

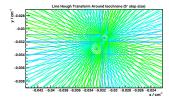
## GPUs: Platform, Programming, Pitfalls GridKa School 2016: Data Science on Modern Architectures

Andreas Herten, Forschungszentrum Jülich, 1 September 2016

## About, Outline

#### Andreas Herten

- Physics in
  - Aachen (Dipl. at CMS)
  - Jülich/Bochum (Dr. at PANDA)



Since then: NVIDIA Application Lab Optimizing scientific applications for/on GPUs

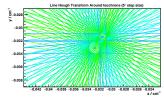




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Motivation Platform Hardware Features Programming Libraries Directives Languages Tools Pitfalls

 Since then: NVIDIA Application Lab Optimizing scientific applications for/on GPUs





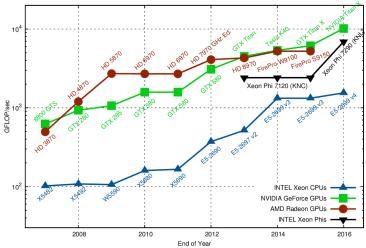




- 1999: General computations with shaders of graphics hardware
- 2001: NVIDIA GeForce 3 with programmable shaders [1]; 2003: DirectX 9 at ATI
- 2016: Top 500:  $1/\!_{10}$  with GPUs, Green 500: 70 % of top 50 with GPUs



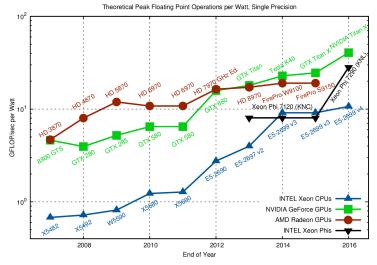




Theoretical Peak Performance, Single Precision

#### **Status Quo** GPU all around





#### **Status Quo** GPU all around



















## Platform

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#4|37

#### **CPU vs. GPU** *A matter of specialties*







#### **CPU vs. GPU** *A matter of specialties*



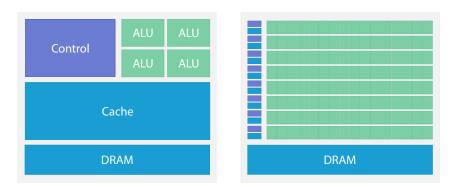


#### Transporting one



Transporting many





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Overview



Aim: Hide Latency *Everything else follows* 

Overview



# Aim: Hide Latency *Everything else follows*

SIMT

Asynchronicity

Memory

Overview



# Aim: Hide Latency *Everything else follows*

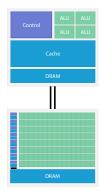
SIMT

Asynchronicity

Memory

- GPU: accelerator / extension card
- ightarrow Separate device from CPU

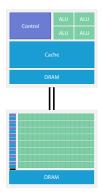






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- → Separate device from CPU Separate memory, but UVA

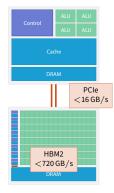






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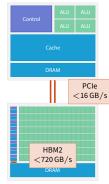


## Memory

#### GPU memory ain't no CPU memory

- GPU: accelerator / extension card
- → Separate device from CPU Separate memory, but UVA
- Memory transfers need special consideration! Do as little as possible!

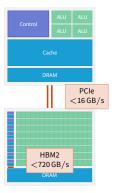






## GPU: accelerator / extension card

- → Separate device from CPU Separate memory, but UVA and UM
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  - Formerly: Explicitly copy data to/from GPU Now: Done automatically (performance...?)

