr\ sequenceA$, we get an Arrow $arr\ [a]\ (Par\ [IVar\ b])$. Then, in order to fetch the results of the different threads, we map over the IVars inside the $Par\ Monad$ with $arr\ (>>= mapM\ get)$ — our intermediary Arrow is of type $arr\ [a]\ (Par\ [b])$. Finally, we execute the computation $Par\ [b]$ by composing with $arr\ runPar$ and get the final Arrow $arr\ [a]\ [b]$.

```
instance (ArrowChoice\ arr) \Rightarrow ArrowParallel\ arr\ a\ b\ (Conf\ b) where parEvalN\ (Conf\ strat)\ fs = evalN\ (map\ (>>> arr\ strat)\ fs) >>> arr\ sequenceA >>> arr\ (>>= mapM\ Control.Monad.Par.get) >>> arr\ runPar
```

Figure 8: Par Monad ArrowParallel instance.

4.2.3 Eden

For both the GpH Haskell and *Par* Monad implementations we could use general instances of *ArrowParallel* that just require the *ArrowChoice* type class. With Eden this is not the case as we can only spawn a list of functions, which we cannot extract from general Arrows. While we could still manage to have only one instance in the module by introducing a type class

```
class (Arrow arr) \Rightarrow ArrowUnwrap arr where unwrap :: arr a b \rightarrow (a \rightarrow b)
```

we avoid doing so for aesthetic reasons. For now, we just implement ArrowParallel for normal functions and the Kleisli type in Fig. 9, where Conf is simply defined as $\mathbf{data}\ Conf = Nil$ since Eden does not have a configurable spawnF variant.

```
instance (Trans a, Trans b) \Rightarrow ArrowParallel (\rightarrow) a b Conf where parEvalN _ = spawnF instance (ArrowParallel (\rightarrow) a (m b) Conf, 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(Kleisli f) \rightarrowf) fs)) >>> Kleisli sequence
```

Figure 9: Eden ArrowParallel instance.

4.2.4 Default configuration instances

While the configurability in the instances of the *ArrowParallel* instances above is nice, users probably would like to have proper default configurations for many parallel programs as well. These can also easily be defined as we can see by the example of the default implementation of *ArrowParallel* for the GpH:

```
instance (NFData b,ArrowChoice arr,ArrowParallel arr a b (Conf b)) \Rightarrow ArrowParallel arr a b () where
parEvalN \_fs = parEvalN (defaultConf fs) fs
defaultConf :: (NFData b) \Rightarrow [arr a b] \rightarrow Conf b
defaultConf fs = stratToConf fs rdeepseq
stratToConf :: [arr a b] \rightarrow Strategy b \rightarrow Conf b
stratToConf \_strat = Conf strat
```

The other backends have similarly structured implementations which we do not discuss here for the sake of brevity. We can, however, only have one instance of $ArrowParallel\,arr\,a\,b\,()$ present at a time, which should not be a problem, though.

Up until now we discussed Arrow operations more in detail, but in the following sections we focus more on the data-flow between the Arrows, now that we have seen that Arrows are capable of expressing parallelism. We do explain new concepts with more details if required for better understanding, though.

4.3 Extending the interface

With the *ArrowParallel* type class in place and implemented, we can now define other parallel interface functions. These are basic algorithmic skeletons that are used to define more sophisticated skeletons.

4.3.1 Lazy parEvalN

The function parEvalN fully traverses the list of passed Arrows as well as their inputs. 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 (Figs. 10, 11) fixes this. It works by first chunking the input from [a] to [[a]]

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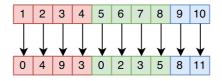


Figure 10: parEvalNLazy depiction.

```
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) >>> \\ evalN \ fchunks >>> \\ arr \ concat \\ \textbf{where} \\ fchunks = map \ (parEvalN \ conf) \ (chunksOf \ chunkSize \ fs)
```

Figure 11: Definition of *parEvalNLazy*.

with the given *chunkSize* in arr (*chunksOf chunkSize*). These chunks are then fed into a list [arr [a] [b]] of chunk-wise parallel Arrows with the help of our lazy and sequential *evalN*. The resulting [[b]] is lastly converted into [b] with arr *concat*.

4.3.2 Heterogeneous tasks

We have only talked about the parallelization of Arrows of the same set of input and output types until now. But sometimes we want to parallelize heterogeneous types as well. We can implement such a parEval2 combinator (Figs. 12, C 12) 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 type class. Here, the general 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 outline parallel Arrow combinator in Fig. 13. In a distributed environment this first evaluates all $[arr\ a\ b]$ in parallel, sends the results back to the master node, rotates



Figure 12: parEval2 depiction.

Arrows for Parallel Computation

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

Figure 13: The outline combinator.

the input once (in the example we require ArrowChoice for this) 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 could be rewritten into a single *parEvalN* call by directly wiring the Arrows together before spawning, it 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 14. This can become a serious bottleneck for a larger amount of data and number of processes (as e.g. Berthold *et al.*, 2009c, showcases).

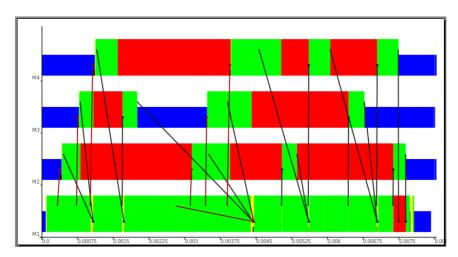


Figure 14: Communication between 4 Eden processes without Futures. All communication goes through the master node. Each bar represents one process. Black lines represent communication. Colours: blue $\hat{=}$ idle, green $\hat{=}$ running, red $\hat{=}$ blocked, yellow $\hat{=}$ suspended.

This is only a problem in distributed memory (in the scope of this paper) and we should allow nodes to communicate directly with each other. Eden already provides 'remote data' that enable this (Alt & Gorlatch, 2003; Dieterle *et al.*, 2010a). But as we want code with our DSL to be implementation agnostic, we have to wrap this concept. We do this with

the *Future* type class (Fig. 15). We require a *conf* parameter here as well, but only so that Haskells type system allows us to have multiple Future implementations imported at once without breaking any dependencies similar to what we did with the *ArrowParallel* type class earlier. Since *RD* is only a type synonym for a communication type that Eden uses

```
class Future fut a conf | a conf \rightarrow fut where put :: (Arrow arr) \Rightarrow conf \rightarrow arr \ a \ (fut \ a) get :: (Arrow arr) \Rightarrow conf \rightarrow arr \ (fut \ a) \ a
```

Figure 15: The *Future* type class.

internally, we have to use some wrapper classes to fit that definition, though, as Fig. C 1 shows. Technical details are in Appendix, in Section C.

For GpH and *Par* Monad, we can simply use *BasicFutures* (Fig. C 2), which are just simple wrappers around the actual data with boiler-plate logic so that the type class is satisfied. This is because the concept of a *Future* does not change anything for shared-memory execution as there are no communication problems to fix. Nevertheless, we require a common interface so the parallel Arrows are portable across backends. The implementation can also be found in Section C.

