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Alpaka Parallel Programming Library In a Nutshell











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alpaka – Abstraction Library for Parallel Kernel Acceleration

Alpaka is...

- A parallel programming library: Accelerate your code by exploiting your hardware's parallelism!
- An abstraction library independent of hardware ecosystem: Create portable code that runs on CPUs and GPUs!
- Free & open-source software





Problem of HPC Systems?

Heterogenous Hardware Ecosystem!

TOP500

- Frontier(USA) 1.194 Exaflop/s, AMD EPYC CPU + AMD Instict GPU
- Aurora(USA) 585 Petaflop/s, Intel Xeon CPU + Intel GPU Max
- Eagle(USA) 561 Petaflop/s, Intel Xeon CPU + NVIDIA H100 GPU
- Fugaku (Japan) 442 Petaflop/s, Fujitsu A64FX CPU
- Lumi (Finland) 380 Petaflop/s, AMD EPYC CPU + AMD Instinct GPU
- Leonardo (Italy) 238.7 Petaflop/s, Intel Xeon CPU + NVIDIA A100 GPU
- MareNostrum (Spain) 183.2 Petaflops/s, Intel Xeon CPU + **NVIDIA H100 GPÚ**
- Summit(USA) 148.3 Petaflops/s, IBM Power9 CPU + NVIDIA Volta **GPU**
- Eos(USA) 121.4 Petaflops/s, Intel Xeon CPU + NVIDIA H100 GPU
- 10 Sierra(USA) 94 Petaflops/s, IBM POWER9 CPU + NVIDIA Tesla V100 **GPU**



THE LIST

www.top500.or



11/2023 Highlights
The 62nd edition of the TOP500 shows five new or upgraded entries in the top 10 but the Frontier system still remains the only true exascale machine with an HPL score of 1.194 Exaflop/s.

The Frontier system at the Oak Ridge National Laboratory, Tennessee, USA remains the No. 1 system on the TOP500 and is still the only system reported with an HPL performance exceeding one Exaflop/s. Frontier brought the pole position back to the USA one year ago on the June 2022 listing and has since been remeasured with an HPL score of 1.194 Exaflop/s.

Frontier is based on the latest HPE Cray EX235a architecture and is equipped with AMD EPYC 64C 2GHz processors. The system has 8,699,904 total cores, a power efficiency rating of 52.59 gigaflops/watt, and relies on HPE's Slingshot 11 network for data transfer.

The Aurora system at the Argonne Leadership Computing Facility, Illinois, USA is currently being commissioned and will at full scale exceed Frontier with a peak performance of 2 Exaflop/s. It was submitted with a measurement on half of the final system achieving 585 Petaflop/s on the HPL benchmark which secured the No. 2 spot on the TOP500.

Aurora is built by Intel based on the HPE Cray EX - Intel Exascale Compute Blade which uses Intel Xeon CPU Max Series processors and Intel Data Center GPU Max Series accelerators which communicate through HPE's Slingshot-11 network interconnect.

The Eagle system installed in the Microsoft Azure cloud in the USA is newly listed as No. 3. This Microsoft NDv5 system is based on Intel Xeon Platinum 8480C processors and NVIDIA accelerators and achieved an HPL score of 561 Pflop/s.

read more »

List Statistics

Vendors System Share

- Frontier HPE Cray EX235a, AMD Optimized 3rd Generation EPYC 64C 2GHz, AMD Instinct MI250X. Slingshot-11, HPE
- Aurora HPE Cray EX Intel Exascale Compute Blade, Xeon CPU Max 9470 52C 2.4GHz, Intel Data Center GPU Max, Slingshot-11, Intel
- Eagle Microsoft NDv5. Xeon Platinum 8480C 48C 2GHz, NVIDIA H100, NVIDIA Infiniband NDR. Microsoft
- Supercomputer Fugaku Supercomputer Fugaku, A64FX 48C 2.2GHz. Tofu interconnect D, Fujitsu
- LUMI HPE Cray EX235a. AMD Optimized 3rd Generation EPYC 64C 2GHz. AMD Instinct MI250X. Slingshot-11, HPE
- Leonardo BullSequana XH2000, Xeon Platinum 8358 32C 2.6GHz, NVIDIA A100 SXM4 64 GB, Quad-rail NVIDIA HDR100 Infiniband, **EVIDEN**
- Summit IBM Power System AC922, IBM POWER9 22C 3.07GHz. NVIDIA Volta GV100. Dual-



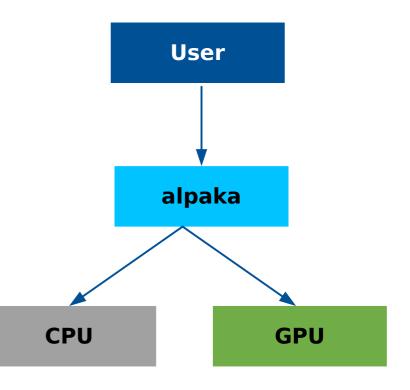


Without alpaka

 Hardware ecosystem is heterogenous, platforms are not inter-operable → parallel programs not easily portable

alpaka: one API to rule them all

- **Abstraction** (not hiding!) of the underlying hardware, compiler and OS platforms
 - No default device, default queue, built-in functions, language extentions
- Easy change of the backend
 - Code needs only minor adjustments to support different accelerators
- Direct usage of vendor APIs, not depend on "unified APIs"
 - GPU Backends: Hip backend uses HIP API, Cuda backend uses Cuda API, SYCL backend can be used for Intel GPUs
 - Cuda profilers and HIP profilers can be used to profile Alaka generated kernel code.
 - CPU Backends: OpenMp, Threads, TbbBlocks
- Easy indexing of threads in kernels
- Alpaka proposes and validates the type of parallelism
- **Heterogenous Programming**: Using different backends in a synchronized manner







Find us on GitHub!

Alpaka library: https://www.github.com/alpaka-group/alpaka

Full source code and many examples, Issue tracker

The documents: https://alpaka.readthedocs.io/en/latest/

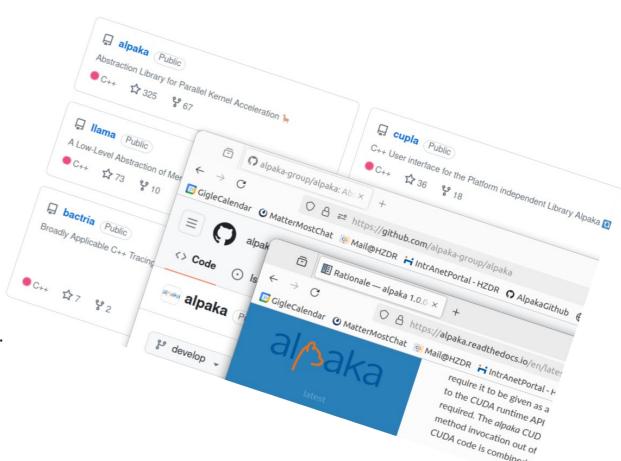
- Installation guide
- Cheatsheet
- Abstraction model and the rationale behind alpaka

Project group: https://www.github.com/alpaka-group

Contains all alpaka-related projects, documentation, samples, ...

