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Natural Environment Research Council







Outline



- OpenMP/ROCm/CUDA interoperability
- Custom data mapper
- Memory types
- Memory allocation
- Asynchronous tasks



Using ROCm in an OpenMP Program



- Sometimes need to call routines from a library that expects a device pointer, not a mapped variable.
- Might be needed if you want to port a computationally heavy kernel to CUDA/HIP to increase performance
- Might be needed if you want to call vendor libraries (rocFFT/cuFFT, cuBLAS,rocBLAS ...)

Using ROCm in an OpenMP Program



- Suppose var_a is a variable already mapped to the device with OpenMP
- Want to avoid re-allocating and transferring data with CUDA/HIP
- Need to obtain the address of the variable mapped on the device

Using rocm in an OpenMP Program



• The clause **use_device_addr** makes all references to **var_a** in the code refer to the variable mapped on the GPU.

```
#pragma omp target data use_device_addr(var_a)
{ . . . . }
C/C++
```

```
!$omp target data use_device_addr(var_a)
...
!$omp end target data
Fortran
```

Using ROCm with OpenMP



Example of a matrix multiplication using rocBLAS

```
!$omp target data map(to:A(1:M,1:K),B(1:K,1:N)) map(tofrom:C(1:M,1:N))
  !$omp target data use_device_addr(A(1:M,1:K), B(1:K,1:N), C(1:M,1:N))
      call rocblas_dgemm( . . . , A, . . . , B, . . . , C, . . .)
  !$omp end target data

!$omp end target data
```



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- Memory allocations on HIP/CUDA
- Want to use OpenMP code that uses memory allocated on the device with HIP/CUDA
- Useful when porting from a HIP/CUDA program to a performance portable implementation with OpenMP



- The clause is <u>device_ptr(a,b, . . .)</u> tells OpenMP a, b, ... are pointers to data available on the device, not on the host.
- Prevents mapping of the pointer variable, allows using the pointer directly on the device.



Ex.: Copying an array b to array c

```
double * b_device, c_device;
hipMalloc( &b_device, n*sizeof(double) ) );
hipMalloc( &c_device, n*sizeof(double) ) );

#pragma omp target teams distribute parallel for
is_device_ptr(c_device, b_device)
for (int i=0;i<n;i++) {
   c_device[i] = b_device[i];
}</pre>
```



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Ex.: Copying an array b to array c

```
integer, parameter :: fp kind = kind(0.0)
type(c ptr) :: cptr b, cptr c
real(fp kind), pointer, dimension(:) :: fptr b => null(), fptr c => null()
integer :: n = 5, i, ierr
ierr = hipMalloc(cptr_b, n * sizeof(fp_kind))
ierr = hipMalloc(cptr c, n * sizeof(fp kind))
call c f pointer(cptr_b, fptr_b, [n])
call c f pointer(cptr c, fptr c, [n])
!$omp target teams distribute parallel do simd is device ptr(fptr b, fptr c)
do i = 1, n
  fptr c(i) = fptr b(i)
end do
                                                                         Fortran
```





- Consider a structure with three arrays: **c**, **d**, **e**.
- Each array is of length n.
- Mapping a variable of type $\bf A$ will correctly map scalar variables (like $\bf n$) but not the value of the arrays $\bf c$, $\bf d$, $\bf e$.

```
struct my_struct {
  int n;
  double * c;
  double * d;
  double * e;
};
```

```
type my_type
  integer :: n
  real, allocatable :: c(:)
  real, allocatable :: d(:)
  real, allocatable :: e(:)
end type
Fortran
```



- All arrays need to be explicitly mapped every time a variable of type my struct is used.
- Can be very verbose and error-prone with complex data structures.

```
my struct a, b;
#pragma omp target data map(tofrom:a,a.c[0:n],a.d[0:n],a.e[0:n]) {
    // Do something with a
#pragma omp target data map(tofrom:b,b.c[0:n],b.d[0:n],b.e[0:n]) {
    // Do something with b
                                                                      C/C++
type (my type) :: a, b
!$omp target data map(tofrom:a,a%c(1:a%n),a%d(1:a%n),a%e(1:a%n))
    ! Do something with a
!$omp end target data
!$omp target data map(tofrom:b,b%c(1:b%n),b%d(1:b%n),b%e(1:b%n))
    ! Do something with b
!$omp end target data
                                                                     Fortran
```