In our communication example we can use this *Future* concept for direct communication between nodes as shown in Fig. 16. In a distributed environment, this gives us a communication scheme with messages going through the master node only if it is needed – similar to what is shown in the trace visualisation in Fig. 17. One especially elegant aspect of the definition in Fig. 15 is that we can specify the type of *Future* to be used per backend with full interoperability between code using different backends, without even requiring to know about the actual type used for communication. We only specify that there has to be a compatible Future and do not care about any specifics as can be seen in Fig. 16. With the PArrows DSL we can also define default instances *Future fut a* () for each backend similar to how *ArrowParallel arr a b* () was defined in Section 4. Details can be found in Section C.

6 Skeletons

Now we have developed Parallel Arrows far enough to define some useful algorithmic skeletons that abstract typical parallel computations. While there are many possible skeletons to implement, we demonstrate the expressive power of PArrows here using four *map*-based and three toplogical skeletons.

6.1 map-based skeletons

The essential differences between the mapping skeletons presented here are in terms of order of evaluation and work distribution but still provide the same semantics as a sequential *map*.

Arrows for Parallel Computation

```
some Combinator :: (Arrow Choice\ arr, \\ Arrow Parallel\ arr\ a\ (fut\ b)\ (), \\ Arrow Parallel\ arr\ (fut\ b)\ c\ (), \\ Future\ fut\ b\ ()) \Rightarrow \\ [arr\ a\ b] \rightarrow [arr\ b\ c] \rightarrow arr\ [a]\ [c] \\ some Combinator\ fsl\ fs2 = \\ par EvalN\ ()\ (map\ (>>>put\ ())\ fsl) >>> \\ right Rotate >>> \\ par EvalN\ ()\ (map\ (get\ ()>>>)\ fs2)
```

Figure 16: The outline combinator in parallel.

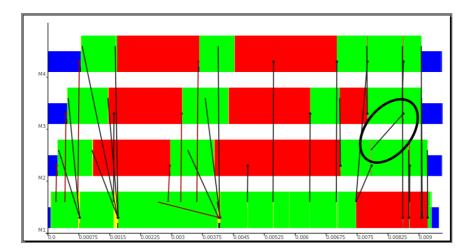


Figure 17: Communication between 4 Eden processes with Futures. Other than in Fig. 14, processes communicate directly (one example message is highlighted) instead of always going through the master node (bottom bar).

Parallel *map* **and laziness.** The *parMap* skeleton (Figs. C 3, C 4) 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 obtain an Arrow *arr* [*a*] [*b*]. Just like *parEvalN*, *parMap* traverses all input Arrows as well as the inputs. Because of this, it has the same restrictions as *parEvalN* as compared to *parEvalNLazy*. So it makes sense to also have a *parMapStream* (Figs. C 5, C 6) which behaves like *parMap*, but uses *parEvalNLazy* instead of *parEvalN*. Implementing these skeletons is straightforward as in Appendix C in Figs.C 4 and C 6.

Statically load-balancing parallel *map*. Our *parMap* spawns every single computation in a new thread (at least for the instances of *ArrowParallel* we presented in this paper). This can be quite wasteful and a statically load-balancing *farm* (Figs. 18, 19) that equally

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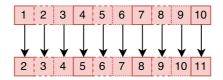


Figure 18: farm depiction.

```
\begin{array}{l} \textit{farm} :: (\textit{ArrowParallel arr a b conf}, \\ \textit{ArrowParallel arr } [a] [b] \textit{conf}, \textit{ArrowChoice arr}) \Rightarrow \\ \textit{conf} \rightarrow \textit{NumCores} \rightarrow \textit{arr a b} \rightarrow \textit{arr } [a] [b] \\ \textit{farm conf numCores} f = \\ \textit{unshuffle numCores} >>> \\ \textit{parEvalN conf (repeat (mapArrf))} >>> \\ \textit{shuffle} \end{array}
```

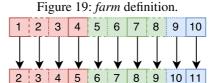


Figure 20: farmChunk depiction.

distributes the workload over *numCores* workers seems useful. The definitions of the helper functions *unshuffle*, *takeEach*, *shuffle* (Fig. C 7) originate from an Eden skeleton⁸.

Since a *farm* is basically just *parMap* with a different work distribution, it has the same restrictions as *parEvalN* and *parMap*. We can, however, define *farmChunk* (Figs. 20, C 10) which uses *parEvalNLazy* instead of *parEvalN*. It is basically the same definition as for *farm*, but with *parEvalNLazy* instead of *parEvalN*.

6.2 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 with more predefined skeletons⁹, among them a *pipe*, a *ring*, and a *torus* implementations (Loogen *et al.*, 2003). These seem like reasonable candidates to be ported to our Arrowbased parallel Haskell. We aim to showcase that we can express more sophisticated skeletons with parallel Arrows as well.

If we used the original definition of *parEvalN*, however, these skeletons would produce an infinite loop with the GpH and *Par* Monad which during runtime would result in the program crash. This materialises with the usage of *loop* of the *ArrowLoop* type class and

⁸ Available on Hackage under https://hackage.haskell.org/package/edenskel-2.1.0.0/docs/src/Control-Parallel-Eden-Map.html.

⁹ Available on Hackage: https://hackage.haskell.org/package/edenskel-2.1.0.0/docs/Control-Parallel-Eden-Topology.html.

we think that this is due to difference of how evaluation is done in these backends when compared to Eden. An investigation of why this difference exists is beyond the scope of this work, we only provide a workaround for these types of skeletons as such they probably are not of much importance outside of a distributed memory environment. However our workaround enables users of the DSL to test their code within a shared memory setting. The idea of the fix is to provide a *ArrowLoopParallel* type class that has two functions – *loopParEvalN* and *postLoopParEvalN*, where the first is to be used inside an *loop* construct while the latter will be used right outside of the *loop*. This way we can delegate to the actual *parEvalN* in the spot where the backend supports it.

```
class ArrowParallel arr a b conf \Rightarrow
ArrowLoopParallel arr a b conf where
loopParEvalN :: conf \rightarrow [arr\ a\ b] \rightarrow arr\ [a]\ [b]
postLoopParEvalN :: conf \rightarrow [arr\ a\ b] \rightarrow arr\ [a]\ [b]
```

As Eden has no problems with the looping skeletons, we use this instance:

```
instance (ArrowChoice arr,ArrowParallel arr a b Conf) ⇒
ArrowLoopParallel arr a b Conf where
loopParEvalN = parEvalN
postLoopParEvalN _ = evalN
```

As *Par* Monad and GpH have problems with *parEvalN* inside of *loop* their respective instances for *ArrowLoopParallel* look like this:

```
instance (ArrowChoice arr, ArrowParallel arr a b (Conf b)) \Rightarrow ArrowLoopParallel arr a b (Conf b) where loopParEvalN \_= evalN postLoopParEvalN = parEvalN
```