alpaka is a free software (MPL 2.0)

moz://a Public License

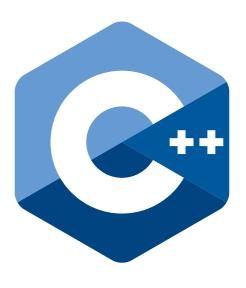






Programming with alpaka

- C++ only!
- Alpaka is written entirely in C++17. Coming soon: C++20.
- Header-only library. No additional runtime dependency. #include <alpaka/alpaka.hpp> is enough!
- Supports a wide range of modern C++ compilers (g++, clang++, Apple LLVM, MS VS)
- Portable across operating systems: Linux, macOS, Windows are supported







Installing alpaka [shorten]

- alpaka requires Boost, Cmake and a modern C++ compiler (q++, clanq++, Visual C++, ...)
- Linux:
 - sudo dnf install boost-devel (RPM)
- sudo apt install libboost-all-dev (DEB)
- brew install boost (Using Homebrew, https://brew.sh)
- sudo port install boost (Using MacPorts, https://macports.org)
- Windows: vcpkg install boost (Using vcpkg, https://github.com/microsoft/vcpkg)
- Depending on your target platform you may need additional packages
- NVIDIA ĞPUs: CUDA Toolkit (https://developer.nvidia.com/cuda-toolkit)
- AMD GPUs: ROCm / HIP (https://rocmdocs.amd.com/en/latest/index.html)
- CMake is the preferred system for building and installing
- Linux: sudo dnf install cmake (RPM) or sudo apt install cmake (DEB)
- macOS and Windows: Download the installer from https://cmake.org/download/
- Download alpaka: git clone -b develop https://github.com/alpaka-group/alpaka.git

Install alpaka's dependencies:

- In the terminal / PowerShell, switch to the downloaded alpaka directory:
 - cd /path/to/alpaka
- · Create a build directory and switch to it:
 - mkdir build
 - cd build
- In the build directory use CMake for configuration (replace the install prefix with an actual path) alpaka_ACC_CPU_B_OMP2_T_SEQ_EN cmake -DCMAKE INSTALL PREFIX=/some/other/path/ ...
- Alpaka determines the installed backends and makes the configuration of Cmake variables according
- Ccmake tool would show the cmake variables configured by the command above. You can change cmake ccmake ../
- Your alpaka installation will later reside in /some/other/path/
- If you are content with the default directories you can omit the install prefix
- This requires administrator privileges!
- This will result in a system-wide installation!
- On Linux and macOS the default directory will be /usr/local
- On Windows the default directory will be C:\Program Files
- Execute the installation command (note the dot):
 - cmake --install .
- You should now have a complete alpaka installation in the directory you chose earlier
- Did you encounter problems or errors?
- Check the documentation: https://alpaka.readthedocs.io/en/latest/install/instructions.html
- Still having problems? Please report them here: https://github.com/alpaka-group/alpaka/issues

```
BUILD TESTING
Boost ATOMIC LIBRARY RELEASE
                                 /usr/lib/x86 64-linux-qnu/libboost atomic.so.
Boost INCLUDE DIR
CMAKE BUILD TYPE
CMAKE CUDA ARCHITECTURES
CMAKE INSTALL PREFIX
FETCHCONTENT BASE DIR
FETCHCONTENT FULLY DISCONNECTE
FETCHCONTENT QUIET
FETCHCONTENT SOURCE DIR MDSPAN
FETCHCONTENT UPDATES DISCONNEC
FETCHCONTENT_UPDATES_DISCONNEC
MDSPAN_CXX_STANDARD
MDSPAN ENABLE BENCHMARKS
MDSPAN ENABLE COMP BENCH
MDSPAN ENABLE CONCEPTS
MDSPAN ENABLE CUDA
MDSPAN ENABLE EXAMPLES
MDSPAN ENABLE OPENMP
MDSPAN ENABLE TESTS
MDSPAN USE SYSTEM GTEST
RT LIBRARY
                                 /usr/lib/x86 64-linux-gnu/librt.a
TBB DIR
                                 /usr/lib/x86 64-linux-qnu/cmake/TBB
alpaka ACC CPU B SEQ T OMP2 EN
alpaka ACC CPU B SEQ T SEQ ENA
alpaka ACC CPU B SEQ T THREADS
alpaka ACC CPU B TBB T SEQ ENA
alpaka ACC CPU DISABLE ATOMIC
alpaka ACC GPU CUDA ENABLE
alpaka ACC GPU CUDA ONLY MODE
alpaka ACC GPU HIP ENABLE
alpaka ACC GPU HIP ONLY MODE
alpaka ACC SYCL ENABLE
alpaka ASSERT ACC ENABLE
alpaka BLOCK SHARED DYN MEMBER
alpaka BUILD BENCHMARKS
alpaka BUILD EXAMPLES
```





Accessing alpaka examples [shorten]

- There are many examples in the example directory! You can build them all at once from your build directory: cmake -Dalpaka BUILD EXAMPLES=ON -DCMAKE BUILD TYPE=Release ... cmake --build . --config Release
- Switch to the vectorAdd example (from the top-level source directory): cd example/vectorAdd mkdir build cd build
- If you installed alpaka to a standard location (= default CMAKE INSTALL PREFIX): cmake -DCMAKE BUILD TYPE=Release ...
- If you installed alpaka to a non-standard location: cmake -DCMAKE_BUILD_TYPE=Release -Dalpaka_R00T=/path/to/installed/alpaka ...
- Alternative: Set the CMAKE PREFIX PATH environment variable (note the missing alpaka):
- Linux / macOS: export CMAKE PREFIX PATH=/path/to/installed
- Windows: set CMAKE PREFIX PATH=C:\path\to\installed
- Switch to the vectorAdd example (from the top-level source directory):
 - cd example/vectorAdd
 - mkdir build
 - cd build
- If you installed alpaka to a standard location (= default CMAKE INSTALL PREFIX):
 - cmake -DCMAKE BUILD TYPE=Release ...
- If you installed alpaka to a non-standard location:
 - cmake -DCMAKE_BUILD_TYPE=Release -Dalpaka_R00T=/path/to/installed/alpaka ...
- · Alternative: Set the CMAKE PREFIX PATH environment variable (note the missing alpaka):
- Linux / macOS: export CMAKE PREFIX PATH=/path/to/installed
- Windows: set CMAKE PREFIX PATH=C:\path\to\installed
- Build the example (note the dot):
- cmake --build . --config Release
- If everything worked correctly you should see the executable somewhere in your build tree:
- Linux / macOS: CMakeCache.txt CMakeFiles cmake_install.cmake Makefile vectorAdd
- Windows: The executable should be in the Release subdirectory of build. cd to the Release directory.
- Execute the example: ./vectorAdd
- Expected output: Execution results correct!





Tenets of Thread-Parallel Programming

- **Massive number of threads.** Large number of threads should run the same code seperately! This code is called "Kernel".
- **Indexing of threads.** Each thread should work on a different data portion or do a specific task. Hence, each thread has an index, inside the kernel this index is accessable.
- **Grouping threads to map them to the cores.** Threads are grouped in blocks to map them to the cores to increase occupancy.

The warpsize of the GPUs and memory(shared memory, registers) usage of kernel affects the performance. Therefore grid of threads are devided into thread blocks.

