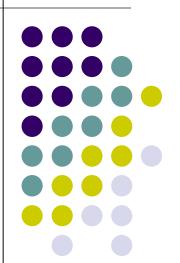
## Practical Parallel Computing (実践的並列コンピューティング)

Part2: GPU (2) May 9, 2022

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- Part 0: Introduction
  - 2 classes
- Part 1: OpenMP for shared memory programming
  - 4 classes
- Part 2: GPU programming
  - 4 classes
     We are here (2/4)
  - OpenACC (1.5 classes) and CUDA (2.5 classes)
- Part 3: MPI for distributed memory programming
  - 3 classes

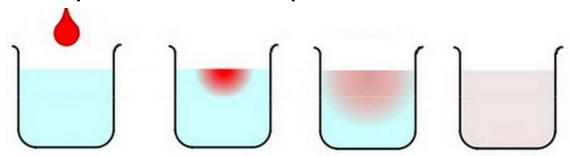
#### Data Region and Kernel Region in OpenACC **GPU** int main() Copy x,y CPU →GPU Α; #pragma acc data copy(x,y) #pragma acc kernels Data Β; Region #pragma acc kernels Kernel regions Copy x,y CPU ←GPU Ε;

- Data movement occurs at beginning and end of data region
- Data region may contain 1 or more kernel regions

### "diffusion" Sample Program related to [G1]



An example of diffusion phenomena:



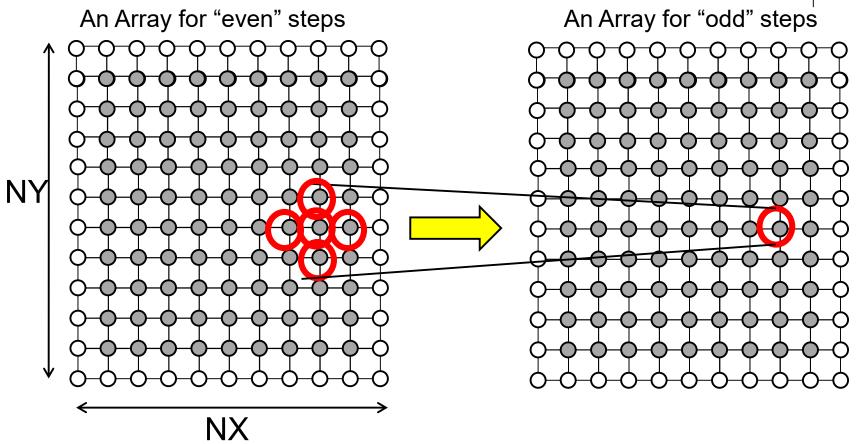
The ink spreads gradually, and finally the density becomes uniform (Figure by Prof. T. Aoki)

Available at /gs/hs1/tga-ppcomp/22/diffusion/ You can use /gs/hs1/tga-ppcomp/22/diffusion-acc/

- Execution:./diffusion [nt]
  - nt: Number of time steps

#### Data Structure in "diffusion"





# Consideration of Parallelizing Diffusion with OpenACC related to [G1]



- x, y loops can be parallelized
  - We can use "#pragma acc loop" twice
- t loop cannot be parallelized

```
[Data transfer from CPU to GPU]
for (t = 0; t < nt; t++) {
  for (y = 1; y < NY-1; y++) {
    for (x = 1; x < NX-1; x++) {
        :
     }
  }
}</pre>
```

Kernel region on GPU Parallel x, y loops

It's better to transfer data *out of* t-loop

[Data transfer from GPU to CPU]

### data Clause for Multi-Dimensional arrays



f loat A[2000][1000]; → an example of a 2-dimension array

```
.... data copy(A)
→ OK, all elements of A are copied
.... data copy(A[0:2000][0:1000])
→ OK, all elements of A are copied
.... data copy(A[500:600][0:1000])
→ OK, rows[500,1100) are copied
.... data copy(A[0:2000][300:400])
→ NG in current OpenACC
```