6.2.1 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 type class 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

```
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 loop to our advantage gives us a first draft of a pipe implementation (Fig. 21) by plugging in the parallel evaluation call loopParEvalN conf fs inside the second argument of &&& and then only picking the first element of the resulting list with arr last.

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

```
\begin{aligned} \textit{pipeSimple} &:: (\textit{ArrowLoop arr, ArrowLoopParallel arr a a conf}) \Rightarrow \\ & \textit{conf} \rightarrow [\textit{arr a a}] \rightarrow \textit{arr a a} \\ & \textit{pipeSimple conf fs} = \\ & \textit{loop (arr snd &&& \\ & (\textit{arr (uncurry (:)} >>>> \textit{lazy}) >>>> \textit{loopParEvalN conf fs)}) >>>> \\ & \textit{arr last} \end{aligned}
```

Figure 21: Simple *pipe* skeleton. The use of *lazy* (Fig. C 8) is essential as without it programs using this definition would never halt. We need to ensure that the evaluation of the input [a] is not forced fully before passing it into *loopParEvalN*.

```
pipe::(ArrowLoop arr,
   ArrowLoopParallel arr (fut a) (fut a) conf,
   Future fut a conf) \Rightarrow
   conf \rightarrow [arr a a] \rightarrow arr a a
pipe conf fs = unliftFut conf (pipeSimple conf (map (liftFut conf) fs))
liftFut::(Arrow arr, Future fut a conf, Future fut b conf) \Rightarrow
   conf \rightarrow arr a b \rightarrow arr (fut a) (fut b)
liftFut conf f = get conf >>> f >>> put conf
unliftFut::(Arrow arr, Future fut a conf, Future fut b conf) \Rightarrow
   conf \rightarrow arr (fut a) (fut b) \rightarrow arr a b
unliftFut conf f = put conf >>> f >>> get conf
```

Figure 22: *pipe* skeleton definition with Futures.

sophisticated version that internally uses Futures and obtain the final definition of *pipe* in Fig. 22.

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 bigger Arrow $arr\ (([a],[b]),[c])\ (([a],[b]),[c])$ suitable for pipe, we can define pipe2 as in Fig. 23.

Extensive use of *pipe2* 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* inside of *evalN*. Nonetheless, we can define a parallel piping operator |>>> |, which is semantically equivalent to >>>> similarly to other parallel syntactic sugar from Appendix D.

6.2.2 Ring skeleton

Eden comes with a ring skeleton¹⁰ (Fig. 24) implementation that allows the computation of a function $[i] \rightarrow [o]$ with a ring of nodes that communicate 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

Available on Hackage: https://hackage.haskell.org/package/edenskel-2.1.0.0/docs/Control-Parallel-Eden-Topology.html

```
pipe2::(ArrowLoop arr, ArrowChoice arr,
   ArrowLoopParallel\ arr\ (fut\ (([a],[b]),[c]))\ (fut\ (([a],[b]),[c]))\ conf,
   Future fut (([a],[b]),[c]) conf) \Rightarrow
      conf \rightarrow arr \ a \ b \rightarrow arr \ b \ c \rightarrow arr \ a \ c
pipe2 conf f g =
   (arr return &&& arr (const [])) &&& arr (const []) >>>
   pipe\ conf\ (replicate\ 2\ (unify\ f\ g)) >>>
   arr snd >>>
   arr head
   where
      unify::(ArrowChoice\ arr) \Rightarrow
         arr\ a\ b \rightarrow arr\ b\ c \rightarrow arr\ (([a],[b]),[c])\ (([a],[b]),[c])
      unify f'g' =
         (mapArr f' *** mapArr g') *** arr (const []) >>>
         arr (\lambda((b,c),a) \rightarrow ((a,b),c))
(|>>>|)::(ArrowLoop arr,ArrowChoice arr,
   ArrowLoopParallel\ arr\ (fut\ (([a],[b]),[c]))\ (fut\ (([a],[b]),[c]))\ (),
   Future fut (([a],[b]),[c])()) \Rightarrow
   arr\ a\ b 
ightarrow arr\ b\ c 
ightarrow arr\ a\ c
(|>>>|) = pipe2()
```

Figure 23: Definition of *pipe2* and (|>>>|), a parallel >>>.

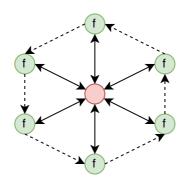


Figure 24: ring skeleton depiction.

the neighbour of each node. This data is sent over direct communication channels, the so called 'remote data'. We depict it in Appendix, Fig. C 11.

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 *loop*'s signature: $arr(a,b)(c,b) \rightarrow arr\ a\ c$. In each iteration we start by rotating the intermediary input from the nodes $[fut\ r]$ with $second\ (rightRotate >>> lazy)$ (Fig. C8). Similarly to the pipe from Section 6.2.1 (Fig. 21), we have to feed the intermediary input into our lazy (Fig. C8) Arrow here, or the evaluation would fail to terminate. The reasoning is explained by Loogen (2012) as a demand problem.

Next, we zip the resulting $([i],[fut\ r])$ to $[(i,fut\ r)]$ with $arr\ (uncurry\ zip)$. We then feed this into our parallel Arrow $arr\ [(i,fut\ r)]\ [(o,fut\ r)]$ obtained by transforming our

```
ring:: (Future fut r conf,
    ArrowLoop arr,
    ArrowLoopParallel arr (i,fut r) (o,fut r) conf,
    ArrowLoopParallel arr o o conf) \Rightarrow
    conf \rightarrow arr (i,r) (o,r) \rightarrow arr [i] [o]
ring conf f =
    loop (second (rightRotate >>> lazy) >>>
    arr (uncurry zip) >>>
    loopParEvalN conf (repeat (second (get conf) >>>> f >>> second (put conf))) >>>
    postLoopParEvalN conf (repeat (arr id))
```

Figure 25: ring skeleton definition.

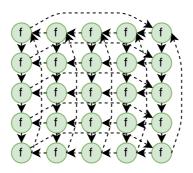


Figure 26: torus skeleton depiction.

input Arrow f :: arr(i,r)(o,r) into arr(i,fut r)(o,fut r) before repeating and lifting it with loopParEvalN. Finally we unzip the output list [(o,fut r)] list into ([o],[fut r]).

Plugging this Arrow $arr([i],[fut\ r])([o],fut\ r)$ into the definition of loop from earlier gives us arr[i][o], our ring Arrow (Fig. 25). To make sure this algorithm has speedup on shared-memory machines as well, we pass the result of this Arrow to $postLoopParEvalN\ conf\ (repeat\ (arr\ id))$. This combinator can, for example, be used to calculate the shortest paths in a graph using Warshall's algorithm.

6.2.3 Torus skeleton

If we take the concept of a *ring* from Section 6.2.2 one dimension further, we obtain a *torus* skeleton (Fig. 26, 27). Every node sends and receives data from horizontal and vertical neighbours in each communication round. With our Parallel Arrows we re-implement the *torus* combinator¹¹ from Eden—yet again with the help of the *ArrowLoop* type class.

Similar to the *ring*, we start by rotating the input (Fig. C 8), but this time not only in one direction, but in two. This means that the intermediary input from the neighbour nodes has

Available on Hackage: https://hackage.haskell.org/package/edenskel-2.1.0.0/docs/Control-Parallel-Eden-Topology.html.

