



# GPU PROGRAMMING 101

## GRIDKA SCHOOL 2019

29 August 2019 | Andreas Herten | Forschungszentrum Jülich *Handout Version*

# About, Outline



## Jülich Supercomputing Centre

- Operation of supercomputers
- Application support
- Research
- Me: *All things GPU*

## Topics

Motivation

Platform

Hardware

Features

Summary

Programming GPUs

Libraries

Directives

Languages

Abstraction Libraries/DSL

Tools

Conclusions

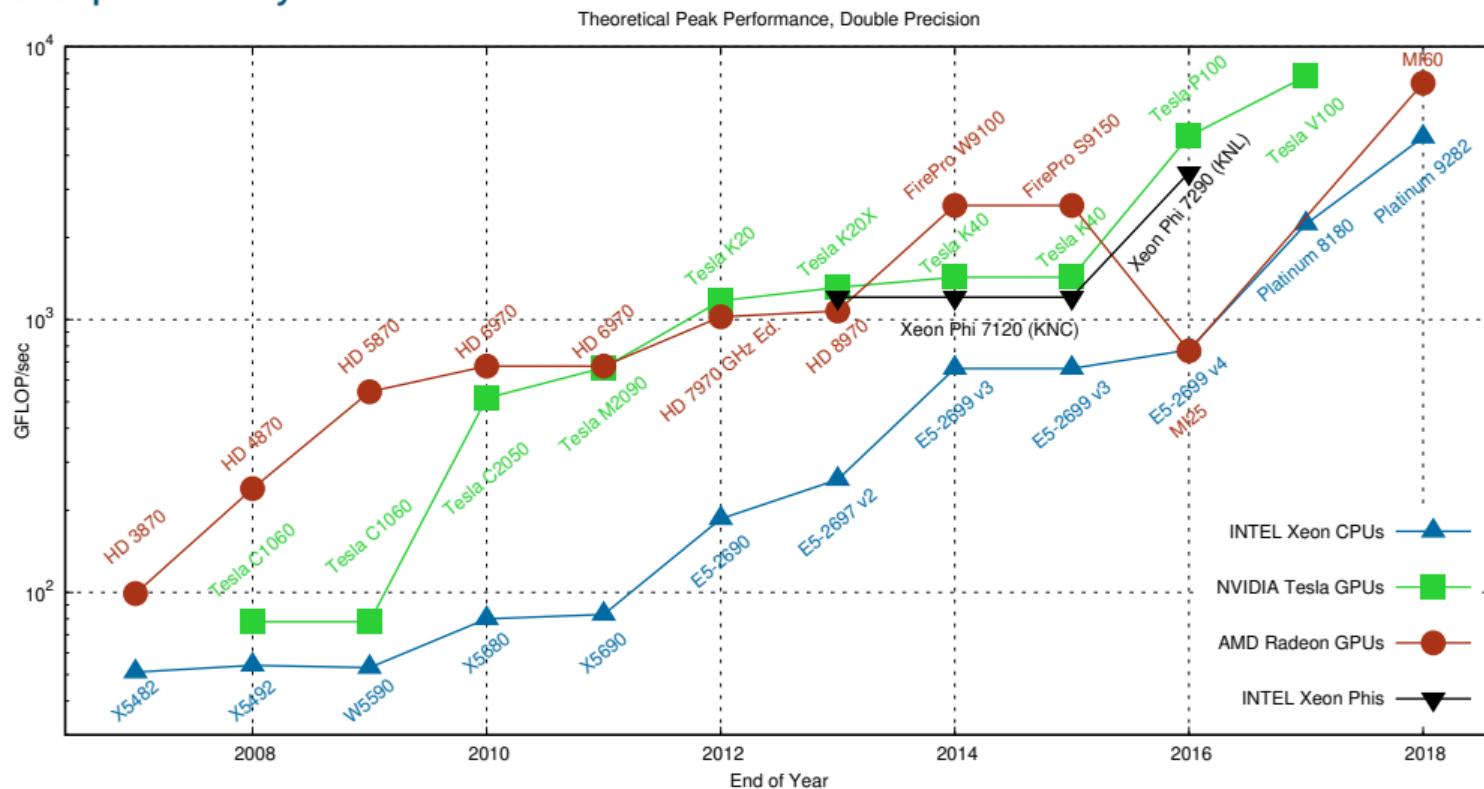
# Status Quo

## A short but parallel story

- 1999 Graphics computation pipeline implemented in dedicated *graphics hardware*  
Computations using OpenGL graphics library [2]  
»GPU« coined by NVIDIA [3]
- 2001 NVIDIA GeForce 3 with *programmable* shaders (instead of fixed pipeline) and floating-point support; 2003: DirectX 9 at ATI
- 2007 CUDA
- 2009 OpenCL
- 2019 Top 500: 25 % with **NVIDIA** GPUs (#1, #2) [4], Green 500: 8 of top 10 with GPUs [5]
- 2021 Aurora: First (?) US exascale supercomputer based on **Intel** GPUs  
Frontier: First (?) US *more-than-exascale* supercomputer based on **AMD** GPUs

