Folding Unfolded Polyglot FP for Fun and Profit Haskell and Scala

Can a left fold ever work over an infinite list? What about a right fold? Find out.

Learn about the other two functions used by functional programmers to implement mathematical induction: iterating and scanning. Learn about the limitations of the accumulator technique and about tupling, a technique that is the dual of the accumulator trick.

Part 4 - through the work of





As promised at the end of **Part 3**, let's now resume looking at **Tony**'s talk. We last saw him using **foldr** to implement the **flatten** function. In the next slide he uses **foldr** to implement the **filter** function.



Tony Morris @dibblego right folds replace constructors

Let's filter a list on predicate

Filter is a little bit ugly. Who's never heard of the filter function? Everyone has heard of it. They are talking about it outside right now. Filter. We are going to take a predicate and keep the elements that match that predicate. It gets a bit tricky.

```
filter a list on predicate (p)
Supposing
list = Cons A (Cons B (Cons C (Cons D Nil)))
```

What are we going to replace *Cons* with? We want to take the **head** at that *Cons*, the **A** or the **B** or the **C** or whatever, and we want to check if it matches a **predicate**. And if it does, we are going to *Cons* that value back on again. If it doesn't, we are just going to leave it off and then move on to the next *Cons* cell.

```
filter a list on predicate (p)
• let Cons = \x -> if p x then Cons x else id
• let Nil = Nil
```

So we want to say, given **x**, the element that I am visiting, if it matches the **predicate**, then **Cons x**, otherwise do nothing, **id**, the **identity function**. And replace **Nil** with **Nil**.



Tony Morris

filter a list on predicate (p)

```
Supposing
applyp = if p x then Cons x else id
list = applyp A (applyp B (applyp C (applyp D Nil)))
filter p list = foldr (\x -> if p x then Cons x else id) Nil list
filter p = foldr (\x -> if p x then Cons x else id) Nil
filter p = foldr (\x -> bool id (Cons x) (p x)) Nil
filter p = foldr (bool id . Cons <*> p) Nil
```

So, if you ever see this code, you'll know that it is filtering a list. Replace *Cons* with $(\x -> if p x then Cons else id)$ and *Nil* with *Nil*.

I happen to have a particular aversion to the **if then else** construct by the way. I wish it would die. It takes three arguments for a start. And for those who attended yesterday, what are my favourite functions? How many arguments do those take? One. So I don't like that one.

There is a function that does if then else that does take one argument. It is called **bool**. I have used it here. And it does if then else with the arguments the right way. There is a right way. It says, if $(p \ x)$, then $(Cons \ x)$, else id.

That takes one argument. The reason I like that function is because I can get down to that point there (last line above), and that point means I don't have to name so many identifiers. That's just my goal in life. That's not really the point. The point is that I can do that. I can look at this code and I can say: replace *Cons* with (bool id . *Cons* <*> p) and *Nil* with *Nil*. What is it going to do? It is going to filter the list on p. OK?

If you found it challenging at all to understand how that last refactoring of the definition of filter works, i.e. what the <*> operator is doing in the refactored definition, then you are not alone.

I am familiar with <*> as Applicative's apply function, which takes a function in a context and the function's argument in that

e.g. if I define function inc x = x + 1, then if inc is in a Maybe context and 3 is also in a Maybe context, I can use <*> to

filter $p = foldr (\langle x - \rangle bool id (Cons x) (p x)) Nil$ filter p = foldr (bool id . Cons <*> p) Nil

apply inc to 3 and return the result 4 in a Maybe context:

context and returns the result of applying the function to its argument, also in that context.

<*> is called apply or ap, but is also known as the advanced tie fighter (|+| being the plain tie fighter), the angry parent, and the sad Pikachu.



@philip schwarz



Tie fighter

[4]

> ((+) <*> inc) 3

e.g.



advanced Tie fighter

instance Functor ((->) a) where > Just inc <*> Just 3 $(a \rightarrow b) \rightarrow (r \rightarrow a) \rightarrow r \rightarrow b$ fmap = (.)Just 4 Similarly for a List context: instance Applicative ((->) a) where **pure** = const $a \rightarrow r \rightarrow a$ (< >) fgz = fx (gx) (r -> (a -> b)) -> (r -> a) -> r -> b > [inc] <*> [3]

Well, it turns out that a function is also an Applicative, so if I have a function f::a - b - c and a function g:a - b then I can use $\langle \rangle$ to create a function **h**:: **a**->**c** which passes its argument **a** to both **f** and **g** as follows: **f a** (**g a**).

> < > :: (r -> (a -> b)) -> (r -> a) -> r -> binc :: Num a => a -> a (+) :: Num a => a -> a -> a

We are using $\langle * \rangle$ to produce the $(\alpha \rightarrow \beta \rightarrow \beta)$ function that we need to pass to **foldr**

foldr :: $(\alpha \rightarrow \beta \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow \beta$

to be able to implement **filter** as follows

```
filter p = foldr (bool id . (:) <*> p) []
```

If we take the signature of <*>

```
<*> ::(r -> (a -> b)) -> (r -> a) -> r -> b
```

and make the following substitutions

a = Bool b = [a] -> [a] r = a

we can understand the signature in our context and see the types that <*> is operating on

```
<*> ::(a -> (Bool -> [a] -> [a])) -> (a -> Bool) -> a -> [a] -> [a]
```

even

> :type **bool id**

bool id . (:)

That's why we are able to refactor

 $x \rightarrow bool id ((:) x) (p x)$

to

```
bool id . (:) <*> p
```

```
foldr:: (\alpha \rightarrow \beta \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow \betafoldr f e []= efoldr f e (x: xs)= f x (foldr f e xs)
```

bool :: $a \rightarrow a \rightarrow Bool \rightarrow a$ Case analysis for the Bool type. bool x y p evaluates to x when p is False and evaluates to y when p is True. This is equivalent to if p then y else x; that is, one can think of it as an if-then-else construct with its arguments reordered. **is operating on is operating operat**

> 3) True

> > > :type (:) (:) :: a -> [a] -> [a]

```
> :type (bool id) . (:)
(bool id) . (:) :: a -> Bool -> [a] -> [a]
```

f

bool id :: $(a \rightarrow a) \rightarrow Bool \rightarrow a \rightarrow a$



For what it is worth, the order of the arguments of a conditional if C then P else Q does not look so 'wrong' to me if I think that in the lambda calculus, true is a binary selector function returning its first argument, false is a binary selector function returning its second argument, and the conditional function is one that takes a truth-valued conditional expression and two other alternative expressions and applies the selector function that is the result of the first expression to the two other expressions, which results in the selector function selecting one of the alternative expressions for evaluation.

Conditional Execution in the Lambda Calculus

true	=	λx.λy.x	true	3	4	β>	3
false	=	λx.λy.y	false	3	4	β>	4

if C then P else Q \rightarrow $\lambda c.\lambda p.\lambda q.cpq$



NOTE: Lazy evaluation is required to stop alternative expressions P and Q from both being evaluated before they are passed to the selector function that is the value of the conditional expression.



Tony's talk now turns to a very interesting subject which is a bit advanced and which we are going to cover separately in an upcoming part of this series, so we are now going to skip the next section of his talk and jump to its final part, in which he uses **foldr** to implement a function called **heador** and also recaps on some of the key things that we have learned in his talk.



Tony Morris

right folds replace constructors

Let's get the head of a list, or a default for no head
:: a -> List[a] -> a

I find this one quite interesting. So what this function does, is it takes an element **a**, and a list of **a**s, and it returns the first element out of the list. The first element. But it could be *Nil*. And if it's *Nil* return the other element that I passed in as an argument. OK, so if this list is *Nil*, return the **a** that is the first argument, otherwise return the first **a** in the list.