```
torus:: (Future fut a conf, Future fut b conf,
  ArrowLoop arr, ArrowChoice arr,
  ArrowLoopParallel\ arr\ (c, fut\ a, fut\ b)\ (d, fut\ a, fut\ b)\ conf,
  ArrowLoopParallel \ arr \ [d] \ [d] \ conf) \Rightarrow
   conf \rightarrow arr(c,a,b)(d,a,b) \rightarrow arr[[c]][[d]]
torus conf f =
   loop (second ((mapArr rightRotate >>> lazy) *** (arr rightRotate >>> lazy)) >>>
     arr (uncurry3 (zipWith3 lazyzip3)) >>>
     arr\ length\ \&\&\&\ (shuffle>>> loopParEvalN\ conf\ (repeat\ (ptorus\ conf\ f)))>>>
     arr (uncurry unshuffle) >>>
     arr (map unzip3) >>> arr unzip3 >>> threetotwo) >>>
  postLoopParEvalN conf (repeat (arr id))
ptorus :: (Arrow arr, Future fut a conf, Future fut b conf) \Rightarrow
   arr(c,a,b)(d,a,b) \rightarrow
  arr(c, fut \ a, fut \ b)(d, fut \ a, fut \ b)
ptorus conf f =
     arr (\lambda \sim (c, a, b) \rightarrow (c, get \ conf \ a, get \ conf \ b)) >>>
     arr (\lambda \sim (d, a, b) \rightarrow (d, put \ conf \ a, put \ conf \ b))
```

Figure 27: *torus* skeleton definition. *lazyzip3*, *uncurry3* and *threetotwo* definitions are in Fig. C 9

to be stored in a tuple ([$[fut\ a]$],[$[fut\ b]$]) in the second argument (loop only allows for two arguments) of our looped Arrow $arr\left([[c]],([[fut\ a]],[[fut\ b]])\right)$ ([[d]],([$[fut\ a]$],[$[fut\ b]$])) and our rotation Arrow becomes

```
second ((mapArr rightRotate >>> lazy) *** (arr rightRotate >>> lazy))
```

instead of the singular rotation in the ring as we rotate $[[fut \ a]]$ horizontally and $[[fut \ b]]$ vertically. Then, we 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 evaluate in parallel.

This, however, is more complicated than in the ring case as we have one more dimension of inputs that needs to be transformed. We first have to *shuffle* all the inputs to then pass them into *loopParEvalN conf* (repeat (ptorus conf f)) to get an output of $[(d,fut\ a,fut\ b)]$. We then unshuffle this list back to its original ordering by feeding it into arr (uncurry unshuffle) which takes the input length we saved one step earlier as additional input to get a result matrix $[[[(d,fut\ a,fut\ b)]]]$. Finally, we unpack this matrix with arr (map unzip3) >>> arr unzip3 >>>> threetotwo to get $([[d]],([[fut\ a]],[[fut\ b]]))$.

This internal looping computation is once again fed into *loop* and we also compose a final *postLoopParEvalN conf* (*repeat* (*arr id*)) for the same reasons as explained for the *ring* skeleton.

As an example of using this skeleton, Loogen *et al.* (2003) showed the matrix multiplication using the Gentleman algorithm (1978). An adapted version can be found in Fig. 28. If

