

# Deterministic parallel programming and data parallelism

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University of Copenhagen

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# Agenda

Content and motivation

Course organisation

Array programming with NumPy

Higher-order array programming

## Content and motivation

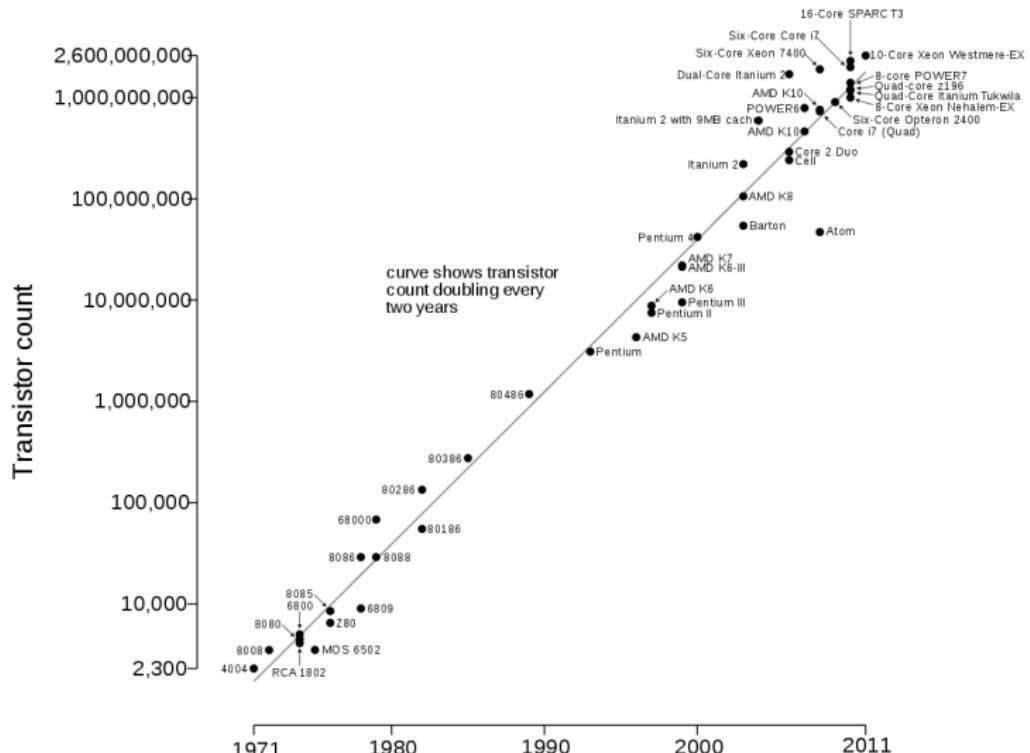
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## The situation so far

Microprocessor transistor counts 1971-2011 & Moore's law



## Moore's Law

*The complexity for minimum component costs has increased at a rate of roughly a factor of two per year. Certainly over the short term this rate can be expected to continue, if not to increase.* –Gordon Moore, 1965

**Note:** not actually a law, and frequently misinterpreted.

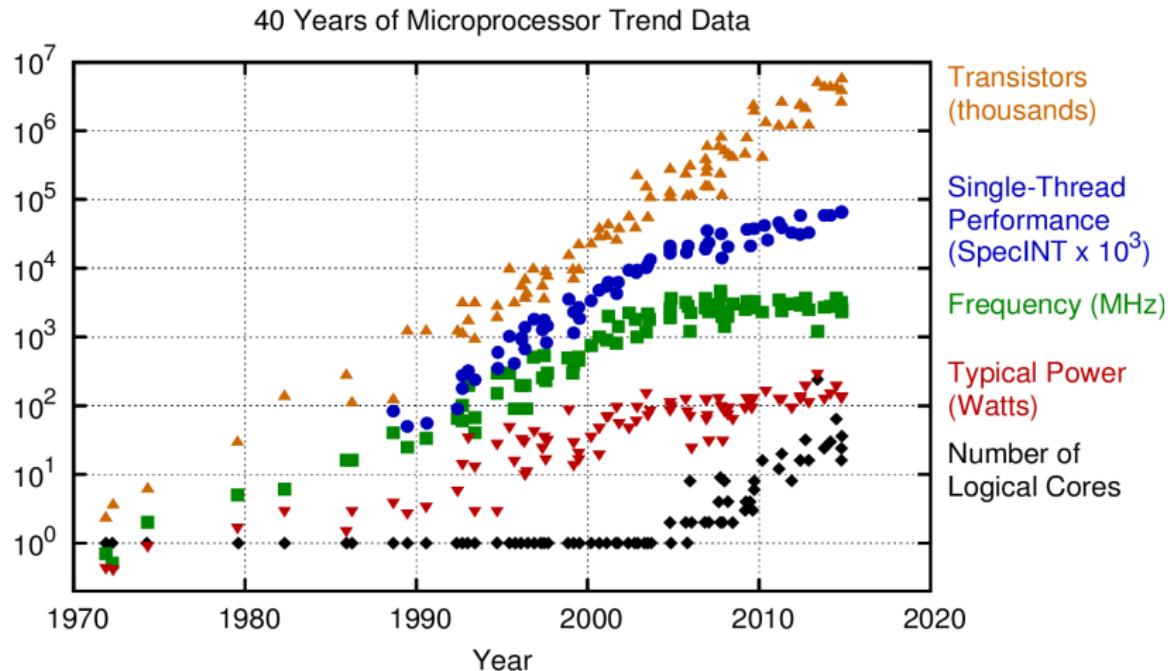
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- Improvements in transistor fabrication technology resulted in smaller transistors (=more transistors for the same chip area).
- This was also translated into higher clock frequencies.