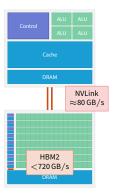






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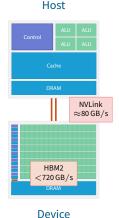






#### GPU: accelerator / extension card

- → 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...?)
  - Values for P100: 16 GB RAM, 720 GB/s





Overview



# Aim: Hide Latency *Everything else follows*

SIMT

Asynchronicity

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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)!
- ightarrow 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

Overview



Aim: Hide Latency *Everything else follows* 

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Overview



# Aim: Hide Latency *Everything else follows*

## SIMT

## Asynchronicity

Memory



#### Scalar

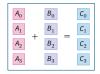


- CPU:
  - Single Instruction, Multiple Data (SIMD)





#### Vector



- CPU:
  - Single Instruction, Multiple Data (SIMD)



- CPU:
  - Single Instruction, Multiple Data (SIMD)
  - Simultaneous Multithreading (SMT)



#### Vector





CPU:

JÜLICH

#### Vector



#### SMT



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Single Instruction, Multiple Data (SIMD)

Simultaneous Multithreading (SMT)

- CPU:
  - Single Instruction, Multiple Data (SIMD)
  - Simultaneous Multithreading (SMT)
- GPU: Single Instruction, Multiple Threads (SIMT)





Vector

SMT



- CPU:
  - Single Instruction, Multiple Data (SIMD)
  - Simultaneous Multithreading (SMT)
- GPU: Single Instruction, Multiple Threads (SIMT)



IÜLICH



SI	M	1	
		_	



SIMT



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



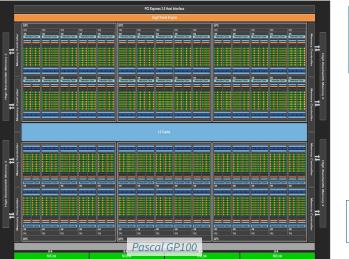
SMT



SIMT







Vector



SMT

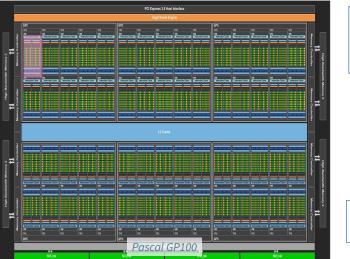


SIMT



#### **SIMT** Of threads and warps





Vector



SMT



SIMT



#### **SIMT** *Of threads and warps*





 $C_0$ 

 $C_1$ 

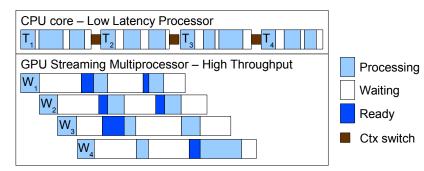
 $C_2$ 

 $C_3$ 



#### • CPU minimizes latency within each thread

GPU hides latency with computations from other thread groups







## **CPU vs. GPU** Low latency vs. high throughput





#### 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

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#### **Preface: CPU** A simple CPU program!



```
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</pre>
```

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





























- GPU-parallel BLAS (all 152 routines)
- 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

#### cuBLAS

Code example

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

```
cublasInit();
```

```
float * d_x, * d_y;
cudaMalloc((void **)&d_x, n * sizeof(x[0]);
cudaMalloc((void **)&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);
cublasShutdown();
```

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#### 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)
- → http://thrust.github.io/ http://docs.nvidia.com/cuda/thrust/



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;

```
x = d_x;
```



## Programming Directives

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Keepin' you portable

Annotate usual source code by directives

#pragma acc loop
for (int i = 0; i < 1; i+\*) {};</pre>



Keepin' you portable

Annotate usual source code by directives

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Also: Generalized functions

acc\_copy();

Compiler interprets directives, creates according instructions



Keepin' you portable

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Also: Generalized 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

- Only few compilers
- Not all the raw power available
- Harder to debug
- Easy to program wrong



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



```
void saxpy acc(int n, float a, float * x, float * y) {
  #pragma acc kernels
  for (int i = 0; i < n; i++)</pre>
    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);
```



```
void saxpy_acc(int n, float a, float * x, float * y) {
    #pragma acc parallel loop copy(y) copyin(x)
    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</pre>
```

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



## Programming Languages

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Finally...

