

Programming Accelerators with OpenMP



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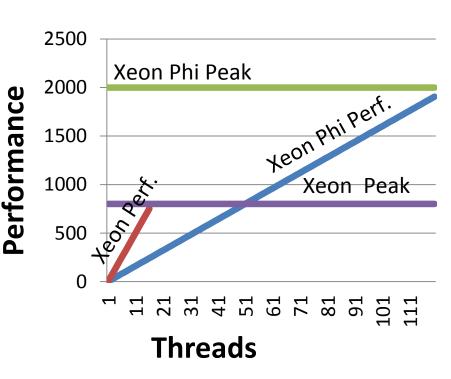


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Motivation



- **Demand for more compute power**
- Reach higher performance with more threads
- **Power consumption: Better** performance / watt ratio
- **GPUs are one alternative, but:** CUDA is hard to learn / program
- Intel Xeon Phi can be programmed with established programming paradigms like OpenMP, MPI, Pthreads



Source: James Reinders, Intel

Architecture (1/2)

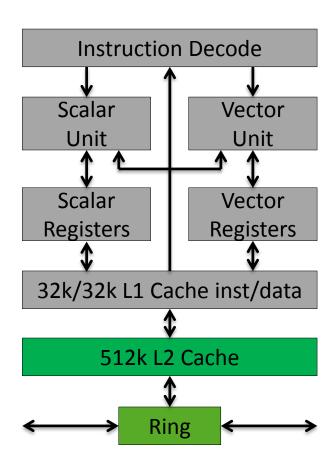




Source: Intel

Intel Xeon Phi Coprocessor

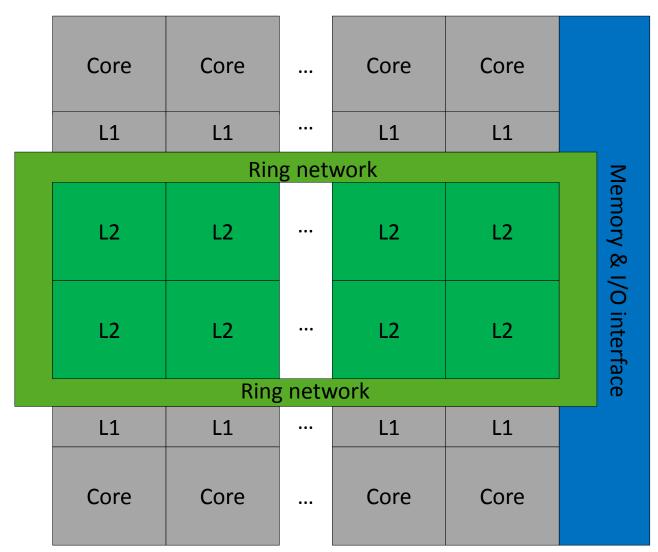
- 1 x Intel Xeon Phi @ 1090 MHz
- 60 Cores (in-order)
- ~ 1 TFLOPS DP Peak
- 4 hardware threads per core (SMT)
- 8 GB GDDR5 memory
- 512-bit SIMD vectors (32 registers)
- Fully-coherent L1 and L2 caches
- Plugged into PCI Express bus

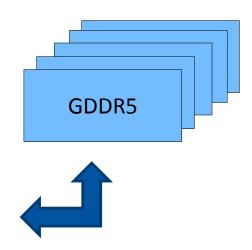


Architecture (2/2)





