An Introduction to (Deterministic) Parallelism in Haskell

Munich Lambda Meetup

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The Haskell Consultants

- Parallelism and Concurrency
- A first example
- Haskell and Effects
- Writing parallel programs
- ► (N Queens)
- Conclusions



- PhD (Utrecht University) 2004 on "Generic Haskell"
- Lecturer at Utrecht University 2007–2010
- Partner at Well-Typed 2010–



- Founded 1998.
- ► Haskell consulting (development, advice, support, training).
- ► Currently ~7 people working full-time in various places.
- Clients mainly in Europe and USA (most work done remotely).
- ► Also helped to set up the Industrial Haskell Group.



Parallelism and Concurrency

Parallelism

Running (parts of) programs in parallel on multiple cores (or nodes), in order to speed up the program.

Concurrency

Language constructs that support structuring a program as if it has many independent threads of control.



Using concurrency to implement parallelism is quite common, but not necessarily a good idea:

- reasoning about threads is difficult,
- communication between threads,
- exceptions,
- potential deadlocks and race conditions.

Often, code we want to parallelise is pure – it involves no side effects at all. So why introduce them just for parallelism?



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In Haskell, function application is free of side effects, and evaluation is non-strict:



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f x

In principle, we can run f in parallel with x:

- f might not need x at all, but no harm is done,
- f might need x immediately, then no harm is done,
- f might not need x immediately, then time is saved!



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- ▶ f might not need x immediately, then time is saved!

(The final case looks particularly attractive if \mathbf{x} produces a data structure lazily that is consumed by \mathbf{f} .)



The enemies of parallelism:

- ► there is overhead in running things in parallel,
- garbage collection is difficult to parallelize,
- ► non-strictness can not only be helpful, but also tricky:
 - ► we might run too many things we don't need,
 - it's unclear how far to evaluate speculatively,
 - we have to make clear how it interacts with GC.



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 - ▶ we have to make clear how it interacts with GC.

Conclusion

Fully automatic parallelism is still a future goal. For now, we need to help the compiler.



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Haskell supports multiple approaches to deterministic parallelism.



A few deterministic approaches:

- nested data parallelism (Data-Parallel Haskell, dph),
- flat data parallelism (repa),
- evaluation strategies (parallel),
- safe dataflow specification (monad-par).



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- flat data parallelism (repa),
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A few non-deterministic approaches:

- concurrency primitives (forkIO, ...),
- dataflow with side effects (monad-par),
- asynchronous computations (async),
- Cloud Haskell (distributed-process).



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 - data parallelism is about doing the same operations for many pieces of data; a particular common form that warrants dedicated support (dph, repa)



- Parallelism is "hot".
- Parallelising programs (even explicitly) is not trivial.
- Different forms of parallelism have different demands:
 - data parallelism is about doing the same operations for many pieces of data; a particular common form that warrants dedicated support (dph, repa)
 - task or control parallelism is about dividing the overall work into many parts – these approaches can be used for data parallelism, too (parallel, monad-par).



Given the lack of time, we have to limit ourselves, and will focus on the Par monad.



A first example

Example



Example

```
\begin{array}{l} \mbox{collatz}:: \mbox{Integer} \rightarrow \mbox{Integer} \\ \mbox{collatz} n \\ | \mbox{ even } n = n \ \mbox{'div'} \ 2 \\ | \ \mbox{odd} \ n = 3 \ast n + 1 \\ \mbox{collatzSeq} :: \mbox{Integer} \rightarrow \mbox{[Integer]} \\ \mbox{collatzSeq} = \mbox{takeWhile} \ (>1) \ . \ \mbox{iterate collatz} \\ \mbox{collatzSteps} :: \mbox{[Int]} \\ \mbox{collatzSteps} = \mbox{map} \ \mbox{(length . collatzSeq)} \ \mbox{[1..]} \end{array}
```

```
GHCi> collatzSeq 9
[9,28,14,7,22,11,34,17,52,26,13,40,20,10,5,16,8,4,2]
GHCi> take 10 collatzSteps
[0,1,7,2,5,8,16,3,19,6]
```

Let's find the maximum number of steps in a given range.