Two solutions:



- Two solutions:
  - 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
    - Different compilers available



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    - CUDA NVIDIA's GPU platform 2007
      - Platform: Drivers, programming language (CUDA C/C++), API, compiler, debuggers, profilers, ...
      - Only NVIDIA GPUs
      - Compilation with nvcc
        - GCC/LLVM solutions on way (slowly)
      - Also: CUDA Fortran



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



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Warp the kernel, it's a thread.

Methods to exploit parallelism:

Warp the kernel, it's a thread.

Methods to exploit parallelism:

Thread





- Methods to exploit parallelism:
  - Threads





- Methods to exploit parallelism:
  - Threads  $\rightarrow$  Block





- Methods to exploit parallelism:
  - Threads  $\rightarrow$  Block
  - Block





- Methods to exploit parallelism:
  - Threads  $\rightarrow$  Block
  - Blocks



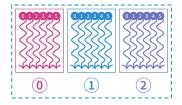


### CUDA C/C++

Warp the kernel, it's a thread.

- Methods to exploit parallelism:
  - Threads  $\rightarrow$  Block
  - Blocks  $\rightarrow$  Grid



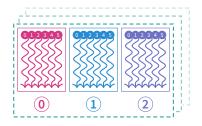


### CUDA C/C++

#### Warp the kernel, it's a thread.

- Methods to exploit parallelism:
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  - $\text{ Blocks} \rightarrow \text{Grid}$
  - All in 3D





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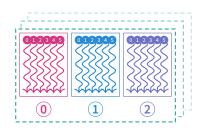
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  - $\text{ Blocks} \rightarrow \text{Grid}$
  - All in 3D
- Execution unit: kernel
  - Function executing in parallel on device

\_\_global\_\_ kernel(int a, float \* b) { }

- Access own ID by global variables threadIdx.x, blockIdx.y,...
- Execution order non-deterministic!
- Only threads in one warp (32 threads of block) can communicate reliably/quickly





### CUDA C/C++

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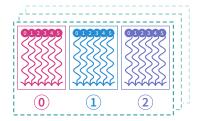
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#### $\Rightarrow$ SAXPY!

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### **CUDA SAXPY**



#### With runtime-managed data transfers

```
global___void saxpy cuda(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, v);
cudaDeviceSynchronize();
```



## Programming Tools

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The helpful helpers helping helpless (and others)

NVIDIA

cuda-gdbGDB-like command line utility for debuggingcuda-memcheckLike Valgrind's memcheck, for checking errors in<br/>memory accessesNsightIDE for GPU developing, based on Eclipse (Linux,<br/>OS X) or Visual Studio (Windows)nvprofCommand line profiler, including detailed<br/>performance countersVisual ProfilerTimeline profiling and annotated performance<br/>experiments

 OpenCL: CodeXL (Open Source, GPUOpen/AMD) – debugging, profiling.