Declare all mapping rules for a custom structure only once.

```
!$omp declare mapper(my_type :: x) map(x, ...)
Fortran
```

• Here **x** is a dummy argument of type **my_struct** or **my_type**.

```
epcc
```

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```
#pragma omp declare mapper(my struct x) \
 map(x, x.c[0:x.n], x.d[0:x.n], x.e[0:x.n])
my struct a, b;
#pragma omp target data map(tofrom:a) {
      // Do something with a
#pragma omp target data map(tofrom:b) {
      // Do something with b
                                        C/C++
```

```
type (my_type) :: a, b
!$omp declare mapper(my type :: x) &
!\$omp map(x, x&c(1:x&n), x&d(1:x&n), &
!$omp x.e(1:x%n))
!$omp target data map(tofrom:a)
 ! Do something with a
!$omp end target data
!$omp target data map(tofrom:b)
 ! Do something with b
!$omp end target data
                                  Fortran
```



Host memory



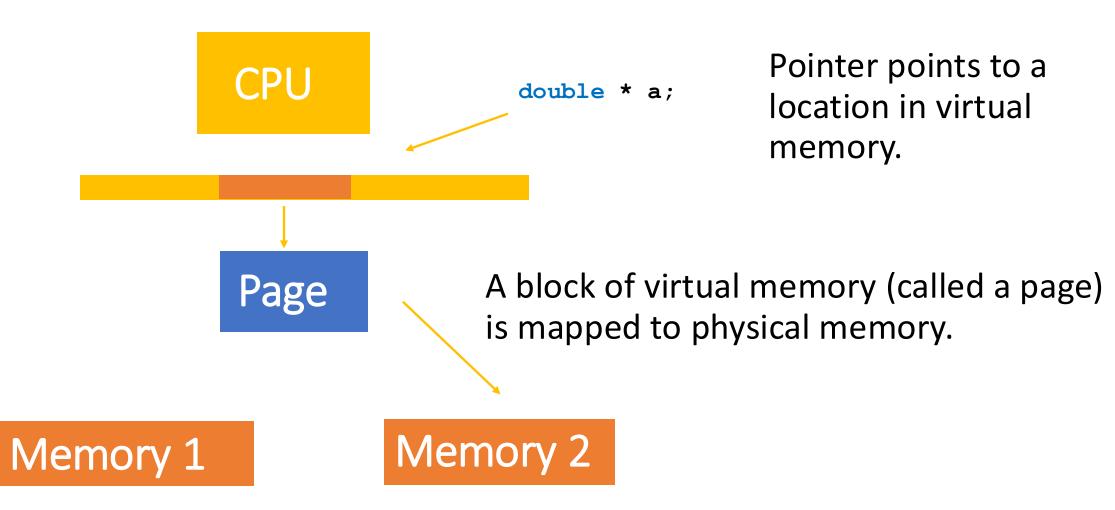
• Pageable memory: Memory allocated in the usual way on the (malloc, new, allocate, etc...). Memory is only available on the host.

 Pinned memory: Memory optimized for transfer to the GPU. Custom allocators are used on the host

Unified memory: Memory is accessible on both the host and the device.