Parallelisation

$\begin{array}{l} \mbox{collatzMax}:: \mbox{Integer} \rightarrow \mbox{Integer} \rightarrow \mbox{Int} \\ \mbox{collatzMax} \mbox{ lo } hi = \mbox{maximum} (\mbox{map} (\mbox{length} \ . \ \mbox{collatzSeq}) \ \mbox{[lo } . \ . \ \mbox{hi}]) \end{array}$



Parallelisation

```
collatzMax :: Integer \rightarrow Integer \rightarrow Int
collatzMax lo hi = maximum (map (length . collatzSeq) [lo . . hi])
```

Binary division:

```
\begin{array}{l} \mbox{parCollatzMax} :: \mbox{Integer} \rightarrow \mbox{Integer} \rightarrow \mbox{Int} \\ \mbox{parCollatzMax} \mbox{ lo hi} = \mbox{ruPar} \\ \mbox{do} \\ \mbox{r1} \leftarrow \mbox{spawnP} \mbox{ (collatzMax lo mi)} \\ \mbox{r2} \leftarrow \mbox{spawnP} \mbox{ (collatzMax (mi + 1) hi)} \\ \mbox{m1} \leftarrow \mbox{get r1} \\ \mbox{m2} \leftarrow \mbox{get r2} \\ \mbox{return} \mbox{ (max m1 m2)} \\ \mbox{where} \\ \mbox{mi} = \mbox{(lo + hi) 'div' 2} \end{array}
```



Compile with:

\$ ghc -02 -threaded -rtsopts Collatz

Run with:

```
$ ./Collatz +RTS -N -s
```

Compared with a plain implementation provides modest speedup.



Compiler flags:

- -02 enables optimisation
- -threaded links in the threaded run-time system
- -rtsopts allows configuration of run-time system at run time
- -eventlog allows eventlog generation for debugging

Run-time system flags:

- -N runs on all available cores
- -s produces run-time statistics
- -1 generates an eventlog for debugging



Haskell and Effects

Effects

Java/C-like

```
int add0 (int x, int y) {
    return x + y;
}
```



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```
int add0 (int x, int y) {
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int add1 (int x, int y) {
    launch_missiles (now);
    return x + y;
}
```



Effects

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```
int add0 (int x, int y) {
    return x + y;
}
int add1 (int x, int y) {
    launch_missiles (now);
    return x + y;
}
```

Both functions have the same type!



Haskell

add0 :: Int \rightarrow Int \rightarrow Int add0 x y = x + y add1 :: Int \rightarrow Int \rightarrow IO Int add1 x y = **do** launch_missiles return (x + y)

Effectful computations are tagged by the type system!



Effects in Haskell's types

We have rather fine-grained control about effects just by looking at the types:

- A some type, no effect
 IO A IO, exceptions, random numbers, concurrency, ...
 Gen A random numbers only
 ST s A mutable variables only
 STM A software transactional memory log variables only
 State s A (persistent) state only
 Error A exceptions only
 Signal A time-changing value
 - All effect types share a common interface (monad; allows sequencing of operations and **do** notation).
 - ► New effect types can be defined. Effects can be combined.



More on **do** notation

```
interaction :: IO String
interaction = do
    putStrLn "Who are you?"
    name ← getLine
    putStrLn ("Hello, " + name + "!")
    return name
```



More on **do** notation

```
interaction :: IO String
interaction = do
putStrLn "Who are you?"
name ← getLine
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Look at the types:



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```

Look at the types:

$$(\gg)$$
 :: IO a \rightarrow (a \rightarrow IO b) \rightarrow IO b



No escape from IO

There's no function of type:

 $\text{IO } a \to a$

Example:

- ► An Int is a constant integer.
- ► An IO Int is an IO action yielding an integer.
- We shouldn't be able to forget about the potential side effects.



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unsafePerformIO :: IO $a \rightarrow a$



The Par monad

- very limited interface
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 $runPar:: Par \ a \to a$

- create an annotated parallel computation of type Par a
- run it with runPar
- obtain a deterministic result of type a



Writing parallel programs