```
 \begin{aligned} \textbf{type} \ \textit{Matrix} &= [[\textit{Int}]] \\ \textit{prMM\_torus} &:: \textit{Int} \rightarrow \textit{Int} \rightarrow \textit{Matrix} \rightarrow \textit{Matrix} \rightarrow \textit{Matrix} \\ \textit{prMM\_torus} \ \textit{numCores} \ \textit{problemSizeVal} \ \textit{ml} \ \textit{m2} = \\ \textit{combine} \$ \ \textit{torus} \ () \ (\textit{mult} \ \textit{torusSize}) \$ \ \textit{zipWith} \ \textit{zip} \ (\textit{split1} \ \textit{m1}) \ (\textit{split2} \ \textit{m2}) \\ \textbf{where} \ \textit{torusSize} &= \ (\textit{floor} \circ \textit{sqrt}) \$ \ \textit{fromIntegral} \$ \ \textit{numCoreCalc} \ \textit{numCores} \\ \textit{combine} \ \textit{x} &= \ \textit{concat} \ (\textit{map} \ (\textit{concat})) \circ \textit{transpose}) \ \textit{x}) \\ \textit{split1} \ \textit{x} &= \ \textit{staggerHorizontally} \ (\textit{splitMatrix} \ (\textit{problemSizeVal} \ \textit{'div'} \ \textit{torusSize}) \ \textit{x}) \\ \textit{split2} \ \textit{x} &= \ \textit{staggerVertically} \ (\textit{splitMatrix} \ (\textit{problemSizeVal} \ \textit{'div'} \ \textit{torusSize}) \ \textit{x}) \\ \textit{--Function performed} \ \textit{by} \ \textit{each} \ \textit{worker} \\ \textit{mult} :: \textit{Int} \rightarrow ((\textit{Matrix}, \textit{Matrix}), [\textit{Matrix}], [\textit{Matrix}]) \rightarrow (\textit{Matrix}, [\textit{Matrix}], [\textit{Matrix}]) \\ \textit{mult} \ \textit{size} \ ((\textit{sm1}, \textit{sm2}), \textit{sm1s}, \textit{sm2s}) = \ (\textit{result}, \textit{toRight}, \textit{toBottom}) \\ \textbf{where} \ \textit{toRight} = \ \textit{take} \ (\textit{size} - 1) \ (\textit{sm1} : \textit{sm1s}) \\ \textit{toBottom} = \ \textit{take} \ (\textit{size} - 1) \ (\textit{sm2} : \textit{sm2s}) \\ \textit{sms} = \ \textit{zipWith} \ \textit{prMM} \ (\textit{sm1} : \textit{sm1s}) \ (\textit{sm2} : \textit{sm2s}) \\ \textit{result} = \ \textit{fold11'} \ \textit{matAdd} \ \textit{sms} \end{aligned}
```

Figure 28: Adapted matrix multiplication in Eden using a the *torus* skeleton. $prMM_torus$ is the parallel matrix multiplication. mult is the function performed by each worker. prMM is the sequential matrix multiplication in the chunks. splitMatrix splits the Matrix into chunks. staggerHorizontally and staggerVertically pre-rotate the matrices. matAdd calculates A + B. Omitted definitions can be found in C 13.

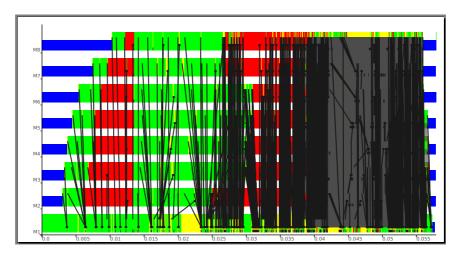


Figure 29: Matrix multiplication with torus (PArrows).

we compare the trace from a call using our Arrow definition of the torus (Fig. 29) with the Eden version (Fig. 30) we can see that the behaviour of the Arrow version and execution times are comparable. We discuss further benchmarks on larger clusters in more detail in the next section.

Arrows for Parallel Computation

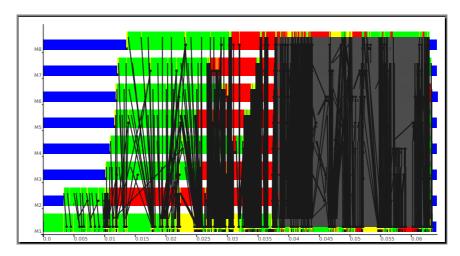


Figure 30: Matrix multiplication with *torus* (Eden).

7 Performance results and discussion

The preceding section has shown that PArrows are expressive. This section evaluates the performance overhead of this compositional abstraction in comparison to GpH and the *Par* Monad on shared memory architectures and Eden on a distributed memory cluster. We describe our measurement platform, the benchmark results – the shared-memory variants (GpH, *Par* Monad and Eden CP) followed by Eden in a distributed memory setting, and conclude that PArrows hold up in terms of performance when compared to the original parallel Haskells.

7.1 Measurement platform

We start by explaining the hardware and software stack and outline the benchmark programs and motivation for choosing them. We also shortly address hyper-threading and why we do not use it in our benchmarks.

7.1.1 Hardware and software

The benchmarks are executed both in a shared and in a distributed memory setting using the Glasgow GPG Beowulf cluster, consisting of 16 machines with 2 Intel[®] Xeon[®] E5-2640 v2 and 64 GB of DDR3 RAM each. Each processor has 8 cores and 16 (hyper-threaded) threads with a base frequency of 2 GHz and a turbo frequency of 2.50 GHz. This results in a total of 256 cores and 512 threads for the whole cluster. The operating system was Ubuntu 14.04 LTS with Kernel 3.19.0-33. Non-surprisingly, we found that hyper-threaded 32 cores do not behave in the same manner as real 16 cores (numbers here for a single machine). We disregarded the hyper-threading ability in most of the cases.

Table 1: The benchmarks we use in this paper.

Name	Area	Туре	Origin	Source
Rabin-Miller test	Mathematics	parMap + reduce	Eden	Lobachev (2012)
Jacobi sum test	Mathematics	workpool + reduce	Eden	Lobachev (2012)
Gentleman	Mathematics	torus	Eden	Loogen et al. (2003)
Sudoku	Puzzle	parMap	Par Monad	Marlow <i>et al.</i> (2011) ¹⁹

Apart from Eden, all benchmarks and libraries were compiled with Stack's 12 lts-7.1 GHC compiler which is equivalent to a standard GHC 8.0.1 with the base package in version 4.9.0.0. Stack itself was used in version 1.3.2. For GpH in its Multicore variant we used the parallel package in version 3.2.1.0¹³, while for the *Par* Monad we used monad-par in version 0.3.4.8¹⁴. For all Eden tests, we used its GHC-Eden compiler in version 7.8.2¹⁵ together with OpenMPI 1.6.5¹⁶.

Furthermore, all benchmarks were done with help of the bench¹⁷ tool in version 1.0.2 which uses criterion (>=1.1.1.0 && < 1.2)¹⁸ internally. All runtime data (mean runtime, max stddev, etc.) was collected with this tool.

We used a single node with 16 real cores as a shared memory test-bed and the whole grid with 256 real cores as a device to test our distributed memory software.

7.1.2 Benchmarks

We measure four benchmarks from different sources. Most of them are parallel mathematical computations, initially implemented in Eden. Table 1 summarises.

Rabin-Miller test is a probabilistic primality test that iterates multiple (here: 32-256) 'subtests'. Should a subtest fail, the input is definitely not a prime. If all n subtest pass, the input is composite with the probability of $1/4^n$.

Jacobi sum test or APRCL is also a primality test, that however, guarantees the correctness of the result. It is probabilistic in the sense that its run time is not certain. Unlike Rabin-Miller test, the subtests of Jacobi sum test have very different durations. Lobachev (2011) discusses some optimisations of parallel APRCL. Generic parallel implementations of Rabin-Miller test and APRCL were presented in Lobachev (2012).

'Gentleman' is a standard Eden test program, developed for their torus skeleton. It implements a Gentleman's algorithm for parallel matrix multiplication (Gentleman, 1978). We ported an Eden-based version (Loogen et al., 2003) to PArrows.