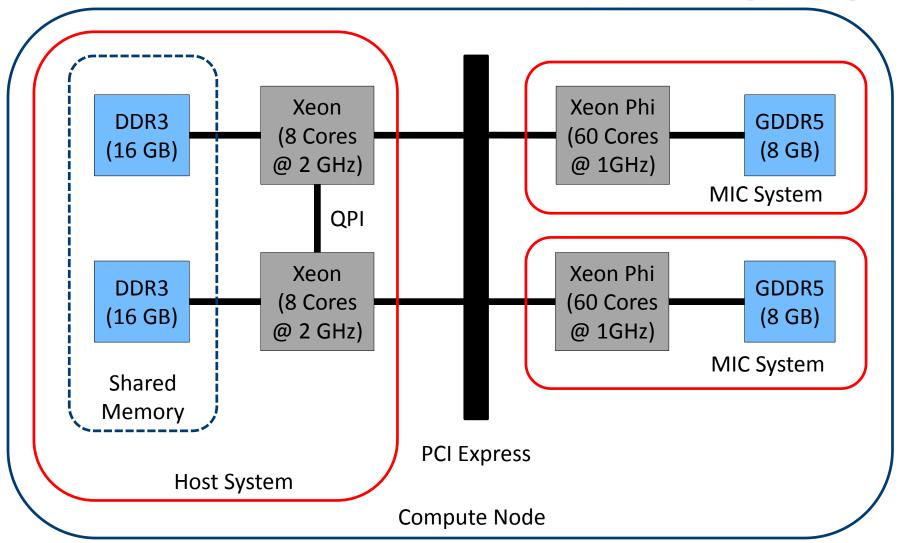




Xeon Phi Nodes at RWTH Aachen



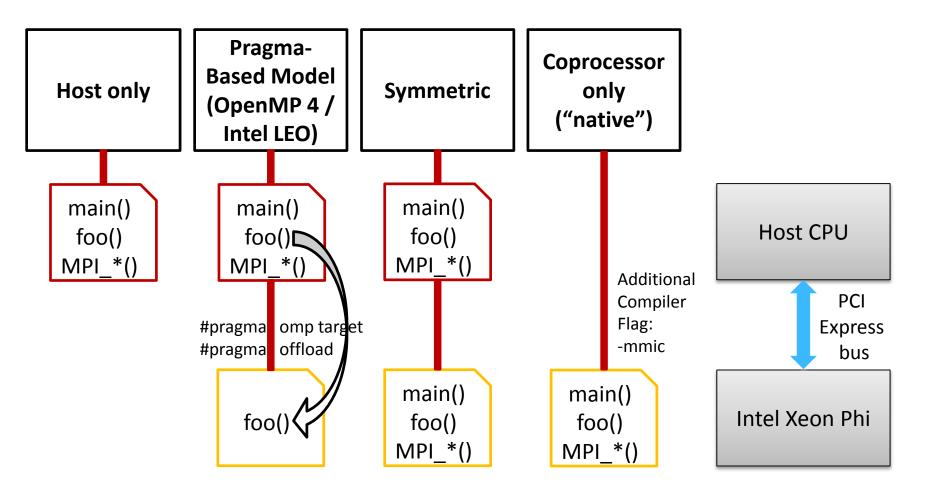




Programming Models







Native/Symmetric Execution



- The Xeon Phi is x86 based, but not binary compatible to the host
 - → recompile the application with "-mmic" to create an executable for Xeon Phi coprocessors
 - → For native execution:
 - →login via ssh
 - → start the executable
 - → For MPI (symmetric) execution:
 - → Compile two executables, one for the host and one for the coprocessor with a certain suffix, e.g. ".mic"
 - → Set "I_MPI_MIC=enabled" and "I_MPI_MIC_POSTFIX=.mic"
 - → Start the executable with mpiexec on host and coprocessor

Performance Expectations





Comparison with our MPI nodes (2 x Intel Westmere CPUs)

	Xeon Phi 5110P	2 x Xeon X5675
Frequency [GHz]	1.09	3.07
Cores	60	12
Peak Performance [GFLOPS]	1171	144
Peak Performance (no SIMD) [GFLOPS]	146	72
Memory Bandwidth (STREAM) [GB/s]	160	40
Memory Bandwidth (single core) [GB/s]	5	12

- → Only scalable, vectorizable codes benefit of the Xeon Phi at the moment
- → Performance is not for free, tuning will be necessary



OpenMP for Accelerators

Agenda



- What is an Accelerator?
- Execution and Data Model
- Target Construct
- Example: SAXPY
- Live Experiment: PI
- Outlook: what will come next?



What is an Accelerator?

in OpenMP

What kind of devices shall be supported?



In how differs an accelerator from just another core?

