# Cost models and advanced Futhark programming

Troels Henriksen (athas@sigkill.dk)
Some material by Martin Elsman

DIKU University of Copenhagen

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#### **Agenda**

Parallel cost models

Prefix sums (scans)

Using scans

**Auxiliary** 

#### Parallel cost models

Prefix sums (scans)

Using scans

Auxiliary

#### The need for cost models

```
Which is better?
import numpy as np

def inc_scalar(x):
   for i in range(len(x)):
      x[i] = x[i] + 1

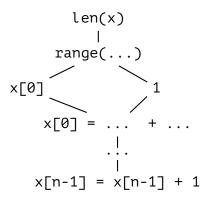
def inc_par(x):
   return x + np.ones(x.shape)
```

#### The need for cost models

```
Which is better?
import numpy as np
def inc scalar(x):
  for i in range(len(x)):
    x[i] = x[i] + 1
def inc par(x):
  return x + np.ones(x.shape)
Intuitively, inc_par is better because it is "more parallel".
```

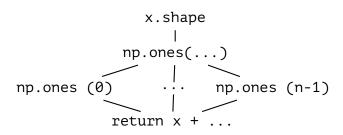
Parallel cost models make this notion precise.

# Dependency DAG for inc\_scalar



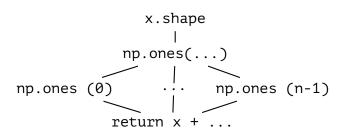
- Total count of nodes is the *work*, W(p).
- Length of longest path from a leaf to the root is the span.
- With an infinite number of processors, if a program p has span k, written S(p) = k, the program can execute in O(k) time.
- Here, W(p) = O(n), S(p) = O(n).

## Dependency DAG for inc\_par



What is the work and span complexity?

## Dependency DAG for inc\_par



#### What is the work and span complexity?

- W(p) = O(n)
- S(p) = O(1)

#### Parallel cost model based on work and span

Instead of giving just a simple cost-model based on the total notion of work carried out by a program, we give instead a *refined* cost model, which aims at providing both:

- a notion of how much total work (W) the program does;
- a notion of the *span*<sup>1</sup> (*S*) of the program, specifying the maximum depth required by the computation.

#### Notice:

- The span is the length of the longest sequence of operations that must be performed sequentially due to data dependencies.
- With an infinite number of processors, if a program p has span k, written S(p) = k, the program can execute in O(k) time.

<sup>&</sup>lt;sup>1</sup>Sometimes also called *depth*.

Writing  $T_i$  for the time taken to execute an algorithm on i processors, Brent's Theorem states that

$$\frac{T_1}{p} \le T_p \le T_\infty + \frac{T_1}{p}$$

**Proof sketch:** At level j of the DAG there are  $M_j$  independent operations, which can clearly be executed by p processors in time

$$\left\lceil \frac{M_j}{p} \right\rceil$$

Sum these for each level of the DAG.

Ramification: We can simulate an "infinitely parallel" machine on a real machine at an overhead proportional to the amount of "missing" hardware parallelism.

#### Language-based cost models

- Tallying up levels in an infinite DAG is impractical for real programs. Instead we prefer a language-based cost model
- E.g. W(x + y) is defined as W(x) + W(y).
- The following slides define work and span cost for a small subset of Futhark.
- Write [e] for the result of evaluating expression e (we are being intuitive about scopes and such).

#### Language-based cost models

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#### Cost model must be implementable

A provable time and space efficient implementation of NESL—Guy Blelloch and John Greiner, 1996

$$W(v) = S(v) = S(e_1 \oplus e_2) = S(e_1 \oplus e_2) = W(\x -> e) = S(\x -> e) =$$

$$W(v) = 1$$
 $S(v) = 1$ 
 $W(e_1 \oplus e_2) =$ 
 $S(e_1 \oplus e_2) =$ 
 $W(\setminus x -> e) =$ 
 $S(\setminus x -> e) =$ 

$$W(v) = 1$$
 $S(v) = 1$ 
 $W(e_1 \oplus e_2) = W(e_1) + W(e_2) + 1$ 
 $S(e_1 \oplus e_2) = S(e_1) + S(e_2) + 1$ 
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$$W([e_1, \dots, e_n]) = S([e_1, \dots, e_n]) = W((e_1, \dots, e_n)) = S((e_1, \dots, e_n)) = S((e_1, \dots, e_n)) = S(e_1, \dots, e_n)$$