#### nvprof

#### Command that line

#### Usage: nvprof ./app

	(Q) (0) (							
🖲 🕘 🔮 🙎 Slides — aherten@JUHYDRA: ~/cudaSamples/NVIDIA_CUDA-7.5_Samples/bin/x86_64/linux/release —linux/release — ssh juhydra								
NOTE: The CUDA Samples are not meant for performance measurements. Results may vary when GPU Boost is enabled.								
==27580== Profiling application: ./matrixMul								
	≔ Profiling							
Time(%)	Time	Calls	Avg	Min	Max	Name		
99.82%	111.33ms	301	369.85us	363.97us	375.62us	<pre>void matrixMulCUDA<int=32>(float*, float*, float*, int, int)</int=32></pre>		
0.11%	124.58us		62.289us	43.393us	81.185us	[CUDA memcpy HtoD]		
0.07%	80.736us		80.736us	80.736us	80.736us	[CUDA memcpy DtoH]		
==27580== API calls:								
Time(%)	Time	Calls	Avg	Min	Max	Name		
	348.27ms		116.09ms	241.59us	347.79ms	cudaMalloc		
	226.68ms		226.68ms	226.68ms	226.68ms	cudaDeviceReset		
	107.40ms		107.40ms	107.40ms	107.40ms	cudaEventSynchronize		
	3.5853ms	301	11.911us	11.045us	34.486us	cudaLaunch		
	2.4915ms	332	7.5040us	196ns	277.14us	cuDeviceGetAttribute		
	1.6478ms			294.19us	539.73us	cuDeviceTotalMem		
	1.3333ms		444.43us	181.15us	813.00us	cudaMemcpy		
0.12%	802.85us			249.19us	299.41us	cudaFree		
0.09%	604.10us		604.10us	604.10us	604.10us	cudaGetDeviceProperties		
	451.30us	1505	299ns	266ns	6.0860us	cudaSetupArgument		
	362.32us			362.32us	362.32us	cudaDeviceSynchronize		
	242.14us		60.534us	56.884us	69.764us	cuDeviceGetName		
	127.99us	301	425ns	384ns	2.4580us	cudaConfigureCall		
	10.920us		5.4600us	4.2100us	6.7100us	cudaEventRecord		
	10.613us		10.613us	10.613us	10.613us	cudaGetDevice		
0.00%	9.4980us		1.1870us	246ns	4.2760us	cuDeviceGet		
	5.7490us		2.8740us		4.5790us	cudaEventCreate		
	5.4630us		5.4630us	5.4630us	5.4630us	cudaEventElapsedTime		
0.00%	3.2900us		1.6450us	1.2160us	2.0740us	cuDeviceGetCount		
_								

#### nvprof



#### Command that line

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

	"void matrixMulCUDA <int=32>(float*,</int=32>				
=27425== Replaying kernel	"void matrixMulCUDA <int=32>(float*,</int=32>	<pre>float*, float*, int, int)" (dor</pre>	ne)		
	"void matrixMulCUDA <int=32>(float*;</int=32>				
	"void matrixMulCUDA <int=32>(float*;</int=32>				
=27425== Replaying kernel	"void matrixMulCUDA <int=32>(float*,</int=32>	<pre>float*, float*, int, int)" (dor</pre>	ne)		
=27425== Replaying kernel	"void matrixMulCUDA <int=32>(float*;</int=32>	<pre>float*, float*, int, int)" (dor</pre>	ıe)		
	"void matrixMulCUDA <int=32>(float*;</int=32>				
	"void matrixMulCUDA <int=32>(float*,</int=32>				
	"void matrixMulCUDA <int=32>(float*,</int=32>				
	"void matrixMulCUDA <int=32>(float*;</int=32>				
	"void matrixMulCUDA <int=32>(float*,</int=32>				
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	"void matrixMulCUDA <int=32>(float*;</int=32>				
	"void matrixMulCUDA <int=32>(float*,</int=32>				
	Time= 40.205 msec, Size= 131072000	Ops, WorkgroupSize= 1024 threads	s/block		
necking computed result fo	r correctness: Result = PASS				
OTE: The CUDA Samples are	not meant for performance measurem	ents. Results may vary when GPU B	Boost is enabled.		
27425== Profiling applica					
=27425=== Profiling result:					
27425== Metric result:					
wocations	Metric Name	Metric Description	on Min	Max	Av
evice "Tesla K40m (0)"					
	MulCUDA <int=32>(float*, float*, flo</int=32>				
301	flop_sp_efficiency	FLOP Efficiency(Peak Single	e) 7.88%	8.19%	8.07
	daSamples/NVIDIA CUDA-7.5 Samples/H				
nvprofmetrics flop sp		11/X00_04/ L1nux/ release [20:58:4			

