## Cost models and advanced Futhark programming

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#### 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

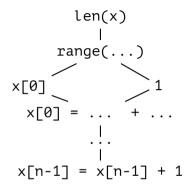
def inc_par(x):
    return x + np.ones(x.shape)
```

### 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".
```

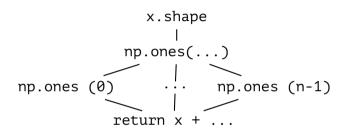
Parallel cost models make this notion precise.

# Dependency DAG for inc\_scalar



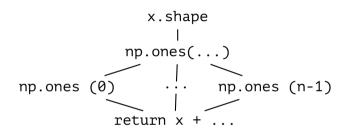
- Total count of nodes is the work, W(p).
- Length of longest path from a leaf to the root is the *span*.
- With an infinite number of processors, if a program p has span k, written S(p) = k, the program can execute in O(k) time.
- Here, W(p) = O(n), S(p) = O(n).

## 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).

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- Write [e] for the result of evaluating expression e (we are being intuitive about scopes and such).

### Cost model must be implementable

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

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

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

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

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

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

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

$$W(\mathtt{iota}\ e) = W(e) + \llbracket e 
rbracket$$
  
 $S(\mathtt{iota}\ e) = S(e) + 1$ 

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$$W($$
let  $x = e$ in  $e') =$  $S($ let  $x = e$ in  $e') =$ 

$$W(\mathtt{iota}\ e) = W(e) + \llbracket e 
rbracket$$
  
 $S(\mathtt{iota}\ e) = S(e) + 1$ 

$$W(\text{let } x = e \text{ in } e') = W(e) + W(e'[x \mapsto [e]]) + 1$$
  
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$$W(e_1 \ e_2) = W(e_1) + W(e_2) + W(e'[x \mapsto [e_2]]) + 1$$
  
where  $[e_1] = \langle x - \rangle e'$   
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# Work and span of map

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

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

# Work and span of 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 \llbracket e_1 \rrbracket = \backslash x \rightarrow e'
   where \llbracket e_2 \rrbracket = [v_1, \cdots, v_n]
```

# **Reduction by contraction**

```
def npow2 (n:i64):i64 =
  loop a = 2 while a < n do 2*a
-- Pad a vector to make its size a power of two
def padpow2 \lceil n \rceil (ne: i32) (v:\lceil n \rceili32) : \lceil \rceili32 =
  concat v (replicate (npow2 n - n) ne)
-- Reduce by contraction
def red (xs : [7i32) : 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]
```

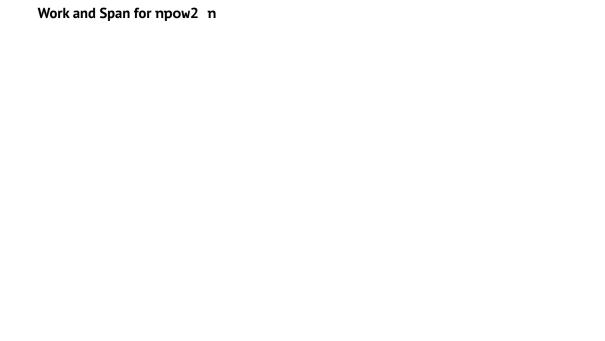
# Work and span of loop

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

$$S({f loop}\ x=e_1\ {f while}\ e_2\ {f do}\ e_3)=$$

## Work and span of loop

```
W(\text{loop } x = e_1 \text{ while } e_2 \text{ do } e_3) = W(e_1) + W(e_2[x \mapsto [e_1]]) +
   if \llbracket e_2 \llbracket x \mapsto \llbracket e_1 \rrbracket \rrbracket \rrbracket = \mathbf{false}
   then 0
   else W(e_3[x \mapsto [e_1]]) +
            W(\text{loop } x = \llbracket e_3[x \mapsto \llbracket e_1 \rrbracket] \rrbracket \text{ while } e_2 \text{ do } e_3)
S(\text{loop } x = e_1 \text{ while } e_2 \text{ do } e_3) = S(e_1) + S(e_2[x \mapsto [e_1]]) +
   if [e_2][x \mapsto [e_1]] = false
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```