Currently, OpenACC does not support non-consecutive transfer



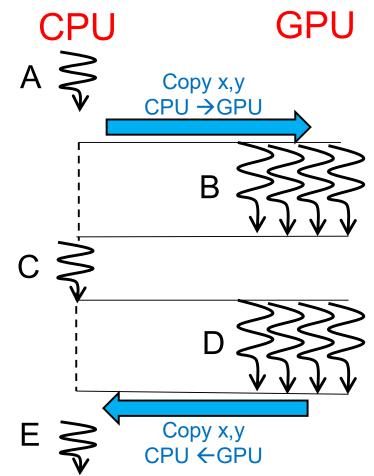


- You will need compiler options different from the diffusion directory for OpenACC
- You can use files in diffusion-acc directory as basis
  - /gs/hs1/tga-ppcomp/22/diffusion-acc/
  - "Makefile" in this directory supports compiler options for OpenACC
  - Don't forget "module load nvhpc" before "make"

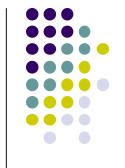
# Data Transfer Costs in GPU Programming Related to [G2]



- In GPU programming, data transfer costs between CPU and GPU have impacts on speed
  - Program speed may be slower than expected ☺



### Speed of GPU Programs: case of mm-acc



In mm-acc, speed in Gflops is computed by

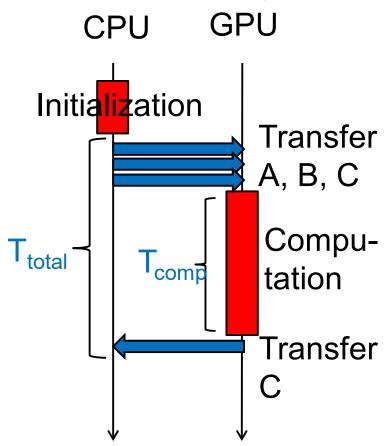
 $S = 2mnk / T_{total}$ 

T<sub>total</sub> includes both computation time and transfer

→ S counts slow-down by transfer

To see the effects, let's try another sample /gs/hs1/tga-ppcomp/22/mm-meas-acc which outputs time for

- copyin (transfer A, B, C)
- computation
- copyout (transfer C)



#### **To Measure Transfer Time**

 Data transfer occurs at the beginning and the end of "data region"
 CPU GPU

```
// A,B,C are on CPU

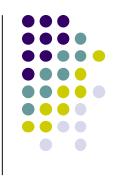
#pragma acc data copyin(A,B) copy(C)
{ // copyin (CPU->GPU) here

#pragma acc kernels
{
:
} //copyout (GPU->CPU) here
```

See mm-meas-acc/mm.c

Also note that gettimeofday() must be called on CPU

### Discussion on Data Transfer Costs



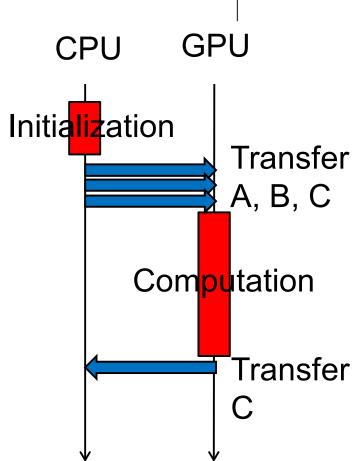
- - M: Data size in bytes
  - B: "Bandwidth" (speed)
  - L: "Latency" (if M is sufficiently large, we can ignore it)
- In a P100 GPU,
  - Theoretical computation speed is 5.3TFlops
  - Theoretical bandwidth B is 16GB/s (2G double values per second)
  - → Transfer of values is much slower than computation 🕾

# Discussion on Computation and Transfer Costs

In mm-acc,

- Computation amount: O(mnk)
- Data transfer amount:
  - A, B, C: CPU  $\rightarrow$  GPU: O(mk+kn+mn)
  - C: GPU → CPU: O(mn)

Transfer costs are relatively smaller with larger m, n, k



In diffusion-acc [G1], how can we reduce transfer costs?





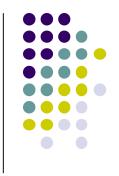
- Calling functions in kernel region is ok, but we need to be careful
  - "acc routine" directive is required by compiler to generate GPU code

### **How about Library Functions?**



Inside kernel regions (#pragma acc kernel),

- Available library functions is very limited
- We cannot use strlen(), memcpy(), fopen(), fflush()...
- We cannot use gettimeofday() <sup>(2)</sup>
- Exceptionally, some mathematical functions are ok ©
  - fabs, sqrt, fmax...
  - #include <math.h> is needed
- Recently, printf() in kernel regions is ok!



