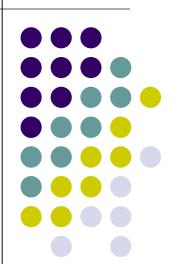
Practical Parallel Computing (実践的並列コンピューティング)

Part3: MPI (2) May 23, 2022

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- Part 0: Introduction
 - 2 classes
- Part 1: OpenMP for shared memory programming
 - 4 classes
- Part 2: GPU programming

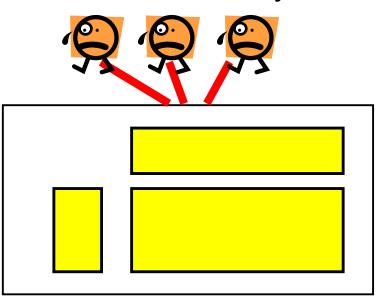
 - 4 classes
 ← We are here (1/4)
 - OpenACC (1.5 classes) and CUDA (2.5 classes)
- Part 3: MPI for distributed memory programming
 - 4 classes

← We are here (2/4)

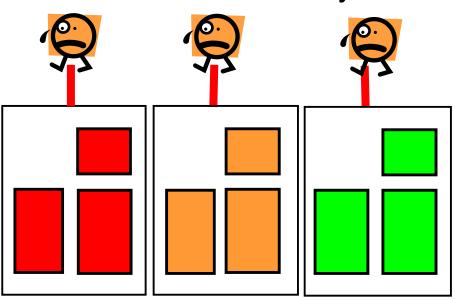
Shared Memory Model and Distributed Memory Model



Shared Memory

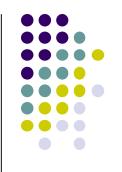


Distributed Memory



- In distributed memory model, a process CANNOT read/write other processes' memory directory
- How can a process access data on others?
- → Message passing (communication) is required





/gs/hs1/tga-ppcomp/22/test-mpi

```
[make sure that you are at a interactive node (r7i7nX)]
module load cuda openmpi [Do once after login]
cd ~/t3workspace [In web-only route]
cp -r /gs/hs1/tga-ppcomp/22/test-mpi
cd test-mpi
make
[An executable file "test" is created]
mpiexec -n 2 ./test
```

This sample is for 2 processes

Basics of Message Passing: Peer-to-peer Communication

Example: /gs/hs1/tga-ppcomp/22/test-mpi/

Rank 0 computes contents of "int a[16]"

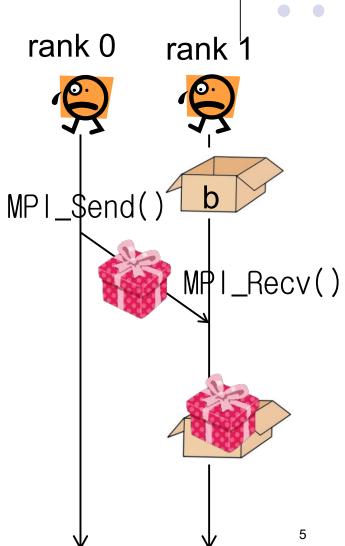
Rank 1 wants to see contents of a!

Rank0:

- Write data to an array a
- MPI_Send(a, 16, MPI_INT, 1, 100, MPI_COMM_WORLD);

Rank1:

- Prepares a memory region (array b here)
- MPI_Recv(b, 16, MPI_INT, 0, 100, MPI_COMM_WORLD, &stat);
- Now b has copy of a!



MPI_Send

```
MPI_Send(a, 16, MPI_INT, 1, 100, MPI_COMM_WORLD);
```

- a: Address of memory region to be sent
- 16: Number of data to be sent
- MPI_INT: Data type of each element
 - MPI_CHAR, MPI_LONG. MPI_DOUBLE, MPI_BYTE
- 1: Destination process of the message
- 100: An integer tag for this message (explained later)
- MPI_COMM_WORLD: Communicator (explained later)



MPI_Recv

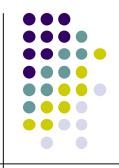


```
MPI_Status stat;
MPI_Recv(b, 16, MPI_INT, 0, 100, MPI_COMM_WORLD, &stat);
```