What am I going to replace *Cons* with? const. const the first argument. I agree. And *Nil*? That value there, that's the thing that we are returning if *Nil*, right? So given an entire list, if you ever see *Nil*, return that value, but if you see *Cons*, then return the value that is sitting right at the head of that *Cons*. That will do exactly what our requirements is for this function. So given this list

The head of a list, or default for no head

```
Supposing
list = Cons A (Cons B (Cons C (Cons D Nil)))
```

I want to return A, that's the first element of that list. Here is a function that given the A, and then the rest of the stuff after the *Cons*, just returns A. Replace *Cons*, with that. And replace *Nil*, with that default value that came in as an argument. That will do what it is we need to do.

The head of a list, or default for no head
• let Cons = \x _ -> x
• let Nil = the default





So there is the code.

The head of a list, or default for no head

```
Supposing
constant x _ = x
list = Cons A (Cons B (Cons C (Cons D Nil)))
heador thedefault list = foldr constant thedefault list
heador thedefault = foldr constant thedefault
heador = foldr constant
```

So who remembers writing **heador** yesterday, in the workshop? We wrote **heador**, right? This is **heador**. And we did it with pattern matching and all of that. We could have just written this: **fold right**, **const**, or **constant**, with that argument.

From https://hoogle.haskell.org/?hoogle=const:

```
const x is a unary function which evaluates to x for all inputs.
```

```
>>> const 42 "hello"
```

42

```
>>> map (const 42) [0..3]
```

```
[42,42,42,42]
```



Tony Morris @dibblego Observations

- There is no *order* specified, however, there is associativity.
- foldr may work on an infinite list.
 - ?

Fold right may work on an infinite list. So if I had the list from zero to infinity, and I said, call that heador function back here

```
heador thedefault list = foldr constant thedefault list
heador thedefault = foldr constant thedefault
heador = foldr constant
```

So here is a list, **zero** to **infinity**, and I say OK, pass some number in here, say four, I should get back **zero**, right? Because that is the first element of the list **zero** to **infinity**. OK? Do you want to see some code? This is an **infinite list** of ones:

\$ infinity = 1 : infinity

You might not believe me, in which case, do you believe me now?

\$ in	fi	nit	у																																							
[1,1	.,1	,1,	1,	1,1	1,1	, 1	, 1 ,	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,	1
,1,1	.,1	,1,	1,	1,1	1,1	, 1	, 1 ,	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,	1
, 1 , 1	, 1	,1,	1,	1,1	1,1	, 1	, 1 ,	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,	Ĺ
, 1 , 1	, 1	,1,	1,	1,1	1,1	, 1	, 1	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,	Ĺ
, 1 , 1	, 1	,1,	1,	1,1	1,1	, 1	, 1 ,	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,	Ĺ
, 1 , 1	, 1	,1,	1,	1,1	1,1	, 1	, 1 ,	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,	Ĺ
, 1 , 1	, 1	, 1,	1,	1,1	1,1	, 1	, 1 ,	,1,	1,	1,	1,	1,	1,1	1,1	.,1	, 1	,1,	1,	1,	1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	,1,	1,	1,	1,1	1,1	1,1	.,1	, 1	, 1	, 1	,1,	1,	1,1	1
etc,	e	tc																																								



Tony Morris
@dibblego

\$ heador a = foldr const a

OK, so, heador 99 infinity.

\$ heador 99 infinity

This will not go all the way to the right of that list, because when it gets there, there is just a *Cons* all the way. So I should get 1. And I do:

\$ heador 99 infinity

Fold right just worked on an **infinite** list. What's the complexity of **heador**? O(1). So, whether or not **fold right** will work on an **infinite list** depends on the **strictness** of the function that we are replacing **Cons** with.

Observations

- foldr may work on an infinite list.
 - There is no *order* specified, however, there is associativity.
 - Depends on the strictness of the given function.
 - •

const, the function I just used, is lazy, it ignores the second argument, and therefore it works on an infinite list.



Tony Morris

Observations

- foldr may work on an infinite list.
 - There is no *order* specified, however, there is associativity.
 - Depends on the strictness of the given function.
 - Replaces the *Nil* constructor *if it ever comes to exist*

It only replaces *Nil* if *Nil* ever comes to exist. It doesn't exist in **infinity**.

Observations

- foldr may work on an infinite list.
 - There is no *order* specified, however, there is associativity.
 - Depends on the strictness of the given function.
 - Replaces the *Nil* constructor *if it ever comes to exist*
- The expression **foldr** *Cons Nil* leaves the list unchanged.
 - In other words, passing the list constructors to **foldr** produces an *identity* function

How about this function: **foldr** *Cons Nil*? Leave the list alone. Replace *Cons* with *Cons* and *Nil* with *Nil*. It does nothing to the list. It's an **identity function**. What's the function that gives me an identity for bool? That's an interesting question? I just showed it to you. I'll let you think about that.



Tony just explained that doing a **right fold** with **Cons** and **Nil** produces the **identity function**.

In **Parts 1** and **3** we saw that doing a **left fold** with a flipped **Cons** and with **Nil** produces the **reverse** function.

The next slide illustrates the above.



Summary



Tony Morris @dibblego the key intuition

- left fold performs a *loop*, just like we are familiar with
- right fold performs constructor replacement

The key intuition is, the thing to take aways is, a **left fold** does a **loop**, and a **right fold** does **constructor replacement**.

If you always remember those two things you'll never go wrong.

from this we derive some observations

- left fold will never work on an infinite list
- right fold *may* work on an infinite list

Left fold will never work on an infinite list. We can see that in the loop. Right fold might. And these are just independent observations. They have nothing to do with programming languages. I have used Haskell as the example. These things are independent of the programming language.

from this we also solve problems

- product
- append
- map
- length

...

```
product = foldl (*) 1
append = flip (foldr Cons)
map f = foldr (Cons . f) Nil
length = foldl (const . (+ 1)) 0
```

So we could solve all these problems, as we just did.

Summary

I just want to be very clear on this: fold left does a loop, fold right does constructor replacement, and there are no footnotes required, this is precise.



• intuitively, this is what list folds do

- foldl performs a loop
- foldr performs constructor replacement
- this intuition is **precise** and requires no footnotes

I don't want you to ring me up next week and say 'you told me it does **constructor replacement** and it didn't this time'. Yes it did. It always will. OK. There are no footnotes required.

Tony Morris @dibblego	The End	Nil	
	Thanks!		

In the last part of his talk, **Tony** explained that whether or not **fold right** will work on an **infinite list** depends on the **strictness** of the function that we are replacing *Cons* with.

What about in **Scala**? Is there a notion of an **infinite list**? Can a **fold right** function be written which, when used to replace *Cons* with a suitable function, is able to handle an **infinite list**?



In **Part 1** I said that we weren't going to be modelling **infinity** in any of the **Scala** code in this series of slide decks. I have changed my mind about that. I do want to look at how to do **right folds** over large and **infinite lists**. And the technique that we'll be using to do that is also used to address limitations in the **accumulator technique**. So I want to use it for that too.

Which limitations in the **accumulator technique** I hear you ask? Let's cover that in this deck. And since we are there, let's cover a technique that is the dual of the **accumulator trick**, i.e. **tupling**.

And how can an **infinite list** be created in **Scala**? We can use the **iterate** function. Let's cover that in this deck too.