By inspection, we have

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

Work and Span for padpow2 ne v

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

#### Work and Span for padpow2 ne v

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

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

$$S(padpow2 ne v) = O(log n)$$

Work and Span for red

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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(2 n) times, we have (where n = l ength v)

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

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## **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 \Gamma100000007i32: 4675.40\mus
Results for reduce.fut:test reduce:
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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]
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#### **Prefix sums are scans**

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

- Scans are a fundamental tool for 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:
    acc = acc + input[i]
    scanned[i] = acc
    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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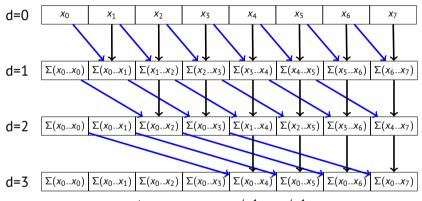
```
Assume S(sum([x_0,\ldots,x_{n-1}])) = log_2(n).

Work: O(\sum_{i < n} i) = O(n^2)

Span: O(max(S(sum([x_0])),\ldots S(sum([x_0,\ldots,x_{n-1}])))) = O(log_2(n))

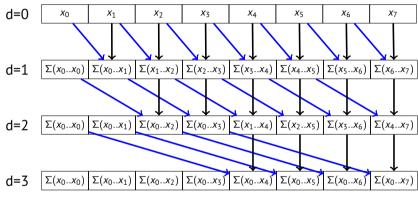
Terrible. The sequential implementation is faster for large n!
```

# Hillis-Steele scan (1986)



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

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

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

Span: log(n)

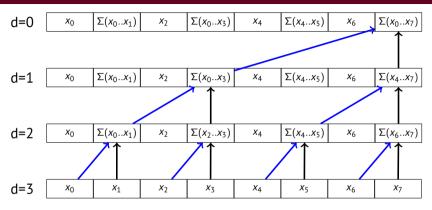
#### **Work-efficient scan**

#### Two passes

Upsweep Build a balanced binary tree of partial sums stored in every other cell. Downsweep Use the partial sums to fill out the missing parts.

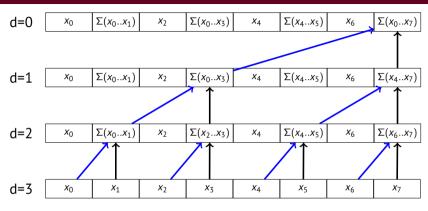
The binary tree does not actually exist as a recursive pointer structure, but is just a communications concept.

# **Upsweep ("reduction phase")**



$$X_i^d = X_{i-2^{m-d-1}}^{d+1} + X_i^{d+1}$$

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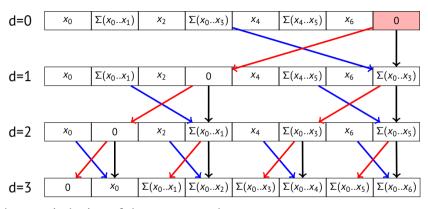


$$x_i^d = x_{i-2^{m-d-1}}^{d+1} + x_i^{d+1}$$

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

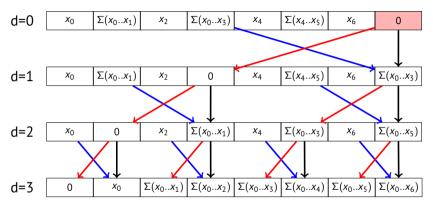
Span: log(n)

# **Downsweep**



Inverse indexing of the upsweep phase.

# Downsweep



Inverse indexing of the upsweep phase.