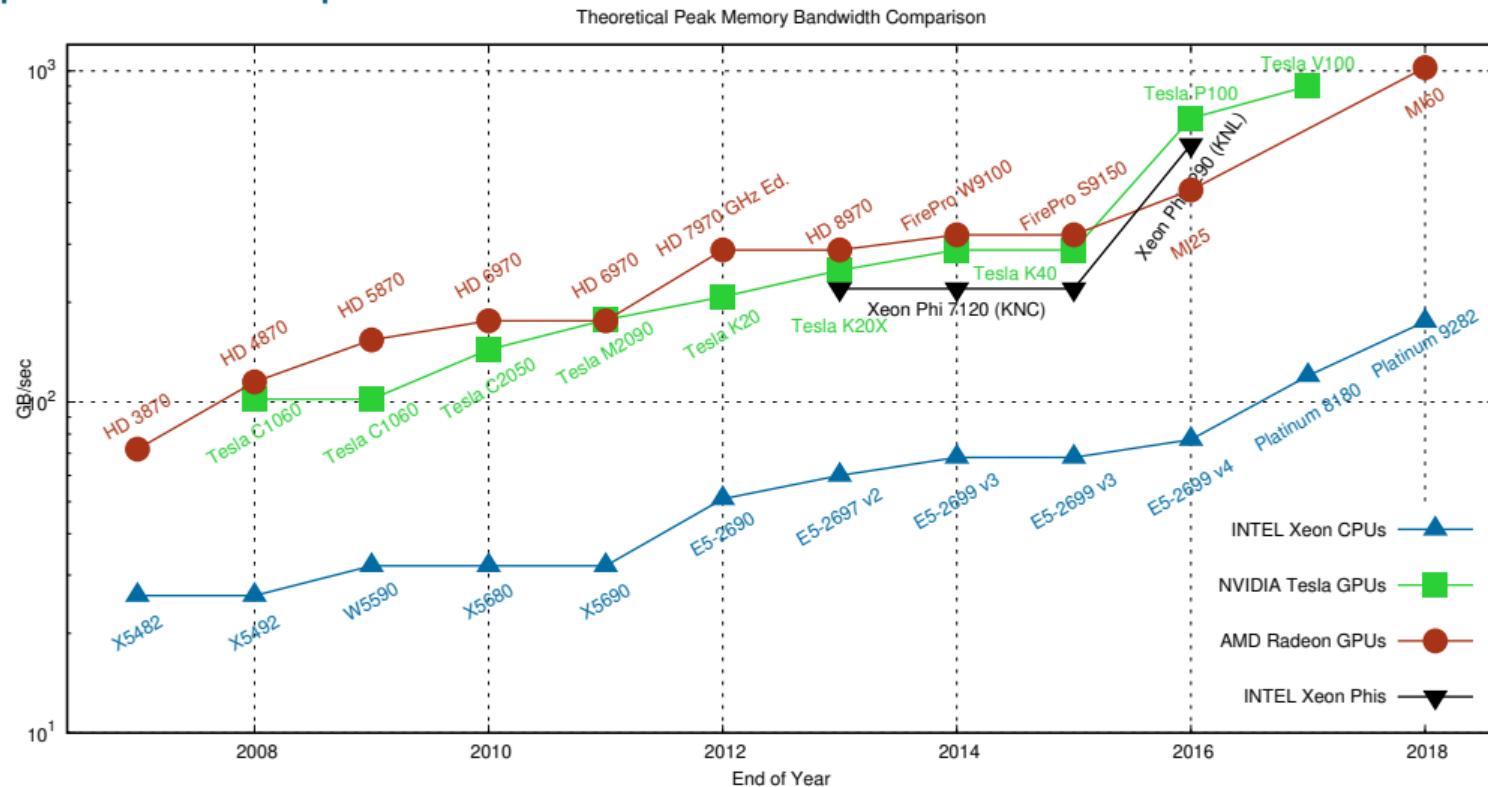
# Status Quo

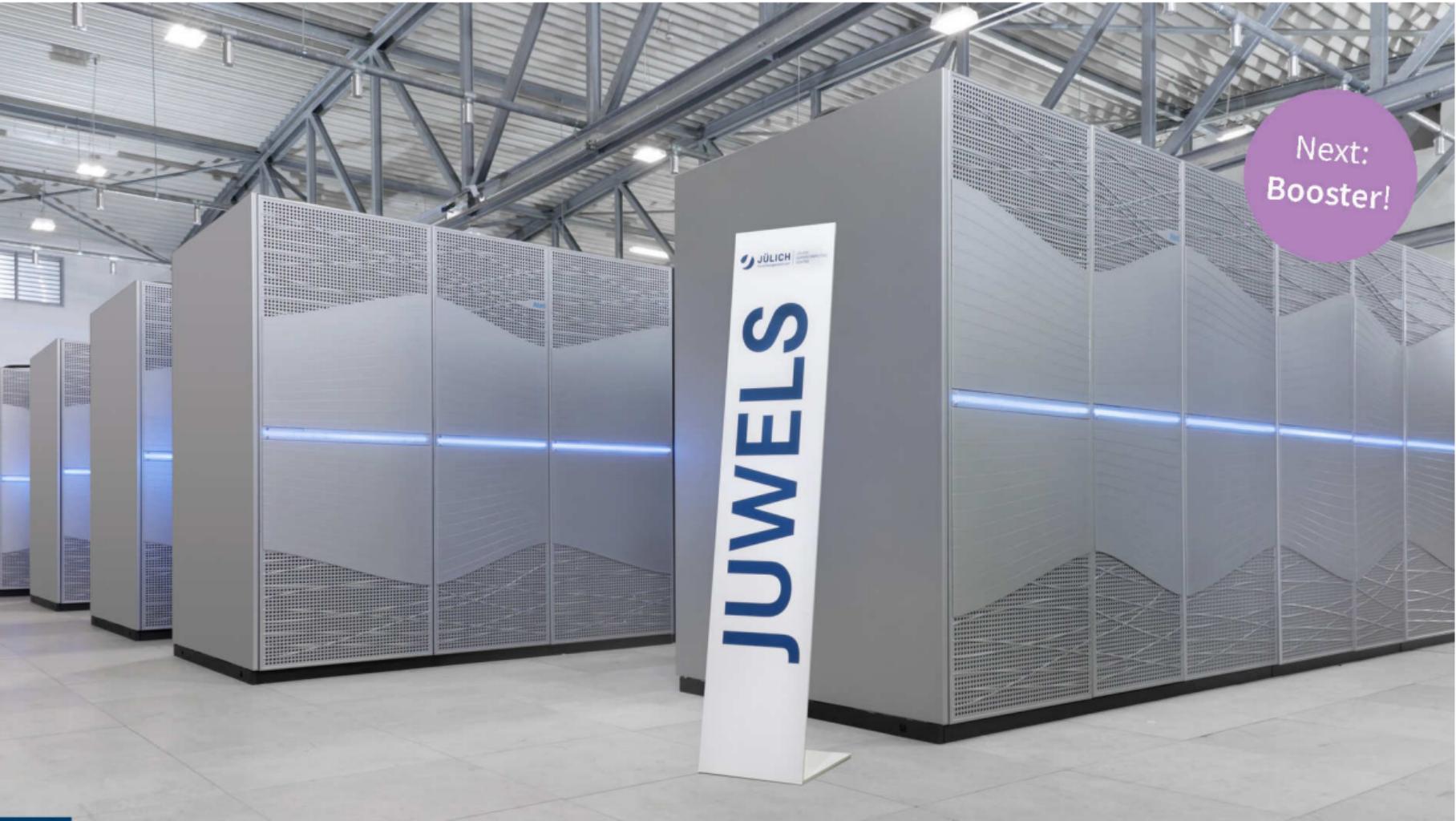
A short but parallel story



# Status Quo

## Peak performance double precision





Next:  
Booster!

**JUWELS**





Next:  
Booster!