Extent: A vector representing the sizes along each dimension. In 3d, a vector of 3 items {height,width,depth}

Dimensions: Set of dimension names. {X-dimension, Y-dimension, Z-dimension}

Number of Dimensions



Grid Block Extent: Vec{3} Block Thread Extent: Vec{4}

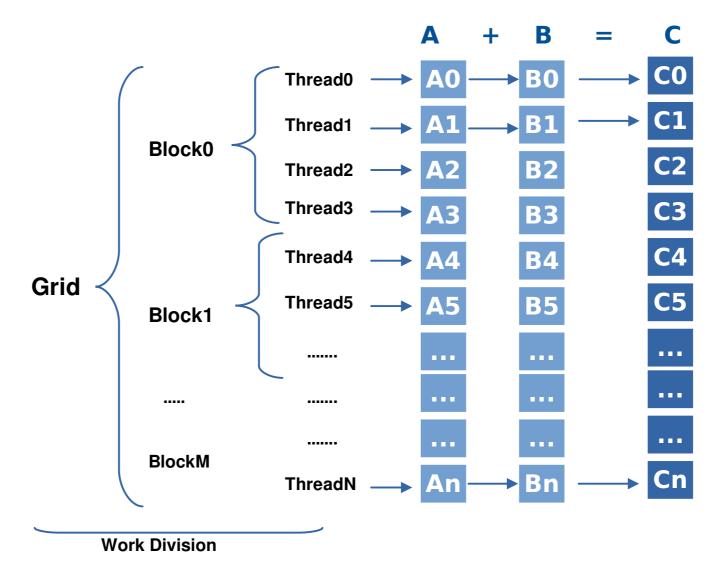


Grid Block Extent: Vec{1,3} Block Thread Extent: Vec{5,5}

Vector Addition Problem



- -Fix the number of data dimensions and the index type
- -Create A and B vectors on Host Use dynamic array alpaka::Buf
- -Create empty C vector on Host Use alpaka::Buf
- -Copy A and B Vectors to Device Use alpaka::memcpy
- -Determine the Paralelism: Work Division. Extents of threads per block, blocks per grids, elements per thread.
- -Create the Queue Use alpaka::Queue
- -Execute the kernel on device with the workdiv on the queue. Pass the A,B,C vectors to the executer. Use alpaka::exec



Parallel vector addition code



- Define kernel
- Select the backend depending on your system, the accelerator!
- Select the device (e.g your specific GPU)
- Decide on proper WorkDivision
- Create input vectors A and B and result vector C on host (e.g CPU)
- •Create input vectors A and B and result vector C on device (e.g. CPU)
- Copy the vectors to the device memory
- Run kernel
- Copy the result vector back to the host vector

```
alpaka::WorkDivMembers<Dim, Idx> const workDiv = alpaka::getValidWorkDiv<Acc>(*
                                                                                                       elementsPerThread,
                                                                                                       false, alpaka::GridBlockExtentSubDivRestrictions::Unrestricted);
class VectorAddKernel
                                                                                                    // Get the host device for allocating memory on the host.
                                                                                                    auto const platformHost = alpaka::PlatformCpu{};
   ALPAKA_NO_HOST_ACC_WARNING
                                                                                                    auto const devHost = alpaka::getDevByIdx(platformHost, 0);
       template<typename <a href="TAcc">TAcc</a>, typename <a href="TIdx">TIdx</a>>
       ALPAKA_FN_ACC auto operator()(
           TAcc const& acc, // the accelerator
                                                                                                    using DevHost = alpaka::DevCpu;
           TElem const* const A,
           TElem const* const B,
                                                                                                    using BufHost = alpaka::Buf<DevHost, DataType, Dim, Idx>;
           TElem* const C,
                                                                                                    BufHost bufHostA(alpaka::allocBuf<DataType, Idx>(devHost, extent));
           TIdx const& numElements) const -> void
                                                                                                    BufHost bufHostB(alpaka::allocBuf<DataType, Idx>(devHost, extent));
                                                                                                    BufHost bufHostC(alpaka::allocBuf<DataType, Idx>(devHost, extent)); **
       TIdx const gridThreadIdx(alpaka::getIdx<alpaka::Grid, alpaka::Threads>(acc)[0u]);
                                                                                                   for(Idx i(0); i < numElements; ++i)</pre>
                                                                                                   { bufHostA[i] = randomA; bufHostB[i] = randomB; bufHostC[i] = 0; }*
       if(gridThreadIdx < numElements)</pre>
           // Use thread index as the data index
                                                                                                    using BufAcc = alpaka::Buf<DevAcc, DataType, Dim, Idx>; ?
           C[gridThreadIdx] = A[gridThreadIdx] + B[gridThreadIdx];
                                                                                                    BufAcc bufAccA(alpaka::allocBuf<DataType, Idx>(devAcc, extent));
                                                                                                    BufAcc bufAccB(alpaka::allocBuf<DataType, Idx>(devAcc, extent));

                                                                                                    BufAcc bufAccC(alpaka::allocBuf<DataType, Idx>(devAcc, extent));
guto main() -> int
                                                                                                    alpaka::Queue<Acc, alpaka::Blocking> queue(devAcc);
   using Dim = alpaka::DimInt<lu>; // Define the index domain ?
   using Idx = std::size_t; // Index type of the threads and buffers
                                                                                                    alpaka::memcpy(queue, bufAccA, bufHostA);
   using DataType = std::uint32_t; // Define the buffer element type
                                                                                                    alpaka::memcpy(queue, bufAccB, bufHostB);
                                                                                                    alpaka::memcpy(queue, bufAccC, bufHostC);
   using Acc = alpaka::AccGpuCudaRt<Dim, Idx>;
                                                                                                    VectorAddKernel kernel;
   using DevAcc = alpaka::Dev<Acc>;
   // Select a device from platform of Acc
   auto const platform = alpaka::Platform<Acc>{};
                                                                                                       kernel, alpaka::getPtrNative(bufAccA), alpaka::getPtrNative(bufAccB),
   quto const devAcc = alpaka::getDevByIdx(platform, 0);
                                                                                                       numElements);
   // Define the work division depending on the data
   Idx const numElements(100000);
                                                                                                    alpaka::memcpy(queue, bufHostC, bufAccC); // bufHostC includes the result!
   Idx const elementsPerThread(1u);
                                                                                                    alpaka::wait(queue);
   alpaka::Vec<Dim, Idx> const extent(numElements);
```





Define the Alpaka Kernel

- Contains the algorithm
- Written on per-data-element basis
- It allows access to the thread index directly
- alpaka Kernels are functors (function-like C++ structs / classes)
- operator() is annotated with ALPAKA FN ACC specifier
- operator() must return void
- operator() must be const
- Arguments can be pointers and trivially copyable types
- Agnostic to device details
- Code piece that is run by each thread

```
// Single header library
#include <alpaka/alpaka.hpp>
#include <iostream>
class VectorAddKernel
    ALPAKA_NO_HOST_ACC_WARNING
    template<typename TAcc, typename TElem, typename TIdx>
    ALPAKA_FN_ACC auto operator()(
        TAcc const& acc, // the accelerator
        TElem const* const A, TElem const* const B, TElem* const C,
        TIdx const& numElements ) const -> void
        static_assert(alpaka::Dim<TAcc>::value == 1, "The kernel expects 1-dimensional indices!");
        // Get thread index
        TIdx const gridThreadIdx(alpaka::getIdx<alpaka::Grid, alpaka::Threads>(acc)[0u]);
        if(gridThreadIdx < numElements)</pre>
                // Use thread index as the data index
                C[gridThreadIdx] = A[gridThreadIdx] + B[gridThreadIdx];
```