#### Implications?

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 $S(e_1 \oplus e_2) = S(e_1) + S(e_2) + 1$ 
 $W(\x -> e) = 1$ 
 $S(\x -> e) = 1$ 

$$W([e_1, \dots, e_n]) = W(e_1) + \dots + W(e_n) + 1$$
  
 $S([e_1, \dots, e_n]) = S(e_1) + \dots + S(e_n) + 1$   
 $W((e_1, \dots, e_n)) = W(e_1) + \dots + W(e_n) + 1$   
 $S((e_1, \dots, e_n)) = S(e_1) + \dots + S(e_n) + 1$ 

#### **Implications?**

$$W(\text{iota } e) = S(\text{iota } e) =$$

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$$W(\text{let } x = e \text{ in } e') = W(e) + W(e'[x \mapsto [\![e]\!]]) + 1$$
  
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 $S(\text{let } x = e \text{ in } e') = S(e) + S(e'[x \mapsto [e]]) + 1$ 

$$W(e_1 \ e_2) = W(e_1) + W(e'[x \mapsto [e_2]]) + 1$$
  
where  $[e_1] = \langle x - \rangle e'$   
 $S(e_1 \ e_2) = S(e_1) + S(e'[x \mapsto [e_2]]) + 1$   
where  $[e_1] = \langle x - \rangle e'$ 

# Work and span for map

$$W(\text{map } e_1 e_2) =$$

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# Work and span for map

```
W(\text{map } e_1 e_2) =
   W(e_1) + W(e_2) + W(e'[x \mapsto v_1]) + \ldots + W(e'[x \mapsto v_n])
   where \llbracket e_1 \rrbracket = \backslash x \rightarrow e'
   where [e_2] = [v_1, \dots, v_n]
S(\text{map } e_1 e_2) =
   S(e_1) + S(e_2) + max(S(e'[x \mapsto v_1]), \dots, S(e'[x \mapsto v_n])) + 1
   where [\![e_1]\!] = \x -> e'
   where \llbracket e_2 \rrbracket = [v_1, \cdots, v_n]
```

## **Reduction by contraction**

```
let npow2 (n:i32):i32 =
 loop a = 2 while a < n do 2*a
-- Pad a vector to make its size a power of two
let padpow2 [n] (ne: i32) (v:[n]i32) : []i32 =
 concat v (replicate (npow2 n - n) ne)
-- Reduce by contraction
let red (xs : []i32) : i32 =
 let xs =
   loop xs = padpow2 0 xs
   while length xs > 1 do
      let n = length xs / 2
      in map2 (+) xs[0:n] xs[n:2*n]
 in xs[0]
```

# Work and span of loop

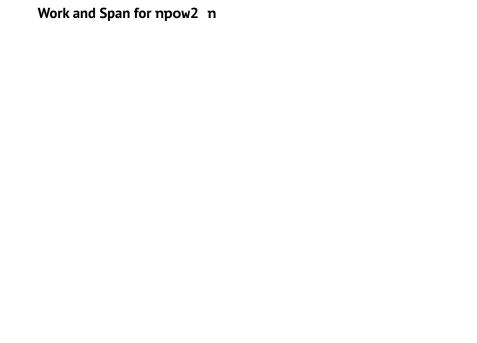
$$W(\operatorname{loop} x = e_1 \operatorname{while} e_2 \operatorname{do} e_3) =$$

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### Work and span of loop

```
W(\mbox{loop}\ x = e_1 \mbox{ while}\ e_2 \mbox{ do}\ e_3) = \mbox{ if } [\![e_2]\!] = \mbox{true} \mbox{ then } 1 + W(e_2) \mbox{ else } 1 + W(e_2) + W(\mbox{loop}\ x = [\![e_3[x \mapsto e_1]]\!] \mbox{ while}\ e_2 \mbox{ do}\ e_3)
```

```
S(\mathbf{loop}\ x = e_1\ \mathbf{while}\ e_2\ \mathbf{do}\ e_3) = 
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# Work and Span for npow2 n

By inspection, we have

$$W(\text{npow2 n}) = S(\text{npow2 n}) = O(\log n)$$

Work and Span for padpow2 ne v

#### Work and Span for npow2 n

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$$W(npow2 n) = S(npow2 n) = O(log n)$$

#### Work and Span for padpow2 ne v

Because npow2  $n \le 2n$ , we have (where n = length v)