```
12 see https://www.haskellstack.org/
```

¹³ see https://hackage.haskell.org/package/parallel-3.2.1.0

see https://hackage.haskell.org/package/monad-par-0.3.4.8

¹⁵ see http://www.mathematik.uni-marburg.de/~eden/?content=build_eden_7_&navi= build

¹⁶ see https://www.open-mpi.org/software/ompi/v1.6/

 $^{^{17}\ {\}rm see\ https://hackage.haskell.org/package/bench}$

see https://hackage.haskell.org/package/criterion-1.1.1.0

actual code from: http://community.haskell.org/~simonmar/par-tutorial.pdf and https://github.com/simonmar/parconc-examples

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A parallel Sudoku solver was used by Marlow et al. (2011) to compare Par Monad to GpH, we ported it to PArrows.

7.1.3 What parallel Haskells run where

The Par monad and GpH – in its multicore version (Marlow et al., 2009) – can be executed on shared memory machines only. Although GpH is available on distributed memory clusters, and newer distributed memory Haskells such as HdpH exist, current support of distributed memory in PArrows is limited to Eden. We used the MPI backend of Eden in a distributed memory setting. However, for shared memory Eden features a 'CP' backend that merely copies the memory blocks between disjoint heaps. In this mode, Eden still operates in the 'nothing shared' setting, but is adapted better to multicore machines. We call this version of Eden 'Eden CP'.

7.1.4 Effect of hyper-threading

In preliminary tests, the PArrows version of Rabin-Miller test on a single node of the Glasgow cluster showed almost linear speedup on up to 16 shared-memory cores (as supplementary materials show). The speedup of 64-task PArrows/Eden at 16 real cores version was 13.65 giving a parallel efficiency of 85.3%. However, if we increased the number of requested cores to 32 – i.e. if we use hyper-threading on 16 real cores – the speedup did not increase that well. It was merely 15.99 for 32 tasks with PArrows/Eden. This was worse for other implementations. As for 64 tasks, we obtained a speedup of 16.12 with PArrows/Eden at 32 hyper-threaded cores and only 13.55 with PArrows/GpH.

While this shows that hyper-threading can be of benefit in scenarios similar to the ones presented in the benchmarks, we only use real cores for the performance measurements in Section 7.2 as the purpose of this paper is to show the performance of PArrows and not to investigate parallel behaviour with hyper-threading.

7.2 Benchmark results

We compare the PArrow performance with direct implementations of the benchmarks in Eden, GpH and the Par Monad. We start with the definition of mean overhead to compare both PArrows-enabled and standard benchmark implementations. We continue comparing speedups and overheads for the shared memory implementations and then study OpenMPI variants of the Eden-enabled PArrows as a representative of a distributed memory backend. We plot all speedup curves and all overhead values in the supplementary materials.

7.2.1 Defining overhead

We compare the mean overhead, i.e. the mean of relative wall-clock run time differences between the PArrow and direct benchmark implementations executed multiple times with the same settings. The error margins of the time measurements, supplied by criterion package²⁰, yield the error margin of the mean overhead.

https://hackage.haskell.org/package/criterion-1.1.1.0

Quite often the zero value lies in the error margin of the mean overhead. This means that even though we have measured some difference (against or even in favour of PArrows), it could be merely the error margin of the measurement and the difference might not be existent. We are mostly interested in the cases where above issue does not persist, we call them *significant*. We often denote the error margin with \pm after the mean overhead value.

7.2.2 Shared memory

Speedup. The Rabin-Miller test benchmark showed almost linear speedup for both 32 and 64 tasks, the performance is slightly better in the latter case: 13.7 at 16 cores for input $2^{11213} - 1$ and 64 tasks in the best case scenario with Eden CP. The performance of the Sudoku benchmark merely reaches a speedup of 9.19 (GpH), 8.78 (Par Monad), 8.14 (Eden CP) for 16 cores and 1000 Sudokus. In contrast to Rabin–Miller, here the *GpH* seems to be the best of all, while Rabin-Miller profited most from Eden CP (i.e. Eden with direct memory copy) implementation of PArrows. Gentleman on shared memory has a plummeting speedup curve with GpH and Par Monad and logarithmically increasing speedup for the Eden-based version. The latter reached a speedup of 6.56 at 16 cores.

Overhead. For the shared memory Rabin-Miller test benchmark, implemented with PArrows using Eden CP, GpH, and Par Monad, the overhead values are within single percents range, but also negative overhead (i.e. PArrows are better) and larger error margins happen. To give a few examples, the overhead for Eden CP with input value $2^{11213} - 1$, 32 tasks, and 16 cores is 1.5%, but the error margin is around 5.2%! Same implementation in the same setting with 64 tasks reaches -0.2% overhead, PArrows apparently fare better than Eden – but the error margin of 1.9% disallows this interpretation. We focus now on significant overhead values. To name a few: $0.41\% \pm 7 \cdot 10^{-2}\%$ for Eden CP and 64 tasks at 4 cores; $4.7\% \pm 0.72\%$ for GpH, 32 tasks, 8 cores; $0.34\% \pm 0.31\%$ for Par Monad at 4 cores with 64 tasks. The worst significant overhead was in case of GpH with $8\% \pm 6.9\%$ at 16 cores with 32 tasks and input value $2^{11213} - 1$. In other words, we notice no major slow-down through PArrows here.

For Sudoku the situation is slightly different. There is a minimal significant (-1.4% \pm 1.2% at 8 cores) speed improvement with PArrows Eden CP version when compared with the base Eden CP benchmark. However, with increasing number of cores the error margin reaches zero again: $-1.6\% \pm 5.0\%$ at 16 cores. The Par Monad shows a similar development, e.g. with $-1.95\% \pm 0.64\%$ at 8 cores. The GpH version shows both a significant speed improvement of $-4.2\% \pm 0.26\%$ (for 16 cores) with PArrows and a minor overhead of 0.87% $\pm 0.70\%$ (4 cores).

The Gentleman multiplication with Eden CP shows a minor significant overhead of $2.6\% \pm 1.0\%$ at 8 cores and an insignificant improvement at 16 cores. Summarising, we observe a low (if significant at all) overhead, induced by PArrows in the shared memory setting.

7.2.3 Distributed memory

Speedup. The speedup of distributed memory Rabin–Miller benchmark with PArrows and Eden showed an almost linear speedup excepting around 192 cores where an unfortunate task distribution reduces performance. As seen in Fig. 31, we reached a speedup of 213.4 with PArrrows at 256 cores (vs. 207.7 with pure Eden). Because of memory limitations, the speedup of Jacobi sum test for large inputs (such as $2^{4253} - 1$) could be measured only in a massively distributed setting. PArrows improved there from 9193 s (at 128 cores) to 1649 s (at 256 cores). A scaled-down version with input $2^{3217} - 1$ stagnates the speedup at about 11 for both PArrows and Eden for more than 64 cores. There is apparently not enough work for that many cores. The Gentleman test with input 4096 had an almost linear speedup first, then plummeted between 128 and 224 cores, and recovered at 256 cores with speedup of 129.

Speedup of distributed Rabin-Miller test '44497 256'

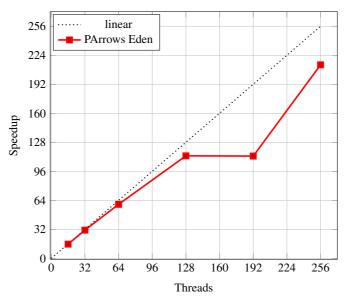


Figure 31: Speedup of the distributed Rabin–Miller test benchmark using PArrows with Eden.

Overhead. We use our mean overhead quality measure and the notion of significance also for distributed memory benchmarks. The mean overhead of Rabin-Miller test in the distributed memory setting ranges from 0.29% to -2.8% (last value in favour of PArrows), but these values are not significant with error margins $\pm 0.8\%$ and $\pm 2.9\%$ correspondingly. A sole significant (by a very low margin) overhead is $0.35\% \pm 0.33\%$ at 64 cores. We measured the mean overhead for Jacobi benchmark for an input of $2^{3217} - 1$ for up to 256 cores. We reach the flattering value $-3.8\% \pm 0.93\%$ at 16 cores in favour of PArrows, it was the sole significant overhead value. The value for 256 cores was $0.31\% \pm 0.39\%$. Mean overhead for distributed Gentleman multiplication was also low. Significant values include

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Table 2: Overhead in the shared memory benchmarks. Bold marks values in favour of PArrows.

Benchmark	Base	Mean of mean overheads	Maximum normalised stdDev	Runtime for 16 cores (s)
Sudoku 1000	Eden CP	-2.1%	5.1%	1.17
	GpH	-0.82%	0.7%	1.11
	Par Monad	-1.3%	2.1%	1.14
Gentleman 512	Eden CP	0.81%	6.8%	1.66
Rabin–Miller test 11213 32	Eden CP	0.79%	5.2%	5.16
	GpH	3.5%	6.9%	5.28
	Par Monad	- 2.5 %	19.0%	5.84
Rabin–Miller test 11213 64	Eden CP	0.21%	1.9%	10.3
	GpH	1.6%	1.3%	10.6
	Par Monad	-4.0%	17.0%	11.4

 $1.23\%\pm1.20\%$ at 64 cores and $2.4\%\pm0.97\%$ at 256 cores. It took PArrows 64.2 seconds at 256 cores to complete the benchmark.

Similar to the shared memory setting, PArrows only imply a very low penalty with distributed memory that lies in lower single-percent digits at most.

7.3 Discussion

PArrows performed in our benchmarks with little to no overhead. Tables 2 and 3 clarify this once more: The PArrows-enabled versions trade blows with their vanilla counterparts when comparing the means over all cores of the mean overheads. If we combine these findings with the usability of our DSL, the minor overhead induced by PArrows is outweighed by their convenience and usefulness to the user.

PArrows are an extendable formalism, they can be easily ported to further parallel Haskells while still maintaining interchangeability. It is straightforward to provide further implementations of algorithmic skeletons in PArrows.

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Table 3: Overhead in the distributed memory benchmarks. Bold marks values in favour of PArrows.

Benchmark	Base	Mean of mean overheads	Maximum normalised stdDev	Runtime for 256 cores (s)
Gentleman 4096	Eden	0.67%	1.5%	110.0
Rabin–Miller test 44497 256	Eden	-0.5%	2.9%	165.0
Jacobi sum test 3217	Eden	-0.74%	1.6%	635.0

8 Conclusion

Arrows are a generic concept that allows for powerful composition combinators. To our knowledge we are first to represent parallel computation with Arrows, and hence to show their usefulness for composing parallel programs. We have shown that for a generic and extensible parallel Haskell, we do not have to restrict ourselves to a monadic interface. We argue that Arrows are better suited to parallelise pure functions than Monads, as the functions are already Arrows and can be used directly in our DSL. Arrows are a better fit to parallelise pure code than a monadic solution as regular functions are already Arrows and can be used with our DSL in a more natural way. We use a non-monadic interface (similar to Eden or GpH) and retain composability. The DSL allows for a direct parallelisation of monadic code via the Kleisli type and additionally allows to parallelise any Arrow type that has an instance for ArrowChoice. (Some skeletons require an additional ArrowLoop instance.)

We have demonstrated the generality of the approach by exhibiting PArrow implementations for Eden, GpH, and the Par Monad. Hence, parallel programs can be ported between task parallel Haskell implementations with little or no effort. We are confident that it will be straightforward to add other task-parallel Haskells. In other words, PArrows greatly increase portability of parallel Haskell programs. Our measurements of four benchmarks on both shared and distributed memory platforms shows that the generality and portability of PArrows has very low performance overheads, i.e. never more than $8\% \pm 6.9\%$ and typically under 2%.

8.1 Future work

Our PArrows DSL can be expanded to other task parallel Haskells, and a specific target is HdpH (Maier et al., 2014). Further Future-aware versions of Arrow combinators can

be defined. Existing combinators could also be improved, for example more specialised versions of >>> and *** combinators are viable.

In ongoing work we are expanding both our skeleton library and the number of skeleton-based parallel programs that use our DSL. It would also be interesting to see a hybrid of PArrows and Accelerate (McDonell *et al.*, 2015). Ports of our approach to other languages such as Frege, Eta, or Java directly are at an early development stage.

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A Utility Arrows

Following are definitions of some utility Arrows used in this paper that have been left out for brevity. We start with the *second* combinator from Hughes (2000), which is a mirrored version of *first*, which is for example used in the definition of ***:

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

Next, we give the definition of *evalN* which also helps us to define *map*, and *zipWith* on Arrows. The *evalN* combinator in Fig. A 1 converts a list of Arrows $[arr\ a\ b]$ into an Arrow $arr\ [a]\ [b]$.

Figure A 1: The definition of evalN

The mapArr combinator (Fig. A 2) lifts any Arrow $arr\ a\ b$ to an Arrow $arr\ [a]\ [b]$. The original inspiration was from Hughes (2005), but the definition as then unified with evalN.