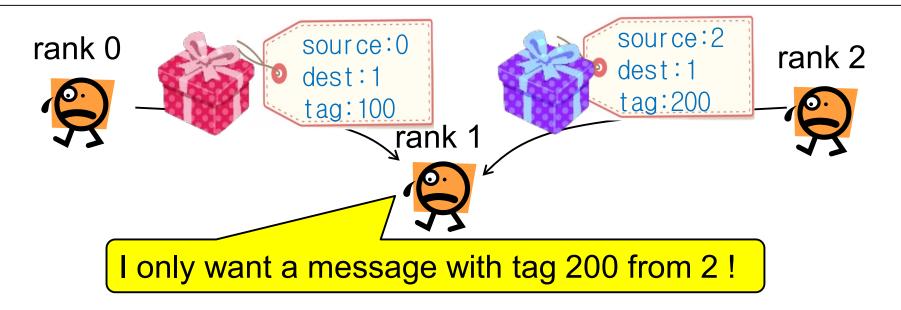
- b: Address of memory region to store incoming message
- 16: Number of data to be received
- MPI_INT: Data type of each element
- 0: Source process of the message
- 100: An integer tag for a message to be received
 - Should be same as one in MPI_Send
- MPI_COMM_WORLD: Communicator (explained later)
- &stat: Some information on the message is stored

Note: MPI_Recv does not return until the message arrives

Notes on MPI_Recv: Message Matching (1)



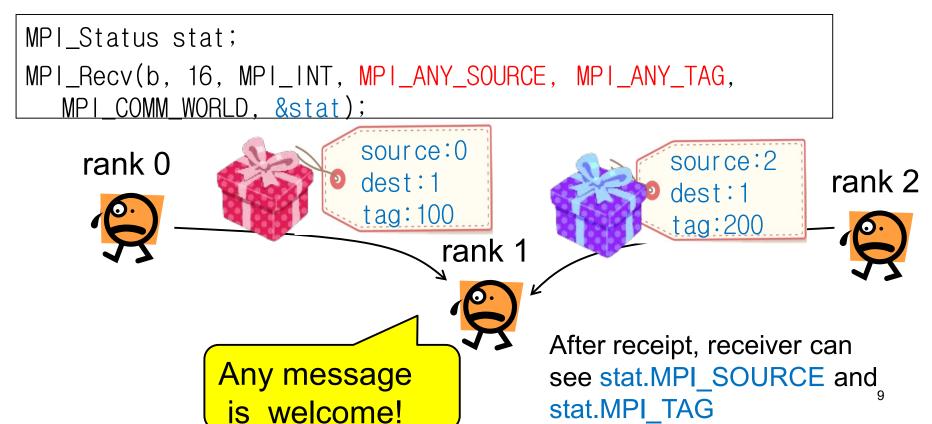
MPI_Recv(b, 16, MPI_INT, 2, 200, MPI_COMM_WORLD, &stat);



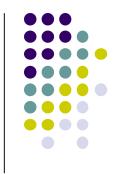
- Receiver specifies "source" and "tag" that it wants to receive
- → The message that matches the condition is delivered
- Other messages should be received by other MPI_Recv calls later

Notes on MPI_Recv: Message Matching (2)

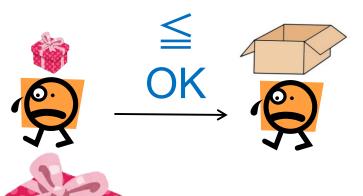
- In some algorithms, the sender may not be known beforehand
 - cf) client-server model
- For such cases, MPI_ANY_SOURCE / MPI_ANY_TAG may be useful



Notes on MPI_Recv: What If Message Size is Unmatched



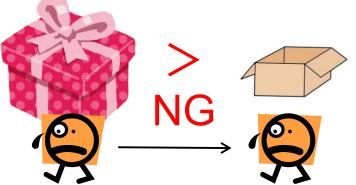
MPI_Recv(b, 16, MPI_INT, 0, 100, MPI_COMM_WORLD, &stat);



If message is smaller than expected, it's ok

→ Receiver can know the actual size by

MPI_Get_Count(&stat, MPI_INT, &s);