Now explanation of OpenACC is finished; we will go to CUDA

### OpenACC and CUDA for GPUs



#### OpenACC

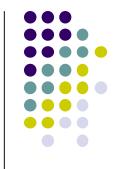
- C/Fortran + directives (#pragma acc ...), Easier programming
- NVIDIA HPC SDK compiler works
  - module load nvhpc
  - pgcc -acc ... XXX.c
- Basically for data parallel programs with for-loops
- → Only for limited types of algorithms ⊗

#### CUDA

- Most popular and suitable for higher performance
- Use "nvcc" command for compile
  - module load cuda
  - nvcc ... XXX.cu

Programming is harder, but more general

### An OpenACC Program Look Like



```
int A[100], B[100];
  int i;
#pragma acc data copy(A,B)
#pragma acc kernels
#pragma_acc_loop_independent
  for (i = 0; i < 100; i++) {
       A[i] += B[i];
```

**Executed on GPU** in parallel

// After kernel region finishes,
CPU can access to A[i],B[i]



### A CUDA Program Look Like

```
int A[100], B[100];
int *DA, *DB;
int i;
cudaMalloc(&DA, sizeof(int)*100);
cudaMalloc(&DB, sizeof(int)*100);
cudaMemcpy(DA,A,sizeof(int)*100,
   cudaMemcpyHostToDevice);
cudaMemcpy(DB,B,sizeof(int)*100,
   cudaMemcpyHostToDevice);
add<<<20, 5>>>(DA, DB);
cudaMemcpy(A,DA,sizeof(int)*100,
   cudaMemcpyDeviceToHost);
```

```
Sample: /gs/hs1/tga-ppcomp/22/add-cuda/
```

```
__global__ void add

(int *DA, int *DB)

{

int i = blockldx.x*blockDim.x

+ threadldx.x;

DA[i] += DB[i];

}
```

Executed on GPU (called a *kernel function*)





```
[make sure that you are at a interactive node (r7i7nX)]
module purge [If you have loaded nvhpc, delete it]
module load cuda [Do once after login]
cd ~/t3workspace [Example in web-only route]
cp -r /gs/hs1/tga-ppcomp/22/add-cuda
cd add-cuda
make
[An executable file "add" is created]
./add
```

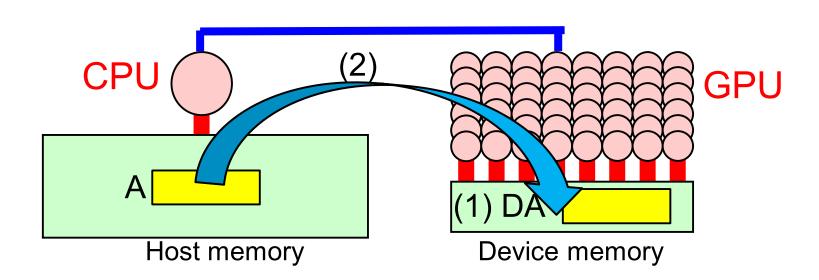
※ [Standard route] A log-in node does not have a GPU

→ You can compile the sample there, but the program does not work!

# Preparing Data on Device Memory



- (1) Allocate a region on device memory
  - cf) cudaMalloc((void\*\*)&DA, size);
- (2) Copy data from host to device
  - cf) cudaMemcpy(DA, A, size, cudaMemcpyDefault);



Note: cudaMalloc and cudaMemcpy must be called on CPU, NOT on GPU

### Comparing OpenACC and CUDA



#### **OpenACC**

Both allocation and copy are done by acc data copyin

One variable name A may represent both

- A on host memory
- A on device memory

```
int A[100]; ✓ on CPU
#pragma acc data copy(A)
#pragma acc kernels

{
    ··· A[i] ···
} on GPU
```