Also, there is something closely related to **folding** that we have not looked at yet, i.e. **scanning**. Let's also cover that in this deck.

So that means the following subjects will be addressed in **Part 5**:

- how to do right folds over large and infinite lists
- how to get around limitations in the applicability of the accumulator trick



In the next four slides we see how Sergei Winitzki describes the iterate function in his book.

2.3 Converting a single value into a sequence

An aggregation converts ("folds") a sequence into a single value; the <u>opposite operation</u> ("unfolding") converts a single value into a sequence. An example of this task is to compute the sequence of decimal digits for a given integer:

```
def digitsOf(x: Int): Seq[Int] = ???
```

```
scala> digitsOf(2405)
res0: Seq[Int] = List(2, 4, 0, 5)
```

We cannot implement digitsOf using map, zip, or foldLeft, because these methods work only if we already have a sequence; but the function digitsOf needs to create a new sequence. We could create a sequence via the expression (1 to n) if the required length of the sequence were known in advance. However, the function digitsOf <u>must produce a</u> <u>sequence</u> <u>whose length</u> <u>is determined by a condition that we cannot easily evaluate in advance</u>.

A general "unfolding" operation needs to build a sequence whose length is not determined in advance. This kind of sequence is called a stream. The elements of a stream are computed only when necessary (unlike the elements of List or Array, which are all computed in advance). The unfolding operation will compute the next element on demand; this creates a stream. We can then apply takeWhile to the stream, in order to stop it when a certain condition holds. Finally, if required, the truncated stream may be converted to a list or another type of sequence. In this way, we can generate a sequence of initially unknown length according to any given requirements.

The Scala library has a general stream-producing function Stream.iterate. This function has two arguments, the initial value and a function that computes the next value from the previous one:

```
scala> Stream.iterate(2) { x => x + 10 }
res0: Stream[Int] = Stream(2, ?)
```

The stream is ready to start computing the next elements of the sequence (so far, only the first element, 2, has been computed).





Sergei Winitzki in sergei-winitzki-11a6431 In order to see the next elements, we need to stop the stream at a finite size and then convert the result to a list:

scala> Stream.iterate(2) { x => x + 10 }.take(6).toList
res1: List[Int] = List(2, 12, 22, 32, 42, 52)

If we try to evaluate toList on a stream without first limiting its size via take or takeWhile, the program will keep producing more elements of the stream until it runs out of memory and crashes.

Streams are similar to sequences, and methods such as map, filter, and flatMap are also defined for streams. For instance, the method drop skips a given number of initial elements:

```
scala> Seq(10, 20, 30, 40, 50).drop(3)
res2: Seq[Int] = List(40, 50)
```

```
scala> Stream.iterate(2) { x => x + 10 }.drop(3)
res3: Stream[Int] = Stream(32, ?)
```

This example shows that in order to evaluate drop(3), the stream had to compute its elements up to 32 (but the subsequent elements are still not computed).

To figure out the code for digitsOf, we first write this function as a mathematical formula. To compute the digits for, say, n = 2405, we need to divide n repeatedly by 10, getting a sequence n_k of intermediate numbers ($n_0 = 2405$, n1 = 240, ...) and the corresponding sequence of last digits, $n_k \mod 10$ (in this example: 5, 0, ...). The sequence n_k is defined using mathematical induction:

- **Base case**: $n_0 = n$, where **n** is the given initial integer.
- Inductive step: $\mathbf{n}_{k+1} = \begin{bmatrix} \mathbf{n}_k \\ 10 \end{bmatrix}$ for $\mathbf{k} = 1, 2, ...$

Here $\left|\frac{\mathbf{n}_k}{10}\right|$ is the mathematical notation for the integer division by 10.



ite bia

Sergei Winitzki

in sergei-winitzki-11a6431

Let us tabulate the evaluation of the sequence n_k for n = 2405:

<i>k</i> =	0	1	2	3	4	5	6
$n_k =$	2405	240	24	2	0	0	0
$n_k \mod 10 =$	5	0	4	2	0	0	0

The numbers n_k will remain all zeros after k = 4. It is clear that the useful part of the sequence is before it becomes all zeros. In this example, the sequence n_k needs to be stopped at k = 4. The sequence of digits then becomes [5, 0, 4, 2], and we need to reverse it to obtain [2, 4, 0, 5]. For reversing a sequence, the Scala library has the standard method reverse. So, a complete implementation for digitsOf is:

```
def digitsOf(n: Int): Seq[Int] =
    if (n == 0) Seq(0) else { // n == 0 is a special case.
        Stream.iterate(n) { nk => nk / 10 }
        .takeWhile { nk => nk != 0 }
        .map { nk => nk % 10 }
        .toList.reverse
}
```

We can shorten the code by using the syntax such as ($\ \% \ 10$) instead of { nk => nk $\% \ 10$ },

```
def digitsOf(n: Int): Seq[Int] =
    if (n == 0) Seq(0) else { // n == 0 is a special case.
        Stream.iterate(n) (_ / 10 )
        .takeWhile ( _ != 0 )
        .map ( _ % 10 )
        .toList.reverse
}
```





The type signature of the method **Stream.iterate** can be written as

```
def iterate[A](init: A)(next: A => A): Stream[A]
```

and shows a close correspondence to a definition by mathematical induction. The base case is the first value, init, and the inductive step is a function, next, that computes the next element from the previous one. It is a general way of creating sequences whose length is not determined in advance.



Before we can look at that example, we are going to do the following:

the dual of the accumulator technique that we have already seen.

• The next two slides are from **Part 2** and remind us of how the accumulator technique can sometimes be used to transform a program so that it becomes tail recursive. Just skip the slides if they are still fresh in your mind.

I want to show you another example of using **iterate**. There is one example that I find quite interesting, which is the computation of **fibonacci numbers** using a technique which is called **tupling** and which is

- The two slides after that, remind us of how the **accumulator technique** can sometimes be used to improve the **efficiency** of a program.
- The subsequent slide briefly explains how the **tupling technique** can sometimes be used to increase the **efficiency** of a program.



def lengthS(s: Seq[Int]): Int =
 if (s.isEmpty) 0
 else 1 + lengthS(s.tail)

```
lengthS(Seq(1, 2, ..., 100000))
= 1 + lengthS(Seq(2, ..., 100000))
= 1 + (1 + lengthS(Seq(3, ..., 100000)))
= ...
```





Sergei Winitzki in sergei-winitzki-11a6431

The function body of lengthS will evaluate the inductive step, that is, the "else" part of the "if/else", about 100_000 times. Each time, the sub-expression with nested computations 1+(1+(...)) will <u>get larger</u>.

This intermediate sub-expression needs to be held somewhere in memory, until at some point the function body goes into the base case and returns a value. When that happens, the entire intermediate sub-expression will contain about 100_000_nested function calls still waiting to be evaluated.

This sub-expression is held in a special area of memory called stack memory, where the not-yet-evaluated nested function calls are held in the order of their calls, as if on a <u>"stack"</u>. Due to the way computer memory is managed, the <u>stack memory has a fixed</u> size and cannot grow automatically. So, when the intermediate expression becomes large enough, it causes an <u>overflow of the</u> <u>stack memory</u> and crashes the program.

A way to avoid stack overflows is to use a trick called <u>tail recursion</u>. Using tail recursion means rewriting the code so that all recursive calls occur at the end positions (at the "tails") of the function body. In other words, each recursive call must be itself the last computation in the function body, rather than placed inside other computations. Here is an example of tail-recursive code:

```
def lengthT(s: Seq[Int], res: Int): Int =
    if (s.isEmpty)
        res
    else
        lengthT(s.tail, 1 + res)
```

In this code, one of the branches of the **if/else** returns a fixed value without doing any **recursive calls**, while the other branch returns the result of a **recursive call** to **lengthT(...)**. In the code of **lengthT**, **recursive calls** never occur within any sub-expressions.