Obtaining the indices and extents in Kernel

- alpaka provides API functions for obtaining indices
 - Index of Thread on the Grid: alpaka::getIdx<alpaka::Grid, alpaka::Threads>(acc)[dim]; Index of Thread on a Block: alpaka::getIdx<alpaka::Block, alpaka::Threads>(acc)[dim];
 - Index of Block on the Grid: alpaka::getIdx<alpaka::Grid, alpaka::Blocks>(acc)[dim];
- alpaka provides API functions for obtaining extents of the grid and the blocks
 - Number of Threads on the Grid: alpaka::getWorkDiv<alpaka::Grid, alpaka::Threads>(acc)[dim];
 - Number of Threads on a Block: alpaka::getWorkDiv<alpaka::Block, alpaka:Threads>(acc)[dim];
 - Number of Blocks on the Grid: alpaka::getWorkDiv<alpaka::Grid, alpaka::Blocks>(acc)[dim];

al/aka in a Nutshell: Accelerator



Select the Accelerator

- alpaka provides a number of pre-defined Accelerators.
- Accelerator is a type that is only instantiated on Device not on the host.
- Accelerator is an abstraction that makes the kernel 'device agnostic'. Gives kernel means the access thread index. Responsible for the thread state, memory management, and synchronization ??)
- For GPUs:
 - AccGpuCudaRt for NVIDIA GPUs
 - AccGpuHipRt for AMD, Intel and NVIDIA GPUs
- For CPUs
 - AccCpuFibers based on Boost.fiber
 - AccCpu0mp2Blocks based on OpenMP 2.x
 - AccCpu0mp4 based on OpenMP 4.x
 - AccCpuTbbBlocks based on TBB
 - AccCpuThreads based on std::thread

```
// Easy change of the backend type!
// Example: CUDA GPU accelerator
using Acc = alpaka::AccGpuCudaRt<Dim, Idx>;
// Switch to AMD: HIP GPU accelerator
using Acc = alpaka::AccGpuHipRt<Dim, Idx>;
// Switch to CPU Omp: CPU accelerator
using Acc = alpaka::AccCpuOmp2Blocks<Dim, Idx>;
```

```
auto main() -> int
   using Dim = alpaka::DimInt<1u>; // Define the index domain ?
   using Idx = std::size_t; // Index type of the threads and buffers
   using DataType = std::uint32_t; // Define the buffer element type
  // Define the accelerator: AccGpuCudaRt, AccGpuHipRt,
   // AccCpuThreads, AccCpuOmp2Threads, AccCpuOmp2Blocks, AccCpuTbbBlocks AccCpuSerial
   using Acc = alpaka::AccGpuCudaRt<Dim, Idx>;
   using DevAcc = alpaka::Dev<Acc>;
   // Select a device from platform of Acc
   quto const platform = alpaka::Platform<Acc>{};
   quto const devAcc = alpaka::getDevByIdx(platform, 0);
   Idx const numElements(100000);
   Idx const elementsPerThread(1u);
   alpaka::Vec<Dim, Idx> const extent(numElements);
   // Let alpaka calculate good block and grid sizes given our full problem extent
   alpaka::WorkDivMembers<Dim, Idx> const workDiv = alpaka::getValidWorkDiv<Acc>(*
       devAcc, // device
       extent, // {length, height, depth} of grid. For 1D only legth of the vector!
       elementsPerThread,
       false, alpaka::GridBlockExtentSubDivRestrictions::Unrestricted);
   // Get the host device for allocating memory on the host.
   quto const platformHost = alpaka::PlatformCpu{};
   auto const devHost = alpaka::getDevByIdx(platformHost, 0);
   using DevHost = alpaka::DevCpu;
   using BufHost = alpaka::Buf<DevHost, DataType, Dim, Idx>;
   VectorAddKernel kernel;
   alpaka::exec<Acc>( // Run the kernel execution task
```



Get device instance

alpaka Devices

- Each alpaka Device instance represents a single physical device
- Determined by programmer's Accelerator choice
- Easy management of physical devices
- Contains device information.

```
using Acc = alpaka::AccGpuCudaRt<Dim, Idx>;
using DevAcc = alpaka::Dev<Acc>; ?
// Select a device
auto const platform = alpaka::Platform<Acc>{};*
auto const devAcc = alpaka::getDevByIdx(platform, 0);
auto const name = alpaka::getName(devAcc);
                                                  // Back-end-defined device name
auto const bytes = alpaka::getMemBytes(devAcc); // Size of device memory
auto const free = alpaka::getFreeMemBytes(devAcc); // Size of available device memory
alpaka::reset(devAcc);
```

```
// Select a device
auto const platform = alpaka::Platform<Acc>{}; ?
auto const devAcc = alpaka::getDevByIdx(platform, 0);
// Define the work division depending on the data
Idx const numElements(100000);
Idx const elementsPerThread(lu);
alpaka::Vec<Dim, Idx> const extent(numElements);
// Let alpaka calculate good block and grid sizes given our full problem extent
alpaka::WorkDivMembers<Dim, Idx> const workDiv = alpaka::getValidWorkDiv<Acc>(*
    devAcc, // device
    extent, // {length, height, depth} of grid. For 1D only legth of the vector!
    elementsPerThread, false, alpaka::GridBlockExtentSubDivRestrictions::Unrestricted);
// Get the host device for allocating memory on the host.
using DevHost = alpaka::DevCpu;
auto const platformHost = alpaka::PlatformCpu{};
auto const devHost = alpaka::getDevByIdx(platformHost, 0);
using BufHost = alpaka::Buf<DevHost, Data, Dim, Idx>;
BufHost bufHostA(alpaka::allocBuf<Data, Idx>(devHost, extent));
BufHost bufHostB(alpaka::allocBuf<Data, Idx>(devHost, extent));
BufHost bufHostC(alpaka::allocBuf<Data, Idx>(devHost, extent));
for(Idx i(0); i < numElements; ++i)</pre>
    bufHostA[i] = randomA; bufHostB[i] = randomB; bufHostC[i] = 0;
// Allocate 3 buffers on the accelerator
using BufAcc = alpaka::Buf<DevAcc, Data, Dim, Idx>; ?
BufAcc bufAccA(alpaka::allocBuf<Data, Idx>(devAcc, extent));
BufAcc bufAccB(alpaka::allocBuf<Data, Idx>(devAcc, extent));
•
BufAcc bufAccC(alpaka::allocBuf<Data, Idx>(devAcc, extent)); **
// choose between Blocking and NonBlocking
using QueueProperty = alpaka::Blocking;
alpaka::Queue<Acc, QueueProperty> queue(devAcc); ?
```

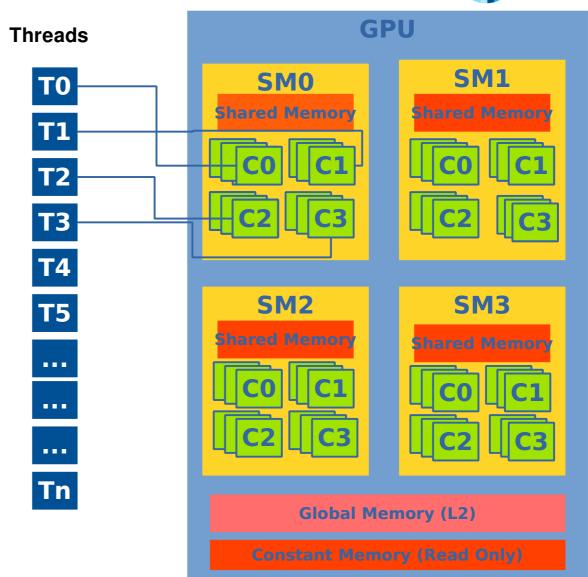




Determine the WorkDivision-I

How to distribute threads among SMs?