```
W(\text{padpow2 ne v})
= W(\text{concat } v \text{ (replicate (npow2 n - n) ne)})
= O(n)
S(\text{padpow2 ne v}) = O(\log n)
```

Work and Span for red

#### Work and Span for npow2 n

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#### Work and Span for padpow2 ne v

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W(\text{padpow2 ne } v)
= W(\text{concat } v \text{ (replicate (npow2 n - n) ne)})
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```

#### Work and Span for red

Each loop iteration in red has span O(1). Because the loop is iterated at-most log(2n) times, we have (where n = length v)

$$W(\text{red } v) = O(n) + O(n/2) + O(n/4) + \cdots + O(1) = O(n)$$
  
 $S(\text{red } v) = O(\log n)$ 

#### Work efficiency

A parallel algorithm is said to be *work efficient* if it has at most the same work as the best sequential algorithm.

Is red work efficient?

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A parallel algorithm is said to be *work efficient* if it has at most the same work as the best sequential algorithm.

Is red work efficient?

**Yes**, because it does O(n) work, which is as good as a sequential summation.

Is it also efficient?

#### Performance Compared to the Built-in Reduction SOAC

```
-- ==
-- entry: test_red test_reduce
-- random input { [10000000]i32 }
entry test_red = red
entry test_reduce = reduce (+) 0
```

### Performance Compared to the Built-in Reduction SOAC

```
-- entry: test red test reduce
-- random input { [10000000]i32 }
entry test red = red
entry test reduce = reduce (+) 0
$ futhark bench --backend=opencl reduce.fut
Compiling reduce.fut...
Results for reduce.fut:test red:
dataset [10000000]i32: 4675.40\mus
Results for reduce.fut:test reduce:
dataset [10000000]i32: 273.80\mu s
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```

If you are not using futhark bench, then you are probably doing it wrong.

Parallel cost models

Prefix sums (scans)

Using scans

Auxiliary

## Inclusive and exclusive prefix sum

### Exclusive prefix sum ("prescan")

Given

[1, 2, 3, 4]

produce

[0, 1, 3, 6]

#### Inclusive prefix sum

Given

[1, 2, 3, 4]

produce

[1, 3, 6, 10]

#### **Prefix sums are scans**

Generalising the addition and zero used by a prefix sum to an arbitrary associative operator  $\oplus$  and neutral element  $0_{\oplus}$ , we get *scan*.

```
-- The scan in Futhark is inclusive.
> scan (+) 0 [1,2,3,4]
[1, 3, 6, 10]
```

#### **Prefix sums are scans**

Generalising the addition and zero used by a prefix sum to an arbitrary associative operator  $\oplus$  and neutral element  $0_{\oplus}$ , we get scan.

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-- The scan in Futhark is inclusive.
> scan (+) 0 [1,2,3,4]
[1, 3, 6, 10]
```

- Scans are a fundamental building block on parallelising seemingly-sequential algorithms.
- Let us see how scans can be computed in parallel.

### Sequential prefix sum

```
acc = 0
for i < n:
  acc = acc + input[i]
  scanned[i] = acc</pre>
```

### Sequential prefix sum

```
acc = 0
for i < n:
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  Work: O(n)
  Span: O(n)</pre>
```

#### **Brute force**

To calculate the prefix sum of  $[x_0, \ldots, x_{n-1}]$ , compute

```
[sum([x_0])

sum([x_0, x_1])

\vdots

sum([x_0, x_1, ..., x_{n-1}])]
```

Assume  $S(sum([x_0,\ldots,x_{n-1}])) = log_2(n)$ .

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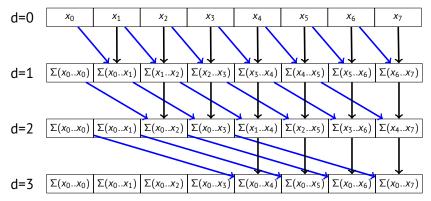
sum([x_0, x_1, ..., x_{n-1}])]
```

Assume 
$$S(sum([x_0, ..., x_{n-1}])) = log_2(n)$$
.