 Memory allocated in the usual way.

Pageable memory





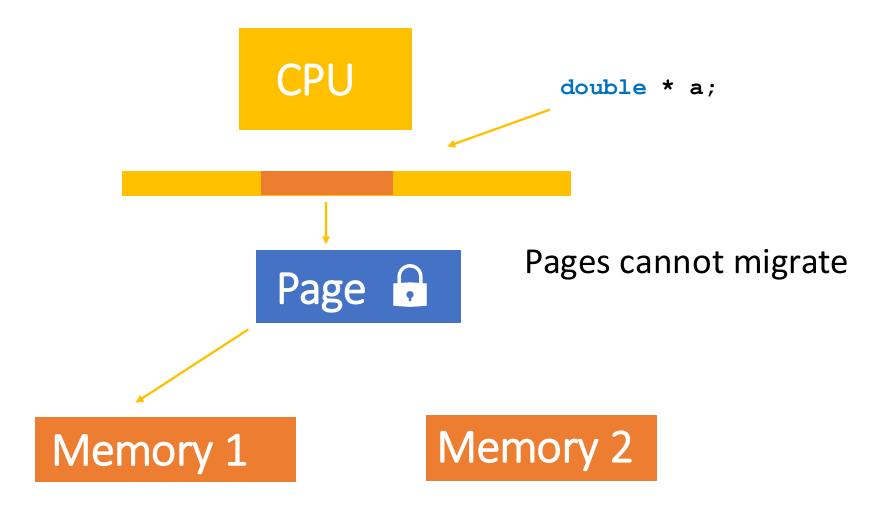
Pageable memory



- Memory is allocated using the usual allocators.
- Memory address is virtual.
- Virtual memory is mapped to physical memory using a page.
- Pages can migrate between different physical devices.
- Memory is only accessible from the host.

Pinned memory





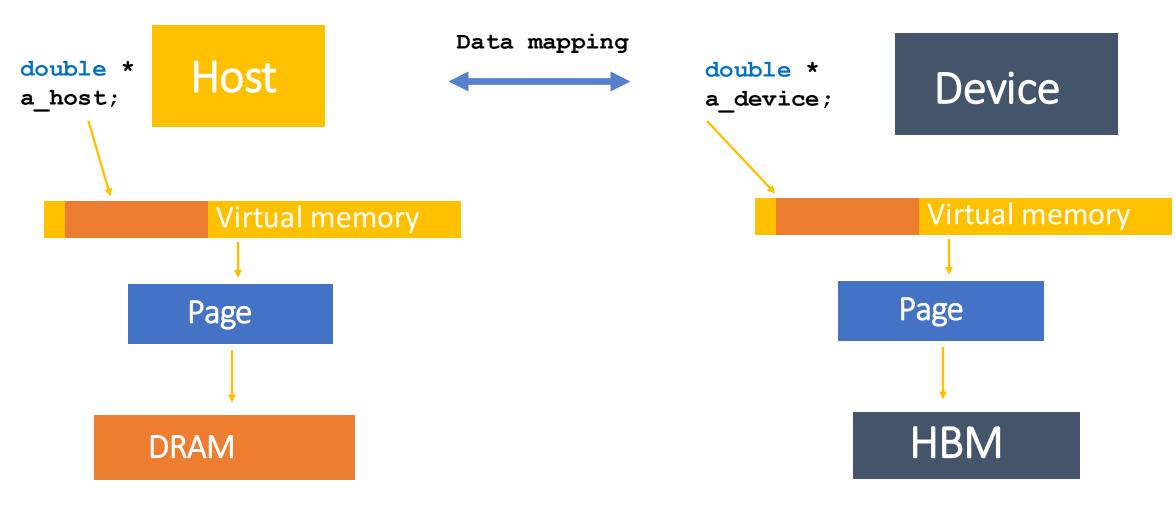
Pinned memory



- Memory is allocated using the usual allocators.
- Memory address is virtual.
- Virtual memory is mapped to physical memory using a page.
- Pages cannot migrate between different physical devices.
- Memory is only accessible from the host.
- Memory transfers are faster than pageable memory.

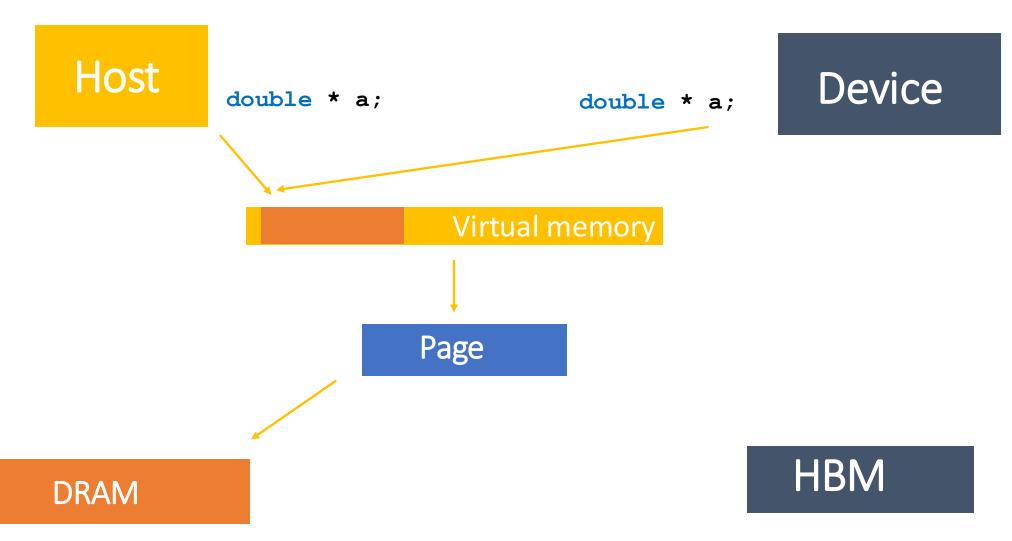
Non-unified memory





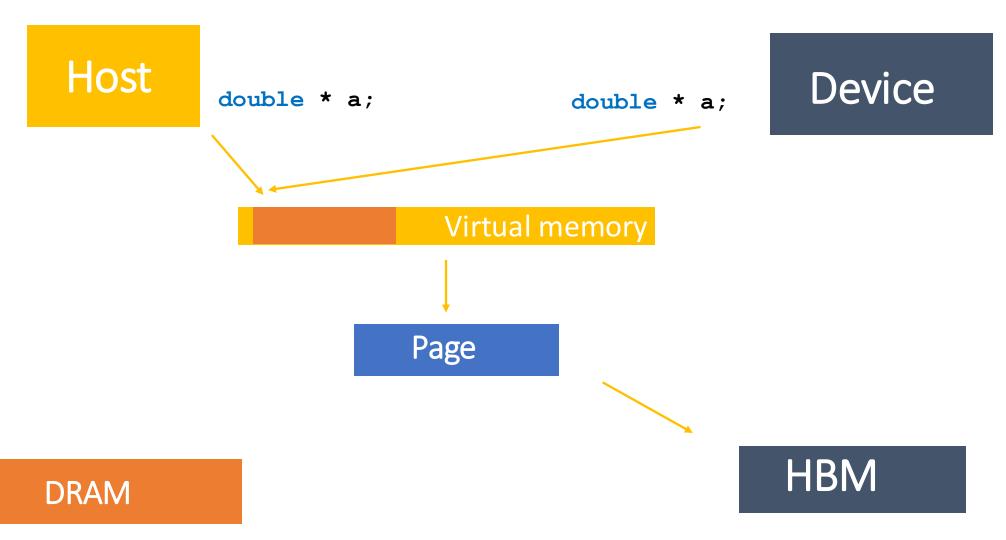
Unified memory





Non-unified memory





Unified memory



- Standard allocators are used, which are not GPU aware.
- Needs to use requires unified_shared_memory
- Host can access pages on the device and vice-versa
- Mapping is optional

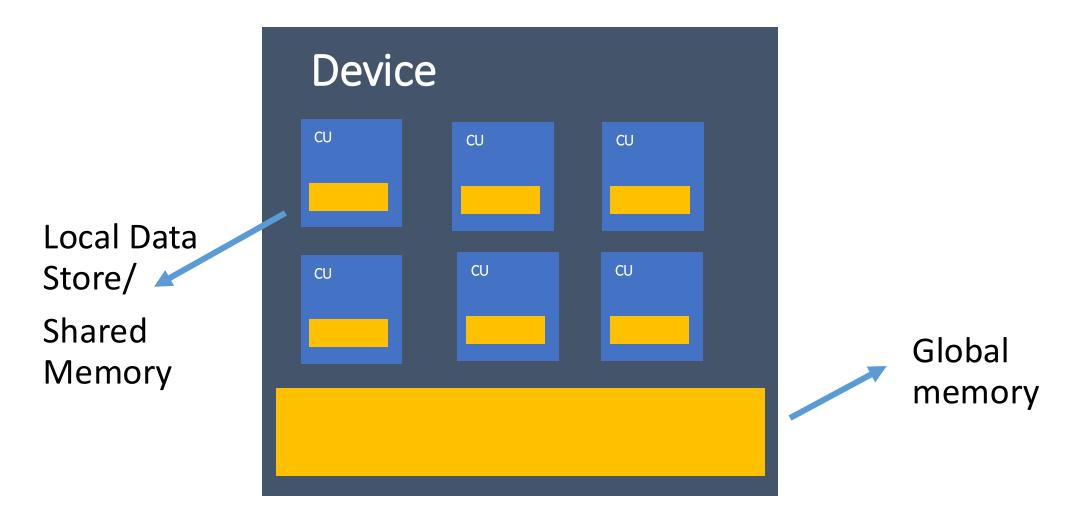
Unified memory



- Does not work on older devices
- Performance is good only on architectures that have fast interconnect between host and device memory (MI300A, Grace-Hopper etc..).

Memory on the Device





Memory



Device		
Global memory	Shared by all threads on GPU	
	Slower than the local data store	
Local Data Store	Shared by threads on a CU/team	
	Fast but small	

Host	
Pageable	Allocated with malloc, new, allocate
	Slow data transfer to GPU
Pinned memory	Fast data transfer
	Requires special allocation calls



OpenMP allocators



• Define how to allocate memory; is a combination of a memory space and an allocator traits.

Memory space: hint to the compiler on where to allocate memory.
 The omp_default_mem_space memory space is sufficient for most applications.

- Allocator traits: hints to the compiler on how the memory should be allocated and how the memory will be accessed.
 - Implementation is undefined.

OpenMP allocator traits



Allowed values	Notes
A positive integer must be a multiple of 2	
all	can be accessed from anywhere.
cgroup	can be accessed by all threads in a contention group.
pteam	only shared by threads in a team.
thread	memory private to each thread.
true , false	
	A positive integer must be a multiple of 2 all cgroup pteam

OpenMP allocation on the host