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

Span: log(n)

#### **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.

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Suppose we wish to remove negative elements from the list

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**let** as = 
$$[-1, 2, -3, 4, 5, -6]$$

For each element, see if we want to keep it:

let keep = map (\a -> if a >= 0 then 1 else 0) as -- 
$$[0, 1, 0, 1, 1, 0]$$

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Suppose we wish to remove negative elements from the list

**let** as = 
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For each element, see if we want to keep it:

**let** offsets1 = scan (+) 0 keep 
$$-- \int 0$$
, 1, 1, 2, 3, 37

Suppose we wish to remove negative elements from the list

**let** as = 
$$[-1, 2, -3, 4, 5, -6]$$

For each element, see if we want to keep it:

offsets[i] now indicates position in filtered list iff

$$keep[1] == 1$$

#### scatter

 $\verb|scatter xs is vs computes equivalent of the imperative pseudocode|\\$ 

```
for j < n:
    xs[is[j]] = vs[j]</pre>
```

- Out-of-bound writes are ignored
- Writing different values to same index is undefined<sup>2</sup>
- Work O(n), span O(1)

Just what we need for filtering!

<sup>&</sup>lt;sup>2</sup>reduce\_by\_index handles conflicts with provided operator.

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#### Just what we need for filtering!

<sup>&</sup>lt;sup>2</sup>reduce\_by\_index handles conflicts with provided operator.

# Implementing filter

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

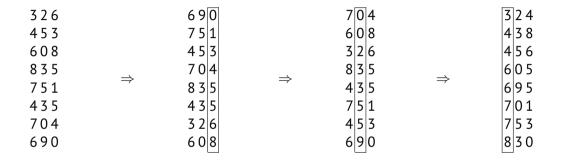
#### Radix sort

- 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	6 9	0	7 C	0 4	[3	3 2 4
453	7 5	1	6 0	8	2	4 3 8
608	4	3	3 2	6	$\rightarrow$	4 5 6
8 3 5	70	4	8 3	5		6 0 5
751	⇒ 83	5  ⇒	4 3	5 5		5 9 5
435	4 3 5 3 2 6	5	7 5	5 1	7	701
704		6	4 5	3	7	7 5 3
690	6 0	8	6 9	90	8	3 0

# Example with radix-10



- 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

```
let bits = map (u32.get_bit b) xs
let bits_neg = map (1-) bits
let offs = reduce (+) 0 bits_neg
```

# Sorting xs:[n]u32 by bit b

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let bits = map (u32.get_bit b) xs
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let offs = reduce (+) 0 bits_neg
```

#### Example

```
b = 0

xs = [0, 1, 2, 3, 4]

bits = [0, 1, 0, 1, 0]

bits_neg = [1, 0, 1, 0, 1]

offs = 3
```

#### Example

```
bits = [0, 1, 0, 1, 0]

bits_neg = [1, 0, 1, 0, 1]

offs = 3

idxs0 = [1, 0, 2, 0, 3]

idxs1 = [0, 4, 0, 5, 0]

map2 (+) idxs0 idxs1 = [1, 4, 2, 5, 3]
```

#### Then scatter as when filtering.

# The whole step

```
def radix sort step [n] (xs: [n]u32) (b: i32): [n]u32 =
  let bits = map (u32.get bit b) xs
  let bits neg = map (1-) bits
  let offs = reduce (+) 0 bits nea
  let idxs0 = map2 (*) bits neg
                   (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**

```
def 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]i64) -> (vs: [n]a) -> *[k]a
Semantics
```

```
for index in 0..k-1:
    i = is[index]
    v = vs[index]
    dest[i] = f(dest[i], v)
```

Futhark uses parallel implementation with GPU *atomics*.

# Proving associativity and neutral elements

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

# argmax: associativity

First, inline definitions:

Then enumerate all possible comparisons between ax, bx, and cx and show that these two expressions are equivalent.

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

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

Similarly, by equational reasoning.

```
(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)
```

## A more calculational approach

https://byorgey.wordpress.com/2020/02/23/what-would-dijkstra-do-proving-the-associativity-of-min/

- Worth a read!
- More elegant and concise, but requires more creative thinking to characterise a useful property of the operator.

## **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.