```
mapArr :: ArrowChoice \ arr \Rightarrow arr \ a \ b \rightarrow arr \ [a] \ [b]

mapArr = evalN \circ repeat
```

Figure A 2: The definition of map over Arrows.

Finally, with the help of mapArr (Fig. A 2), we can define zipWithArr (Fig. A 3) that lifts any Arrow arr(a,b)c to an Arrow arr([a],[b])[c].

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

Figure A 3: zipWith over Arrows.

These combinators make use of the ArrowChoice type class which provides the \parallel combinator. It takes two Arrows $arr\ a\ c$ and $arr\ b\ c$ and combines them into a new Arrow $arr\ (Either\ a\ b)\ c$ which pipes all $Left\ a$'s to the first Arrow and all $Right\ b$'s to the second Arrow:

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

B Profunctor Arrows

In Fig. B 1 we show how specific Profunctors can be turned into Arrows. This works because Arrows are strong Monads in the bicategory *Prof* of Profunctors as shown by Asada (2010). In Standard GHC (>>>>) has the type (>>>>):: *Category cat* \Rightarrow *cat* a b \rightarrow *cat* b c \rightarrow *cat* a c and is therefore not part of the *Arrow* type class like presented in this paper.²¹

For additional information on the type classes used, see: https://hackage.haskell.org/package/profunctors-5.2.1/docs/Data-Profunctor.html and https://hackage.haskell.org/package/base-4.9.1.0/docs/Control-Category.html.

Arrows for Parallel Computation

```
instance (Category p, Strong p) ⇒ Arrow p where
    arr f = dimap id f id
    first = first'
instance (Category p, Strong p, Costrong p) ⇒ ArrowLoop p where
    loop = loop'
instance (Category p, Strong p, Choice p) ⇒ ArrowChoice p where
    left = left'
```

Figure B 1: Profunctors as Arrows.

C Additional function definitions

We have omitted some function definitions in the main text for brevity, and redeem this here. We begin with warping Eden's build-in *RemoteData* to *Future* in Figure C 1