- Main determinants of mapping threads to the SMs:
 - Number of cores SM,
 - Warpsize,
 - Register and local memory (lm) usage of each thread,
 - Shared memory usage of each thread,
 - Threads per SM, Threads per Block, Blocks per Device
- Memory latencies: Global Memory and Constant memory has different latencies.
- Memory sizes: Size of shared memory used by threads in a block or blocks assigned to an SM







Determine the workdivision-II

- WorkDivision data structure consists 3 vectors:
 - Grid block extent.

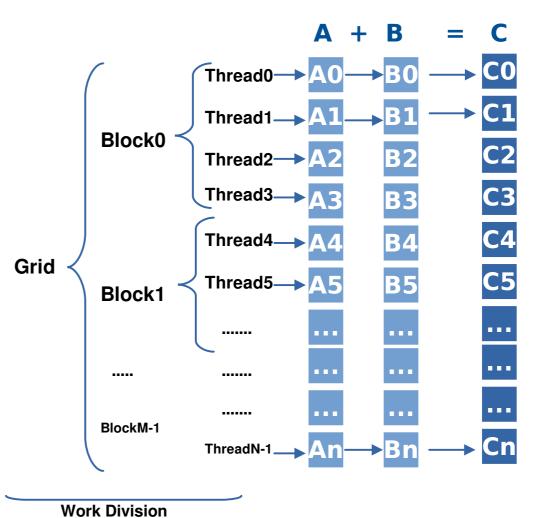
Vec{M} or Vec{1, 1, M} depending on the number of dimensions.

Block thread extent.

 $Vec{4}$ or $Vec{1,1,4}$.

- Elements per thread
- Setting work-div manually

```
using Dim1D = alpaka::DimInt<1>;//Set number of dims to 1
using Vec1D = alpaka::Vec<Dim1D, Idx>;//Define alias
auto workDiv1D = alpaka::WorkDivMembers(Vec1D{M}, Vec1D{4u}, Vec1D{1u});
// alternatively
using Dim3D = alpaka::DimInt<3>;//Set number of dims to 3
using Vec3D = alpaka::Vec<Dim3D, Idx>; //Define alias
auto workDiv3D = alpaka::WorkDivMembers(Vec3D{1,1,M}, Vec3D{1,1,4u}, Vec3D{1,1,1u});
```







1D WorkDivision!

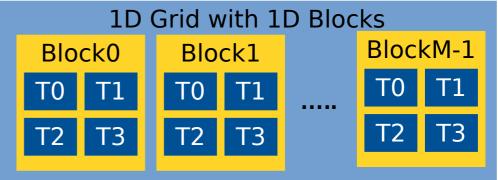
Grid Block Extent 1D vector = Vec{M} Block Thread Extent 1D vector = Vec{4u} Thread Elem Extent 1D vector = Vec{1u} Grid Thread Extent 1D vector =Vec{4*M} or Vec{N}

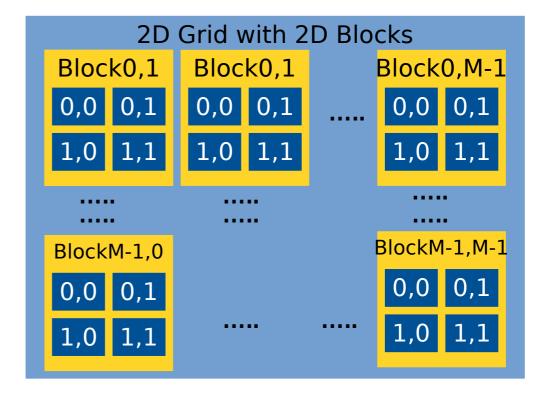
1D WorkDivision {{M},{4u},{1u}} // if Dim is 3 then fill with 1u 1D WorkDivision {{1u,1u,M},{1u,1u,4},{1u,1u,1u}}

2D WorkDivision!

Grid Block Extent 2D vector = Vec{M,M} Block Thread Extent 2D vector = Vec{2u,2u} Thread Elem Extent 2D vector = Vec{1u,1u} Grid Thread Extent 2D vector = Vec{4u*M,4u*M} or Vec{N,N}

2D WorkDivision {{M,M},{2u,2u},{1u,1u}} 2D WorkDivision {{1u,M,M},{1u,2u,2u},{1u,1u,1u}}









Determine the WorkDivision-III

- Use getValidWorkDiv function
- Inputs are
 - Full grid-thread extent. Hence we guarantee total number of threads needed.
 - Elements per thread extent
- getValidWorkDiv generates devides the full grid-thread extent to blocks of given dimension.

```
auto const platform = alpaka::Platform<Acc>{};
auto const devAcc = alpaka::getDevByIdx(platform, 0);
// Define the work division depending on the data
Idx const numElements(100000);
Idx const elementsPerThread(1u);
alpaka::Vec<Dim, Idx> const extent(numElements);
// Let alpaka calculate good block and grid sizes given our full problem extent
alpaka::WorkDivMembers<Dim, Idx> const workDiv = alpaka::getValidWorkDiv<Acc>(*
    devAcc, // device
    elementsPerThread,
    false, alpaka::GridBlockExtentSubDivRestrictions::Unrestricted);
VectorAddKernel kernel;
alpaka::exec<Acc>( // Run the kernel execution task
    queue,
    workDiv,
   kernel, alpaka::getPtrNative(bufAccA), alpaka::getPtrNative(bufAccB),
    alpaka::getPtrNative(bufAccC),
    numElements);
alpaka::memcpy(queue, bufHostC, bufAccC); // bufHostC includes the result!
alpaka::wait(queue);
```