# CPU trends over 40 years



Original data up to the year 2010 collected and plotted by M. Horowitz, F. Labonte, O. Shacham, K. Olukotun, L. Hammond, and C. Batten  
New plot and data collected for 2010-2015 by K. Rupp

# CPU trends over 40 years

- Physical limitations stopped most increases in CPU clock frequency around 2005. Instead, the still-continuing increases in transistor counts was used to make multicore CPUs.
- We had to program with a little more parallelism, but within each thread, we still had a standard CPU model.
- Now even Moore's Law is beginning to end, and we must make better use of the transistors we have—we won't get that many more.
- This has made exotic non-CPU architectures more attractive.

# Two kinds of parallelism

## Task parallelism

Simultaneously performing *different* operations on potentially different pieces of data. Often nondeterministic.

# Two kinds of parallelism

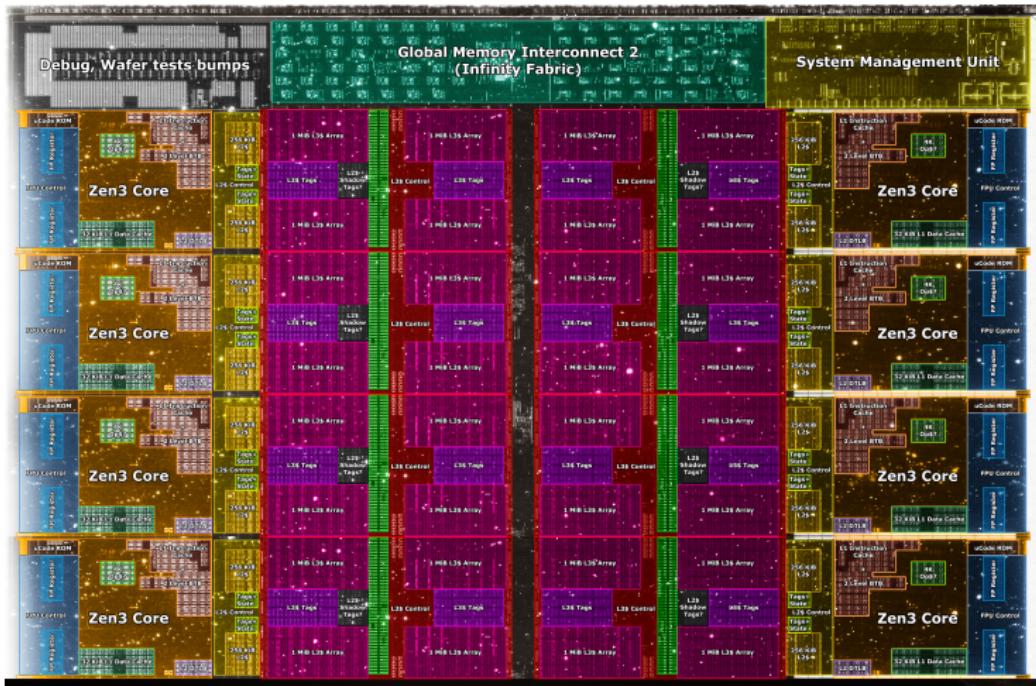
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## Data parallelism