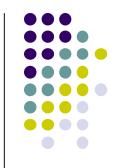
#### **CUDA**

cudaMalloc and cudaMemcpy are separated

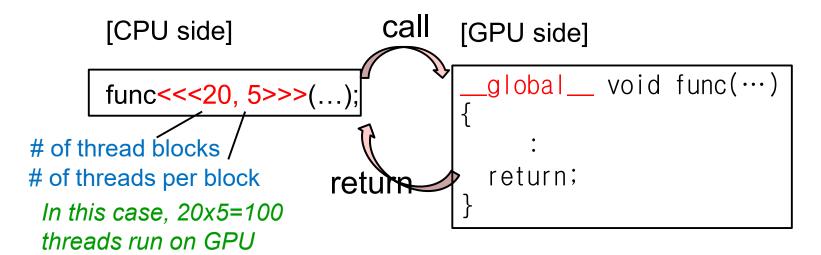
Programmer have to prepare two pointers, such as A and DA

```
int A[100];
int *DA;
cudaMalloc(&DA, ...);
cudaMemcpy(DA, A, ..., ...);
// Here CPU cannot access DA[i]
func<<<..., ...>>>(DA, ...);
22
```

# Calling A GPU Kernel Function from CPU



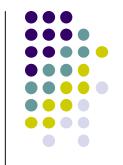
- A region executed by GPU must be a distinct function
  - called a GPU kernel function

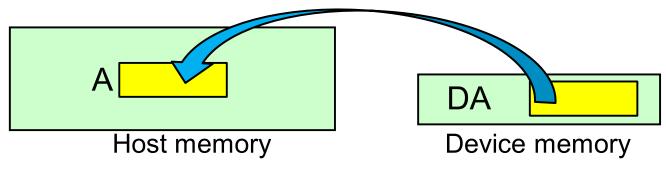


A GPU kernel function (called from CPU)

- needs \_\_global\_\_ keyword
- can take parameters
- can NOT return value; return type must be void



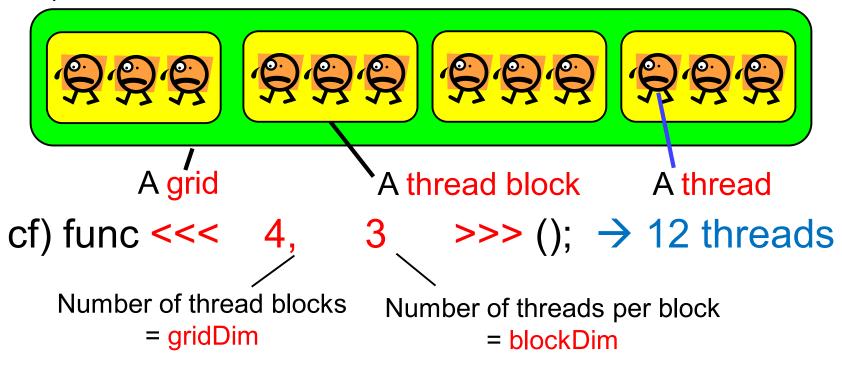




- Copy data using cudaMemcpy
  - cf) cudaMemcpy(A, DA, size, cudaMemcpyDefault);
  - 4<sup>th</sup> argument is one of
    - cudaMemcpyHostToDevice, cudaMemcpyDeviceToHost
    - cudaMemcpyDeviceToDevice, cudaMemcpyHostToHost
    - cudaMemcpyDefault ← Detect memory type automatically ☺
- When a memory area is unnecessary, free it
  - cf) cudaFree(DA);

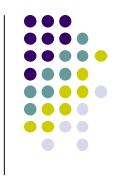
#### Threads in CUDA

When calling a GPU kernel function, specify 2 numbers (at least) for number of threads



The reason is related to GPU hardware
Thread block ⇔ SMX, Thread ⇔ CUDA core

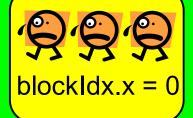
#### To See Who am I



- By reading the following special variables, each thread can see its thread ID in GPU kernel function
- My ID
  - blockldx.x: Index of the block the thread belong to (≥0)
  - threadIdx.x: Index of the thread (inside the block) (≥0)
- Number of thread/blocks
  - gridDim.x: How many blocks are running
  - blockDim.x: How many threads (per block) are running

### Thread Block ID, Thread ID

blockldx.x = 1threadIdx.x = 0 blockldx.x = 2threadIdx.x = 2





blockldx.x = 1



blockldx.x = 2



blockldx.x = 3

A grid

A thread block

A thread

For every thread, gridDim.x = 4, blockDim.x = 3

Note: In order to see the entire sequential ID, we should compute blockldx.x \* blockDim.x + threadldx.x