*But why?!*

*Let's find out!*

# Platform

# CPU vs. GPU

A matter of specialties



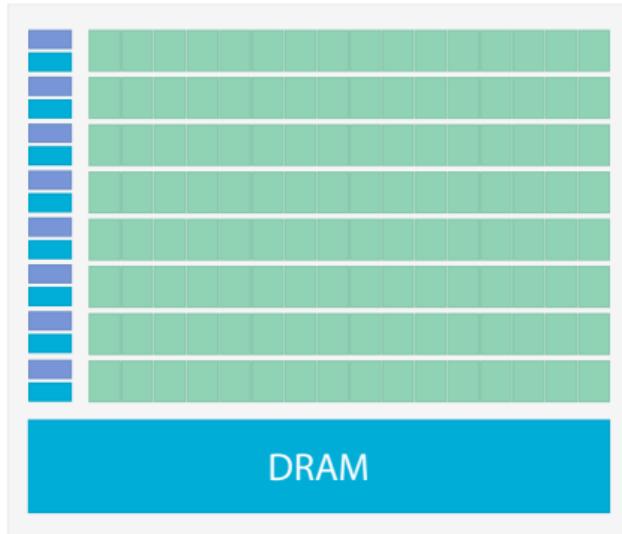
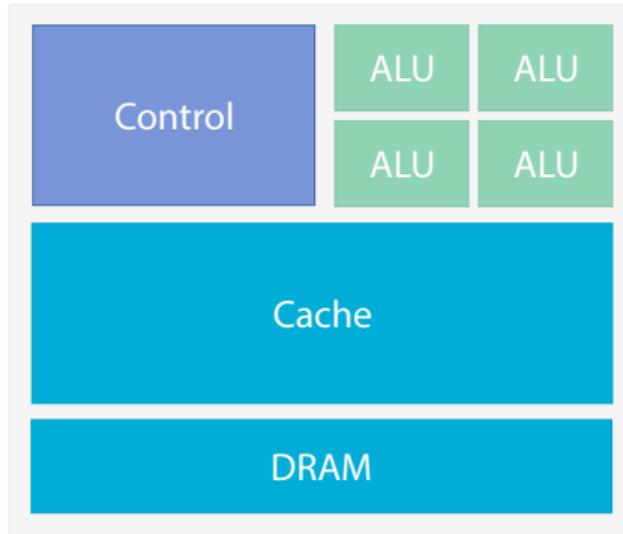
Transporting one



Transporting many

# CPU vs. GPU

## Chip



# GPU Architecture

## Overview

Aim: Hide Latency  
*Everything else follows*

SIMT

Asynchronicity

Memory

# GPU Architecture

## Overview

Aim: Hide Latency  
*Everything else follows*

SIMT

Asynchronicity

Memory

# Memory

GPU memory ain't no CPU memory

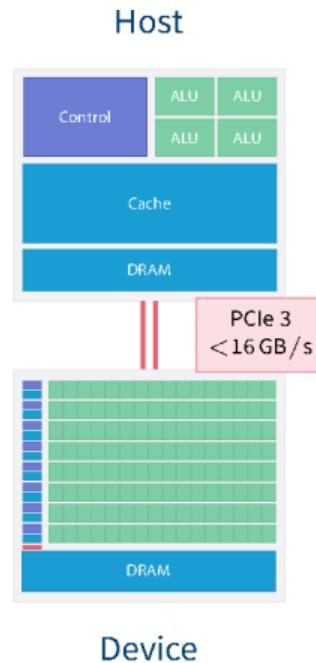
Unified Virtual Addressing

- GPU: accelerator / extension card
- Separate device from CPU
- Separate memory, but UVA**
- Memory transfers need special consideration!  
*Do as little as possible!*
- Formerly: Explicitly copy data to/from GPU  
Now: Done automatically (performance...?)

P100  
16 GB RAM, 720 GB/s

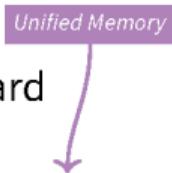


V100  
32 GB RAM, 900 GB/s



# Memory

GPU memory ain't no CPU memory

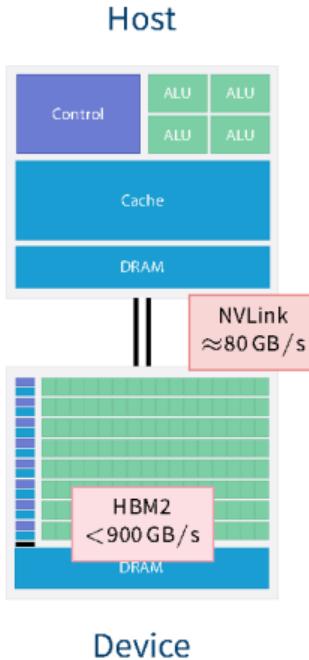
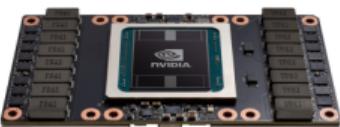


- GPU: accelerator / extension card  
→ Separate device from CPU
- Separate memory, but UVA and UM**
- Memory transfers need special consideration!  
*Do as little as possible!*
- Formerly: Explicitly copy data to/from GPU  
Now: Done automatically (performance...?)

P100  
16 GB RAM, 720 GB/s



V100  
32 GB RAM, 900 GB/s



# GPU Architecture

## Overview

Aim: Hide Latency  
*Everything else follows*

SIMT

Asynchronicity

Memory

# Async

## Following different streams

- Problem: Memory transfer is comparably slow  
Solution: Do something else in meantime (**computation**)!  
→ Overlap tasks
- Copy and compute engines run separately (*streams*)



- GPU needs to be fed: Schedule many computations
- CPU can do other work while GPU computes; synchronization
- Also: Fast switching of contexts to keep GPU busy (*KGB*)

# GPU Architecture

## Overview

Aim: Hide Latency  
*Everything else follows*

**SIMT**

Asynchronicity

Memory

# SIMT

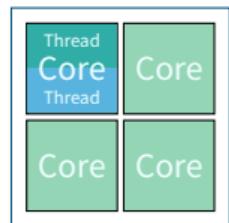
## Of threads and warps

- CPU:
  - Single Instruction, Multiple Data (**SIMD**)
  - Simultaneous Multithreading (**SMT**)
- GPU: Single Instruction, Multiple Threads (**SIMT**)
  - CPU core  $\approx$  GPU multiprocessor (**SM**)
  - Working unit: set of threads (32, a *warp*)
  - Fast switching of threads (large register file)  $\rightarrow$  **hide latency**
  - Branching    

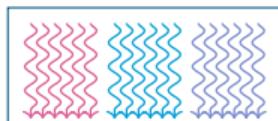
Vector

$$\begin{array}{ccc} A_0 & B_0 & C_0 \\ A_1 & B_1 & C_1 \\ A_2 & B_2 & C_2 \\ A_3 & B_3 & C_3 \end{array} + \begin{array}{ccc} & & \\ & & \end{array} = \begin{array}{ccc} & & \\ & & \end{array}$$