Work: 
$$O(\sum_{i < n} i) = O(n^2)$$
  
Span:  $O(max(S(sum([x_0])), ... S(sum([x_0, ..., x_{n-1}])))) = O(log_2(n))$ 

**Terrible.** The sequential implementation is faster for large n!

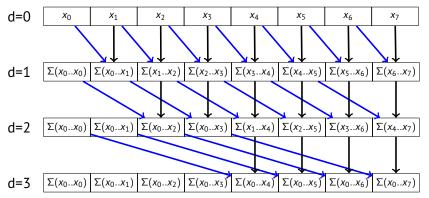
## Hillis—Steele scan (1986)<sup>2</sup>



For each d, element  $x_i^d$  is updated by  $x_{i-2^i}^{d-1} + x_i^{d-1}$ .

<sup>&</sup>lt;sup>2</sup>This slide took 60 minutes to make.

# Hillis-Steele scan (1986)<sup>2</sup>



For each d, element  $x_i^d$  is updated by  $x_{i-2}^{d-1} + x_i^{d-1}$ .

Work: For 
$$n = 2^m$$
,  $O(\sum_{i < m} 2^m - 2^i) = O(n \log(n))$ 

Span: log(n)

<sup>&</sup>lt;sup>2</sup>This slide took 60 minutes to make.

#### **Work-efficient scan**

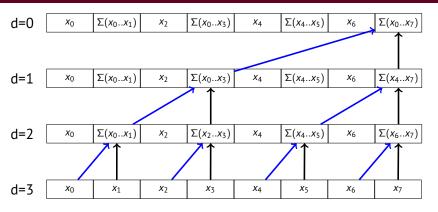
#### Two passes

Up-sweep Build a balanced binary tree of partial sums stored in every other cell.

Down-sweep Use the partial sums to fill out the missing parts.

The binary tree does not actually exist as a recursive pointer structure, but is just a communications concept.

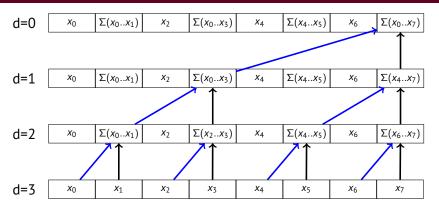
# Up-sweep ("reduction phase")<sup>3</sup>



$$x_i^d = x_{i-2^{m-d-1}}^{d+1} + x_i^{d+1}$$

<sup>&</sup>lt;sup>3</sup>This slide took 30 minutes to make.

# **Up-sweep ("reduction phase")**<sup>3</sup>



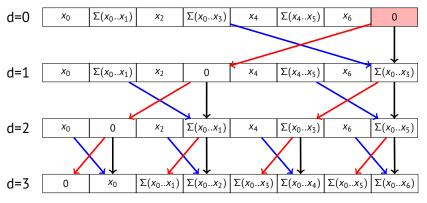
$$x_i^d = x_{i-2^{m-d-1}}^{d+1} + x_i^{d+1}$$

Work: For  $n = 2^m$ ,  $O(\sum_{i=1}^m 2^i) = O(n)$ 

Span: log(n)

<sup>&</sup>lt;sup>3</sup>This slide took 30 minutes to make.

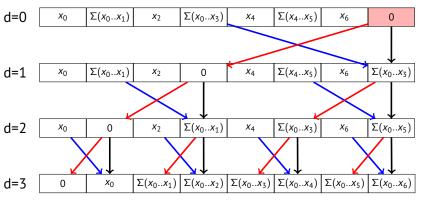
### Down-sweep<sup>4</sup>



Inverse indexing of the up-sweep phase.

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## Down-sweep<sup>4</sup>



Inverse indexing of the up-sweep phase.

Work: For 
$$n = 2^m$$
,  $O(\sum_{i < m} 2^i) = O(n)$ 

Span: log(n)

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#### Work efficient scan

#### Complexity of scan on size-n input

Work: O(n)Span:  $\log(n)$ 

- Optimal, as reduce is the same.
- Can now depend on scan as a relatively cheap building block.