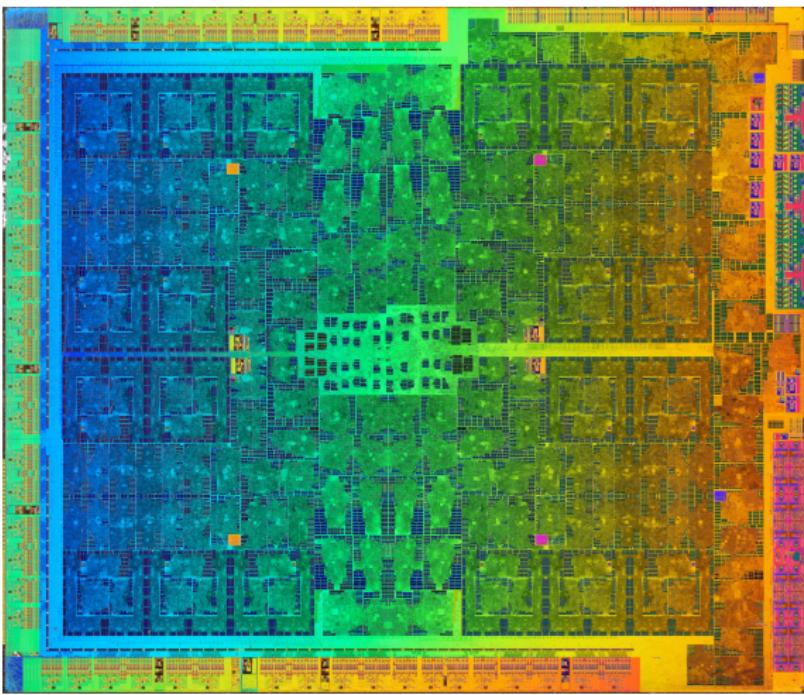
Simultaneously performing *the same* operation on different pieces of the same data. Almost always deterministic.

# AMD Ryzen 3



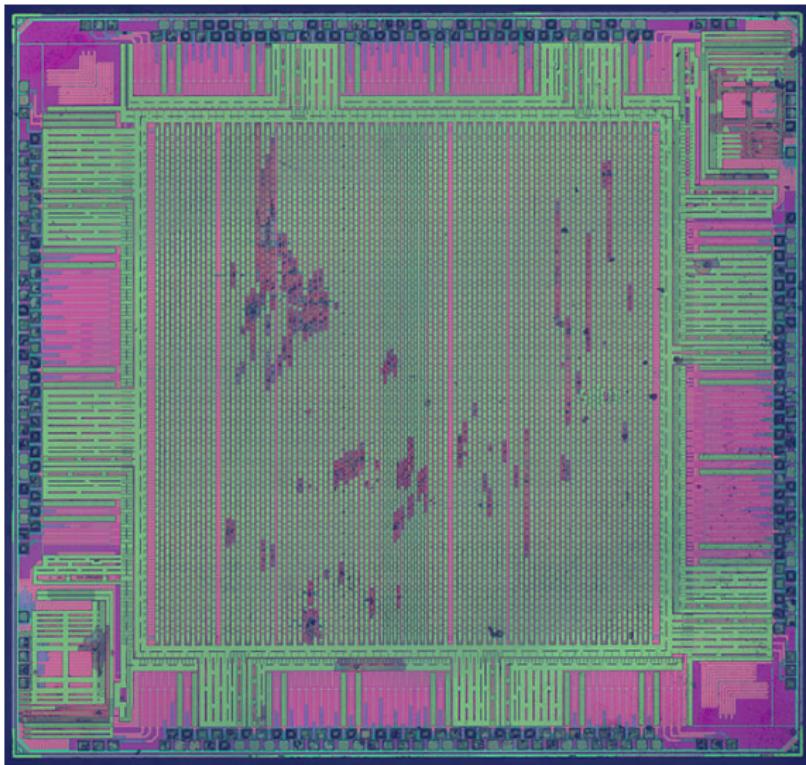
- Irregular layout with many parts.
- Most transistors spent on OOO execution, speculation, caches, branch prediction, etc.
- Few transistors spent on actual computation.
- Task parallel. Fairly easy to program.
- <https://www.reddit.com/r/Amd/comments/jqjg8e>

# NVIDIA Pascal GPU



- Layout consists of replicated components.
- Most transistors spent on computation.
- Data parallel. Quite hard to program.