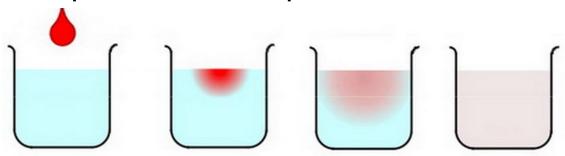
If message is larger than expected, it's an error (the program aborts)

If the message size is UNKNOWN beforehand, the receiver should prepare enough memory

Case of "diffusion" Sample related to [M1]



An example of diffusion phenomena:

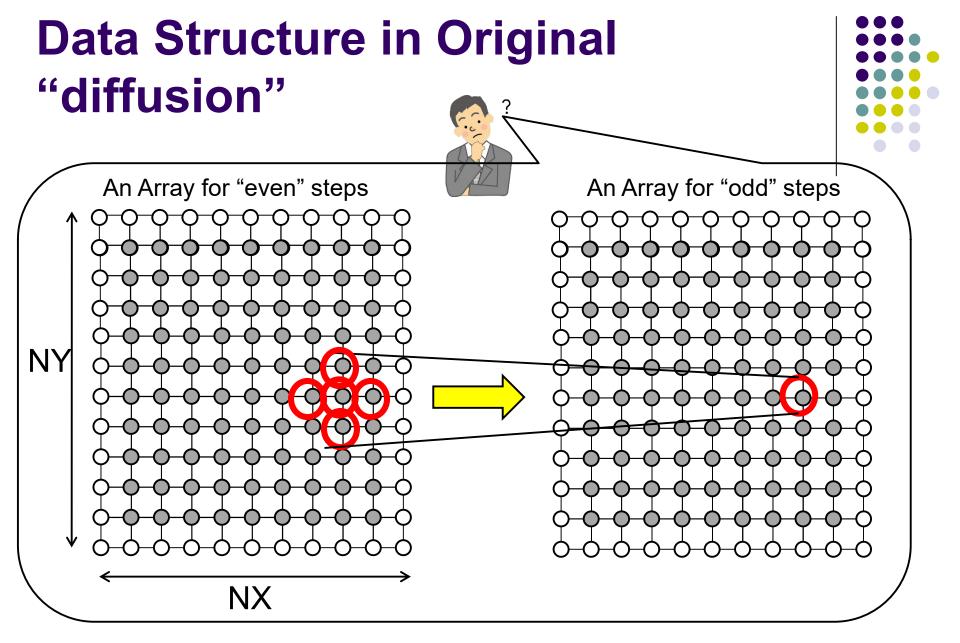


The ink spreads gradually, and finally the density becomes uniform (Figure by Prof. T. Aoki)

Available at /gs/hs1/tga-ppcomp/22/diffusion/

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

You can use /gs/hs1/tga-ppcomp/22/diffusion-mpi/as a base. Makefile uses mpicc



How Do We Parallelize "diffusion" Sample?



On OpenMP:

[Algorithm] Parallelize spatial (Y or X) for-loop

- Each thread computes its part in the space
- Time (T) loop cannot be parallelized, due to dependency

[Data] Data structure is same as original:

2 x 2D arrays → float data[2][NY][NX];

On MPI:

[Algorithm] Same as above

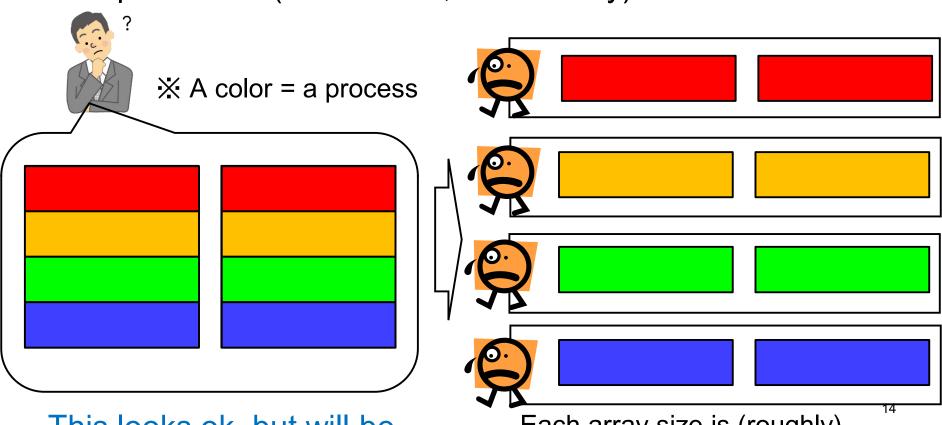
Each process computes its part in the space