Real-world scan implementations are often very different for technical reasons, but we can depend on these asymptotics when analysing and designing parallel algorithms. Parallel cost models

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Suppose we wish to remove negative elements from the list **let** as =  $\begin{bmatrix} -1, 2, -3, 4, 5, -6 \end{bmatrix}$ 

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For each element, see if we want to keep it: **let** keep = map (\a -> **if** a >= 0 **then** 1 **else** 0) as --  $\begin{bmatrix} 0, 1, 0, 1, 1 \end{bmatrix}$ 

```
Suppose we wish to remove negative elements from the list let as = [-1, 2, -3, 4, 5, -6]
```

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For each element, see if we want to keep it: **let** keep = map (\a -> **if** a >= 0 **then** 1 **else** 0) as -- [0, 1, 0, 1, 1, 0]

**let** offsets1 = scan (+) 0 keep -- [ 0, 1, 1, 2, 3, 3]

let offsets = map (\x -> x - 1) offsets1
-- [-1, 0, 0, 1, 2, 2]

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-- [ 0, 1, 1, 2, 3, 3]

let offsets = map (\x -> x - 1) offsets1
-- [-1, 0, 0, 1, 2, 2]

offsets[i] now indicates position in filtered list iff keep[1]
== 1.

#### scatter

scatter xs is vs computes equivalent of the imperative
pseudocode
for j < n:
 xs[is[j]] = vs[j]</pre>

- Out-of-bound writes are ignored
- Writing different values to same index is undefined<sup>5</sup>
- Work O(n), span O(1)

#### Just what we need for filtering!

<sup>&</sup>lt;sup>5</sup>reduce\_by\_index handles conflicts with provided operator.

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#### Just what we need for filtering!

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### Implementing filter

```
let filter 'a (p: a -> bool) (as: []a): []a =
  let keep = map (a \rightarrow if p a then 1 else 0) as
  let offsets1 = scan (+) 0 keep
  let num to keep = reduce (+) 0 keep
  in if num to keep == 0
     then []
     else scatter (replicate num to keep as[0])
                  (map2 (\i k ->
                             if k == 1
                             then i-1
                             else -1)
                         offsets1 keep)
                   as
```

#### Radix sort

- Many classical sorting algorithms are a poor fit for data parallelism, but radix sort works well.
- Radix-2 sort works by repeatedly partitioning elements according to one bit at a time, while preserving the ordering of the previous steps.

### **Example with radix-10**

326	69	0	7	04	ļ.	3	24
453	7 5	5   1	6	0 8	3	4	38
608	4 5	3	3	2 6		4	5 6
8 3 5	70	7 0 4	8	3 5	$\Rightarrow$	6	0 5
751	⇒ 83	5	⇒ 4	3 5	<i>→</i>	6	9 5
435	4 3	5	7	5 1	-	7	01
704	3 2	6	4	5 3	5	7	5 3
690	60	8	6	90	)	8	30

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- Radix sort is not as general as a comparison-based sort.
- Assumes sorting key can be decomposed into "digits".

## Sorting xs:[n]u32 by bit b

```
-- 1 if bit b set.

let check_bit b x =

(i32.u32 (x >> u32.i32 b)) & 1
```

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let bits_neg = map (1-) bits
let offs = reduce (+) 0 bits_neg
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```

```
xs = [0, 1, 2, 3, 4]
bits = [0, 1, 0, 1, 0]
bits_neg = [1, 0, 1, 0, 1]
offs = 3
```

**let** idxs0 = map2 (\*)

### Example

Then scatter as when filtering.

### The whole step

```
let check bit b x =
  (i32.u32 (x >> u32.i32 b)) & 1
let radix sort step [n]
    (xs: \lceil n \rceil u 32) (b: i 32): \lceil n \rceil u 32 =
  let bits = map (check bit b) xs
  let bits neg = map (1-) bits
  let offs = reduce (+) 0 bits neg
  let idxs0 = map2 (*) bits neq
                    (scan (+) 0 bits neg)
  let idxs1 = map2 (*) bits
                    (map (+offs) (scan (+) 0 bits))
  let idxs2 = map2 (+) idxs0 idxs1
  let idxs = map (x->x-1) idxs2
  let xs' = scatter (copy xs) idxs xs
  in xs'
```

#### **Radix sort in Futhark**

```
let radix_sort [n] (xs: [n]u32): [n]u32 =
  loop xs for i < 32 do radix_sort_step xs i

See worked example at https:
//futhark-lang.org/examples/radix-sort.html</pre>
```

## **Segmented scan**

```
val segmented_scan [n] 't
    : (op: t -> t -> t) -> (ne: t)
    -> (flags: [n]bool) -> (as: [n]t)
    -> [n]t
```

true starts a segment and false continues a segment.