# Altera Cyclone 1 FPGA



- Extremely regular layout.
- Not cleanly task- or data parallel. Extremely hard to program.

# The simplified tradeoff

**For a fixed transistor budget, a hardware designer can choose to spend transistors on**

1. Making the chip more flexible and easier to program.
2. Improving the peak computational capacity.

Modern CPUs are heavily invested in (1).

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A silver bullet?

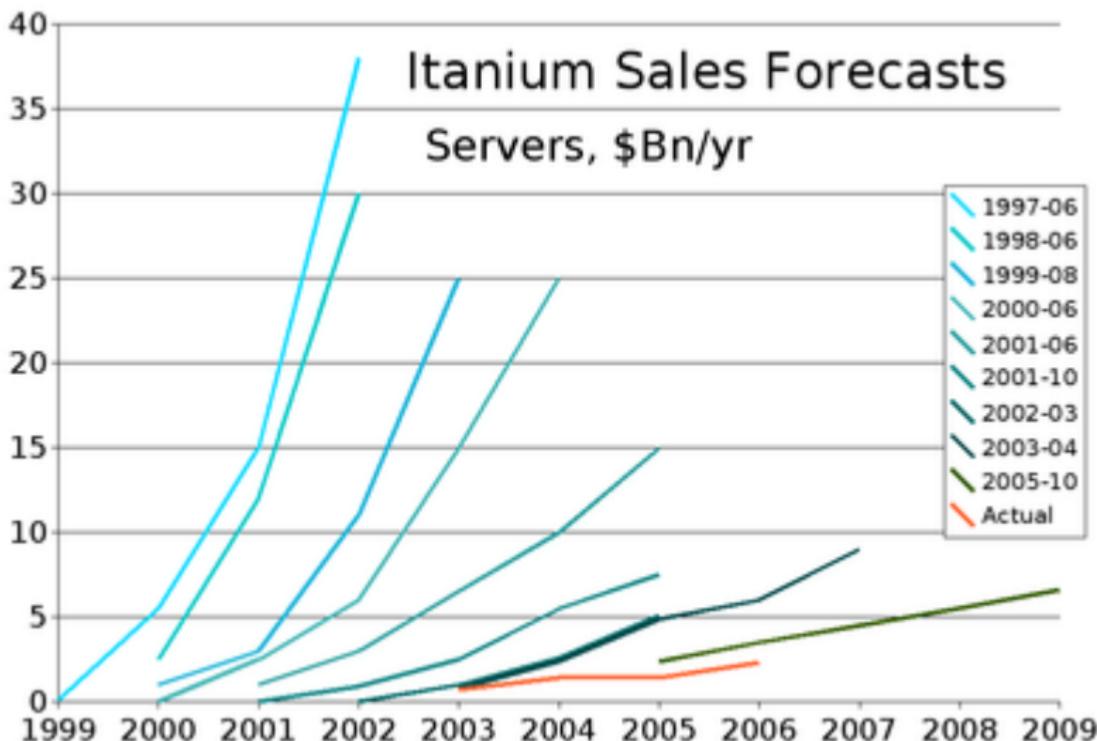
Can we build hardware that is fast and inflexible, and paper over the programming difficulty with a Sufficiently Smart Compiler?

# The Itanium

- Intel's *Itanium* was a VLIW 64-bit replacement for x86.
- Assuming sufficient compilers, performed well.

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1. Bridging the gap from conventional C programs to Itanium's explicitly parallel EPIC architecture was very hard.
2. Increasing transistor counts meant CPUs could just do their own dynamic superscalar scheduling in hardware.
3. Itanium was outcompeted by ever more complex (soon multicore) x86-64 CPUs that handled legacy code just fine.

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## Important lesson

We cannot expect compilers to map legacy models of programming to new hardware.

# The usual compiler disappointments

Probstring's Law

Compiler Advances Double Computing Power Every 18 Years

# The usual compiler disappointments

## Probstring's Law

Compiler Advances Double Computing Power Every 18 Years

- Empirically true in isolation.
- *Probably not true* in a wider sense, where we use compiler technology to make exotic hardware more accessible.

# The purpose of this course

## What we want

Hardware that maximises useful work per transistor and watt.

## What we need

High-level human-friendly programming models with *sequential semantics* that can be efficiently compiled to the efficient hardware.

## What we investigate in Parallel Functional Programming

What should such a programming model look like, and how is it mapped to hardware?