### The Case of add-cuda Sample

- /gs/hs1/tga-ppcomp/22/add-cuda
- We want to do

```
for (i = 0; i < 100; i++) { DA[i] += DB[i]; }

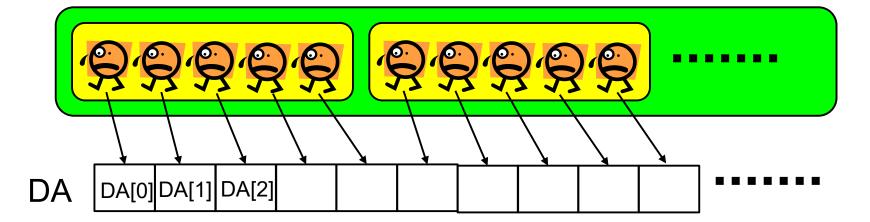
[CPU side]

[add<<<20, 5>>>(...);

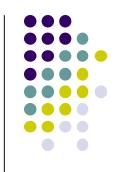
20x5=100 threads
will execute add function

[GPU side]

[global__ void add(int *DA, int *DB)
{
    int i = blockldx.x * blockDim.x + threadldx.x;
    DA[i] += DB[i];
    return;
}
```



# How is Number of Threads Determined? (1)



On CUDA, a different strategy is required from on OpenMP

- On OpenMP, number of threads (OMP\_NUM\_THREADS) should be ≤ CPU cores (or hyper threads)
  - The number is basically determined by hardware
  - $\leq$  14 on q\_node node,  $\leq$  56 on f\_node
- On CUDA, it is better to use number of thread ≥ GPU cores
  - ■ 3584 on TSUBAME3's P100 GPU
  - You can use >1,000,000 threads!

# How is Number of Threads Determined? (2)



We have to deicide 2 numbers <<<br/>block number, block size>>>

A better way would be

- (1)We decide total number of threads P
- (2)We tune each block size BS
  - Good candidates are 32, 64, ... 1024
- (3) Then block number is P/BS
  - We consider indivisible cases later

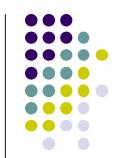


#### Comparing OpenMP/OpenACC/CUDA



	OpenMP	OpenACC	CUDA
Processors	CPU	CPU+GPU	
File extension	.C, .CC		.cu
To start parallel (GPU) region	#pragma omp parallel	#pragma acc kernels	func<<<,>>>()
To specify # of threads	export OMP_NUM _THREADS=	(num_gangs, vector_length etc)	
Desirable # of threads	# of CPU cores or less	# of GPU cores or "more"	
To get thread ID	omp_thread_num()	-	blockldx, threadldx
Parallel for loop	#pragma omp for	#pragma acc loop	-
Task parallel	#pragma omp task	-	-
To allocate device memory	-	#pragma acc data	cudaMalloc()
To copy to/from device memory	-	#pragma acc data #pragma acc update	cudaMemcpy()
Functions on GPU	-	#pragma acc routine	global,_device

# **Assignments in GPU Part** (Abstract)

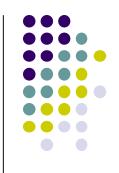


Choose one of [G1]—[G3], and submit a report

Due date: May 26 (Thursday)

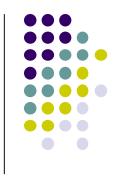
- [G1] Parallelize "diffusion" sample program by OpenACC or CUDA
- [G2] Evaluate speed of "mm-acc" or "mm-cuda" in detail
- [G3] (Freestyle) Parallelize any program by OpenACC or CUDA.

### **Notes in Report Submission (1)**



- Submit the followings via T2SCHOLA
  - (1) A report document
    - PDF, MS-Word or text file
    - 2 pages or more
    - in English or Japanese (日本語もok)
  - (2) Source code files of your program
  - Try "zip" to submit multiple files

### **Notes in Report Submission (2)**



#### The report document should include:

- Which problem you have chosen
  - Also, which you have used, OpenACC or CUDA
- How you parallelized
  - It is even better if you mention efforts for high performance or new functions
- Performance evaluation on TSUBAME
  - With varying number of threads
  - With varying problem sizes
  - Discussion with your findings
  - Other machines than TSUBAME are ok, if available



- GPU Programming (3) on May 12
  - mm sample on CUDA
- Also please note due date of OpenMP assignment is May 12