SMT



SIMT



# SIMT

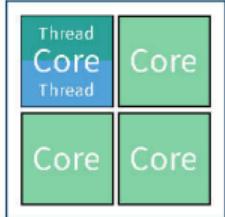
of t



Vector

$$\begin{matrix} A_0 \\ A_1 \\ A_2 \\ A_3 \end{matrix} + \begin{matrix} B_0 \\ B_1 \\ B_2 \\ B_3 \end{matrix} = \begin{matrix} C_0 \\ C_1 \\ C_2 \\ C_3 \end{matrix}$$

SMT



SIMT



# SIMT

of



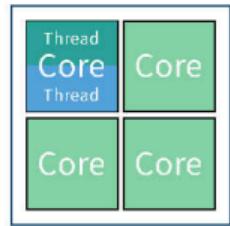
# Multiprocessor



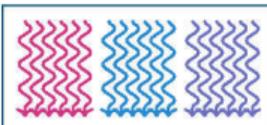
# Vector

$$\begin{matrix} A_0 \\ A_1 \\ A_2 \\ A_3 \end{matrix} + \begin{matrix} B_0 \\ B_1 \\ B_2 \\ B_3 \end{matrix} = \begin{matrix} C_0 \\ C_1 \\ C_2 \\ C_3 \end{matrix}$$

# SMT



# SIMT



# New: Tensor Cores

## New in Volta

- 8 Tensor Cores per Streaming Multiprocessor (SM) (640 total for V100)
  - Performance: 125 TFLOP/s (half precision)
  - Calculate  $\mathbf{A} \times \mathbf{B} + \mathbf{C} = \mathbf{D}$  ( $4 \times 4$  matrices;  $\mathbf{A}, \mathbf{B}$ : half precision)
- 64 floating-point FMA operations per clock (mixed precision)

$$\begin{array}{c} \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \end{array} \times \begin{array}{c} \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \end{array} + \begin{array}{c} \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \end{array} = \begin{array}{c} \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \\ \bullet \bullet \bullet \bullet \end{array}$$

FP16      FP32      FP16      FP32      FP16      FP32      FP16      FP32

FP16  
FP32

# CPU vs. GPU

Let's summarize this!



Optimized for **low latency**

- + Large main memory
- + Fast clock rate
- + Large caches
- + Branch prediction
- + Powerful ALU
- Relatively low memory bandwidth
- Cache misses costly
- Low performance per watt



Optimized for **high throughput**

- + High bandwidth main memory
- + Latency tolerant (parallelism)
- + More compute resources
- + High performance per watt
- Limited memory capacity
- Low per-thread performance
- Extension card

# Programming GPUs

# Preface: CPU

A simple CPU program as reference!

SAXPY:  $\vec{y} = a\vec{x} + \vec{y}$ , with single precision

Part of LAPACK BLAS Level 1

```
void saxpy(int n, float a, float * x, float * y) {
    for (int i = 0; i < n; i++)
        y[i] = a * x[i] + y[i];
}

int a = 42;
int n = 10;
float x[n], y[n];
// fill x, y

saxpy(n, a, x, y);
```

# Libraries

Programming GPUs is easy: Just don't!

*Use applications & libraries*



# Libraries

Programming GPUs is easy: Just don't!

*Use applications & libraries*



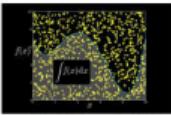
cuSPARSE  
rocSPARSE



cuDNN  
rocDNN



cuFFT  
rocFFT



cuRAND  
rocRAND



{ } ARRAYFIRE

Numba

theano



CUDA Math

# BLAS on GPU

Parallel algebra



## cuBLAS

- GPU-parallel BLAS (all 152 routines) by NVIDIA
- Single, double, complex data types
- Constant competition with Intel's MKL
- Multi-GPU support

→ <https://developer.nvidia.com/cublas>  
<http://docs.nvidia.com/cuda/cublas>

## rocBLAS

- AMD BLAS implementation

→ <https://github.com/ROCmSoftwarePlatform/rocBLAS>  
<https://rocm.readthedocs.io/en/latest/>

# cuBLAS

## Code example

```
int a = 42;  int n = 10;
float x[n], y[n];
// fill x, y

cublasHandle_t handle;
cublasCreate(&handle);

float * d_x, * d_y;
cudaMallocManaged(&d_x, n * sizeof(x[0]));
cudaMallocManaged(&d_y, n * sizeof(y[0]));
cublasSetVector(n, sizeof(x[0]), x, 1, d_x, 1);
cublasSetVector(n, sizeof(y[0]), y, 1, d_y, 1);

cublasSaxpy(n, a, d_x, 1, d_y, 1);

cublasGetVector(n, sizeof(y[0]), d_y, 1, y, 1);

cudaFree(d_x); cudaFree(d_y);
cublasDestroy(handle);
```

# cuBLAS

## Code example

```
int a = 42;  int n = 10;  
float x[n], y[n];  
// fill x, y
```

```
cublasHandle_t handle;  
cublasCreate(&handle);
```

Initialize

```
float * d_x, * d_y;
```

```
cudaMallocManaged(&d_x, n * sizeof(x[0]));
```

Allocate GPU memory

```
cudaMallocManaged(&d_y, n * sizeof(y[0]));
```

Copy data to GPU

```
cublasSetVector(n, sizeof(x[0]), x, 1, d_x, 1);  
cublasSetVector(n, sizeof(y[0]), y, 1, d_y, 1);
```

```
cublasSaxpy(n, a, d_x, 1, d_y, 1);
```

Call BLAS routine

```
cublasGetVector(n, sizeof(y[0]), d_y, 1, y, 1);
```

Copy result to host

```
cudaFree(d_x); cudaFree(d_y);  
cublasDestroy(handle);
```

Finalize

# Libraries

Programming GPUs is easy: Just don't!

*Use applications & libraries*



cuSPARSE  
rocSPARSE



cuDNN  
rocDNN



cuFFT  
rocFFT



cuRAND  
rocRAND



Numba

theano



# Thrust

Iterators! Iterators everywhere! 