# The two themes in this course

## Properties of programming models

- How should we express parallel programs?
- How can we characterise their performance in a portable way?
- How do we make them composable?

## Implementation of programming models

- Mapping to specific hardware?
- Which features are particularly hard to handle?
- Automatically optimising programs?

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- Mapping to specific hardware?
- Which features are particularly hard to handle?
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**What are the tradeoffs of different parallel programming models?**

Content and motivation

Course organisation

Array programming with NumPy

Higher-order array programming

# When and where?

All information always available at

<https://github.com/diku-dk/pfp-e2020-pub>

## This Wednesday

- 13:00–15:00, **Lecture** in Kursussal 3 at Zoo
- 15:00–17:00, **Lab** in Kursussal 3 at Zoo

## Generally

- Monday 13:00–15:00, **Lecture** in *cyberspace*
- Wednesday 10:00–12:00, **Lecture** in *cyberspace*
- Wednesday 13:00–17:00, **Lab** in Kursussal 3 at Zoo
- **Office hours?**

# Tentative Lecture/Lab schedule

Cosmin and Troels will give lectures, conduct the Lab, and correct assignments.

Continuous-evaluation assessment:

- four individual weekly assignments: 40% of final grade
- one group project + final presentation and discussion: 60% of final grade.
  - ▶ you may chose from multiple possible projects
  - ▶ suggestions presented during Lab in December
  - ▶ or discuss your own project with us
  - ▶ projects can be practical programming in a parallel programming language, or more theoretical

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If  $x$  and  $y$  are vectors, then  $x + y$  is a data parallel operation.

- **You already know data parallel programming!**
  - ▶ NumPy, R, Matlab, Julia, SQL are all data parallel models.
- Completely sequential and deterministic semantics.
- Parallel *execution* straightforward.
- **Distinguish between parallel execution and parallel potential.**
  - ▶ Implementations can always be parallelised as long as the code is written in a parallel style.

# NumPy

- Python library that provides  $n$ -dimensional arrays.
- Not usually parallel in practice, but mostly implemented in efficient C and Fortran.
- *Bulk operations* on entire arrays.

```
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)
```

- Very often essentially purely functional.
- *First order*. All parallel operations do just one thing. No map.

## The good news

- NumPy is extremely widely used. This proves empirically that parallel programming does not have to be hard! Let's try some parallel programming.

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- NumPy is extremely widely used. This proves empirically that parallel programming does not have to be hard! Let's try some parallel programming.

```
def dotprod_seq(x, y):  
    acc = 0  
    for i in range(len(x)):  
        acc += x[i] * y[i]  
    return acc
```

Can we write a potentially parallel dotprod with NumPy?

# Just to get it out the way

```
dotprod = np.dot
```

## Just to get it out the way

```
dotprod = np.dot
```

Yes, in practice do it this way, but let's see how we could built it from *primitives*.

# Parallel dot product

```
def dotprod(x, y):
    products = x * y
    return np.sum(products)
```

x	=	[ 1, 2, -3]
y	=	[ 4, 5, 6]
products	=	[ 4, 10, -18]
np.sum(products)	=	-4.0

What about that np.sum?

# Parallel summation

```
def sum_pow2(x):
    while len(x) > 1:
        x_first = x[0:len(x)/2]
        x_last = x[len(x)/2:]
        x = x_first + x_last
    return x[0]
```

# Parallel summation

```
def sum_pow2(x):
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        x = x_first + x_last
    return x[0]
```

x	=	[1, 2, 3, 4, 5, 6, 7, 8]
x_first	=	[1, 2, 3, 4]
x_last	=	[5, 6, 7, 8]
<hr/>		
x	=	[6, 8, 10, 12]
x_first	=	[ 6, 8]
x_last	=	[10, 12]
<hr/>		
x	=	[16, 20]
x_first	=	[16]
x_last	=	[20]
<hr/>		
x	=	[36]

# Wait, is that correct?

In our summation, we changed

$$((((((x_0 + x_1) + x_2) + x_3) + x_4) + x_5) + x_6) + x_7)$$

to

$$((x_0 + x_4) + (x_2 + x_6)) + ((x_1 + x_5) + (x_3 + x_7))$$

Is that valid?

# Wait, is that correct?

In our summation, we changed

$$((((((x_0 \oplus x_1) \oplus x_2) \oplus x_3) \oplus x_4) \oplus x_5) \oplus x_6) \oplus x_7)$$

to

$$((x_0 \oplus x_4) \oplus (x_2 \oplus x_6)) \oplus ((x_1 \oplus x_5) \oplus (x_3 \oplus x_7))$$

Is that valid?