It is not a problem that the **recursive call** to **lengthT** has some sub-expressions such as 1 + **res** as its arguments, because all these sub-expressions will be computed before **lengthT** is **recursively called**.

The recursive call to **lengthT** is the last computation performed by this branch of the **if/else**. A **tail-recursive** function can have many **if/else** or **match/case** branches, with or without recursive calls; but all recursive calls must be always the last expressions returned.

The Scala compiler has a feature for checking automatically that a function's code is tail-recursive : the @tailrec annotation. If a function with a @tailrec annotation is not tail-recursive, or is not recursive at all, the program will not compile.

```
@tailrec def lengthT(s: Seq[Int], res: Int): Int =
if (s.isEmpty) res
else lengthT(s.tail, 1 + res)
```

Let us trace the evaluation of this function on an example:

```
lengthT(Seq(1,2,3), 0)
= lengthT(Seq(2,3), 1 + 0) // = lengthT(Seq(2,3), 1)
= lengthT(Seq(3), 1 + 1) // = lengthT(Seq(3), 2)
= lengthT(Seq(), 1 + 2) // = lengthT(Seq(), 3)
= 3
```

All sub-expressions such as 1 + 1 and 1 + 2 are computed before recursive calls to lengthT. Because of that, sub-expressions do not grow within the stack memory. This is the main benefit of tail recursion.

How did we rewrite the code of lengthS to obtain the tail-recursive code of lengthT? An important difference between lengthS and lengthT is the additional argument, res, called the accumulator argument. This argument is equal to an intermediate result of the computation. The next intermediate result (1 + res) is computed and passed on to the next recursive call via the accumulator argument. In the base case of the recursion, the function now returns the accumulated result, res, rather than 0, because at that time the computation is finished. Rewriting code by adding an accumulator argument to achieve tail recursion is called the accumulator technique or the "accumulator trick".

def lengthS(s: Seq[Int]): Int =
 if (s.isEmpty) 0
 else 1 + lengthS(s.tail)





Sergei Winitzki in sergei-winitzki-11a6431



That was a reminder of how the **accumulator technique** can sometimes be used to transform a program so that it becomes **tail recursive**.

The next two slides remind us of how the **accumulator technique** can sometimes be used to improve the **efficiency** of a program.



If we rewrite a recursive function using the accumulator technique then we end up with a tail recursive function, i.e. one that is stack-safe. In **Part 3** we saw that sometimes the version using the accumulator is more efficient than the version that doesn't, but at other times it can be less efficient, or even just as efficient.

Here is how Richard Bird expresses the case where it is more efficient: "By adding an extra parameter to a function we can sometimes improve the running time. The most important use of this idea is to eliminate possibly expensive # operations from a program"

$T(reverse)(n) = \Theta(n^2)$	$T(reverse')(n) = \Theta(n)$	
reverse:: $[\alpha] \rightarrow [\alpha]$ reverse []=[]=reverse (x : xs)=reverse (x : xs)=	$reverse' :: [\alpha] \rightarrow [\alpha]$ reverse' xs = accum [] xs	the accumulator version
$\frac{1}{1} \frac{1}{1} \frac{1}$	accum ws [] = ws accum ws (x : xs) = accum (x : ws) xs	is more efficient
$T(concat)(m,n) = \Theta(mn)$	$T(concat')(m,n) = \Theta(m^2n)$	
concat:: $[[\alpha]] \rightarrow [\alpha]$ concat[]=concat(xs:xss)=xs + concat xss	$concat' :: [[\alpha]] \rightarrow [\alpha]$ concat' xss = accum [] xss	the accumulator version
# = append	accum ws [] = ws accum ws (xs : xss) = accum (ws # xs) xss	
$T(append)(m,n) = \Theta(m)$	$T(append')(m,n) = \Theta(m)$	
append:: $[\alpha] \rightarrow [\alpha] \rightarrow [\alpha]$ append[]=append(x : xs)ys = x : (append xs ys)	$append'$ $:: [\alpha] \rightarrow [\alpha] \rightarrow [\alpha]$ $append'$ xs ys $accum$ ys $[] = ys$ $accum$ ys $[] = accum$ $(x : xs) = accum$ $(x : ys)$	the accumulator version is just as efficient

Since we saw, also in **Part 3**, that eliminating **foldr** leads to a **recursive** definition of a function and eliminating **foldl** leads to a **tail-recursive** definition of the function, one that uses an **accumulator**, this slide is simply a restatement of the previous one in which the **recursive** definition has been replaced by a definition using **foldr** and the **tail-recursive** definition (using an **accumulator**) has been replaced by a definition using **foldl**.

 $T(reverse)(n) = \Theta(n^2)$

 $\begin{array}{ll} reverse & :: & [\alpha] \rightarrow [\alpha] \\ reverse & = & foldr \ snoc \ [\] \\ & where \ snoc \ x \ xs = append \ xs \ [x] \end{array}$

 $T(reverse')(n) = \Theta(n)$

$$\begin{array}{rcl} reverse' & :: & [\alpha] \rightarrow [\alpha] \\ reverse' & = & foldl \ prefix \ [\] \\ & where \ prefix \ xs \ x = x \ : \ xs \end{array}$$

the *foldl* version is **more efficient**

$T(concat)(m,n) = \Theta(mn)$								
concat	::	$[[\alpha]] \to [\alpha]$						

 $concat :: [[\alpha]] \rightarrow [\alpha]$ concat = foldr append []

 $T(append)(m,n) = \Theta(m)$

append	::	$[\alpha] \to [\alpha] \to [\alpha]$
append xs ys	=	foldr (:) ys xs

 $T(concat')(m,n) = \Theta(m^2n)$ $concat' :: [[\alpha]] \rightarrow [\alpha]$ concat' = foldl append []

the *foldl* version is **less efficient**

 $T(append')(m,n) = \Theta(m)$

append' ::	:	$[\alpha] \to [\alpha] \to [\alpha]$
append'xs ys =	=	foldl scon ys (reverse'xs)
		where scon $xs x = x : xs$





That was a reminder of how the **accumulator technique** can sometimes be used to improve the **efficiency** of a program.

The next slide briefly explains how the **tupling technique** can sometimes be used to increase the **efficiency** of a program.

7.4 Tupling

...

The technique of program optimisation known as tupling is dual to that of accumulating parameters: a function is generalised, not by including an extra argument, but by including an extra result. Our aim in this section is to illustrate this important technique through a number of instructive examples.

7.4.2 Fibonacci function

Another example where **tupling** can improve the **order of growth** of the **time complexity** of a program is provided by the **Fibonacci** function.

The time to evaluate fib n by these equations is given by T(fib)(n), where

T(fib)(0) = O(1) T(fib)(1) = O(1)T(fib)(n+2) = T(fib)(n) + T(fib)(n+1) + O(1)

The timing function T(fib) therefore satisfies equations very like that of *fib* itself. It is easy to check by induction that $T(fib)(n) = \Theta(fib n)$, so the time to compute *fib* is proportional to the size of the result. Since $fib(n) = \Theta(\phi^n)$, where ϕ is the **golden ratio** $\phi = (1 + \sqrt{5})/2$, the time is therefore **exponential** in *n*. Now consider the function *fibtwo* defined by

fibtwo n = (fib n, fib (n + 1))

Clearly, fib n = fst(fibtwo n). Synthesis of a **recursive** program for fibtwo yields

 $\begin{array}{l} fibtwo \ 0 &= (0,1) \\ fibtwo \ (n+1) = (b,a+b), \ \text{where} \ (a,b) = fibtwo \ n \end{array}$

It is clear that this program takes linear time. In this example the tupling strategy leads to a dramatic increase in efficiency, from exponential to linear.





