

# Message Passing Interface (MPI)

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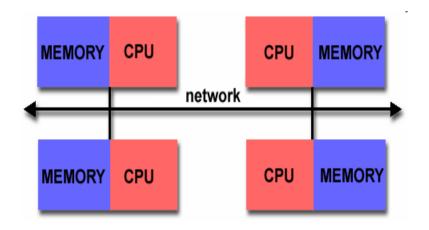
## Introduction

#### Ph.D. course on **High Performance Computing**

## **Distributed Memory**

Distributed memory requires a communication network for the information exchange.

Each processor has its own local memory. Each memory has a separate, independent address space.



Read/write operations are local, for which there are not problems of cache coherence. To allow a task to access remote data, the programmer must explicitly manage the communication among tasks.

The corresponding programming model is called **message passing**. It dictates that each task can directly access only its local memory and must use communicate with remote tasks to access remote data.



#### What is MPI?

A message-passing library specification

- message-passing model
- not a compiler specification
- not a specific product

For parallel computers, clusters, and heterogeneous networks

Full-featured

Designed to permit the development of parallel software libraries

Designed to provide access to advanced parallel hardware for

- end users
- library writers
- tool developers



#### What is MPI?

MPI is a specification for a library interface, not an implementation; there are multiple implementations of MPI.

MPI is not a language, and all MPI operations are expressed as functions, subroutines, or methods, according to the appropriate language bindings, which for C, C++, Fortran-77, and Fortran-95, are part of the MPI standard.

The standard has been defined through an open process by a community of parallel computing vendors, computer scientists, and application developers.



## MPI goals

Design an application programming interface (not necessarily for compilers or a system implementation library).

Allow efficient communication: Avoid memory-to-memory copying, allow overlap of computation and communication, and offload to communication co-processor, where available.

Allow for implementations that can be used in a heterogeneous environment.

Allow convenient C, C++, Fortran-77, and Fortran-95 bindings for the interface.

Assume a reliable communication interface: the user need not cope with communication failures. Such failures are dealt with by the underlying communication subsystem.

Define an interface that can be implemented on many vendor's platforms, with no significant changes in the underlying communication and system software.

Semantics of the interface should be language independent.

The interface should be designed to allow for thread safety.



## MPI history

Version 1.0: May, 1994

Version 1.1: June, 1995

Version 1.2: July 18, 1997

Version 2.0: July 18, 1997

Version 1.3: May 30, 2008

Version 2.1: June 23, 2008

Version 2.2: September 4, 2009

Version 3.0: September 21, 2012

Version 3.1: June 4, 2015

Version 4.0: work in progress

#### **MPI Forum**

http://www.mpi-forum.org



#### What is included in the MPI 3.0 standard?

- Point-to-point communication
- Datatypes
- Collective operations
- Process groups
- Communication contexts
- Process topologies
- Environmental Management and inquiry

- The info object
- Process creation and management
- One-sided communication
- External interfaces
- Parallel file I/O
- Language Bindings for Fortran, C and C++
- Profiling interface



#### What is new in the MPI 3.0 standard?

MPI-3.0 is a major update to the MPI standard:

- extension of collective operations to include nonblocking versions
- extensions to the one-sided operations
- a new Fortran 2008 binding

In addition, the deprecated C++ bindings have been removed, as well as many of the deprecated routines and MPI objects (such as the MPI\_UB datatype).

**MPI** 



#### What is new in the MPI 3.1 standard?

The MPI 3.1 document contains mostly corrections and clarifications to the MPI-3.0 document.

The largest change is a correction to the Fortran bindings introduced in MPI-3.0.

Additionally, new functions added include routines to manipulate MPI\_Aint values in a portable manner, nonblocking collective I/O routines, and routines to get the index value by name for MPI\_T performance and control variables.



#### What is not included in the MPI standard?

- Operations that require more operating system support than is currently standard (for example, interrupt-driven receives, remote execution, or active messages)
- Program constructions tools
- Debugging facilities



## **MPI** implementations

Several MPI implementations are available (e.g., MPICH and Open MPI).

- **MPICH** (C/C++, fully supports MPI-3) <a href="https://www.mpich.org/">https://www.mpich.org/</a>
- Open MPI (C/C++, fully supports MPI-3)
   <a href="https://www.open-mpi.org/">https://www.open-mpi.org/</a>
- **Intel MPI Library** (C/C++, fully supports MPI-3) <a href="https://software.intel.com/en-us/intel-mpi-library">https://software.intel.com/en-us/intel-mpi-library</a>



## Open MPI

With respect to MPICH, it is more efficient and portable.

#### Compiling:

mpicc -o myprog myprog.c



#### Running:

#### mpirun [-np <num>] [-hostfile <hostfile>] <execfile>

where <num> is the number of processes, <hostfile> is the name of a file that lists the IP addresses of the machines, <execfile> is the name of the program to execute



#### Hostfiles

Hostfiles are simple text files with hosts specified, one per line. Each host can also specify a default a maximum number of slots to be used on that host (i.e., the number of available processors on that host). Comments are also supported, and blank lines are ignored.