What about now?

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Is that valid?

What about now?

**In general, what must hold for an operator  $\oplus$  for such a rewrite to be valid?**

# Commutativity and associativity

A binary operator  $\oplus : \alpha \rightarrow \alpha \rightarrow \alpha$  is said to be *commutative* if

$$x \oplus y = y \oplus x$$

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It is *associative* if

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If we are a little more clever, we can do a parallel “sum” with any associative operator.

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If we are a little more clever, we can do a parallel “sum” with any associative operator.

For convenience, we also tend to require a *neutral element*  $0_{\oplus}$ :

$$x \oplus 0_{\oplus} = 0_{\oplus} \oplus x = x$$

# Monoidal summation

An associative binary operator  $\oplus : \alpha \rightarrow \alpha \rightarrow \alpha$  with a neutral element  $0_{\oplus}$  is called a *monoid* and written  $(\oplus, 0_{\oplus})^1$ . Examples:

- $(+, 0)$
- $(*, 1)$
- $(++, [])$

---

<sup>1</sup>A monoid without a neutral element is called a *semigroup*

# Monoidal summation

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- $(+, 0)$
- $(*, 1)$
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For any monoid  $(\oplus, 0_{\oplus})$  we can construct a function `sum` with *semantics*

$$\text{sum}([x_0, \dots, x_{n-1}]) = x_0 \oplus \dots \oplus x_{n-1}$$

but which can *operationally* be executed in parallel.

---

<sup>1</sup>A monoid without a neutral element is called a *semigroup*

# Monoidal reduction

We can abstract the notion of summation into a higher-order function typically called `reduce`:

$$\text{reduce} : (\alpha \rightarrow \alpha \rightarrow \alpha) \rightarrow \alpha \rightarrow []\alpha \rightarrow \alpha$$

Then

$$\text{sum} = \text{reduce } (+) \ 0$$

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Sadly, NumPy is a first-order programming model, and we cannot define an efficient general `reduce`.

**But first order programming models have even bigger problems.**

## The problem with control flow

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$$y[i] = \begin{cases} \sqrt{x[i]} & \text{if } 0 \leq x[i] \\ x[i] & \text{otherwise} \end{cases}$$

```
def sqrt_when_pos_0(x):
    x = x.copy()
    for i in range(len(x)):
        if x[i] >= 0:
            x[i] = np.sqrt(x[i])
    return x
```

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    return x
```

Runtime for  $n = 10^6$ : 14s

## Using flags, filters, and scatters

```
def sqrt_when_pos_1(x):
    x = x.copy()
    x_nonneg = x >= 0
    x[x_nonneg] = np.sqrt(x[x_nonneg])
    return x
```

# Using flags, filters, and scatters

```
def sqrt_when_pos_1(x):
    x = x.copy()
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```

```
x           [ 1,  2, -3]
x_nonneg    [ True, True, False]
x[x_nonneg] [1, 2]
```