We have actually already been introduced numbers more efficiently by using tupling, but @philip_schwarz	(in Part 1) to the idea of computing Fibonacci at in a less formal way. See below for a reminder.
$ \begin{array}{cccc} fib & :: & Nat \rightarrow Nat \\ fib Zero & = & Zero \end{array} $ introduce tur	$fib :: Nat \rightarrow Nat$ $fib = fst \cdot foldn \ g \ (Zero, Succ Zero)$ where $g(m, n) = (n, m + n)$
$\begin{aligned} fib (Succ Zero) &= Succ Zero \\ fib (Succ (Succ n)) &= fib (Succ n) + fib n \end{aligned}$	$ \begin{aligned} foldn & :: & (\alpha \to \alpha) \to \alpha \to Nat \to \alpha \\ foldn \ h \ c \ Zero & = & c \\ foldn \ h \ c \ (Succ \ n) & = & h \ (foldn \ h \ c \ n) \end{aligned} $
	<pre>def fib(n: Nat): Nat = { def fst(pair: (Nat, Nat)): Nat = pair match { case (n,_) => n }</pre>
<pre>val fib: Nat => Nat = { case Zero => Zero case Succ(Zero) => Succ(Zero) introduce tug </pre>	<pre>def g(pair: (Nat, Nat)): (Nat, Nat) = pair match { case (m,n) => (n, m + n) }</pre>

case Succ(Succ(n)) => fib(Succ(n)) + fib(n)

introduce tupling fst(foldn(g, (Zero, Succ(Zero)), n))

```
def foldn[A](h: A => A, c: A, n: Nat): A =
 n match {
    case Zero => c
    case Succ(n) => h(foldn(h,c,n))
```

On the previous slide we saw a definition of the **fibonacci** function that uses **tupling** in conjunction with **foldn**, which operates on the **Nat** type, i.e. **Zero**, **Succ(Zero)**, **Succ(Succ(Zero)**), etc.

What if we want to use **Int** rather than **Nat**? Just for curiosity: does it make any sense to try and use **foldLeft** to write a definition of the **fibonacci** function that uses **tupling**?

It does make some sense. Here is how we can do it:

```
def fibFoldLeft(n: Int): Long =
  (1 to n) foldLeft ((0L, 1L)){
    case ((f1, f2), _) => (f2, f1 + f2)
  }._1
```

The odd thing about the above function is that while it creates a **sequence** 1 to n, it uses the sequence purely to determine how many times to **iterate**: it doesn't actually use the values in the **sequence** – e.g. it would work just as well if the **sequence** were 101 to 100 + n rather than 1 to n.

Still, it works. Let's compare the time taken to compute a large-ish **Fibonacci** number using both **fibFoldLeft** and the traditional definition of the **Fibonacci** function:

```
def fibonacci(n: Int): Long = n match {
   case 0 => 0L
   case 1 => 1L
   case n =>
     fibonacci(n - 1) + fibonacci(n - 2)
}
```







Eleven slides ago I said I wanted to show you a second example of using the **iterate** function. Now that we have gone through the following we are finally in a position to look at that example:

- a refresher of the accumulator technique
- a brief introduction to the **tupling technique**
- a look at how to implement a Fibonacci function using tupling

The example consists of implementing the Fibonacci function using iterate. We are going to use the same tupling technique that we have just used in fibFoldLeft. When Sergei Winitzki introduced the iterate function, we saw that it is a function that returns a Stream, e.g. Stream.iterate(2) { x => x + 10 }. Since Scala 2.13 however, the Stream class is deprecated in favour of LazyList, so we are going to use LazyList.iterate, whose signature is otherwise identical to that of Stream.iterate. The fibIterate function below is similar to fibFoldLeft from the previous slide in that it uses the same tupling approach, but it is simpler because it doesn't have to create a sequence whose data it doesn't even need. It also performs the same way.

/** An infinite LazyList that repeatedly applies a given function
 * to a start value.

```
* @param start the start value of the LazyList
* @param f the function that's repeatedly applied
* @return the Stream returning the infinite sequence
* of values `start, f(start), f(f(start)), ...`
*/
```

def iterate[A](start: A)(f: A => A): LazyList[A] = ...

https://www.scala-lang.org/blog/2018/06/13/scala-213-collections.html

LazyList Is Preferred Over Stream

Stream is deprecated in favor of LazyList. As its name suggests, a LazyList is a linked list whose elements are lazily evaluated. An important semantic difference with Stream is that in LazyList both the head and the tail are lazy, whereas in Stream only the tail is lazy.



Remember how the naive Scala definition of the fibonacci function was taking over 80 seconds to compute the 50 th Fibonacci number? Well the Haskell version seems to be even slower! Here is the Haskell version fibIterate. Just like the Scala version, it is blazingly fast. FibIterate n = fst (fib2 !! n) where fib2 = iterate f (0,1) where f (a,b) = (b,a+b)	<pre>> fib 10 55 > fib 20 6765 > fib 30 832040 > fib 40 I got bored of waiting so I interrupted it</pre>
<pre>> fibIterate 50</pre>	
12586269025	
> TIDILEFALE 10000 > 226/476/876/217822666216120051075/22102021/84606800620065647600746800814421666622681555055126227240	255920652226909261502727
33044704070431703200021012003107545310302140400000005900504705974000001442100002500155555555756 347904838652682630408974630564318873545443695598274916066020998841839338646527313000888302692356736	233020033320000301333737 1313511757979743785 <i>44</i> 137
5775070500520020507005270505075100755754545055550277510000020550040527515000000502052550750	182989697916131278562650
321902203043477010022047305109005270900351343001595025072750025075100512005754207054552155151050700	.025721143349611800230912
082870460889239623288354615057765832712525460935911282039252853934346209042452489294039017062338889	910858410651831733604374
707370085526317673257330037128710375877768077700263058370657728301616377089601787263786272128352581	128205163702080803320090
057079200643674262023897831114700540749984592503606335609338838319233867830561364353518921332797329	081337376476576339897639
07075200045074202025057051114700540745504552505000555005550050515255007050501504555510521552757525	647767428472748674114620
22/2340/0029201//9535005/09950910491/54/000095104105014052255021/40505/52124022050509210529/7010400	6877532834823740024114550
19731397776773679108710679780787718035379131176778977659089938635759377897789776777061977703376386	740040213303432974969020
28228145022418826817682802072002624705622117102101201052160704607622722525777580252526722552750427884245	040677155557700564504420
265261455554166266176656556726656547556251171651612515551657546676527575652555567725525755457664545	085871350167603170719031
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/20000340012903213010//2020/03531000240112/024535/200532505305//59504500/251//445150515590009051000	032203431032220400032400
2/376825907/1903063726508/1/7/081036007852730/5030707165//3156/7137815768800805072030005554/451500	7052022256/8/6/05106112/
002003133/007/200302040/0001320430/0030/40113120//322/40021330042330/1123044110202200103/1033130032	
\0113134132341332413300\17305\113\133535\0303133300020503020500500500505541005030205430\01\343\01\11515	



Next, let's look at some of the things that Richard Bird has to say about the iterate function.