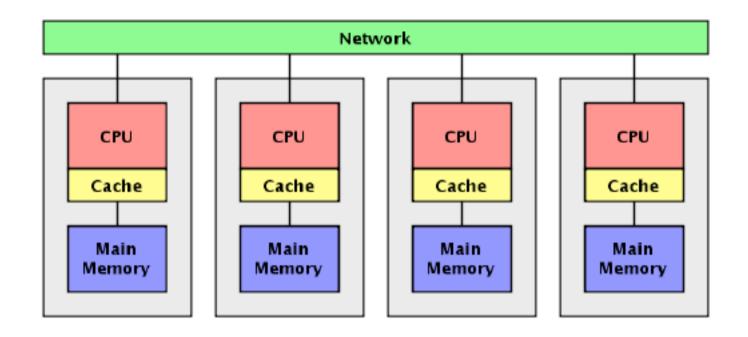
#### For example:

```
# This is an example hostfile. Comments begin with #
#
# The following node is a single processor machine:
foo.example.com
# The following node is a dual-processor machine:
bar.example.com slots=2
# The following node is a quad-processor machine, and we absolutely
# want to disallow over-subscribing it:
yow.example.com slots=4 max-slots=4
```

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## Message passing with MPI

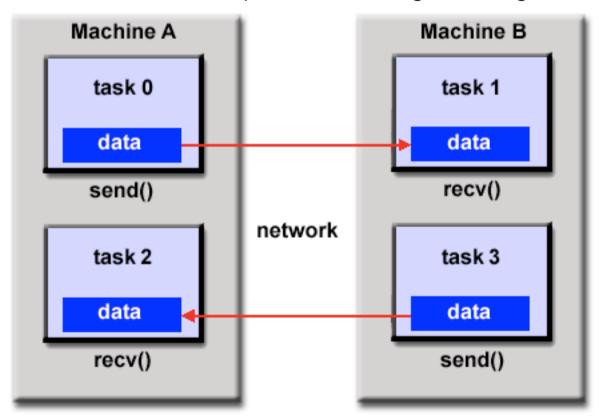


Task = Process = Instance of running program



## Message passing with MPI

Cooperation among processes is based on **explicit communication**. A **sender** process and a **receiver** process exchange messages.





## Message passing with MPI

Each process is an instance of running sub-program. Usually, the same sub-program is executed over different data sets (each execution being a process).

#### SPMD (single program, multiple data)

Each process is identified by an integer number, called *rank*, ranging from 0 to n-1, where n (size) is the total number of processes.

MPMD can be emulated with SPMD.



## Messages

Messages are composed by two parts:

#### **Envelope**

source: rank of the sender

destination: rank of the receiver

tag: ID of the message (from 0 to MPI\_TAG\_UB)

communicator: context of the communication

#### **Body**

type: MPI datatype

length: number of elements

buffer: array of elements

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## Messages

- basic datatypes
   (corresponding to standard types of C or Fortran)
- derived datatypes
   (built from basic or other derived datatypes)

## **MPI Basic Datatypes - C**

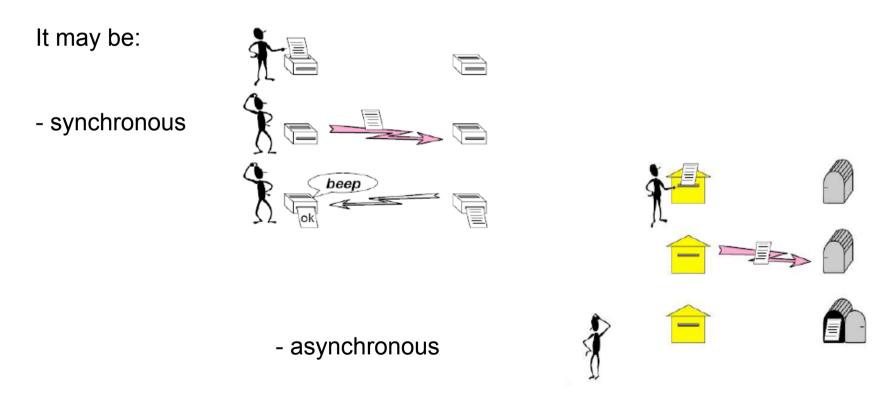
MPI Datatype	C Datatype
MPI_CHAR	signed char
MPI_SHORT	signed short int
MPI_INT	signed int
MPI_LONG	signed long int
MPI_UNSIGNED_CHAR	unsigned char
MPI_UNSIGNED_SHORT	unsigned short int
MPI_UNSIGNED	unsigned int
MPI_UNSIGNED_LONG	unsigned long int
MPI_FLOAT	float
MPI_DOUBLE	double
MPI_LONG_DOUBLE	long double
MPI_BYTE	
MPI_PACKED	

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## Point-to-point communication