# Using flags, filters, and scatters

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def sqrt_when_pos_1(x):
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x           [ 1, 2, -3]
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x[x_nonneg] [1, 2]
```

- Parallel filtering is not so simple (you'll see in a later lecture).
- Runtime for  $n = 10^6$ : 0.18s.

# Arithmetic control flow

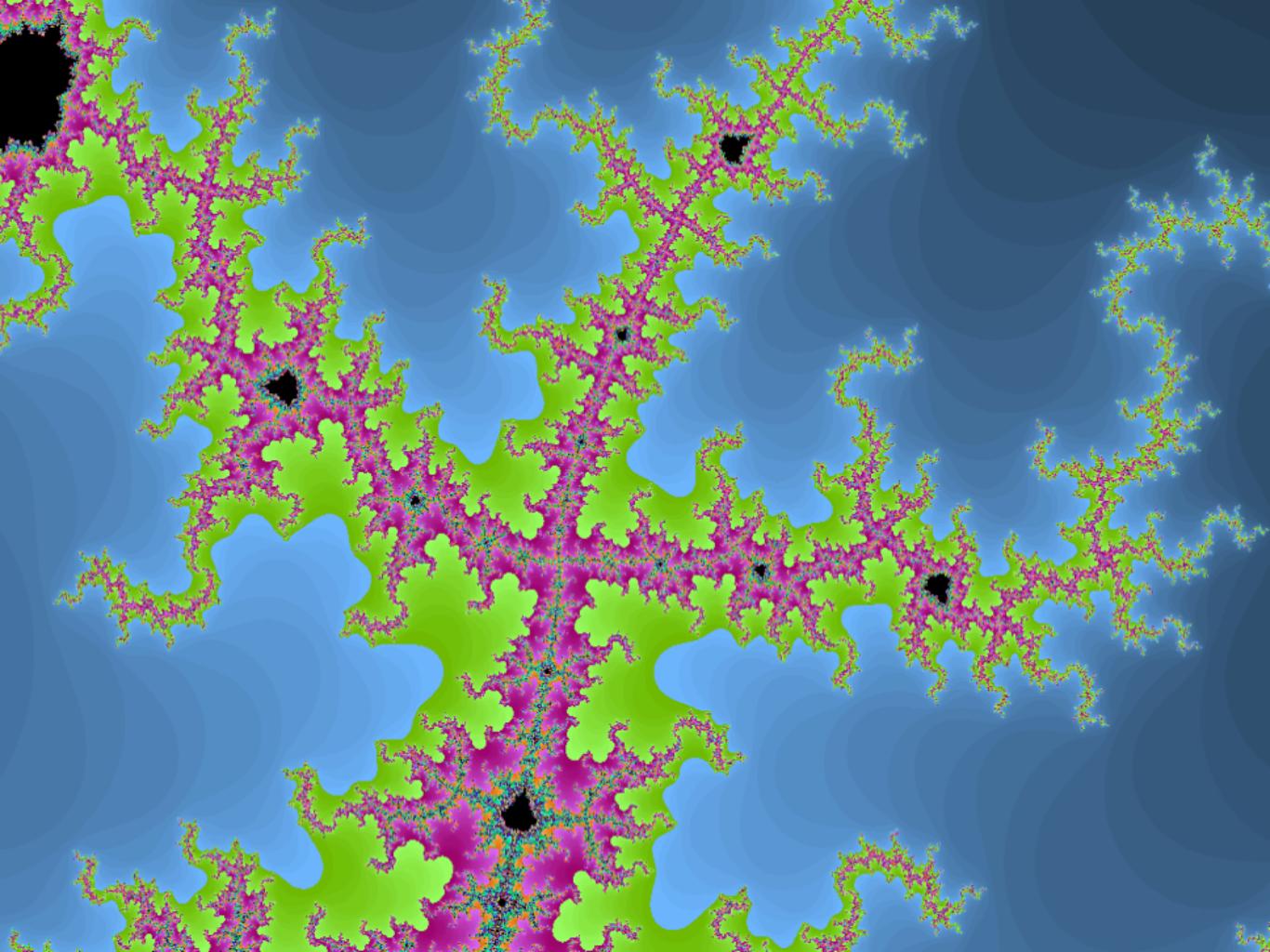
```
def sqrt_when_pos_2(x):
    x_nonneg = x >= 0
    x_neg = x < 0
    x_zero_when_neg = x * x_nonneg.astype(int)
    x_zero_when_nonneg = x * x_neg.astype(int)
    x_sqrt_or_zero = np.sqrt(x_zero_when_neg)
    return x_sqrt_or_zero + x_zero_when_nonneg

x           [ 1,      2,     -3]
x_nonneg    [ True,    True, False]
x_neg       [False,   False, True]
x_zero_when_neg [ 1,      2,     0]
x_zero_when_nonneg [ 0,      0,     -3]
x_sqrt_or_zero [ 1, 1.4142135,  0]
return      [ 1, 1.4142135, -3]
```

Runtime for  $n = 10^6$ : 0.10s.

## It gets worse

That's for a simple conditional. What if we need a *loop*?



# Computing the Mandelbrot Set

The root of those pretty visuals is calling this function (pseudocode) with a bunch of complex numbers:

```
def divergence(c, d):
    i = 0
    z = c
    while i < d and dot(z) < 4.0:
        z = c + z * z
        i = i + 1
    return i
```

# Mandelbrot in NumPy<sup>2</sup>