[Data] 2 x 2D arrays are divided among processes

Each process has its own part of arrays



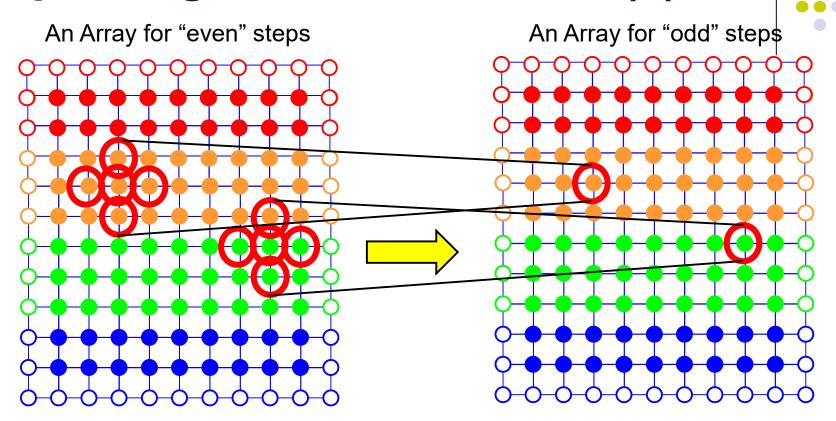
2 x 2D arrays are divided among P processes (in this case, horizontally)



This looks ok, but will be improved next

Each array size is (roughly) NX x (NY/P)

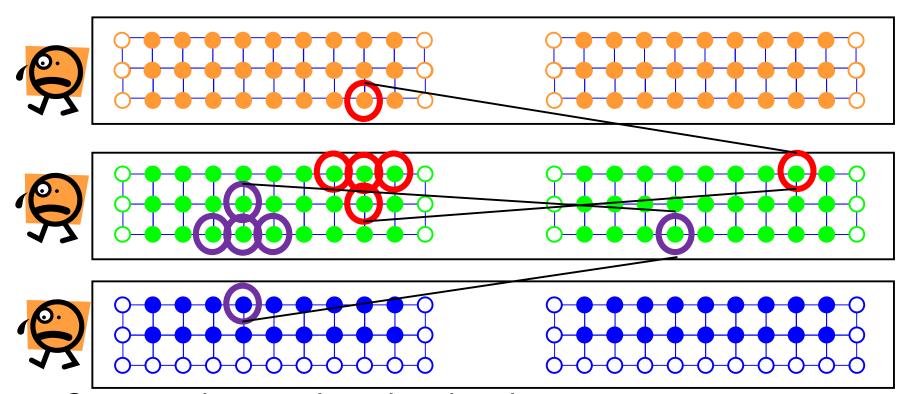
Improving Data Distribution (1)



- Let's remember computation of each point
- → 5 points are read and 1 point is written

Improving Data Distribution (2)

What's wrong with the simple distribution?



Computation requires data in other processes

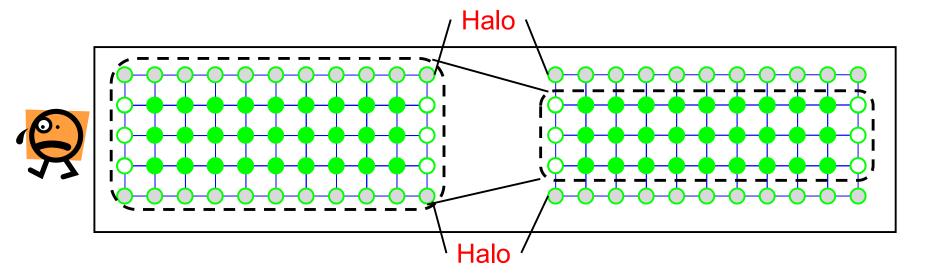
→ Message passing is required

We need memory region for received data!