### **Visual Profiler**



#### Your new favorite tool

💺 *NewSession1 🖾							
	0,3 s	0,35 s	0,4 s	0,45 s	0,5 s	0,55 s	0,6 s
Process "matrixMul" (18924)				·			
<ul> <li>Thread 39720768</li> </ul>							
Runtime API			cudaEventS	ynchronize			
Driver API							
Profiling Overhead							
<ul> <li>[0] Tesla K40m</li> </ul>							
<ul> <li>Context 1 (CUDA)</li> </ul>							
- T MemCpy (HtoD)							
- 🍸 MemCpy (DtoH)							
<ul> <li>Compute</li> </ul>							
- 🍸 100,0% void mat							
Streams							
Default							
			_				
🖬 Analysis 🗔 Details 📮 Console	🖾 📑 Settings		- X X	🗟 🔊 🕫 🛃		E Properties 🖾	- [
terminated> matrixMul on juhydra						Default	
[Matrix Multiply Using CUDA] - Si						▼ Duration	
SPU Device 0: "Tesla K40m" with o	compute capability	3.5				* Duration Session	
MatrixA(320,320), MatrixB(640,320	9)					Session	
Computing result using CUDA Kerne	i						
done Performance= 351.01 GFlop/s, Time	- 0 373 mean Siz	- 121972999 00	r WorkerounSize	- 1924 threads /h	lock		
Checking computed result for corr			s, workgroupsize	= 1024 cm euus/b	LOCK		



## **Pitfalls**

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## Pitfalls; Caveats; Tips



There are mistakes to be made, opportunities to be missed

- Try to use a library if possible; let others do the hard work
- Profile! Don't trust your gut!
- Gradually improve and specialize when porting and optimizing
- Expose enough parallelism! The GPU wants to be fed
- Express data locality
- Study your data transfers, can you reduce it?
- Unified Memory is a good start, but explicit transfers might be fast
- Use specialized memory: constant memory, shared memory! Pinned host memory is sometimes a very easy performance booster
- Overlap computation and transfer
- Does your code really need double precision? Is single precision sufficient? Or, maybe, even half precision?
- The number of threads and blocks is a tunable parameter; 128 is a good start

#### **Omitted** There's so much more!



What I did not talk about

- Atomic operations
- Shared memory
- Pinned memory
- How debugging works
- Overlapping streams
- Cross-compilation for heterogeneous systems



### **Summary & Conclusion**



- GPUs can improve your performance many-fold
- For a fitting, parallelizable application
- Libraries are easiest
- Direct programming (plain CUDA) is most powerful
- OpenACC is somewhere in between (and portable)
- There are many tools helping the programmer
- → Felice will surely give you more details in today's GPU tutorial!

### **Summary & Conclusion**



- GPUs can improve your performance many-fold
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#### Appendix Further Reading & Links Pascal Performances Glossary References



# Further Reading & Links



- A discussion of SIMD, SIMT, SMT by Y. Kreinin.
- NVIDIA's documentation: docs.nvidia.com
- NVIDIA's Parallel For All blog

### **Pascal Performance**



Tesla Products	Tesla K40	Tesla M40	Tesla P100
GPU	GK110 (Kepler)	GM200 (Maxwell)	GP100 (Pascal)
SMs	15	24	56
TPCs	15	24	28
FP32 CUDA Cores / SM	192	128	64
FP32 CUDA Cores / GPU	2880	3072	3584
FP64 CUDA Cores / SM	64	4	32
FP64 CUDA Cores / GPU	960	96	1792
Base Clock	745 MHz	948 MHz	1328 MHz
GPU Boost Clock	810/875 MHz	1114 MHz	1480 MHz
Peak FP32 GFLOPs1	5040	6840	10600
Peak FP64 GFLOPs <sup>1</sup>	1680	210	5300
Texture Units	240	192	224
Memory Interface	384-bit GDDR5	384-bit GDDR5	4096-bit HBM2
Memory Size	Up to 12 GB	Up to 24 GB	16 GB
L2 Cache Size	1536 KB	3072 KB	4096 KB
Register File Size / SM	256 KB	256 KB	256 KB
Register File Size / GPU	3840 KB	6144 KB	14336 KB
TDP	235 Watts	250 Watts	300 Watts
Transistors	7.1 billion	8 billion	15.3 billion
GPU Die Size	551 mm²	601 mm²	610 mm <sup>2</sup>
Manufacturing Process	28-nm	28-nm	16-nm FinFET

#### Figure: Tesla P100 performance characteristics in comparison [5]

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### **Glossary** I



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