@philip_schwarz

...there are three kinds of list:

- A *finite* list, which is built from (:) and []; for example, 1:2:3:[]
- A *partial* list, which is built from (:) and **undefined**; for example, the list **filter** (<4) [1..] is the **partial** list 1:2:3:**undefined**. We know there is no integer after 3 that is less than 4, but **Haskell** is an evaluator, not a theorem prover, so it ploughs away without success looking for more answers.
- An *infinite* list, which is built from (:) alone; for example, [1..] is the *infinite* list of nonnegative integers.

All three kinds of list arise in everyday programming. Chapter 9 is devoted to exploring the world of **infinite lists** and their uses. For example, the prelude function **iterate** returns an **infinite list**:

```
iterate :: (a \rightarrow a) \rightarrow a \rightarrow [a]
iterate f x = x : iterate f (f x)
```

In particular, *iterate* (+1) 1 is an **infinite** list of the positive integers, a value we can also write as [1..].

```
The type of 1 `div` 0 is an integral number.
ghci> :type 1 `div` 0
1 `div` 0 :: Integral a => a
```

The expression 1 `div` 0 is therefore well-formed and possesses a value.

```
ghci> 1 `div` 0
*** Exception: divide by zero
```

GHCi returns an error message. So what is the type of $1 \circ div$ 0? The answer is that it is a special value, written mathematically as \bot and pronounced 'bottom'. In fact, **Haskell** provides a predefined name for this value, except that it is called **undefined**, not \perp .

ghci> :type undefined undefined :: a

ghci> undefined
*** Exception: Prelude.undefined

Haskell is not expected to produce the value ⊥. It may return with an error message, or remain perpetually silent, computing an infinite loop, until we interrupt the computation. It may even cause GHCi to crash.





Richard Bird

...consider the following three definitions of the standard prelude function *iterate*:

iterate1 f x = x: iterate1 f (f x)iterate2 f x = xs where xs = x: map f xsiterate3 f x = x: map f (iterate3 f x)

All three functions have type $(a \rightarrow a) \rightarrow a \rightarrow [a]$ and produce an **infinite list** of **iterates** of f applied to x. The three functions are equal... The first definition is the one used in the standard prelude, but it does not create a **cyclic** list. The second definition does, and the third is obtained from the second by eliminating the *where* clause.

Assuming f x can be computed in constant time, the first definition takes $\Theta(n)$ steps to compute the first n elements of the result, but the third takes $\Theta(n^2)$ steps:

iterate3 (2 *) 1= 1 : map (2 *)(iterate3 (2 * 1)) = 1 : 2 : map (2 *) (map (2 *) (iterate3 (2 * 1))) = 1 : 2 : 4 : map (2 *) (map (2 *) (map (2 *) (iterate3 (2 * 1))))

Evaluating the n^{th} element requires n applications of (2 *), so it takes $\Theta(n^2)$ to produce the first n elements.

That leaves the second definition. Does it take **linear** or **quadratic** time? The evaluation of *iterate* 2 (2 *) 1 proceeds as follows:

 xs
 where xs = 1 : map (2 *) xs

 1: ys
 where ys = map (2 *) (1: ys)

 1: 2: zs
 where zs = map (2 *) (2: zs)

 1: 2: 4: ts
 where ts = map (2 *) (4: zs)

Each element of the result is produced in constant time, so *iterate* 2 (2 *) 1 takes $\Theta(n)$ steps to produce *n* elements.





Richard Bird



Now that we have covered the **iterate** function and the **tupling** technique, let's turn to the subject of **scanning**.

In the next two slides we see how Sergei Winitzki describes scanning in his book.

2.4 Transforming a sequence into another sequence

We have seen methods such as map and zip that transform sequences into sequences. However, <u>these methods cannot express a</u> general transformation where the elements of the new sequence are defined by induction and depend on previous elements.

An example of this kind is computing the **partial sums** of a given sequence x_i , say $b_k = \sum_{i=0}^{k-1} x_i$. This formula defines $b_0 = 0$, $b_1 = x_0$, $b_2 = x_0 + x_1$, $b_3 = x_0 + x_1 + x_2$, etc. A definition via mathematical induction may be written like this:

- Base case: $b_0 = 0$
- Inductive step: Given b_k , we define $b_{k+1} = b_k + x_k$ for k = 0,1,2,...

The **Scala** library method **scanLeft** implements a general, **sequence** to **sequence** transformation defined in this way. The code implementing the partial sums is

```
def partialSums(xs: Seq[Int]): Seq[Int] =
    xs.scanLeft(0){ (x, y) => x + y }
```

scala> partialSums(Seq(1, 2, 3, 4))
val res0: Seq[Int] = List(0, 1, 3, 6, 10)

The first argument of scanLeft is the base case, and the second argument is an updater function describing the inductive step. In general, the type of elements of the second sequence is different from that of the first sequence. The updater function takes an element of the first sequence and a previous element of the second sequence, and returns the next element of the second sequence. Note that the result of scanLeft is one element longer than the original sequence, because the base case provides an initial value.

Until now, we have seen that **foldLeft** is sufficient to re-implement almost every method that works on sequences, such as map, filter, or flatten. Let us show, as an illustration, how to implement the method scanLeft via foldLeft.

The Science of Functional Programming

A tutorial, with examples in Scala

Sergei Winitzki



Sergei Winitzki in sergei-winitzki-11a6431 In the implementation, the accumulator contains the previous element of the second sequence together with a growing fragment of that sequence, which is updated as we iterate over the first sequence. The code is

```
def scanLeft[A, B](xs: Seq[A])(b0: B)(next: (B, A) => B): Seq[B] = {
  val init: (B, Seq[B]) = (b0, Seq(b0))
  val (_, result) = xs.foldLeft(init) {
    case ((b, seq), x) =>
      val newB = next(b, x)
      (newB, seq :+ newB)
  }
  result
}
```

To implement the (nameless) **updater function** for **foldLeft** we used the **Scala** feature that makes it easier to define functions with several arguments containing tuples. In our case, the **updater function** in **foldLeft** has two arguments: the first is a tuple (B, **Seq**[B]), the second is a value of type A. The pattern expression **case** ((b, seq), x) => appears to match a nested tuple. In reality, this expression matches the two arguments of the **updater function** and, at the same time, destructures the tuple argument as (b, seq).



The Science of Functional Programming A tutorial, with examples in Scala





@philip_schwarz

While in his explanation of the scanLeft function, Sergei showed us (for illustration only) that it is possible to implement scanLeft using foldLeft, I was a bit surprised by the fact that he did not mention that the meaning of the scanLeft function is very closely related to that of foldLeft.

That close relationship is one of the things we are going to learn in the next five slides, in which **Richard Bird** explains left and right scans.