Allocate data vectors on host and device. Copy data vectors to the Device

- alpaka::Buf is multi dimensional dynamic array for all devices, copyable among host and accelerator devices.
- alpaka::allocBuf allocates memory to the given device.
- alpaka::memcpy copies the data from one buffer to another. User should not have to specify which buffer resides on which device.
- alpaka::Buf is device-type bound. (Host device or Acc device)
- Memory copy needs to be synchronized , therefore needs a queue

```
// Define the work division depending on the data
 Idx const numElements(100000);
 Idx const elementsPerThread(lu);
 alpaka::Vec<<u>Dim</u>, Idx> const extent(numElements);
 // Let alpaka calculate good block and grid sizes given our full problem extent
 alpaka::WorkDivMembers<Dim, Idx> const workDiv = alpaka::getValidWorkDiv<Acc>(*
    extent, // {length, height, depth} of grid. For 1D only legth of the vector!
    elementsPerThread, false, alpaka::GridBlockExtentSubDivRestrictions::Unrestricted);
 // Get the host device for allocating memory on the host.
 auto const platformHost = alpaka::PlatformCpu{};
 auto const devHost = alpaka::getDevByIdx(platformHost, 0);
// Host device type is needed, because it is not known (for the backend it is known in Acc)
 using DevHost = alpaka::DevCpu;
 using BufHost = alpaka::Buf<DevHost, DataType, Dim, Idx>;
 BufHost bufHostA(alpaka::allocBuf<DataType, Idx>(devHost, extent));
 BufHost bufHostB(alpaka::allocBuf<DataType, Idx>(devHost, extent));*
 BufHost bufHostC(alpaka::allocBuf<DataType, Idx>(devHost, extent));*
// Fill the buffers
for(Idx i(0); i < numElements; ++i)</pre>
{ bufHostA[i] = randomA; bufHostB[i] = randomB; bufHostC[i] = 0; }♥
using BufAcc = alpaka::Buf<DevAcc, DataType, Dim, Idx>; ?
 BufAcc bufAccA(alpaka::allocBuf<DataType, Idx>(devAcc, extent));
 BufAcc bufAccB(alpaka::allocBuf<DataType, Idx>(devAcc, extent));*
 BufAcc bufAccC(alpaka::allocBuf<DataType, Idx>(devAcc, extent)); **
 // Create a queue on the device, define the synchronization behaviour
 alpaka::Queue<Acc, alpaka::Blocking> queue(devAcc);
 alpaka::memcpy(queue, bufAccA, bufHostA);
 alpaka::memcpy(queue, bufAccB, bufHostB);
 alpaka::memcpy(queue, bufAccC, bufHostC);
```



Create the Queue for memcpy and kernel executions

- Alpaka::Queue provides communication between the host and the device
- Queue of tasks on the device
- Queue is always FIFO
- Create alpaka::Queue using the device instance and the accelerator type (e.g GPU)
- Two queue types: blocking and non-blocking
- Blocking queues block the Host until Deviceside command returns.
- Non-blocking queues return control to Host immediately, Device-side command runs asynchronously



```
using BufHost = alpaka::Buf<DevHost, DataType, Dim, Idx>;
 BufHost bufHostA(alpaka::allocBuf<DataType, Idx>(devHost, extent));
 BufHost bufHostB(alpaka::allocBuf<DataType, Idx>(devHost, extent)); **
 BufHost bufHostC(alpaka::allocBuf<DataType, Idx>(devHost, extent)); **
for(Idx i(0); i < numElements; ++i)</pre>
{ bufHostA[i] = randomA; bufHostB[i] = randomB; bufHostC[i] = 0; } •
// Allocate 3 buffers on the accelerator
using BufAcc = alpaka::Buf<DevAcc, DataType, Dim, Idx>; **
 BufAcc bufAccA(alpaka::allocBuf<DataType, Idx>(devAcc, extent));
 BufAcc bufAccB(alpaka::allocBuf<DataType, Idx>(devAcc, extent));*
 BufAcc bufAccC(alpaka::allocBuf<DataType, Idx>(devAcc, extent)); **
 // Create a queue on the device, define the synchronization behaviour
 alpaka::Queue<Acc, alpaka::Blocking> queue(devAcc);
 // Copy from Host to Acc
 alpaka::memcpy(queue, bufAccA, bufHostA);
 alpaka::memcpy(queue, bufAccB, bufHostB);
 alpaka::memcpy(queue, bufAccC, bufHostC);
 VectorAddKernel kernel;
 alpaka::exec<Acc>( // Run the kernel execution task
     queue,
     workDiv,
     kernel, alpaka::getPtrNative(bufAccA), alpaka::getPtrNative(bufAccB),
     alpaka::getPtrNative(bufAccC),
     numElements);
 alpaka::memcpy(queue, bufHostC, bufAccC); // bufHostC includes the result!
 alpaka::wait(queue);
 // bufHostC is filled with the result!
```



Execute the kernel using the workdiv, copy result back to the Host

- alpaka::exec function is used
- Result is stored in an alpaka::Buf
- Copy the result in device to the host
- If the queue not blocking then we can use **alpaka::wait** on the queue for sychronization.



```
// Allocate 3 buffers on the accelerator
using BufAcc = alpaka::Buf<DevAcc, DataType, Dim, Idx>; ?
BufAcc bufAccA(alpaka::allocBuf<DataType, Idx>(devAcc, extent));
BufAcc bufAccB(alpaka::allocBuf<DataType, Idx>(devAcc, extent)); **
BufAcc bufAccC(alpaka::allocBuf<DataType, Idx>(devAcc, extent));
// Create a queue on the device, define the synchronization behaviour
alpaka::Queue<Acc, alpaka::Blocking> queue(devAcc);
// Copy from Host to Acc
alpaka::memcpy(queue, bufAccA, bufHostA);
alpaka::memcpy(queue, bufAccB, bufHostB);
alpaka::memcpy(queue, bufAccC, bufHostC);
// Instantiate the kernel function object
VectorAddKernel kernel;
alpaka::exec<Acc>( // Run the kernel execution task
     queue,
    workDiv,
    kernel.
    alpaka::getPtrNative(bufAccA),
    alpaka::getPtrNative(bufAccB),
    alpaka::getPtrNative(bufAccC),
    numElements);
// Copy back the result
alpaka::memcpy(queue, bufHostC, bufAccC); // bufHostC includes the result!
alpaka::wait(queue);
```





On Tasks and Events

- Device-side operations (kernels, memory operations) are called Tasks
- Tasks on the same queue are executed in order (FIFO principle)

```
alpaka::enqueue(queueA, task1);
alpaka::enqueue(queueA, task2); // task2 starts after task1 has finished
```

- Order of tasks in different queues is unspecified
 - alpaka::enqueue(queueA, task1); alpaka::enqueue(queueB, task2); // task2 starts before, after or in parallel to task1
- For easier synchronization, alpaka Events can be inserted before, after or between Tasks:

```
auto myEvent = alpaka::Event<alpaka::Queue>(myDev);
alpaka::enqueue(queueA, myEvent);
alpaka::wait(queueB, myEvent); // queueB will only resume after queueA reached
myEvent
```





Programming Heterogeneous Systems-I

Using multiple Platforms Synchronously

- Alpaka enables easy heterogeneous programming!
- Create one Accelerator per back-end
- Acquire at least one Device per Accelerator
- Create one Queue per Device

```
// Define Accelerators
using AccCpu = alpaka::AccCpu0mp2Blocks<Dim, Idx>;
using AccGpu = alpaka::AccGpuCudaRt<Dim, Idx>;
// Acquire Devices
auto devCpu = alpaka::getDevByIdx<AccCpu>(0u);
auto devGpu = alpaka::getDevByIdx<AccGpu>(0u);
// Create Oueues
using QueueProperty = alpaka::NonBlocking;
using QueueCpu = alpaka::Queue<AccCpu, QueueProperty>;
using QueueGpu = alpaka::Queue<AccGpu, QueueProperty>;
auto queueCpu = QueueCpu{devCpu};
auto queueGpu = QueueGpu{devGpu};
```