```
\begin{array}{l} \mbox{parCollatzMax} :: \mbox{Integer} \rightarrow \mbox{Integer} \rightarrow \mbox{Int} \\ \mbox{parCollatzMax} \mbox{ lo hi} = \mbox{ruPar} \mbox{\$} \\ \mbox{do} \\ \mbox{r1} \leftarrow \mbox{spawnP} \mbox{ (collatzMax lo mi)} \\ \mbox{r2} \leftarrow \mbox{spawnP} \mbox{ (collatzMax lo mi)} \\ \mbox{m1} \leftarrow \mbox{get r1} \\ \mbox{m2} \leftarrow \mbox{get r2} \\ \mbox{return} \mbox{ (max m1 m2)} \\ \mbox{where} \\ \mbox{mi} = \mbox{(lo + hi) 'div' 2} \end{array}
```



A more recent approach to deterministic parallel programming:

- ► an interface with explicit forking of subcomputations,
- communication via write-once variables ensured deterministic results,
- ► reading a variable blocks until the result is available.



From Control.Monad.Par :

```
\begin{array}{rcl} \mbox{data} \mbox{ Par } a & -- \mbox{ abstract} \\ \mbox{instance } \mbox{Monad } \mbox{Par} \\ \mbox{data } \mbox{IVar } a & -- \mbox{ abstract} \\ \mbox{spawn} & :: \mbox{NFData } a \Rightarrow \mbox{Par } a \to \mbox{Par } (\mbox{IVar } a) \\ \mbox{spawnP} & :: \mbox{NFData } a \Rightarrow a \to \mbox{Par } (\mbox{IVar } a) \\ \mbox{get} & :: \mbox{IVar } a \to \mbox{Par } a \\ \mbox{runPar } & :: \mbox{Par } a \to a \end{array}
```

Let's ignore NFData for a moment.



```
\begin{array}{l} \mathsf{new}::\mathsf{Par}\;(\mathsf{IVar}\;a)\\ \mathsf{put}\;::\mathsf{NFData}\;a\Rightarrow\mathsf{IVar}\;a\to\mathsf{a}\to\mathsf{Par}\;()\\ \mathsf{get}\;::\mathsf{IVar}\;a\to\mathsf{Par}\;a\\ \mathsf{fork}\;::\mathsf{Par}\;()\to\mathsf{Par}\;() \end{array}
```

- The functions spawn and spawnP can be implemented in terms of the functions above.
- ► Writing twice to an IVar is an error.



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Here, we want to make sure that the result is fully computed before it is communicated to the consuming computation:

- the NFData type class contains functions for fully evaluating terms of a given type;
- it is used in spawn and put to make sure that results are fully evaluated before they're written to a write-once variable.



Static partitioning is bad:

- fixed number of tasks, limited speedup on many cores;
- difficult to balance the load;
- difficult to control granularity.

Let's create parallel tasks depending on the problem size.



```
\begin{array}{l} \mbox{parMap}:: \mbox{NFData} \ b \Rightarrow (a \rightarrow b) \rightarrow [a] \rightarrow \mbox{Par} \ [b] \\ \mbox{parMap} \ f \ xs = \mbox{do} \\ \ vs \leftarrow \mbox{mapM} \ (spawnP \ . \ f) \ xs \\ \ mapM \ get \ vs \end{array}
```



```
\begin{array}{l} \mbox{parMap}:: \mbox{NFData} \ b \Rightarrow (a \rightarrow b) \rightarrow [a] \rightarrow \mbox{Par} \ [b] \\ \mbox{parMap} \ f \ xs = \mbox{do} \\ \ vs \leftarrow \mbox{mapM} \ (spawnP \ . \ f) \ xs \\ \ mapM \ get \ vs \end{array}
```