```
! Use the default memory space.
integer(omp_memspace_handle_kind) :: c_memspace = omp_default_mem_space
! Create an arry of 2 traits.
type(omp_alloctrait) :: c_traits(2)
! Create a custom allocator.
integer(omp_allocator_handle_kind) :: c_alloc
c_traits(1) = omp_alloctrait(omp_atk_pinned, .TRUE.)
c_traits(2) = omp_alloctrait(omp_atk_alignment, 128)
c_alloc = omp_init_allocator(c_memspace, 2, c_traits)

Fortran
```

OpenMP allocation on the host



- The allocate directive provides a hint to the compiler.
 It is not guaranteed that the compiler will use the traits you provide.
- You can use the clause allocate (my_allocator: <item-list>)
- The clause needs to be associated with an allocation statement

OpenMP allocation on the host



```
#pragma omp allocate(my_alloc:b)
b = new double [n];
C/C++
```

```
!$omp allocate(my_alloc:b)
allocate (b(n))
Fortran
```

OpenMP allocation on the device



• On the device you must specify which allocators can be used on the host using the uses allocators clause

```
#pragma omp target uses_allocators(my_allocator)
    map(alloc:c[0:BLOCK_SIZE]) allocate(my_allocator:c)
```

```
real, allocatable, dimension(:) :: c
!$omp target teams num_teams(4) & Fortran
!$omp private(c) &
!$omp uses_allocators(my_allocator) &
!$omp map(alloc:c(1:BLOCK_SIZE)) &
!$omp allocate(my_allocator:c)
```

OpenMP allocation on the LDS



- OpenMP does not (yet) provide a mechanism for allocating on the local store.
- However, you can provide hints to the compiler.
- You can use also compiler extensions, but they are not portable between compilers

OpenMP allocation on the LDS



• Statically allocate memory inside a team.

```
#pragma omp target teams num_teams(4) {
    double c[BLOCK_SIZE];
    // do something with a team
    #pragma omp parallel for
    for(int i=0; i<BLOCK_SIZE; i++) {
        c[i] = i;
    }
}</pre>
```

```
program main
  implicit none
  !$omp target teams num teams(4)
    call loop()
  !$omp end target teams
end program main
subroutine loop
  integer, parameter :: N = 1000
  integer :: i
  real, dimension(N) :: c
  !$omp parallel do
  do i = 1, N
    c(i) = i
  end do
end subroutine loop
                                 Fortran
```

OpenMP allocation on the LDS



- Use the default allocator omp_pteam_mem_alloc
- On some compilers you need to use omp_cgroup_mem_alloc
- On some compilers neither is implemented

```
double c[BLOCK_SIZE];

#pragma omp target teams
    num_teams(4) private(c)
    uses_allocators(omp_pteam_mem_alloc)
    map(alloc:c[0:BLOCK_SIZE])
    allocate(omp_pteam_mem_alloc:c)