```
def mandelbrot_numpy(c, d):
    output = np.zeros(c.shape)
    z = np.zeros(c.shape, np.complex32)
    for it in range(d):
        notdone =
            (z.real*z.real + z.imag*z.imag) < 4.0
        output[notdone] = it
        z[notdone] = z[notdone]**2 + c[notdone]
    return output
```

## Problems

- Control flow obscured.
- Always runs for *maxiter* iterations.
- *Lots* of memory traffic.

---

<sup>2</sup>[https://www.ibm.com/developerworks/community/blogs/jfp/entry/How\\_To\\_Compute\\_Mandelbrodt\\_Set\\_Quickly](https://www.ibm.com/developerworks/community/blogs/jfp/entry/How_To_Compute_Mandelbrodt_Set_Quickly)

# The root of the problem

- NumPy is a *first-order parallel language*—all operations are parameterised over simple values, not functions.
- Easy way to design and implement a library, but inflexible.

Now we will look at a *higher-order parallel language*: **Futhark**.

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Higher-order array programming

# Futhark is a high-level language!

- Futhark is *not* a “GPU language”—it is a hardware-agnostic parallel language.
- However, we have written a Futhark *compiler* that can generate good GPU code.
- The compiler is *not* a “parallelising compiler”. The parallelism is *explicitly given* by the programmer. The compiler’s job is to figure out what to do with it. It’s more of a *sequentialising compiler*.
- *Co-design* between compiler and language, inspired by hand-written code for target hardware.

# Futhark at a Glance

## ■ Size-dependent types

An  $n$  by  $m$  integer matrix has type  $[n][m]\text{i32}$ .

- ▶ In the full language, there is an entire type theory to deal with existentially quantified sizes.
- ▶ For this lecture, we assume that the sizes are already known symbolically.

# Futhark at a Glance

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## ■ Nested Parallelism

```
let add_two [n] (a: [n]i32): [n]i32 = map (+2) a
let sum [n] (a: [n]i32): i32 = reduce (+) 0 a
let sumrows [n] (as: [n][m]i32): [n]i32 = map sum as
```

# Futhark at a Glance

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```

## ■ Parallelism must be *regular*; this is not supported:

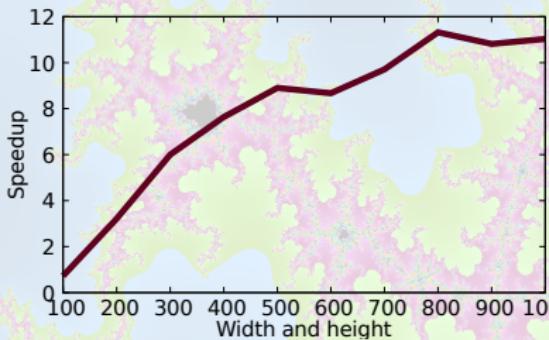
```
let f [n] (a: [n]i32) =
    map (\x -> reduce (+) 0 (iota x)) a
```

# Mandelbrot in Futhark

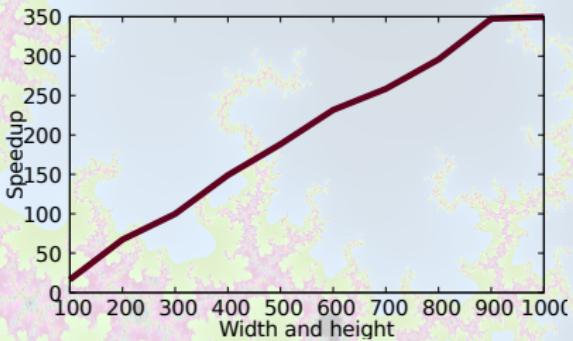
```
let divergence (c: complex) (d: i32): i32 =  
    let (_ , i') =  
        loop (z, i) = (c, 0)  
        while i < d && dot(z) < 4.0 do  
            (addComplex c (multComplex z z),  
             i + 1)  
    in i'  
  
let mandelbrot [n][m] (css: [n][m]complex)  
                  (d: i32) : [n][m]i32 =  
    map (\cs -> map (\c -> divergence c d) cs) css
```

- Only one array written, at the end.
- while loop terminates when the element diverges.

# Mandelbrot speedup on GPU compared to sequential implementation in C



NumPy-style



Futhark-style

**The vectorised style can sacrifice a lot of potential performance.**

# Summary

- Parallel programming is a physical necessity.
- Good language models combined with good compilers can be used to make hardware more efficient.
- Deterministic data parallel programming is demonstrably both accessible to humans and efficient to execute.
- A first-order model is insufficient.
- Sheer parallelism is not enough to make it fast.

**Go install Futhark; you will need it for the assignments:**

<https://futhark-lang.org>

**Then go solve these exercises:**

<https://github.com/diku-dk/pfp-e2020-pub/blob/master/bootstrap-exercises.md>