- $\frac{\text{Thrust}}{\text{CUDA}} = \frac{\text{STL}}{\text{C++}}$
  - Template library
  - Based on iterators
  - Data-parallel primitives (`scan()`, `sort()`, `reduce()`, ...)
  - Fully compatible with plain CUDA C (comes with **CUDA Toolkit**)
  - Great with `[](){}` lambdas!
- <http://thrust.github.io/>  
<http://docs.nvidia.com/cuda/thrust/>
- AMD backend available: <https://github.com/ROCmSoftwarePlatform/Thrust>

# Thrust

## Code example

```
int a = 42;
int n = 10;
thrust::host_vector<float> x(n), y(n);
// fill x, y

thrust::device_vector d_x = x, d_y = y;

using namespace thrust::placeholders;
thrust::transform(d_x.begin(), d_x.end(), d_y.begin(), d_y.begin(), a * _1 + _2);

x = d_x;
```

*One more example with classical lambdas in appendix!*

# Programming GPUs

## Directives

# GPU Programming with Directives

Keepin' you portable

- Annotate usual source code by directives

```
#pragma acc loop  
for (int i = 0; i < 1; i++) {};
```

- Also: Generalized API functions
- acc\_copy();
- Compiler interprets directives, creates according instructions

## Pro

- Portability
  - Other compiler? No problem! To it, it's a serial program
  - Different target architectures from same code
- Easy to program

## Con

- Compilers support limited
- Raw power hidden
- Somewhat harder to debug

# GPU Programming with Directives

The power of... two.

OpenMP Standard for multithread programming on CPU, GPU since 4.0, better since 4.5

```
#pragma omp target map(tofrom:y), map(to:x)
#pragma omp teams num_teams(10) num_threads(10)
#pragma omp distribute
for ( ) {
    #pragma omp parallel for
    for ( ) {
        // ...
    }
}
```

OpenACC Similar to OpenMP, but more specifically for GPUs

Might eventually be re-merged into OpenMP standard

# OpenACC

## Code example

```
void saxpy_acc(int n, float a, float * x, float * y) {  
    #pragma acc kernels  
    for (int i = 0; i < n; i++)  
        y[i] = a * x[i] + y[i];  
}  
  
int a = 42;  
int n = 10;  
float x[n], y[n];  
// fill x, y  
  
saxpy_acc(n, a, x, y);
```

# OpenACC

## Code example

```
void saxpy_acc(int n, float a, float * x, float * y) {  
    #pragma acc kernels  
    for (int i = 0; i < n; i++)  
        y[i] = a * x[i] + y[i];  
}  
  
int a = 42;  
int n = 10;  
float x[n], y[n];  
// fill x, y  
  
saxpy_acc(n, a, x, y);
```

GPU tutorial  
this afternoon!

# Programming GPUs

## Languages, finally

# Programming GPU Directly

Finally...

**OpenCL** Open Computing Language by Khronos Group (Apple, IBM, NVIDIA, ...) 2009

- Platform: Programming language (OpenCL C/C++), API, and compiler
- Targets CPUs, GPUs, FPGAs, and other many-core machines
- Fully open source

**CUDA** NVIDIA's GPU platform 2007

- Platform: Drivers, programming language (CUDA C/C++), API, compiler, tools, ...
- Only NVIDIA GPUs
- Compilation with nvcc (free, but not open)  
clang has CUDA support, but CUDA needed for last step
- Also: CUDA Fortran

**HIP** AMD's new unified programming model for AMD (via ROCm) and NVIDIA GPUs 2016+

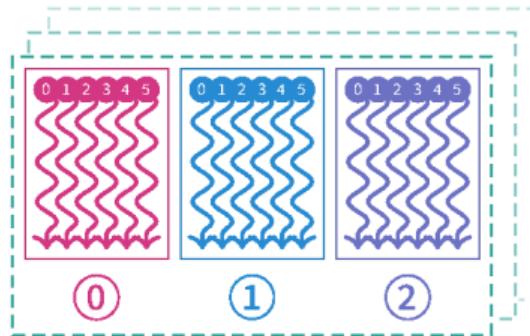
- Choose what flavor you like, what colleagues/collaboration is using
- **Hardest: Come up with parallelized algorithm**

# CUDA Threading Model

Warp the kernel, it's a thread!

- Methods to exploit parallelism:

- Thread → Block
- Block → Grid
- Threads & blocks in 3D



- Parallel function: **kernel**

- **\_\_global\_\_** kernel(**int** a, **float** \* b) { }
- Access own ID by global variables **threadIdx.x**, **blockIdx.y**, ...

- Execution entity: **threads**

- Lightweight → fast switching!
- 1000s threads execute simultaneously → order non-deterministic!

⇒ SAXPY!

# CUDA SAXPY

With runtime-managed data transfers

```
__global__ void saxpy(int n, float a, float * x, float * y) {  
    int i = blockIdx.x * blockDim.x + threadIdx.x;  
    if (i < n)  
        y[i] = a * x[i] + y[i];  
}
```

Specify kernel

ID variables

Guard against  
too many threads

```
int a = 42;  
int n = 10;  
float x[n], y[n];  
// fill x, y  
cudaMallocManaged(&x, n * sizeof(float));  
cudaMallocManaged(&y, n * sizeof(float));  
  
saxpy_cuda<<<2, 5>>>(n, a, x, y);
```

Allocate GPU-capable  
memory

Call kernel  
2 blocks, each 5 threads

```
cudaDeviceSynchronize();
```

Wait for  
kernel to finish

# HIP SAXPY

From CUDA to HIP

```
#include <cuda_runtime.h>
__global__ void saxpy(int n, float a, float * x, float * y) {
    int i = blockIdx.x * blockDim.x + threadIdx.x;
    if (i < n)
        y[i] = a * x[i] + y[i];
}
```

```
int a = 42;
int n = 10;
float x[n], y[n];
// fill x, y
cudaMallocManaged(&x, n * sizeof(float));
cudaMallocManaged(&y, n * sizeof(float));

saxpy_cuda<<<2, 5>>>(n, a, x, y);

cudaDeviceSynchronize();
```

Works on AMD and NVIDIA GPUs!

# HIP SAXPY

From CUDA to HIP

```
#include <hip/hip_runtime.h>
__global__ void saxpy(int n, float a, float * x, float * y) {
    int i = blockIdx.x * blockDim.x + threadIdx.x;
    if (i < n)
        y[i] = a * x[i] + y[i];
}
```

Works on AMD and NVIDIA GPUs!

```
int a = 42;
int n = 10;
float x[n], y[n];
// fill x, y
hipMallocManaged(&x, n * sizeof(float));
hipMallocManaged(&y, n * sizeof(float));

hipLaunchKernelGGL(saxpy, 2, 5, 0, 0, n, a, x, y);

hipDeviceSynchronize();
```

# Programming GPUs

## Abstraction Libraries/DSL

# Abstraction Libraries & DSLs

- Libraries with ready-programmed abstractions; partly compiler/transpiler necessary
- Have different backends to choose from for targeted accelerator
- Between Thrust, OpenACC, and CUDA
- Examples: **SYCL**, **Kokkos**, **Alpaka**, **Futhark**, **C++AMP**, ...