A Technique in Stencil: Introducing "Halo" Region



- In stencil computation, it is a good idea to make additional rows to arrays
- → called "Halo" region



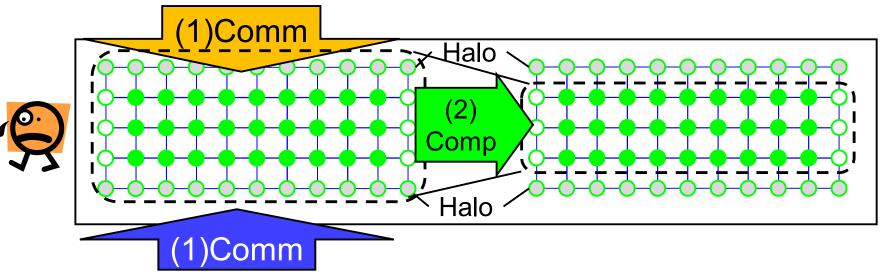
Each array size is (roughly) NX x (NY/P + 2)

Halo regions are used to receive outside border data from neighbor processes



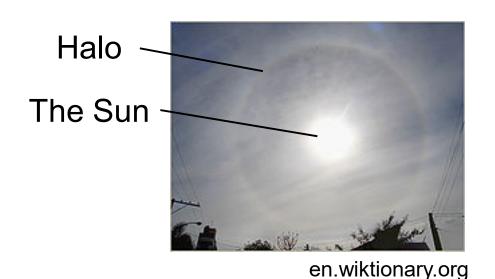
Each time step consists of:

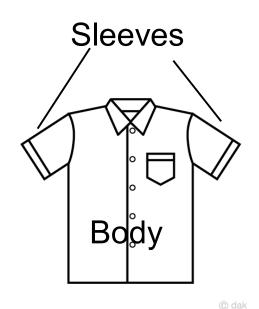
- (1) Communication: Recv data and store into "halo" region
 - Also neighbor processes need "my" data
- (2) Computation: Old data at time t (including "halo")
 - → New data at time t+1



The name of "Halo" Region

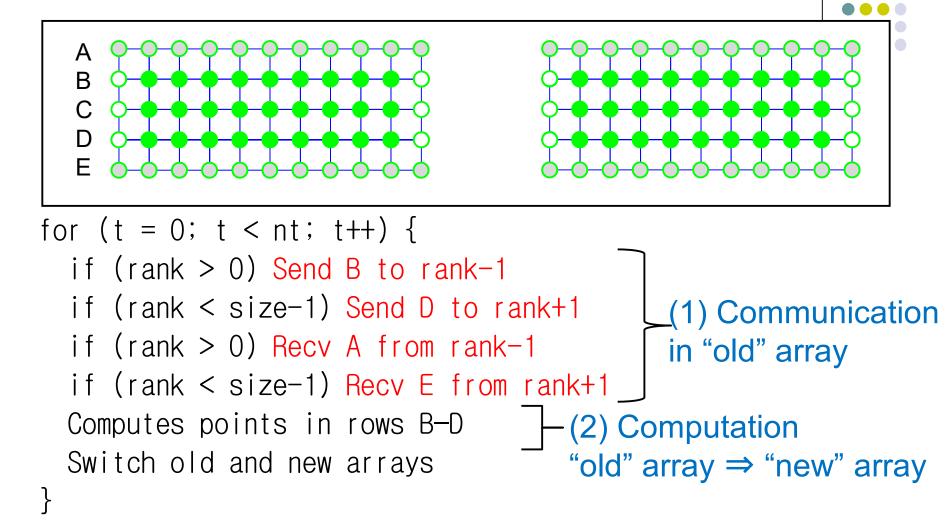






"Halo regions" are sometimes called "sleeve regions" or "overlap regions"

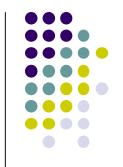
Overview of MPI "diffusion"



This version is still unsafe, for possibility of deadlock

→ Explained next

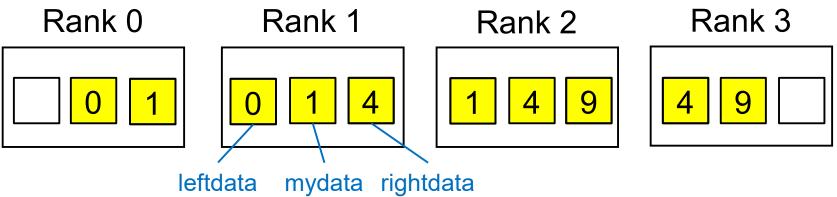
A Sample with Neighbor Communication



When considering neighbor communication, we have to avoid deadlock (a serious bug)!