#pragma omp parallel for shared(c)
for(int i=0;i<BLOCK_SIZE;i++) {
    // do something with c
}</pre>
```

```
real, allocatable, dimension(:) :: c
!$omp target teams num_teams(4) &
!$omp private(c) &
!$omp uses_allocators(omp_pteam_mem_alloc) &
!$omp map(alloc:c(1:BLOCK_SIZE)) &
!$omp allocate(omp_pteam_mem_alloc:c)

!$omp parallel do shared(c)
do i = 1, BLOCK_SIZE
   ! do something with c
end do
!$omp end target teams
Fortran
```



Asynchronous offloading



- map directives are blocking: the CPU and all GPU kernel launches will stall until the data is transferred to the GPU.
- kernels offloaded to the device are blocking: the CPU will wait until the kernel finishes executing on the GPU. Only one kernel at a time is executed.
- You can use non blocking calls by adding the nowait clause.
- Synchronize tasks launched from a CPU thread using the taskwait directive.
- Synchronize tasks and multiple CPU threads using the barrier directive.
- Specify dependencies using the depend clause.

Asynchronous data transfer



 Add a nowait clause to map directives to overlap memory transfers and computations

```
#pragma omp target enter data map(to:xsi[0:m],ysi[0:m]) nowait
depend(out:xsi[0:m],ysi[0:m])

my_host_computation()

#pragma omp taskwait

C/C++
```

```
!$omp target enter data map(alloc:xsi(1:m),ysi(1:m)) nowait
depend(out:xsi(1:m),ysi(1:m))

call my_host_computation()

!$omp taskwait
Fortran
```

Asynchronous execution



Add a nowait clause to overlap computation on the host and device.

```
!$omp target teams distribute parallel do nowait
do j = 1, m
   ! computation A on device
end do
do j = 1, m
   ! independent computation B on host
end do
!$omp taskwait
Fortran
```

Parallel execution on the device



- Add a nowait clause to allow running multiple kernels on the device at the same time.
- Compiler might choose to run the kernels serially on the device or in parallel.
- Only advantageous if the offloaded region does not exhaust the resources on the device.

Parallel execution on the device



```
#pragma omp target teams distribute parallel for nowait
for (int j=0; j<m; j++) {
          // computation A
}
#pragma omp target teams distribute parallel for nowait
for (int j=0; j<m; j++) {
          // independent computation B
}
#pragma omp taskwait</pre>
C/C++
```

```
!$omp target teams distribute parallel do nowait
do j = 1, m
  ! computation A on device
end do
!$omp target teams distribute parallel do nowait
do j = 1, m
  ! independent computation B
end do
!$omp taskwait
Fortran
```

Dependencies



- If your kernels depend on data produced by other kernels you can tell that the kernel has a data dependency.
- You use the depend clause to mark data dependencies.

depend(out:x)

Mark that other tasks depend on the variable **x**, which might be modified by the current task.

depend(in:x)

This kernel depend on the variable **x**, modified by other kernels that specify the out modifier.

depend(inout:x) Specify both input and output dependencies.

Dependencies



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```
// Initiate data transfer. Mark x as a dependency of a future task.
#pragma omp target enter data map(to:x) nowait depend(out: x[0:m])
// Once x is available on the device start computation on the device.
#pragma omp target teams distribute parallel for nowait depend(inout: x[0:m])
for (int j=0; j<m; j++) {
    // update the array x
// Once x is updated from the previous kernel start this kernel.
#pragma omp target teams distribute parallel for nowait depend(in: x[0:m])
for (int j=0; j<m; j++) {
    // use the x array for my computation
```

Dependencies



```
! Initiate data transfer. Mark x as a dependency of a future task.
!$omp target enter data map(to:x) nowait depend(out: x(1:m))
! Once x is available on the device start computation on the device.
!$omp target teams distribute parallel do nowait depend(inout: x(1:m))
do j = 1, m
    ! update the array x
end do
! Once x is updated from the previous kernel start this kernel.
!$omp target teams distribute parallel do nowait depend(in: x(1:m))
do j = 1, m
    ! use the x array for my computation
end do
                                                                    Fortran
```

Summary

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- OpenMP/ROCm/CUDA interoperability
- Custom data mapper
- Memory types
- Memory allocation
- Asynchronous tasks