```
data RemoteData\ a = RD\ \{rd::RD\ a\}
put'::(Arrow\ arr)\Rightarrow arr\ a\ (BasicFuture\ a)
put'=arr\ BF
get'::(Arrow\ arr)\Rightarrow arr\ (BasicFuture\ a)\ a
get'=arr\ (\lambda(\sim(BF\ a))\rightarrow a)
instance NFData\ (RemoteData\ a) where
rnf=rnf\circ rd
instance (Trans\ a)\Rightarrow Future\ RemoteData\ a\ Conf\ where
put\ _=put'
get\ _=get'
instance (Trans\ a)\Rightarrow Future\ RemoteData\ a\ () where
put\ _=put'
get\ _=get'
```

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

Next, we have the definition of *BasicFuture* in Fig. C 2 and the corresponding *Future* instances.

Figures C 3–C 6 show the definitions and a visualisations of two parallel *map* variants, defined using *parEvalN* and its lazy counterpart.

Arrow versions of Eden's *shuffle*, *unshuffle* and the definition of *takeEach* are in Figure C 7. Similarly, Figure C 8 contains the definition of Arrow versions of Eden's *lazy* and *rightRotate* utility functions. Fig. C 9 contains Eden's definition of *lazyzip3* together with the utility functions *uncurry3* and *threetotwo*. The full definition of *farmChunk* is in Figure C 10. Eden definition of *ring* skeleton is in Figure C 11. It follows Loogen (2012).

The parEval2 skeleton is defined in Figure C 12. 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

```
data BasicFuture\ a = BF\ a

put'::(Arrow\ arr)\Rightarrow arr\ a\ (BasicFuture\ a)

put'=arr\ BF

get'::(Arrow\ arr)\Rightarrow arr\ (BasicFuture\ a)\ a

get'=arr\ (\lambda(\sim(BF\ a))\rightarrow a)

instance NFD at a\Rightarrow NFD at a (BasicFuture\ a) where rnf\ (BF\ a)=rnf\ a

instance F uture\ BasicF uture\ a\ (Conf\ a) where put\ _=put'

get\ _=get'

instance\ F uture\ BasicF uture\ a\ () where put\ _=put'

get\ _=get'
```

Figure C 2: BasicFuture type and its Future instance for the Par Monad and GpH.

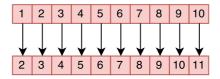


Figure C 3: *parMap* depiction.

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

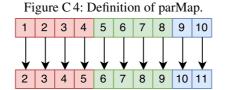


Figure C 5: parMapStream depiction.

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

Figure C 6: parMapStream definition.

f + ++ g as described in the paper. After the calculation is finished, we convert the 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. Furthermore, Fig. C 13 contains the omitted definitions of prMM (sequential matrix multiplication), splitMatrix (which splits the a matrix into chunks), staggerHorizontally

```
shuffle::(Arrow\ arr)\Rightarrow arr\ [[a]]\ [a] shuffle=arr\ (concat\circ transpose) unshuffle::(Arrow\ arr)\Rightarrow Int\rightarrow arr\ [a]\ [[a]] unshuffle\ n=arr\ (\lambda xs\rightarrow [takeEach\ n\ (drop\ i\ xs)\ |\ i\leftarrow [0\mathinner{.\,.} n-1]]) takeEach::Int\rightarrow [a]\rightarrow [a] takeEach\ n\ ([]=[] takeEach\ n\ (x:xs)=x:takeEach\ n\ (drop\ (n-1)\ xs)
```

Figure C7: shuffle, unshuffle, takeEach definition.

```
\begin{aligned} &lazy :: (Arrow \ arr) \Rightarrow arr \ [a] \ [a] \\ &lazy = arr \ (\lambda \sim (x : xs) \rightarrow x : lazy \ xs) \\ &rightRotate :: (Arrow \ arr) \Rightarrow arr \ [a] \ [a] \\ &rightRotate = arr \$ \lambda list \rightarrow \mathbf{case} \ list \ \mathbf{of} \\ &[] \rightarrow [] \\ &xs \rightarrow last \ xs : init \ xs \end{aligned}
```

Figure C 8: lazy and rightRotate definitions.

and staggerVertically (to pre-rotate the matrices), and lastly matAdd, that calculates A + B for two matrices A and B.

D Syntactic sugar

Finally, we also give the definitions for some syntactic sugar for PArrows, namely |***| and |&&&|. For basic Arrows, we have the *** combinator (Fig. 3) 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. 3):

```
(|\&\&\&|):: (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
```

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```
\begin{aligned} &lazyzip3::[a] \rightarrow [b] \rightarrow [c] \rightarrow [(a,b,c)] \\ &lazyzip3 \ as \ bs \ cs = zip3 \ as \ (lazy \ bs) \ (lazy \ cs) \\ &uncurry3::(a \rightarrow b \rightarrow c \rightarrow d) \rightarrow (a,(b,c)) \rightarrow d \\ &uncurry3 \ f \ (a,(b,c)) = f \ a \ b \ c \\ &threetotwo::(Arrow \ arr) \Rightarrow arr \ (a,b,c) \ (a,(b,c)) \\ &threetotwo = arr \$ \lambda \sim (a,b,c) \rightarrow (a,(b,c)) \end{aligned}
```

Figure C9: lazyzip3, uncurry3 and threetotwo definitions.

```
farmChunk :: (ArrowParallel \ arr \ a \ b \ conf \ , ArrowParallel \ arr \ [a] \ [b] \ conf, ArrowChoice \ arr \ , ArrowApply \ arr) \Rightarrow conf \rightarrow ChunkSize \rightarrow NumCores \rightarrow arr \ a \ b \rightarrow arr \ [a] \ [b] farmChunk \ conf \ chunkSize \ numCores \ f = unshuffle \ numCores \ >>> parEvalNLazy \ conf \ chunkSize \ (repeat \ (mapArr \ f)) >>> shuffle
```

Figure C 10: farmChunk definition.

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

Figure C 11: Eden's definition of the *ring* skeleton.

```
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 C 12: parEval2 definition.

Arrows for Parallel Computation

```
 prMM :: Matrix \rightarrow Matrix \rightarrow Matrix \\ prMM m1 m2 = prMMTr m1 (transpose m2) \\ \textbf{where} \\ prMMTr m1' m2' = [[sum (zipWith (*) row col) | col \leftarrow m2'] | row \leftarrow m1'] \\ splitMatrix :: Int \rightarrow Matrix \rightarrow [[Matrix]] \\ splitMatrix size matrix = map (transpose <math>\circ map (chunksOf size)) \circ chunksOf size \circ matrix staggerHorizontally :: [a] \rightarrow [a] \\ staggerHorizontally matrix = zipWith leftRotate [0..] matrix staggerVertically :: [a] \rightarrow [a] \\ staggerVertically :: [a] \rightarrow [a] \\ leftRotate :: Int \rightarrow [a] \rightarrow [a] \\ leftRotate :: Int \rightarrow [a] \rightarrow [a] \\ leftRotate :: xs = xs2 + xs1 \\ \textbf{where} \\ (xs1,xs2) = splitAt : xs \\ matAdd = chunksOf (dimX x) <math>\circ zipWith (+) (concat x) (concat y)
```

Figure C13: prMMTr, splitMatrix, staggerHorizontally, staggerVertically and matAdd definition.