```
\begin{array}{ll} \mathsf{mapM} :: (a \to \mathsf{Par} \ b) \to [a] \to \mathsf{Par} \ [b] \\ \mathsf{mapM} \ f \ [] &= \mathsf{return} \ [] \\ \mathsf{mapM} \ f \ (x : xs) = \textbf{do} \\ \mathsf{r} &\leftarrow \mathsf{f} \ x \\ \mathsf{rs} \leftarrow \mathsf{mapM} \ \mathsf{f} \ \mathsf{xs} \\ \mathsf{return} \ (\mathsf{r} : \mathsf{rs}) \end{array}
```



Sequential version:

collatzMax :: Integer \rightarrow Integer \rightarrow Int collatzMax lo hi = maximum (map (length . collatzSeq) [lo . . hi])

Parallel version:

 $parCollatzMax :: Integer \rightarrow Integer \rightarrow Int$ parCollatzMax lo hi =maximum (runPar (parMap (length . collatzSeq) [lo . . hi]))



Spawning a computation for each list element causes a granularity problem:

- too many too small computations are spawned too fast;
- ► we still get some speedup, but not as much as we'd like.



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A common solution is to chunk the list:

```
type ChunkSize = Int
chunk :: Int \rightarrow [a] \rightarrow [[a]]
chunk n xs = case splitAt n xs of
(ys,[]) \rightarrow [ys]
(ys,zs) \rightarrow ys : chunk n zs
```

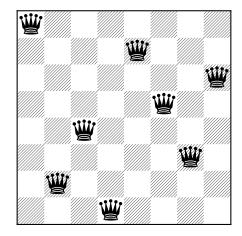


```
parCollatzMax :: ChunkSize \rightarrow Integer \rightarrow Integer \rightarrow Integer
parCollatzMax cs lo hi =
  maximum (
     concat (
        runPar (
          parMap
             (map (length . collatzSeq))
             (chunk cs [lo..hi])
```



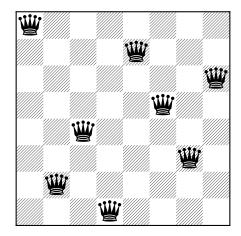
N Queens

The N Queens problem





The N Queens problem



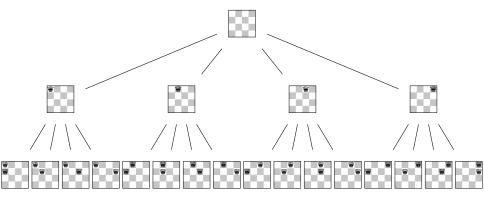
How many solutions for a given board size?



- Pick queens row by row.
- Generate a tree of all possible choices.
- Remove illegal choices from the tree.
- Traverse the tree, counting the number of valid solutions.

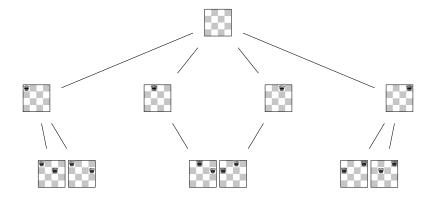


Generate, filter, explore





Generate, filter, explore







Conclusions

- Annotating a program for parallelisation is (relatively) easy.
- We can build domain-specific abstractions such as parMap.
- Deterministic results are guaranteed no deadlocks, no race conditions.
- ► We can focus on achieving speedup.



- The ParlO monad combines IO with Par at the price of determinism.
- The Async monad is similar to IO, but for concurrent applications.



- Time for some exercises now.
- Lots of online material.
- Simon Marlow's book.



Exercises

Exercises

- ► Try to write twice to a single IVar.
- Reproduce the Collatz example.
- Replace the Collatz function by the Fibonacci function what changes?
- Try to abstract and define a function

```
parMapChunked :: NFData b \Rightarrow
ChunkSize \rightarrow (a \rightarrow b) \rightarrow [a] \rightarrow Par [b]
```

- Try to abstract and define a "skeleton" for map-reduce.
- ► Reproduce the N Queens example.
- From Simon Marlow's materials: try Sudoku solving, k-means, conference timetable scheduling; all using the Par monad.