A sample is available at /gs/hs1/tga-ppcomp/22/neicomm-mpi Execution: mpiexec -n [P] ./neicomm

- (1) Each process prepares its local data
- (2) Each process receives data from its neighbors, rank-1 and rank+1



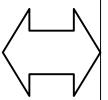
Behavior of neicomm-mpi Sample



Unsafe version ⊗

When neicomm_unsafe() is used

Send to rank-1
Send to rank+1
Recv from rank-1
Recv from rank+1



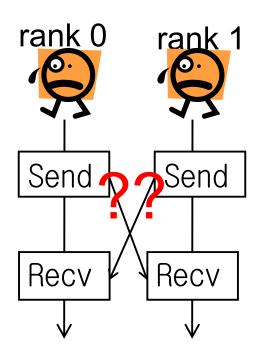
Safe version ©

When neicomm_safe() is used

Start to recv from rank-1
Start to recv from rank+1
Sent to rank-1
Sent to rank+1
Finish to recv from rank-1
Finish to recv from rank+1

Deadlock in MPI

This case "deadlocks" with 2 processes. Why?



One of reasons is MPI_Send and MPI_Recv uses blocking communication

Blocking: a process waits until "some event"

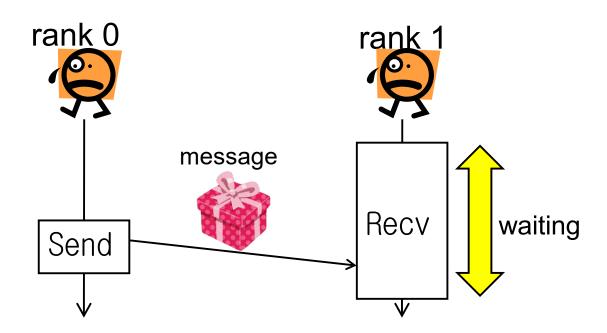




- MPI_Send is called by rank0, and MPI_Recv is called on rank1
 - Processes are running independently

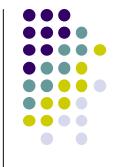
If MPI Recv is called earlier,

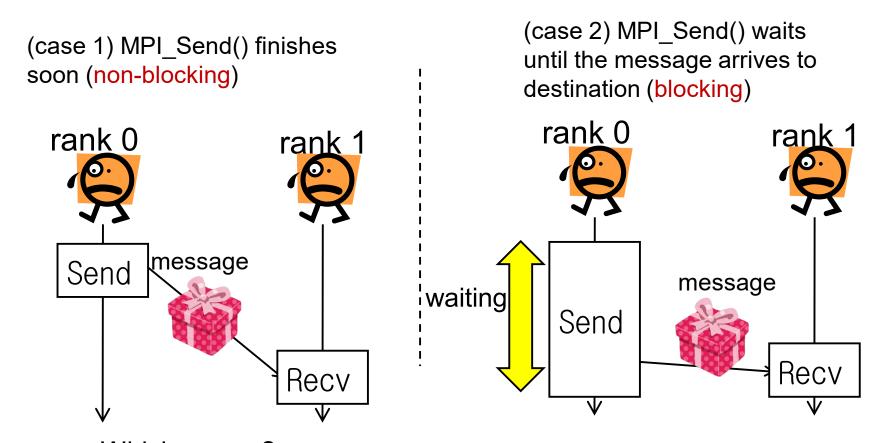
→ MPI_Recv() waits until the message arrives (blocking)



Behavior of MPI_Send()

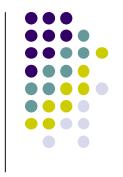
If MPI_Send is called earlier, there are two possibilities

