

The Par Monad: Dataflow Parallelism

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In Control.Monad.Par,

```
newtype Par a = ...  
instance Applicative Par ...  
instance Monad Par ...  
  
runPar :: Par a -> a  
fork :: Par () -> Par ()  
  
data IVar a = ...  
new :: Par (IVar a)  
put :: NFData =>  
      IVar a -> a -> Par ()  
get :: IVar a -> Par a
```

put forces evaluation of its argument (NFData)

```
runPar $ do -- parmonad.hs  
  i <- new           -- Create IVar  
  j <- new  
  fork (put i (fib n)) -- Write result  
  fork (put j (fib m))  
  a <- get i         -- Wait for result  
  b <- get j  
  return (a+b)
```

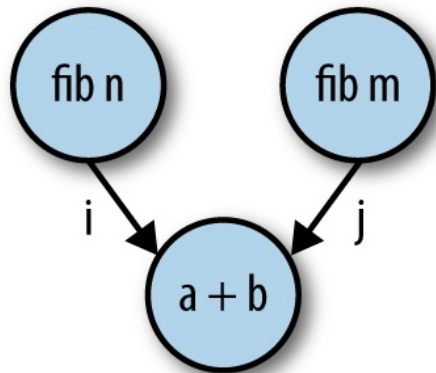
An IVar is a write-once variable

get waits for data to be *put*

Multiple *puts* to the same IVar cause a runtime error

Restrict each IVar to a single Par

The Par Monad: Dataflow Parallelism



Marlow, Fig. 4-1

```
runPar $ do -- parmonad.hs
  i <- new           -- Create IVar
  j <- new
  fork (put i (fib n)) -- Write result
  fork (put j (fib m))
  a <- get i         -- Wait for result
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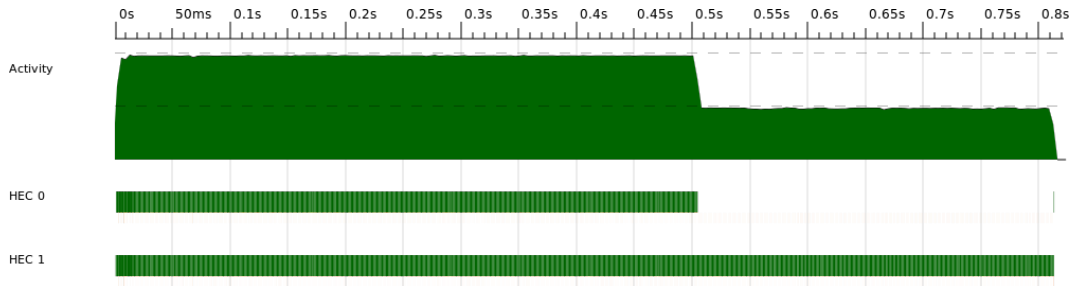
Restrict each IVar to a single Par

Running Marlow's parmonad.hs Example

```
$ stack ghc -- -O2 -threaded -rtsopts -eventlog parmonad.hs
```

```
$ ./parmonad 34 35 +RTS -N2 -ls
```

```
24157817
```



Works OK with -N2; obvious load-balancing problem. -N8 slows it down

Control.Monad.Par.{spawn,spawnP}: Fork and Return New IVar

```
spawn :: NFData a => Par a -> Par (IVar a) -- Spawn a process
spawn p = do i <- new                      -- Create a new IVar i
              fork $ do x <- p              -- Run p
                      put i x              -- Put the result in i
              return i                      -- Return the IVar i
```

```
spawnP :: NFData a -> a -> Par (IVar a) -- Evaluate pure expression
spawnP = spawn . return
```

```
runPar $ do
  i <- spawnP (fib n) -- Start fib n in parallel with
  j <- spawnP (fib m) -- Start fib m
  a <- get i          -- Wait for fib n to finish
  b <- get j          -- Wait for fib m to finish
  return (a+b)
```

Control.Monad.Par.{parMapM,parMap}

parMapM applies a function that works in the monad:

```
parMapM :: NFData b => (a -> Par b) -> [a] -> Par [b]
parMapM f as = do
  ibs <- mapM (spawn . f) as  -- Run each in parallel
  mapM get ibs               -- Wait for all list elements
```

parMap is similar but applies a pure function:

```
parMap :: NFData b => (a -> b) -> [a] -> Par [b]
parMap f as = do
  ibs <- mapM (spawn . return . f) as
  mapM get ibs
```

Actual implementations in Control.Monad.Par are more general

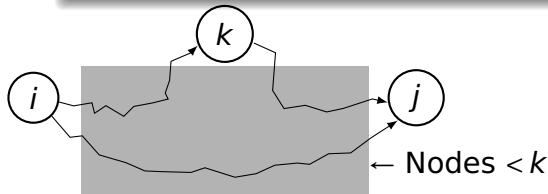
The Floyd-Warshall Shortest Paths Algorithm

The edge in g from i to j has weight $g[i][j]$

Vertices are numbered $0 \dots n$

In pseudocode,

```
shortestPath :: Graph -> Vertex -> Vertex -> Vertex -> Weight
shortestPath g i j 0 = weight g i j
shortestPath g i j k = min (shortestPath g i j (k-1))
                           (shortestPath g i k (k-1) +
                            shortestPath g k j (k-1))
```



Like *Fibonacci*, a recursive definition that should be implemented bottom-up with results recorded. $O(n^3)$ overall

Sparse Graph Representation: Maps of Maps

Marlow's code in `fwsparse/SparseGraph.hs`

```
import qualified Data.IntMap.Strict as Map

type Vertex = Int
type Weight = Int

type Graph = IntMap (IntMap Weight)

weight :: Graph -> Vertex -> Vertex -> Maybe Weight
weight g i j = do
    jmap <- Map.lookup i g
    Map.lookup j jmap
```

`IntMap` is tuned to better work with `Int` keys

$O(n^3)$ Sequential Implementation

```
shortestPaths :: [Vertex] -> Graph -> Graph
shortestPaths vs g = foldl' update g vs where -- For each vertex k
  update g k = Map.mapWithKey shortmap g where -- For each vertex i
    shortmap i jmap = foldr shortest Map.empty vs -- Shortest from i
    where
      shortest j m = case (old,new) of -- Update path from i to j via k
        (Nothing, Nothing) -> m -- No path
        (Nothing, Just w ) -> Map.insert j w m -- Found a new path
        (Just w, Nothing) -> Map.insert j w m -- Existing path only
        (Just w1, Just w2) -> Map.insert j (min w1 w2) m -- Best
      where
        old = Map.lookup j jmap -- Previous i → j path
        new = do w1 <- weight g i k -- i → k
                  w2 <- weight g k j -- k → j
                  return (w1+w2)
```

Running Sequential Floyd-Warshall

Random graph with 1000 vertices and 800 nodes:

```
$ stack ghc -- -O2 -rtsopts fwspare.hs  
$ ./fwspare 1000 800 +RTS -s  
Total time 14.531s ( 14.575s elapsed)
```

Fundamentally, three nested loops:

```
shortestPaths vs g = foldl' update g vs where  
  update g k = Map.mapWithKey shortmap g where  
    shortmap i jmap = foldr shortest Map.empty vs
```

Two are folds, which are difficult to parallelize unless operation is associative

However, mapWithKey is a map

Parallelizing Floyd-Warshall

mapWithKey is an unusual map over an *IntMap*.

We need a map that runs in the *Par* monad. Fortunately, *IntMap* provides

```
traverseWithKey :: Applicative t =>
  (Key -> a -> t b) -> IntMap a -> t (IntMap b)
```

and in *Traversable*,

```
traverse :: (Traversable t, Applicative f) =>
  (a -> f b) -> t a -> f (t b)
```

So we can update our *update* function to spawn *shortmap* in parallel:

```
update g k = runPar $ do
  m <- Map.traverseWithKey
    (\i jmap -> spawnP (shortmap i jmap)) g
  traverse get m -- get each IVar in the IntMap
```

Running Parallel Floyd-Warshall

```
$ stack ghc -- -O2 -threaded -rtsopts -eventlog fwparse1.hs
$ ./fwparse1 1000 800 +RTS -s -N1
    Total    time    6.091s  ( 6.150s elapsed)
$ ./fwparse1 1000 800 +RTS -s -N8
    Total    time   12.071s  ( 1.832s elapsed)
```

A $3.4\times$ speedup on 8 cores, but we beat the sequential version (?)

Note the total time increased substantially (parallel overhead), but the elapsed time decreased anyway