Programming Heterogeneous Systems-II

Communication by Buffers

- Buffers are defined and created per Device
- Buffers can be copied between different Devices / Queues
- Not restricted to a single platform!
- **Restriction**: CPU to GPU copies (and vice versa) require GPU queue

```
// Allocate buffers
auto bufCpu = alpaka::allocBuf<float, Idx>(devCpu, extent);
auto bufGpu = alpaka::allocBuf<float, Idx>(devGpu, extent);
/* Initialization ... */
// Copy buffer from CPU to GPU - destination comes first
alpaka::memcopy(gpuQueue, bufGpu, bufCpu, extent);
// Execute GPU kernel
alpaka::enqueue(qpuQueue, someKernelTask);
// Copy results back to CPU and wait for completion
alpaka::memcopy(qpuQueue, bufCpu, bufGpu, extent);
// Wait for GPU, then execute CPU kernel
alpaka::wait(cpuQueue, gpuQueue);
alpaka::enqueue(cpuQueue, anotherKernelTask);
```





Tips

- If you want to pass pointers of views generated from data with numdim 2d or larger, you need to consider pitch
- If there are unused number of dimentions in workdiv use 1 for that dimension
- CPU to GPU copies require GPU Queue
- 2 queues synchorinization





Summary of Alpaka Structures

- Accelerator provides abstract view of all capable physical devices
- Device represents a single physical device
- Queue enables communication between the host and a single Device
- Platform is a union of Accelerator, Device and Kernel
- **Task** is a device-side operation (e.g kernel, memory operation)
- Others: **Event, Buffer** (dynamic array), **Vector** (static array)
- **Question**: How is portability between back-ends achieved?





[ToBeChanged] CUPLA: **Converting CUDA code to Alpaka**

- Change the suffix *.cu of the CUDA source files to *.cpp
- Remove cuda specific includes on top of your header and source files
- Add #include <cuda to cupla.hpp> remove #include <cuda runtime.h>
- Tansform kernels (global functions) to functors
- The functor's operator() must be qualified as const
- . Add the function prefix ALPAKA FN ACC to the operator() const
- Add as first (templated) kernel parameter the accelerator with the name acc (it is important that the accelerator is named acc because all cupla-to-alpaka replacements use the variable acc)
- If the kernel calls other functions you must pass the accelerator acc to each call
- Add the qualifier const to each parameter which is not changed inside the kernel

CUDA kernel

```
template< int blockSize >
 _global__ void fooKernel( int * ptr, float value )
```

Cupla kernel

```
template< int blockSize >
struct fooKernel
 template< typename T Acc >
 ALPAKA FN ACC
 void operator()( T_Acc const & acc, int * const ptr, float const value) const
```

CUDA kernel call at host

```
dim3 gridSize(42,1,1);
            dim3 blockSize(256,1,1);
            // extern shared memory and stream is optional
            fooKernel< 16 ><<< gridSize, blockSize, 0, 0 >>>( ptr. 23 ):
CUPLA kernel call at host:
            dim3 gridSize(42,1,1);
            dim3 blockSize(256,1,1);
            // extern shared memory and stream is optional
            CUPLA KERNEL(fooKernel< 16 >)( gridSize, blockSize, 0, 0)( ptr, 23);
```

CUDA shared memory (in kernel)

shared int foo;

```
shared int fooCArray[32];
            __shared__ int fooCArray2D[4][32];
            // extern shared memory (size was defined during the host side kernel call)
            extern shared float fooPtr[];
            int bar = fooCArray2D[ threadIdx.x ][ 0 ];
CUPLA shared memory (in kernel)
            sharedMem( foo, int );
            sharedMem(fooCArray, cupla::Array<int, 32 > );
            sharedMem( fooCArray, cupla::Array< cupla::Array< int, 4 >, 32 > );
            sharedMemExtern( fooPtr, float );
            int bar = fooCArray2D[ threadIdx.x ][ 0 ]
```

Use ALPAKA FN ACC in device function definitions and add an acc

(Note that to be exact the acc parameter is only necessary when alpaka functions like blockldx or atomicMax etc are used.)

CUDA Device Function

```
template< typename T Elem >
  device int deviceFunction( T Elem x )
CUPLA
template< typename T Acc, typename T Elem >
ALPAKA_FN_ACC int deviceFunction( T_Acc const & acc, T_Elem x )
```

CUDA Device Function Call

```
auto result = deviceFunction(x);
CUPLA
     auto result = deviceFunction( acc, x );
```





How to start using Alpaka

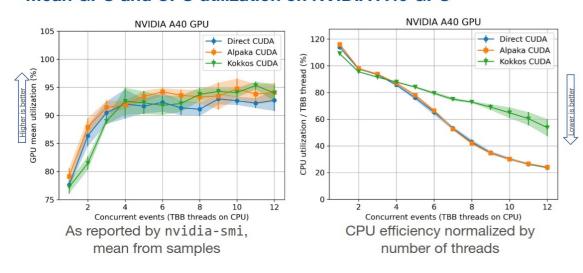
- Use Alpaka directly using examples and the cheat-sheet
- Don't write code initially on Cuda because alpaka is already low level, like Cuda.
- If you have a codebase in cuda, converting to Cupla can be a fast solution to benefit from alpaka features!



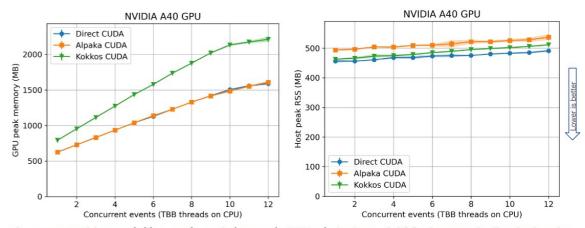


Alpaka Performance

Mean GPU and CPU utilization on NVIDIA A40 GPU



Peak memory usage on NVIDIA A40 GPU



As reported by nvidia-smi and /proc/<PID>/status. A100 shows similar behavior.

Evaluating Performance Portability with the CMS Source: **Heterogeneous Pixel Reconstruction code**

N. Andriotis¹, A. Bocci², E. Cano², L. Cappelli³, M. Dewing⁴, T. Di Pilato^{5,6}, J. Esseiva⁷, L. Ferragina⁸, G. Hugo², M. Kortelainen⁹, M. Kwok⁹, J. J. Olivera Loyola¹⁰, F. Pantaleo², A. Perego¹¹, W. Redjeb^{2,12} ¹BSC ²CERN ³INFN Bologna ⁴ANL ⁵CASUS ⁶University of Geneva ⁷LBNL ⁸University of Bologna ⁹FNAL ¹⁰ITESM ¹¹University of Milano Bicocca ¹² RWTH

CHEP 2023

https://indico.jlab.org/event/459/contributions/11824/attachments/9281/14171/20230511-CHEaP23 CMSPortability.pdf



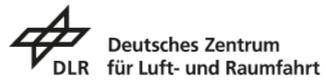


Community and Long Term Support

Partners using and contributing to Alpaka









• Alpaka is a part of Helmholtz Roadmap 2027-2034





Alpaka in the wild...