Before we do that, I am just going to have a go at a **Haskell** version of **Sergei's** implementation of **scanLeft** in terms of **foldLeft** (again, for illustration only).

```
def scanLeft[A, B](xs: Seq[A])(b0: B)(next: (B, A) => B): Seq[B] = {
  val init: (B, Seq[B]) = (b0, Seq(b0))
  val (_, result) = xs.foldLeft(init) {
    case ((b, seq), x) =>
      val newB = next(b, x)
      (newB, seq :+ newB)
  }
  result
}
```

scala> scanleft(List(1,2,3,4))(0)(_+_)
val res0: Seq[Int] = List(0, 1, 3, 6, 10)
scala>



4.5.2 Scan left

Sometimes it is convenient to apply a *foldl* operation to every initial segment of a list. This is done by a function *scanl* pronounced 'scan left'. For example,

 $scanl (\bigoplus) e [x_0, x_1, x_2] = [e, e \bigoplus x_0, (e \bigoplus x_0) \bigoplus x_1, ((e \bigoplus x_0) \bigoplus x_1) \bigoplus x_2]$

In particular, *scanl* (+) 0 computes the list of accumulated sums of a list of numbers, and *scanl* (×) 1 [1..*n*] computes a list of the first *n* factorial numbers. ... We will give two programs for *scanl*; the first is the clearest, while the second is more efficient. For the first program we will need the function *inits* that returns the list of all initial segments of a list. For Example,

inits $[x_0, x_1, x_2] = [[], [x_0], [x_0, x_1], [x_0, x_1, x_2]]$

The **empty list** has only one **segment**, namely the **empty list** itself; A list (x: xs) has the **empty list** as its shortest **initial segment**, and all the other **initial segments** begin with x and are followed by an **initial segment** of xs. Hence

inits :: $[\alpha] \rightarrow [[\alpha]]$ *inits* [] = [[]]*inits* (x:xs) = []:map(x:)(inits xs)

The function *inits* can be defined more succinctly as an instance of *foldr* :

inits = foldr f[[]] where f x xss = []: map(x:) xss

Now we define

```
scanl :: (\beta \rightarrow \alpha \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow [\beta]
scanl f e = map (fold f e).inits
```

This is the clearest definition of *scanl* but it leads to an **inefficient program**. The function *f* is applied *k* times in the evaluation of





Richard Bird

foldl f e on a list of length k and, since the **initial segments** of a list of length n are lists with lengths 0,1,...,n, the function f is applied about $n^2/2$ times in total.

Let us now synthesise a more efficient program. The synthesis is by an **induction argument** on xs so we lay out the calculation in the same way.

<....not shown...>

In summary, we have derived

 $\begin{array}{ll} scanl & :: (\beta \rightarrow \alpha \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow [\beta] \\ scanl \ f \ e \ [\] & = [e] \\ scanl \ f \ e \ (x: xs) & = e: \ scanl \ f \ (f \ e \ x) \ xs \end{array}$

This program is **more efficient** in that function f is applied exactly n times on a list of length n.



Note the similarities and differences between *scanl* and *foldl*, e.g. the left hand sides of their equations are the same, and their signatures are very similar, but *scanl* returns [β] rather than β and while *foldl* is **tail recursive**, *scanl* isn't.

$foldl (\bigoplus) e [x_0, x_1, x_2]$	scanl $(\bigoplus) e [x_0, x_1, x_2]$
Ļ	Ļ
$((e \oplus x_0) \oplus x_1) \oplus x_2$	$[e, e \oplus x_0, (e \oplus x_0) \oplus x_1, ((e \oplus x_0) \oplus x_1) \oplus x_2]$
foldl :: $(\beta \rightarrow \alpha \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow \beta$	<i>scanl</i> :: $(\beta \rightarrow \alpha \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow [\beta]$
foldl f e [] = e	scanl $f e [] = [e]$
foldl $f e(x:xs) = foldl f(f e x) xs$	scanl $f e(x:xs) = e:$ scanl $f(f e x) xs$





Richard Bird



By the way, just for completeness, the appendix of **Richard Bird's** book, contains the following additional definition of *scanl*

$$scanl :: (\beta \to \alpha \to \beta) \to \beta \to [\alpha] \to [\beta]$$

$$scanl f e xs = e : scanl' f e xs$$

where scanl' f e [] = []

$$scanl' f e (y; ys) = scanl f (f e y) ys$$

And just for comparison, here are the other definitions of *scanl*

$$scanl (\bigoplus) e [x_0, x_1, x_2]$$

$$\downarrow$$

$$[e, e \bigoplus x_0, (e \bigoplus x_0) \bigoplus x_1, ((e \bigoplus x_0) \bigoplus x_1) \bigoplus x_2]$$



 $scanl :: (\beta \to \alpha \to \beta) \to \beta \to [\alpha] \to [\beta]$ scanl f e [] = [e]scanl f e (x:xs) = e : scanl f (f e x) xs

 $scanl :: (\beta \to \alpha \to \beta) \to \beta \to [\alpha] \to [\beta]$ scanl f e = map (foldl f e) . inits

 $inits :: [\alpha] \rightarrow [[\alpha]]$ inits [] = [[]]inits (x:xs) = []: map (x:)(inits xs)

4.5.3 Scan right

The dual computation is given by *scanr*.

 $\begin{array}{ll} scanr & :: \ (\alpha \to \beta \to \beta) \to \beta \to [\alpha] \to [\beta] \\ scanr \ f \ e &= map \ (foldr \ f \ e) \ . \ tails \end{array}$

The function *tails* returns the tail segments of a list. For example,

tails
$$[x_0, x_1, x_2] = [[x_0, x_1, x_2], [x_1, x_2], [x_2], []]$$

$$\begin{array}{ll} scanl & :: \ (\beta \to \alpha \to \beta) \to \beta \to [\alpha] \to [\beta] \\ scanl \ f \ e &= map \ (foldl \ f \ e) \ . \ inits \end{array}$$

inits $[x_0, x_1, x_2] = [[], [x_0], [x_0, x_1], [x_0, x_1, x_2]]$

Note that while *inits* produces a list of **initial segments** in **increasing** order of length, tails produces the **tail segments** in **decreasing** order of length. We can define *tails* by

tails:: $[\alpha] \rightarrow [[\alpha]]$ tails []= [[]]tails (x:xs) = (x:xs) : tails xs

The corresponding efficient program for *scanr* is given by

 $\begin{array}{ll} scanr & :: (\alpha \rightarrow \beta \rightarrow \beta) \rightarrow \beta \rightarrow [\alpha] \rightarrow [\beta] \\ scanr f e [] & = [e] \\ scanr f e (x:xs) & = f x (head ys) : ys \\ & & \text{where } ys = scanr f e xs \end{array}$

for reference: a reminder of what *scanr* and *scanl* do

$$scanr (\bigoplus) e [x_0, x_1, x_2]$$

$$\downarrow$$

$$[(x_0 \bigoplus (x_1 \bigoplus (x_2 \bigoplus e))), x_1 \bigoplus (x_2 \bigoplus e), x_2 \bigoplus e, e]$$

$$inits :: [\alpha] \rightarrow [[\alpha]]$$

$$inits [] = [[]]$$

$$inits (x: xs) = []: map (x:)(inits xs)$$

$$scanl :: (\beta \to \alpha \to \beta) \to \beta \to [\alpha] \to [\beta]$$

$$scanl f e [] = [e]$$

$$scanl f e (x:xs) = e : scanl f (f e x) xs$$

scanl $(\bigoplus) e [x_0, x_1, x_2]$ $[e, e \oplus x_0, (e \oplus x_0) \oplus x_1, ((e \oplus x_0) \oplus x_1) \oplus x_2]$