# An Alternative: Kokkos

From Sandia National Laboratories

- C++ library for *performance* portability
- Data-parallel patterns, architecture-aware memory layouts, ...

```
Kokkos::View<double*> x("X", length);
Kokkos::View<double*> y("Y", length);
double a = 2.0;

// Fill x, y

Kokkos::parallel_for(length, KOKKOS_LAMBDA (const int& i) {
    x(i) = a*x(i) + y(i);
});
```

→ <https://github.com/kokkos/kokkos/>

# Another Alternative: SYCL

- Extension of/upon OpenCL
- With buffers, queues, accessors, lambdas, ...
- Main programming model for Aurora's Intel GPUs

→ [khronos.org/sycl/](http://khronos.org/sycl/)

```
class mySaxpy;  
  
std::vector<double> h_x(length), h_y(length);  
// Fill x, y  
cl::sycl::buffer<double, 1> d_x(h_x), d_y(h_y);  
  
cl::sycl::queue queue;  
  
queue.submit([&] (cl::sycl::handler& cgh) {  
    auto x_acc = d_x.get_access<cl::sycl::access::mode::read>(cgh);  
    auto y_acc = d_y.get_access<cl::sycl::access::mode::read>(cgh);  
  
    cgh.parallel_for<class mySaxpy>(length,  
        [=] (cl::sycl::id<1> idx) {  
            y_acc[idx] = a * x_acc[idx] + y_acc[idx];  
        });  
});
```

# Programming GPUs

## Tools

# GPU Tools

The helpful helpers helping helpless (and others)

- NVIDIA

- [cuda-gdb](#) GDB-like command line utility for debugging
- [cuda-memcheck](#) Like Valgrind's memcheck, for checking errors in memory accesses
- [Nsight](#) IDE for GPU developing, based on Eclipse (Linux, OS X) or Visual Studio (Windows)
- [nvprof](#) Command line profiler, including detailed performance counters
- [Visual Profiler](#) Timeline profiling and annotated performance experiments
- New** [Nsight Systems](#) (timeline), [Nsight Compute](#) (kernel analysis)

- OpenCL/HIP:

- [CodeXL](#) Debugging, profiling.
- [ROCmGDB](#) AMD's GDB symbolic debugger
- [RadeonComputeProfiler](#) Profiler for OpenCL and ROCm

# nvprof

## Command that line

Usage: nvprof ./app

```
$ nvprof ./matrixMul -wA=1024 -hA=1024 -wB=1024 -hB=1024
==37064== Profiling application: ./matrixMul -wA=1024 -hA=1024 -wB=1024 -hB=1024
==37064== Profiling result:
Time(%)      Time      Calls      Avg      Min      Max  Name
 99.19%  262.43ms      301  871.86us  863.88us  882.44us void matrixMulCUDA<int=32>(float*, float*, float*, int, int)
   0.58%  1.5428ms       2  771.39us  764.65us  778.12us [CUDA memcpy HtoD]
   0.23%  599.40us       1  599.40us  599.40us  599.40us [CUDA memcpy DtoH]

==37064== API calls:
Time(%)      Time      Calls      Avg      Min      Max  Name
 61.26%  258.38ms       1  258.38ms  258.38ms  258.38ms cudaEventSynchronize
 35.68%  150.49ms       3  50.164ms  914.97us  148.65ms cudaMalloc
   0.73%  3.0774ms       3  1.0258ms  1.0097ms  1.0565ms cudaMemcpy
   0.62%  2.6287ms       4  657.17us  655.12us  660.56us cuDeviceTotalMem
   0.56%  2.3408ms      301  7.7760us  7.3810us  53.103us cudaLaunch
   0.48%  2.0111ms      364  5.5250us   235ns  201.63us cuDeviceGetAttribute
   0.21%  872.52us       1  872.52us  872.52us  872.52us cudaDeviceSynchronize
```

# nvprof

## Command that line

With metrics: nvprof --metrics flop\_sp\_efficiency ./app

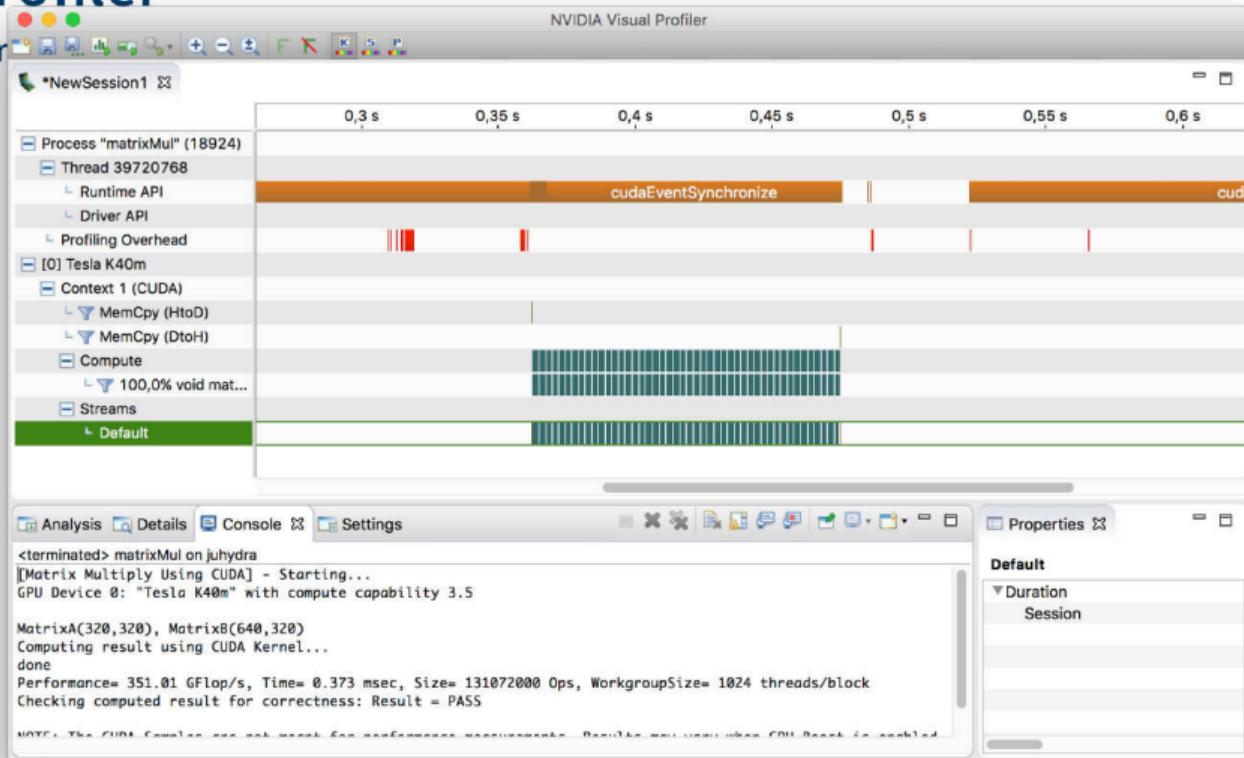


```
$ nvprof --metrics flop_sp_efficiency ./matrixMul -wA=1024 -hA=1024 -wB=1024 -hB=1024
[Matrix Multiply Using CUDA] - Starting...
==37122== NVPROF is profiling process 37122, command: ./matrixMul -wA=1024 -hA=1024 -wB=1024 -hB=1024
GPU Device 0: "Tesla P100-SXM2-16GB" with compute capability 6.0

MatrixA(1024,1024), MatrixB(1024,1024)
Computing result using CUDA Kernel...
==37122== Some kernel(s) will be replayed on device 0 in order to collect all events/metrics.
done122== Replaying kernel "void matrixMulCUDA<int=32>(float*, float*, float*, int, int)" (0 of 2)...
Performance= 26.61 GFlop/s, Time= 80.697 msec, Size= 2147483648 Ops, WorkgroupSize= 1024 threads/block
Checking computed result for correctness: Result = PASS
==37122== Profiling application: ./matrixMul -wA=1024 -hA=1024 -wB=1024 -hB=1024
==37122== Profiling result:
==37122== Metric result:
      Invocations          Metric Name          Metric Description      Min      Max      Avg
Device "Tesla P100-SXM2-16GB (0)"
      Kernel: void matrixMulCUDA<int=32>(float*, float*, float*, int, int)
            301                flop_sp_efficiency    FLOP Efficiency(Peak Single)  22.96%  23.40%  23.15%
```