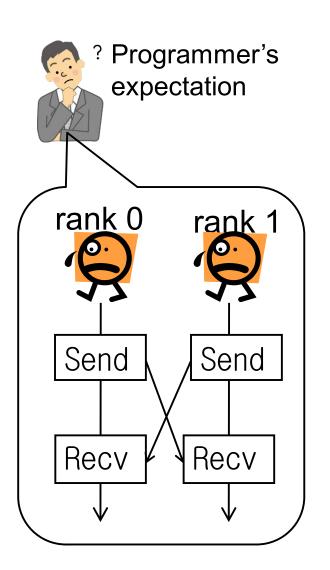




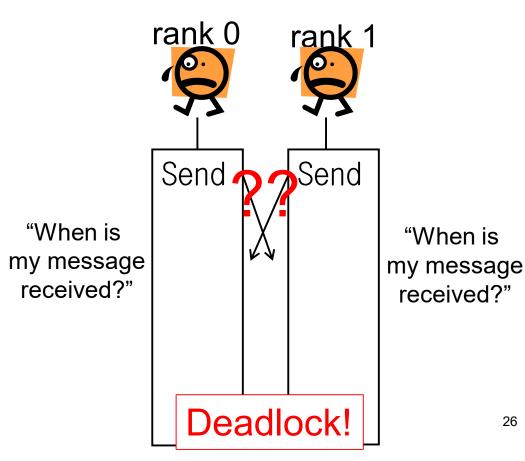
Which occurs?
It depends on MPI library, message size, etc. → Unknown

Deadlock Happens

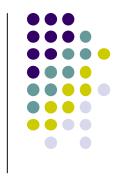




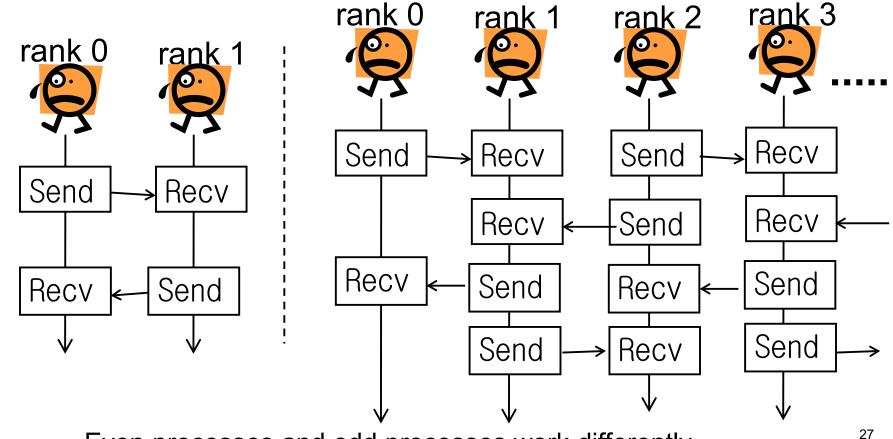
If MPI_Send is blocked until arrival in destination ...



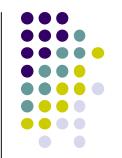
To Avoid Deadlock: An Approach



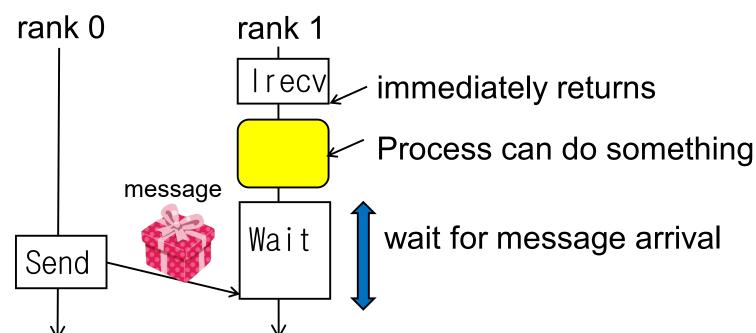
Change order of MPI_Send and MPI_Recv



Another Approach: Using Non-Blocking Communication



- Non-blocking communication: starts a communication (send or receive), but does not wait for its completion
 - MPI_Recv is blocking communication, since it waits for message arrival
- Program must wait for its completion later: MPI_Wait()