Richard Bird

Let's try out <i>scanl</i> and <i>scanr</i> with (+) and 0	$inits :: [\alpha] \rightarrow [[\alpha]]$ $inits [] = [[]]$ $inits (x:xs) = []: map (x:) (inits xs)$	$\begin{array}{ll} tails & :: \ [\alpha] \rightarrow [[\alpha]] \\ tails \left[\right] & = \left[\left[\right] \right] \\ tails (x:xs) = (x:xs) : tails xs \end{array}$
	<i>inits</i> $[x_0, x_1, x_2] = [[], [x_0], [x_0, x_1], [x_0, x_1, x_2]]$	<i>tails</i> $[x_0, x_1, x_2] = [[x_0, x_1, x_2], [x_1, x_2], [x_2], []]$
@philip_schwarz scanl	$scanl :: (\beta \to \alpha \to \beta) \to \beta \to [\alpha] \to [\beta]$ scanl f e = map (foldl f e) . inits	$scanr :: (\alpha \to \beta \to \beta) \to \beta \to [\alpha] \to [\beta]$ scanr f e = map (foldr f e) . tails
applies <i>foldl</i> to every initial segment of a list	$scanl (\bigoplus) e [x_0, x_1, x_2]$ \downarrow $[e, e \bigoplus x_0, (e \bigoplus x_0) \bigoplus x_1, ((e \bigoplus x_0) \bigoplus x_1) \bigoplus x_2]$	$scanr (\bigoplus) e [x_0, x_1, x_2]$ \downarrow $[(x_0 \bigoplus (x_1 \bigoplus (x_2 \bigoplus e))), x_1 \bigoplus (x_2 \bigoplus e), x_2 \bigoplus e, e]$
scanr		
applies <i>foldr</i> to every tail segment of a list	<i>inits</i> $[2,3,4] = [[], [2], [2,3], [2,3,4]]$	<i>tails</i> [2,3,4] = [[2,3,4], [3,4], [4], []]
	$ \begin{array}{l} foldl (+) 0 [] = 0 \\ foldl (+) 0 [2] = 2 \\ foldl (+) 0 [2,3] = 5 \\ foldl (+) 0 [2,3,4] = 9 \end{array} $	foldr (+) 0 [2,3,4] = 9 foldr (+) 0 [3,4] = 7 foldr (+) 0 [4] = 4 foldr (+) 0 [] = 0
	<i>scanl</i> (+) 0 [2,3,4] = [0, 2, 5, 9]	<i>scanr</i> (+) 0 [2,3,4] = [9,7,4,0]
	<i>scanl</i> (+) 0 [2,3,4] ↓	<i>scanr</i> (+) 0 [2,3,4] ↓
	[0, 0+2, (0+2)+3, ((0+2)+3)+4]	[(2 + (3 + (4 + 0))), 3 + (4 + 0), 4 + 0, 0]



Now that we have seen the **iterate** function and the **scanning** functions, let's see how **Sergei Winitzki** describes that fact that **folding**, **iterating** and **scanning** are what we use in **functional programming** to implement **mathematical induction**.

2.5 Summary

We have seen a broad overview of translating mathematical induction into Scala code.

What problems can we solve now?

- Compute mathematical expressions involving arbitrary recursion.
- Use the accumulator trick to enforce tail recursion.
- Implement functions with type parameters.
- Use arbitrary inductive (i.e., recursive) formulas to:
 - convert sequences to single values (aggregation or "folding");
 - create new sequences from single values ("unfolding");
 - transform existing sequences into new sequences.

Definition by induction	Scala code example
$f([]) = b; f(s \leftrightarrow [x]) = g(f(s), x)$	<pre>f(xs) = xs.foldLeft(b)(g)</pre>
$x_0 = b$; $x_{k+1} = g(x_k)$	<pre>xs = Stream.iterate(b)(g)</pre>
$y_0 = b; y_{k+1} = g(y_k, x_k)$	<pre>ys = xs.scanLeft(b)(g)</pre>

Table 2.1: Implementing mathematical induction

Table 2.1 shows Scala code implementing those tasks. <u>Iterative calculations are implemented by translating mathematical induction directly into code</u>. In the functional programming paradigm, the programmer does not need to write any loops or use array indices. Instead, the programmer reasons about sequences as mathematical values: "Starting from this value, we get that sequence, then transform it into this other sequence," etc. This is a powerful way of working with sequences, dictionaries, and sets. Many kinds of programming errors (such as an incorrect array index) are avoided from the outset, and the code is shorter and easier to read than conventional code written using loops.



Sergei Winitzki in sergei-winitzki-11a6431



The next slide is a reminder of the **Functional Programming triad** of **folding**, **scanning** and **iteration**, with some very simple examples of their usage in **Scala** and **Haskell**.

Definition by induction	Scala code example
$f([]) = b; f(s \leftrightarrow [x]) = g(f(s), x)$	<pre>f(xs) = xs.foldLeft(b)(g)</pre>
$x_0 = b$; $x_{k+1} = g(x_k)$	<pre>xs = Stream.iterate(b)(g)</pre>
$y_0 = b$; $y_{k+1} = g(y_k, x_k)$	ys = xs.scanLeft(b)(g)

Implementing mathematical induction

>>Haskell

Haskell> foldl (+) 0 [1,2,3,4]

```
10
```

```
Haskell> take 4 (iterate (+ 1) 1)
[1,2,3,4]
```

```
Haskell> scanl (+) 0 [1,2,3,4]
[0,1,3,6,10]
```

Haskell>

5Scala

```
scala> List(1,2,3,4).foldLeft(0)(_+_)
val res0: Int = 10
scala> Stream.iterate(1)(_ + 1).take(4).toList
val res1: List[Int] = List(1, 2, 3, 4)
scala> List(1,2,3,4).scanLeft(0)(_+_)
val res2: List[Int] = List(0, 1, 3, 6, 10)
scala>
```



Remember when we tried to compute the ten-thousandth Fibonacci number using the traditional recursive definition of the fibonacci function, and we got a stack overflow error?

```
def fibonacci(n: Int): Long = n match {
  case 0 => 0L
  case 1 => 1L
  case n =>
    fibonacci(n - 1) + fibonacci(n - 2)
}
```

scala> fibonacci(10_000)
java.lang.StackOverflowError
 at fibonacci(<console>:1)
 at fibonacci(<console>:4)



With a function like **Factorial**, which makes a single **recursive** call, we can address the problem simply by using the **accumulator technique** to write a **tail recursive** version of the function.

```
def factorial(n: Int): Long = {
    if (n == 0) 1L
    else n * factorial(n - 1L)
}
```

```
faciter 0 acc = acc
faciter n acc = faciter (n - 1) (n * acc)
fac n = faciter n 1
> map fac [0..10]
[1,1,2,6,24,120,720,5040,40320,362880,3628800]
```

But sometimes it is not possible to do that. E.g. if a function makes more than one **recursive** call to itself, it is not possible to rewrite those two calls as one.

In the particular case of the **Fibonacci** function, which calls itself **recursively** twice, we can get round the problem by complementing the **accumulator technique** with **tupling**:

fibiter 0 (a,b) = a
fibiter n (a,b) = fibiter (n-1) (b,a+b)

fib n = fibiter n (0,1)

> map fib [0..10]
[0,1,1,2,3,5,8,13,21,34,55]

On the next slide we conclude this part with Sergei Winitzki describing the limitations of the accumulator technique.



Sergei Winitzki in sergei-winitzki-11a6431 We cannot implement a non-tail-recursive function without stack overflow (i.e., without unlimited growth of intermediate expressions).

The accumulator trick does not always work! In some cases, it is impossible to implement tail recursion in a given recursive computation. An example of such a computation is the "merge-sort" algorithm where the function body must contain two recursive calls within a single expression. (It is impossible to rewrite two recursive calls as one.)

What if our recursive code cannot be transformed into tail-recursive code via the accumulator trick, but the recursion depth is so large that stack overflows occur? There exist special tricks (e.g., "continuations"" and "trampolines") that convert non-tail-recursive code into iterative code without stack overflows. Those techniques are beyond the scope of this chapter.



Sergei Winitzki