- → different functionality, i.e. optimized for something special
- → different (possibly limited) instruction set
- → heterogeneous device

Assumptions used as design goals for OpenMP 4.0:

- → every accelerator device is attached to one host device
- → it is probably heterogeneous
- → it may not be programmable in the same language as the host, or it may not implement all operations available on the host
- → it may or may not share memory with the host device
- → some accelerators are specialized for loop nests



Execution and Data Model

Execution Model



- Host-centric: the execution of an OpenMP program starts on the host device and it may offload target regions to target devices
 - → In principle, a target region also begins as a single thread of execution: when a target construct is encountered, the target region is executed by the implicit device thread and the encountering thread/task [on the host] waits at the construct until the execution of the region completes
- If a target device is not present, or not supported, or not available, the target region is executed by the host device
- If a construct creates a data environment, the data environment is created at the time the construct is encountered

Data Model



- When an OpenMP program begins, each device has an initial device data environment
- Directives accepting data-mapping attribute clauses determine how an original variable is mapped to a corresponding variable in a device data environment
 - →original: the variable on the host
 - → corresponding: the variable on the device
 - → the corresponding variable in the device data environment may share storage with the original variable (danger of data races)
- If a corresponding variable is present in the enclosing device data environment, the new device data environment inherits the corresponding variable from the enclosing device

Execution + Data Model



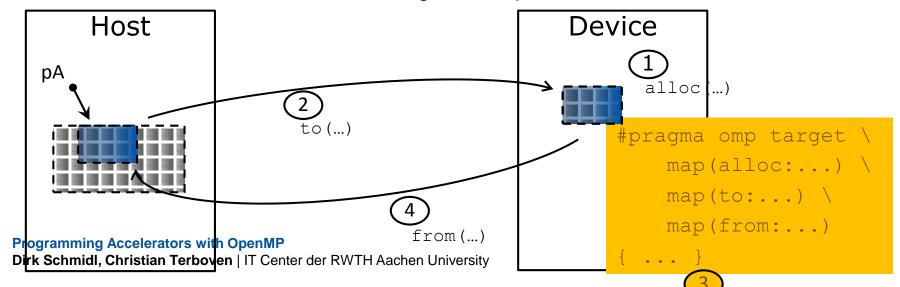


Data environment is lexically scoped

- → Data environment is destroyed at closing curly brace
- → Allocated buffers/data are automatically released

Use target construct to

- → Transfer control from the host to the device
- → Establish a data environment (if not yet done)
- → Host thread waits until offloaded region completed



Device constructs



```
data region
#pragma omp target data device(0) map(alloc:tmp[:N])
                   map(to:input[:N]) map(from:res)
                                                      shaping and
                                                         slicing
#pragma omp target_device(0)
#pragma omp parallel for
                                           offload
    for (i=0; i< N; i++)
      tmp[i] = some computation(input[i], i);
    do some other stuff on host();
#pragma omp target device(0)
#pragma omp parallel for reduction(+:res)
    for (i=0; i< N; i++)
      res += final computation(tmp[i], i)
```

This is a very simple example, but you see functionality similar to OpenACC 1.0. Device support has to come with the implementation.

Example: Execution and Data Model



Environment Variable OMP_DEFAULT_DEVICE=<int>: sets the device number to use in target constructs

```
double B[N] = ...; // some initialization
#pragma omp target map(tofrom:B)
#pragma omp parallel for
for (i=0; i<N; i++)
   B[i] += sin(B[i]);</pre>
```

→ map variable B to device, then execute parallel region on device, works probably pretty well on Intel Xeon Phi

Comparing OpenMP with OpenACC





OpenMP 4.0 – for Intel Xeon Phi:

```
#pragma omp target device(0) map(tofrom:B)
#pragma omp parallel for
for (i=0; i<N; i++)
   B[i] += sin(B[i]);</pre>
```

OpenMP 4.0 – for NVIDIA GPGPU:

OpenACAlsopossible (and equivalent) in OpenMP 4.0:



Target Construct

target data construct



Creates a device data environment for the extent of the region

- → when a target data construct is encountered, a new device data environment is created, and the encountering task executes the target data region
- → when an if clause is present and the if-expression evaluates to false, the device is the host

C/C++:

```
The syntax of the target data construct is as follows:

#pragma omp target data [clause[[,] clause],...] new-line
structured-block

where clause is one of the following:

device( integer-expression )
map( [map-type : ] list )
if( scalar-expression )
```

map clause



- Map a variable from the current task's data environment to the device data environment associated with the construct
 - → the list items that appear in a map clause may include array sections
 - → alloc-type: each new corresponding list item has an undefined initial value
 - → to-type: each new corresponding list item is initialized with the original lit item's value
 - → **from**-type: declares that on exit from the region the corresponding list item's value is assigned to the original list item
 - → *tofrom*-type: the default, combination of to and from