#### Example

```
segmented_scan (+) 0
  [true, false, true, false, false, true]
  [0, 1, 2, 3, 4, 5]
== scan (+) 0 [0,1] ++
    scan (+) 0 [2,3,4] ++
    scan (+) 0 [5]
== [0, 1, 2, 5, 9, 5]
```

### **Segmented reduction**

```
val segmented_reduce [n] 't
    : (op: t -> t -> t) -> (ne: t)
    -> (flags: [n]bool) -> (as: [n]t)
    -> []t
```

#### Example

```
segmented_reduce (+) 0
  [true, false, true, false, false, true]
  [0, 1, 2, 3, 4, 5]
== reduce (+) 0 [0,1] ++
  reduce (+) 0 [2,3,4] ++
  reduce (+) 0 [5]
== [1, 9, 5]
```

## **Generalised histograms**

Like scatter, but uses a provided reduce-like operator to handle multiple writes to same index.

#### Type

```
val reduce_by_index [k] [n] 'a :
        (dest: *[k]a)
   -> (f: a -> a -> a) -> (ne: a)
   -> (is: [n]i32) -> (vs: [n]a) -> *[k]a
```

#### **Semantics**

### Proving associativity and neutral elements

- Is op associative?
- Is (i32.smallest, -1) a neutral element?

### argmax: associativity

```
(a 'op' b) 'op' c
== ((ax, ai) 'op' (bx, bi)) 'op' (cx, ci)
== let (x, i) = if ax < bx then (bx, bi)
                           else (ax, ai)
   in if x < cx then (cx, ci)
                else (x, i)
   a 'op' (b 'op' c)
== (ax, ai) 'op' ((bx, bi) 'op' (cx, ci))
== let (x, i) = if bx < cx then (cx, ci)
                           else (bx, bi)
   in if ax < x then (x, i)
                else (ax, ai)
```

Enumerate all possible comparisons between ax, bx, and cx and show associativity for each.

# E.g. for !(ax < bx) && bx < cx && cx < ax

```
let (x, i) = if ax < bx then <math>(bx, bi)
                            else (ax. bx)
   in if x < cx then (cx, ci)
                else (x, i)
== if ax < cx then (cx, ci)
              else (ax, ai)
== (ax, ai)
   let (x, i) = if bx < cx then (cx, ci)
                            else (bx, bi)
   in if ax < x then (x, i)
                else (ax, ai)
== if ax < cx then (cx, ci)
              else (ax, ai)
== (ax, ai)
```

### argmax: neutral element

```
(a 'op' (i32.smallest, -1))
== ((x, i) 'op' (i32.smallest, -1))
== if x < i32. smallest then (i32.smallest, -1)
                       else (x, i)
== (x, i)
   ((i32.smallest, -1) 'op' a)
== ((i32.smallest, -1) 'op' (x, i))
== if i32.smallest < x then (x, i)
                       else (i32.smallest, -1)
== (x, i)
```

### **Commutativity?**

**Exercise for home:** The argmax operator is not commutative. Try to come up with a counterexample, and see if you can change its definition such that it becomes commutative.

### **Commutativity?**

**Exercise for home:** The argmax operator is not commutative. Try to come up with a counterexample, and see if you can change its definition such that it becomes commutative.

#### Commutative reductions

Futhark has a reduce\_comm function that can be used for commutative operators. This runs faster than normal reduce. Not necessary for built-in operators.

### **Summary**

- Work measures the total number of operations, span measures the longest chain of dependencies.
- Language-based cost models let us reason about program performance in a hardware-agnostic and composable way.
- Scans are a useful building block in advanced data parallel algorithms, but an efficient implementation is not straightforward.