Non-Blocking Receive

```
MPI_Status stat;
MPI_Recv(buf, n, type, src, tag, comm, &stat);

MPI_Status stat;
MPI_Request req;
MPI_Irecv(buf, n, type, src, tag, comm, &req);←start recv
: (Do domething)
MPI_Wait(&req, &stat); ←wait for completion
```

MPI_Irecv: starts receiving, but it returns Immediately

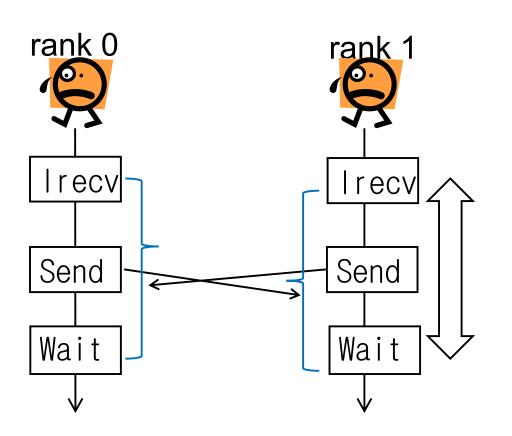
MPI_Wait: wait for message arrival

MPI Request is like a "ticket" for the communication

Algorithm Avoiding Deadlock with Non-Blocking Communication



On each process, Recv is divided into Irecv & Wait

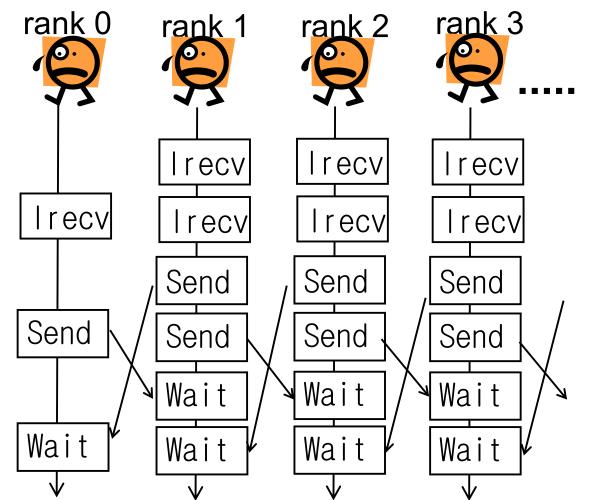


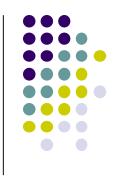
What's difference than before?

A message can be (internally) received during Irecv and Wait → MPI_Send can finish in finite time



See neicomm_safe() in neicomm.c





Each Irecv has to use distinct MPI_Request

Functions Related to Nonblocking Communication



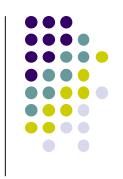
- MPI_Isend(buf, n, type, dest, tag, comm, &req); ←start send
 - MPI_Isend must be followed by MPI_Wait (or alternatives)
- MPI_Wait(&req, &stat); ←wait for completion of one communication
- MPI_Test(&req, &flag, &stat); ←check completion of one communication
- MPI_Waitall, MPI_Waitany, MPI_Testall, MPI_Testany...

Algorithm Avoiding Deadlock with Non-Blocking Communication (2)



- The following patterns are also Ok
- Each process does
 - Irecv, Irecv, Send, Send, Wait, Wait
 - neicomm_safe()
 - Isend, Isend, Recv, Recv, Wait, Wait
 - Isend, Isend, Irecv, Irecv, Wait, Wait, Wait, Wait
 - 4 MPI_Request required
 - Irecv, Irecv, Send, Send, Wait, Wait, Wait,
 - 4 MPI_Request required

Assignments in MPI Part (Abstract)



Choose <u>one of [M1]—[M3]</u>, and submit a report

Due date: June 9 (Thursday)

[M1] Parallelize "diffusion" sample program by MPI.

[M2] Improve mm-mpi sample in order to reduce memory consumption.

[M3] (Freestyle) Parallelize any program by MPI.

For more detail, please see May 19 slides

Next Class

- MPI (3)
 - Improvement of "matrix multiply" sample
 - Related to [M2]
 - Group Communication