C/C++:

The syntax of the **map** clause is as follows:

map([map-type:]list)

target construct



- Creates a device data environment and execute the construct on the same device
 - → superset of the target data constructs in addition, the target construct specifies that the region is executed by a device and the encountering task waits for the device to complete the target region

C/C++:

```
The syntax of the target construct is as follows:

#pragma omp target [clause[[,] clause],...] new-line
structured-block

where clause is one of the following:

device( integer-expression )
map( [map-type : ] list )
if( scalar-expression )
```

Note: map clause does not work correctly with Intel 14.0 compilers

target update construct



Makes the corresponding list items in the device data environment consistent with their original list items, according to the specified motion clauses

C/C++:

```
The syntax of the target update construct is as follows:

#pragma omp target update motion-clause[, clause[,] clause],...] new-line
where motion-clause is one of the following:

to( list )
from( list )
and where clause is one of the following:

device( integer-expression )
if( scalar-expression )
```

declare target directive



- Specifies that [static] variables, functions (C, C++ and Fortran) and subroutines (Fortran) are mapped to a device
 - → if a list item is a function or subroutine then a device-specific version of the routines is created that can be called from a target region
 - → if a list item is a variable then the original variable is mapped to a corresponding variable in the initial device data environment for all devices (if the variable is initialized it is mapped with the same value)
 - all declarations and definitions for a function must have a declare target directive

C/C++:

The syntax of the **declare target** directive is as follows:

#pragma omp declare target new-line declarations-definition-seq #pragma omp end declare target new-line

Example: Target Construct





```
#pragma omp target device(0)
#pragma omp parallel for
for (i=0; i<N; i++) ...</pre>
```

```
#pragma omp target
#pragma omp teams num_teams(8) num_threads(4)
#pragma omp distribute
  for ( k = 0; k < NUM_K; k++ )
   {
      #pragma omp parallel for
      for ( j = 0; j < NUM_J; j++ )
      {
         ...
      }
    }
}</pre>
```

teams construct (1/2)



- Creates a league of thread teams where the master thread of each team executes the region
 - → the number of teams is determined by the num_teams clause, the number of threads in each team is determined by the num_threads clause, within a team region team numbers uniquely identify each team (omp_get_team_num())
 - → once created, the number of teams remeinas constant for the duration of the teams region
- The teams region is executed by the master thread of each team
- The threads other than the master thread to not begin execution until the master thread encounteres a parallel region
- Only the following constructs can be closely nested in the team region: distribute, parallel, parallel loop/for, parallel sections and parallel workshare

distribute construct



- Specifies that the iteration of one or more loops will be executed by the thread teams, the iterations are distributed across the master threads of all teams
 - → there is no implicit barrier at the end of a distribute construct
 - → a distribute construct must be closely nested in a teams region

C/C++:

```
#pragma omp distribute [clause[[,] clause],...] new-line for-loops

Where clause is one of the following:

private( list )
firstprivate( list )
collapse( n )
dist_schedule( kind[, chunk_size] )

All associated for-loops must have the canonical form described in Section 2.5.
```

teams construct (2/2)



- A teams construct must be contained within a target construct, which must not contain any statements or directives outside of the teams construct
- After the teams have completed execution of the teams region, the encountering thread resumes execution of the enclosing target region

C/C++:

```
#pragma omp teams [clause[[,] clause],...] new-line structured-block
where clause is one of the following:
num_teams( integer-expression )
num_threads( integer-expression )
default(shared | none)
private( list )
firstprivate( list )
shared( list )
reduction( operator : list )
```



Example: SAXPY

SAXPY: Serial (Host)



```
int main(int argc, const char* argv[]) {
 int n = 10240; float a = 2.0f; float b = 3.0f;
 float *x = (float*) malloc(n * sizeof(float));
 float *y = (float*) malloc(n * sizeof(float));
  // Initialize x, y
 // Run SAXPY TWICE
 for (int i = 0; i < n; ++i) {
       y[i] = a*x[i] + y[i];
  // y is needed and modified on the host here
 for (int i = 0; i < n; ++i) {
       y[i] = b*x[i] + y[i];
 free(x); free(y); return 0;
```