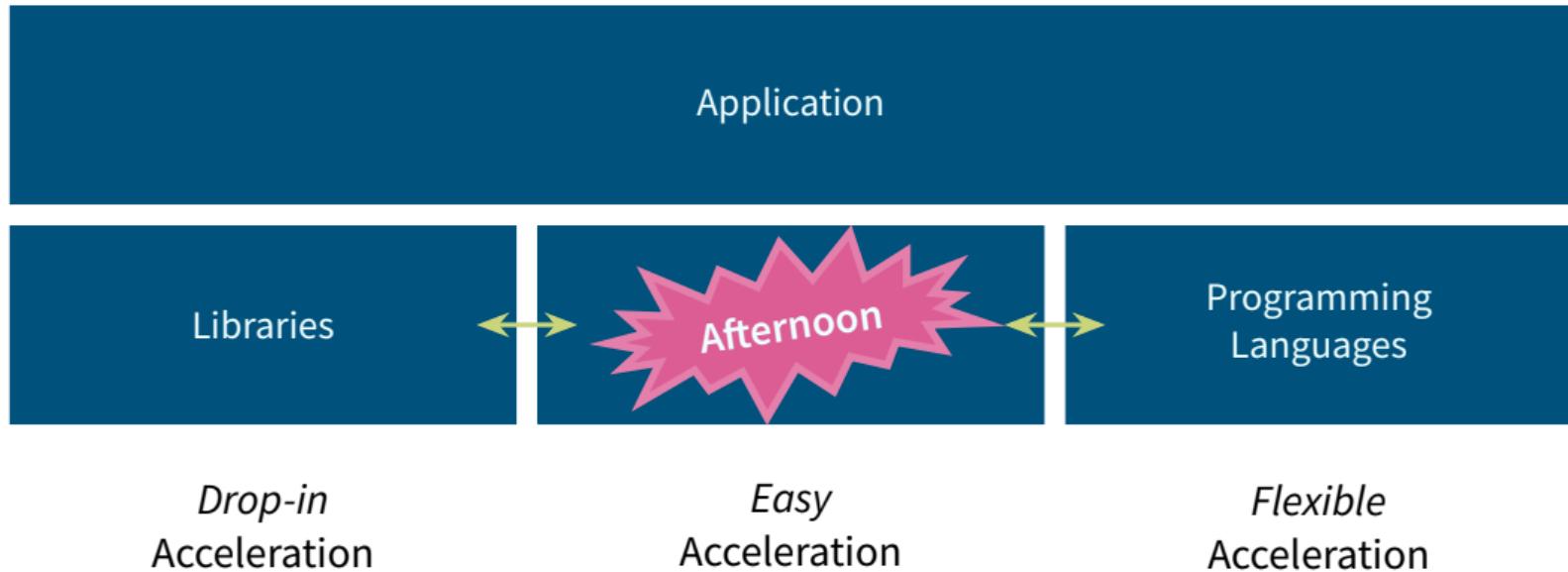
# Visual Profiler

Your new favor



# Conclusions

# Summary of Acceleration Possibilities

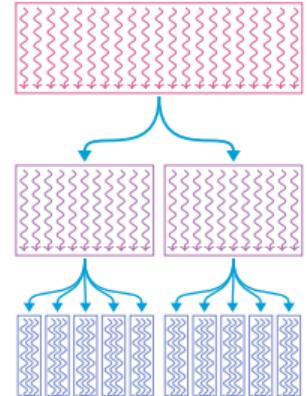


# Omitted

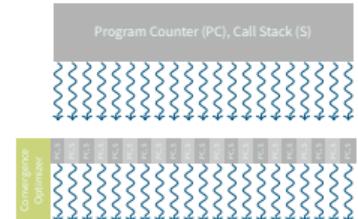
There's so much more!

What I did not talk about

- Atomic  operations
- Shared memory
- Pinned memory
- Managed memory
- Debugging
- Overlapping streams
- Multi-GPU programming (intra-node; [MPI](#))
- Cooperative groups
- Independent thread progress
- Half precision FP16
- ...



Cooperative Groups



Independent Thread Progress

# Summary & Conclusion

- GPUs can improve your performance many-fold
  - For a fitting, parallelizable application
  - Libraries are easiest
  - Direct programming (plain CUDA, HIP) is most powerful
  - OpenACC/OpenMP is somewhere in between (and portable)
  - Many abstraction layers available (mostly using C++)
  - There are many tools helping the programmer
- See it in action this afternoon at [OpenACC tutorial](#)

Thank you  
for your attention!  
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# APPENDIX

## Appendix

[Further Reading & Links](#)

[GPU Performances](#)

[Supplemental: Thrust](#)

[Glossary](#)

[References](#)

# Further Reading & Links

More!