PIConGPU: https://github.com/ComputationalRadiationPhysics/picongpu

- Fully relativistic, manycore, 3D3V particle-in-cell (PIC) code
- Central algorithm in plasma physics
- Scalable to more than 18,000 GPUs
- Developed at Helmholtz-Zentrum Dresden-Rossendorf







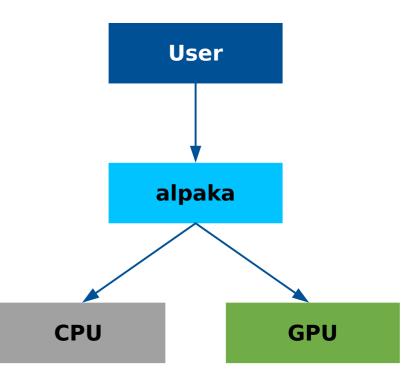
As a summary

Without alpaka

- Multiple hardware types are available from different vendors (CPUs, GPUs, ...)
- Increasingly heterogeneous hardware configurations available
- Platforms not inter-operable → parallel programs not easily portable

alpaka: one API to rule them all

- **Abstraction** (not hiding!) of the underlying hardware & software platforms
 - AMD, Nvidia, Intel GPUs, Different CPU parallelisations like TbbBlocks, OpenMP, Threads
- Easy change of the backend
 - Code needs only minor adjustments to support different accelerators
- Easy indexing of threads in kernels
- Easy setup of the type of parallelism by WorkDivision (Block sizes in grid, Thread sizes in block
- Heterogenous Programming: Using different backends in a synchronized manner







Thank you! If you use alpaka for your research, please cite one of the following publications:

Matthes A., Widera R., Zenker E., Worpitz B., Huebl A., Bussmann M. (2017): Tuning and Optimization for a Variety of Many-Core Architectures Without Changing a Single Line of Implementation Code Using the Alpaká Library. In: Kunkel J., Yokota R., Taufer M., Shalf J. (eds) High Performance Computing. ISC High Performance 2017. Lecture Notes in Computer Science, vol 10524. Springer, Cham, DOI: 10.1007/978-3-319-67630-2 36

E. Zenker et al., "Alpaka - An Abstraction Library for Parallel Kernel Acceleration", 2016 IEEE International Parallel and Distributed Processing Symposium Workshops (IPDPSW), Chicago, IL, 2016, pp. 631 – 640, DOI: 10.1109/IPDPSW.2016.50.

Worpitz, B. (2015, September 28). Investigating performance portability of a highly scalable particle-in-cell simulation code on various multi-core architectures. Zenodo. DOI: 10.5281/zenodo.49768.

al Nutshell



Appendix





[Appendix1] WorkDivision Example

- Determines the number of kernel instantiations
- Determines the type of parallelism
 - Dimensions of a grid in terms of blocks,
 - Dimensions of a block in terms of threads
 - Elements per thread

```
// Define the work division
// The workdiv is divided in three levels of parallelization:
// - grid-blocks:
                      The number of blocks in the grid
                      The number of threads per block (parallel, synchronizable).
// - thread-elements: The number of elements per thread (sequential, not synchronizable)
                       Each kernel has to execute its elements sequentially.
using Vec = alpaka::Vec<Dim, Idx>;
auto const elementsPerThread = Vec::all(static_cast<Idx>(1));
auto const threadsPerGrid = Vec{4, 2, 4};
using WorkDiv = alpaka::WorkDivMembers<Dim, Idx>;
WorkDiv const workDiv = alpaka::getValidWorkDiv<Acc>( devAcc, threadsPerGrid,
    elementsPerThread, false, alpaka::GridBlockExtentSubDivRestrictions::Unrestricted);
// Instantiate the kernel function object
HelloWorldKernel helloWorldKernel;
// Run the kernel
// To execute the kernel, you have to provide the
// work division as well as the additional kernel function parameters.
alpaka::exec<Acc>(queue,workDiv,helloWorldKernel/* put kernel arguments here */);
```





[Appendix2] Matrix Vector Multiplication Example

al/aka Accelerator Details



APPENDIX-5

- Accelerator chosen by the programmer and hides hardware specifics behind alpaka's abstract API using Acc = acc::AccGpuCudaRt<Dim, Idx>;
- Inside Kernel: contains thread state, provides access to alpaka's device-side API
 - The Accelerator provides the means to access to the indices

```
// get thread index on the grid
auto gridThreadIdx = alpaka::getIdx<Grid, Threads>(acc);
// get block index on the grid
auto gridBlockIdx = alpaka::getIdx<Grid, Blocks>(acc);
```

The Accelerator gives access to alpaka's shared memory (for threads inside the same block)

```
// allocate a variable in block shared static memory
auto & mySharedVar = block::shared::st::allocVar<int, COUNTER >(acc);
// get pointer to the block shared dynamic memory
float * mySharedBuffer = block::shared::dyn::getMem<float>(acc);
```

It also enables synchronization on the block level

```
// synchronize all threads within the block
block::sync::syncBlockThreads(acc);
```

- Internally, the accelerator maps all device-side functions to their native counterparts
 - Device-side functions require the accelerator as first argument: math::sqrt(acc, /* ... */); time::clock(acc); atomic::atomicOp<atomic::op::Or>(acc, /* ... */, hierarchy::Grids); (Atomics)
- On Host: Meta-parameter for choosing correct physical device and dependent types





APPENDIX-6 Physical device information and management by "alpaka Device"

- Each alpaka Device represents a single physical device;
- Contains device information:

```
auto const name = alpaka::getName(myDev);
                                                   // Back-end-defined device
 name
```

- auto const bytes = alpaka::getMemBytes(myDev); // Size of device memory
- auto const free = alpaka::getFreeMemBytes(myDev); // Size of available device memory
- Provides the means for device management:

```
alpaka::reset(myDev);
                                                   // Reset GPU device state
```

- Encapsulates back-end device:
 - auto nativeDevice = alpaka::getDev(myDev); // nativeDevice is not portable!





APPENDIX-7 Queue operations

- Queues execute Tasks (see next slide):
 - alpaka::enqueue(myQueue, taskRunKernel);
- Check for completion:
 - bool done = alpaka::empty(myQueue);
- Wait for completion, Events (see next slide), or other Queues:
 - alpaka::wait(myQueue); // blocks caller until all operations have completed
 - alpaka::wait(myQueue, myEvent); // blocks myQueue until myEvent has been reached
 - alpaka::wait(myQueue, otherQueue); // blocks myQueue until otherQueue's ops have completed





Appendix8 - Changing the target platform

```
using namespace alpaka;
using Dim = dim::DimInt<1u>;
using Idx = std::size t;
/*** BEFORE ***/
using Acc = alpaka::AccCpuOmp2Blocks<Dim, Idx>;
/*** AFTER ***/
using Acc = alpaka::AccGpuHipRt<Dim, Idx>;
/* No change required - dependent types and variables are automatically changed */
auto myDev = alpaka::getDevByIdx<Acc>(0u);
using Queue = alpaka::Queue<Acc, queue::NonBlocking>;
auto myQueue = Queue{myDev};
```



APPENDIX 9

Acc versus Device

Acc is a host side compile time construct but instantiated in Kernel Abstraction that make the kernel device agnostic