SAXPY: OpenACC v2 (NVIDIA GPGPU)



```
int main(int argc, const char* argv[]) {
  int n = 10240; float a = 2.0f; float b = 3.0f;
  float *x = (float*) malloc(n * sizeof(float));
 float *v = (float*) malloc(n * sizeof(float));
  // Initialize x, y
 // Run SAXPY TWICE
#pragma acc data copyin(x[0:n])
#pragma acc parallel copy(y[0:n])
#pragma acc loop
  for (int i = 0; i < n; ++i) {
        y[i] = a*x[i] + y[i];
  // y is needed and modified on the host here
#pragma acc parallel copy(y[0:n])
#pragma acc loop
  for (int i = 0; i < n; ++i) {
        y[i] = b*x[i] + y[i];
 free(x); free(y); return 0;
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```

SAXPY: OpenMP 4.0 (Intel MIC)





```
int main(int argc, const char* argv[]) {
 int n = 10240; float a = 2.0f; float b = 3.0f;
 float *x = (float*) malloc(n * sizeof(float));
 float *v = (float*) malloc(n * sizeof(float));
  // Initialize x, y
 // Run SAXPY TWICE
#pragma omp target data map(to:x[0:n])
#pragma omp target map(tofrom:y[0:n])
#pragma omp parallel for
for (int i = 0; i < n; ++i) {
       y[i] = a*x[i] + y[i];
  // y is needed and modified on the host here
#pragma omp target map(tofrom:y[0:n])
#pragma omp parallel for
 for (int i = 0; i < n; ++i) {
       y[i] = b*x[i] + y[i];
 free(x); free(y); return 0;
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```

SAXPY: OpenMP 4.0 (NVIDIA GPGPU)





```
int main(int argc, const char* argv[]) {
  int n = 10240; float a = 2.0f; float b = 3.0f;
  float *x = (float*) malloc(n * sizeof(float));
 float *y = (float*) malloc(n * sizeof(float));
  // Initialize x, y
  // Run SAXPY TWICE
#pragma omp target data map(to:x[0:n])
#pragma omp target map(tofrom:y[0:n])
#pragma omp teams
#pragma omp distribute
#pragma omp parallel for
for (int i = 0; i < n; ++i) {
        y[i] = a*x[i] + y[i];
  // y is needed and modified on the host here
#pragma omp target map(tofrom:y[0:n])
#pragma omp teams
#pragma omp distribute
#pragma omp parallel for
  for (int i = 0; i < n; ++i) {
        y[i] = b*x[i] + y[i];
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 free(x); free(y); return 0;
```



Live Experiment: PI



Host version

```
double f(double x) {
   return (double) 4.0 / ((double) 1.0 + (x*x));
void computePi() {
   double h = (double) 1.0 / (double) iNumIntervals;
   double sum = 0, x;
#pragma omp parallel for private(x) reduction(+:sum)
   for (int i = 1; i <= iNumIntervals; i++) {
      x = h * ((double)i - (double)0.5);
      sum += f(x);
  myPi = h * sum;
```



Accelerator version (Intel Xeon Phi)

```
#pragma omp declare target
double f(double x) {
   return (double) 4.0 / ((double) 1.0 + (x*x));
void computePi() {
   double h = (double) 1.0 / (double) iNumIntervals;
   double sum = 0, x;
#pragma omp target
#pragma omp parallel for private(x) reduction(+:sum)
   for (int i = 1; i <= iNumIntervals; i++) {
      x = h * ((double)i - (double)0.5);
      sum += f(x);
   myPi = h * sum;
```



PI on Host and Accelerator



Outlook

Accelerators – current work



- OpenMP 4.0 introduced the target construct for device programming, i.e. compute accelerators some aspects were missing / undefined:
 - → combinations of constructs, #pragma omp target teams distribute parallel for simd
 - → concurrent mapping of the same variable by two threads still undefined

Current Plans for OpenMP 4.1:

- → async execution: currently the host waits for target operation to complete
 - → task target construct: asyn offload, host will continue execution
 - -> dependencies between host and device: adding target will maintain them
- → unstructured data movement: sometimes data regions are too static
 - → target enter data / exit data constructs to push / pull at encountering
- → open discussion on support for multiple devices (i.e. work distribution)



Questions?