- A discussion of SIMD, SIMT, SMT by Y. Kreinin.
- NVIDIA's documentation: [docs.nvidia.com](https://docs.nvidia.com)
- NVIDIA's [Parallel For All blog](#)
- SYCL Hello World, SYCL Vector Addition

# Volta Performance

9-@ &gt;+3/2	9-@ CDO	9-@ E DO	9-@ &lt;100	9-@ V100
GPU	GK180 (Kepler)	GM200 (Maxwell)	GP100 (Pascal)	GV100 (Volta)
SMs	15	24	56	80
TPCs	15	24	28	40
FP32 Cores / SM	1B2	128	64	64
FP32 Cores / GPU	2880	30C2	3584	5120
FP64 Cores / SM	64	4	32	32
FP64 Cores / GPU	B50	B6	1CB2	2560
Tensor Cores / SM	EF	EF	EF	8
Tensor Cores / GPU	EF	EF	EF	640
GPU Boost Clock	810/8C5 MHz	1114 MHz	1480 MHz	1462 MHz
PeakTF32 TFLPS <sup>1</sup>	5	6M	10M	15
PeakTF64 TFLPS <sup>1</sup>	1N8	M1	5M	ON
PeakTensor TFLPS <sup>1</sup>	EF	EF	EF	120
Texture UDR	240	1B2	224	320
Memory Bandwidth	384 TB/s	384 TB/s	40B6 TUDI GM2	40B6 TUDI GM2
Memory QDR	Up to 12 GG	Up to 24 GG	16 GG	16 GG
L1 Cache Size	1536 KB	30C2 KB	40B6 KB	6144 KB
System Memory / SM	16 KB/32 KB/48 KB	B6 KB	64 KB	CoD2NaUe Np to B6 KB
WZG per FLOPs / SM	256 KB	256 KB	256 KB	256KB
WZG per FLOPs / GPU	3840 KB	6144 KB	14336 KB	20480 KB
TFP	235 f atts	250 f atts	300 f atts	300 f atts
Tensor cores	ON UDD	8 UDD	15M UDD	21M UDD
GPU VRAM	551 PPV	601 PPV	610 PPV	815 PPV
Memory Bandwidth Process	28 DP	28 DP	16 DP FIFTA	12 DP FFE

<sup>1</sup> PeakTFLPS rates are based on GPU Boost Clock

Figure: Tesla V100 performance characteristics in comparison [9]

# Thrust

## Code example with lambdas

```
#include <thrust/for_each.h>
#include <thrust/execution_policy.h>
constexpr int gGpuThreshold = 10000;
void saxpy(float *x, float *y, float a, int N) {
    auto r = thrust::counting_iterator<int>(0);

    auto lambda = [=] __host__ __device__ (int i) {
        y[i] = a * x[i] + y[i];};

    if(N > gGpuThreshold)
        thrust::for_each(thrust::device, r, r+N, lambda);
    else
        thrust::for_each(thrust::host, r, r+N, lambda);}

```

Source

# Appendix

## Glossary & References

# Glossary I

- AMD** Manufacturer of **CPUs** and **GPUs**. [3](#), [27](#), [31](#), [39](#), [42](#), [43](#), [49](#), [64](#), [65](#)
- API** A programmatic interface to software by well-defined functions. Short for application programming interface. [34](#), [39](#), [65](#)
- ATI** Canada-based **GPUs** manufacturing company; bought by AMD in 2006. [3](#)
- CUDA** Computing platform for **GPUs** from NVIDIA. Provides, among others, CUDA C/C++. [3](#), [31](#), [39](#), [40](#), [41](#), [42](#), [43](#), [45](#), [56](#), [64](#)
- DSL** A Domain-Specific Language is a specialization of a more general language to a specific domain. [2](#), [44](#), [45](#)

## Glossary II

**HIP** GPU programming model by AMD to target their own and NVIDIA GPUs with one combined language. Short for Heterogeneous-compute Interface for Portability.  
[39](#), [42](#), [43](#), [49](#), [56](#)

**MPI** The Message Passing Interface, a API definition for multi-node computing. [55](#)

**NVIDIA** US technology company creating **GPUs**. [3](#), [27](#), [39](#), [42](#), [43](#), [49](#), [59](#), [63](#), [64](#), [66](#)

**OpenACC** Directive-based programming, primarily for many-core machines. [35](#), [36](#), [37](#), [45](#), [56](#)

**OpenCL** The *Open Computing Language*. Framework for writing code for heterogeneous architectures (**CPU**, **GPU**, DSP, FPGA). The alternative to **CUDA**. [3](#), [39](#), [47](#), [49](#)

# Glossary III

- OpenGL** The *Open Graphics Library*, an API for rendering graphics across different hardware architectures. [3](#)
- OpenMP** Directive-based programming, primarily for multi-threaded machines. [35](#), [56](#)
- POWER** CPU architecture from IBM, earlier: PowerPC. See also POWER8. [65](#)
- POWER8** Version 8 of IBM's **POWER**processor, available also under the OpenPOWER Foundation. [65](#)
- ROCM** AMD software stack and platform to program AMD GPUs. Short for Radeon Open Compute (*Radeon* is the GPU product line of AMD). [39](#), [49](#)
- SAXPY** Single-precision  $A \times X + Y$ . A simple code example of scaling a vector and adding an offset. [24](#), [40](#), [41](#)

# Glossary IV

**Thrust** A parallel algorithms library for (among others) GPUs. See  
<https://thrust.github.io/>. 31

**Volta** GPU architecture from [NVIDIA](#) (announced 2017). 21

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- [2] Kenneth E. Hoff III et al. “Fast Computation of Generalized Voronoi Diagrams Using Graphics Hardware”. In: *Proceedings of the 26th Annual Conference on Computer Graphics and Interactive Techniques*. SIGGRAPH ’99. New York, NY, USA: ACM Press/Addison-Wesley Publishing Co., 1999, pp. 277–286. ISBN: 0-201-48560-5. DOI: [10.1145/311535.311567](https://doi.org/10.1145/311535.311567). URL: <http://dx.doi.org/10.1145/311535.311567> (page 3).
- [3] Chris McClanahan. “History and Evolution of GPU Architecture”. In: *A Survey Paper* (2010). URL: <http://mcclanahoochie.com/blog/wp-content/uploads/2011/03/gpu-hist-paper.pdf> (page 3).
- [4] Jack Dongarra et al. *TOP500*. June 2019. URL: <https://www.top500.org/lists/2019/06/> (page 3).

## References II

- [5] Jack Dongarra et al. *Green500*. June 2019. URL:  
<https://www.top500.org/green500/lists/2019/06/> (page 3).
- [6] Karl Rupp. *Pictures: CPU/GPU Performance Comparison*. URL:  
<https://www.karlrupp.net/2013/06/cpu-gpu-and-mic-hardware-characteristics-over-time/> (pages 4, 5).
- [10] Wes Breazell. *Picture: Wizard*. URL:  
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