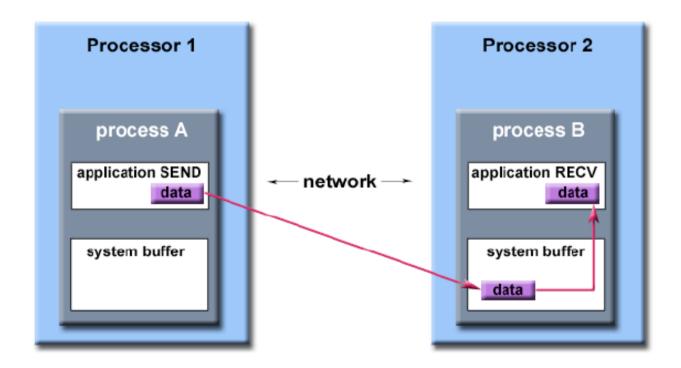
It is the most simple type of communication, involving only 2 processes (sender and receiver).





## **Buffering**

Buffering is implemented differently by each MPI library (the standard does not provides any related specification).



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## **Buffering**

Addressing spaces managed by programmers, for allocating variables, are called **application buffers**.

#### The **system buffer**:

- may exist at both sides: sender and receiver
- is usually limited
- has usually unpredictable behavior (not well documented)
- cannot be controlled by the application programmer

MPI also provides for a user managed **transfer buffer** (buffered send).

## Operation modes

Some operations may cause the blocking of the caller.

Nonblocking operations allow the process to continue, immediately after the call.

The process may **test** or **wait** for remote process completion (test), right after the nonblocking call.

Nonblocking op. + wait = blocking op.



## Operation modes

#### **Blocking**:

- synchronous send
- asynchronous
  - buffered send
  - standard send
- ready send
- recv
- sendrecv

**Non Blocking**: same primitives, but the sender never blocks - it is necessary to check when buffers are reusable, when the communication has completed, etc. using

- test
- wait



## Collective operations

This type of communication involves more than 2 processes (usually all).

**Barrier** (for synchronization)

**Data Movement** (collective communications)

- Broadcast
- Scatter
- Gather

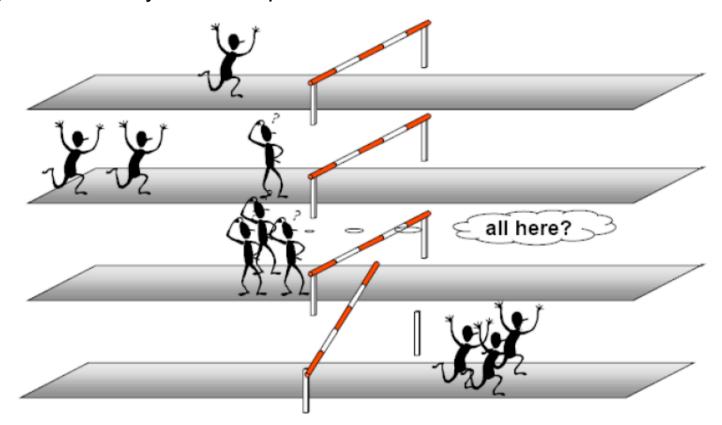
**Reduction** (collective computations)

- Minimum, Maximum
- Sum
- Logical OR, AND, etc.
- User-defined



## Collective operations

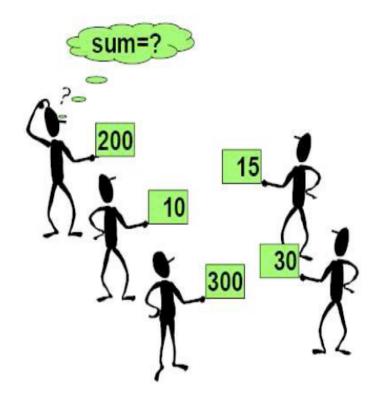
Example: using a barrier to synchronize processes





#### Reductions

Combine data from several processes to obtain a single result.





## Programming with MPI



#### **Functions in MPI**

```
error = MPI_Xxxxx(parameter, ...);
MPI_Xxxxx(parameter, ...);
```

**MPI**\_ is a reserved namespace for MPI constants and routines.

After the prefix, only the first letter is capitalized.

All MPI functions return an integer code.

Constant names are completely capitalized.

#### MPI header file

#include <mpi.h>

mpi.h is the standard header file for C. mpif.h is the standard header file for Fortran.

The header file contains definitions, macros and function prototypes that are necessary for compiling MPI programs.

MPI handles → communication

MPI datatypes → standard datatypes



#### Communicators

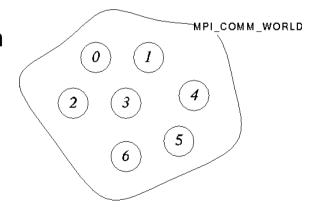
A communicator is a set of processes that can communicate with each other.

- has a name
- has a size (number of processes)
- each process can be univocally identified
- processes are equals

Two processes can communicate iff they belong to the same communicator.

The default communicator is  $MPI\_COMM\_WORLD$ , which includes n processes (having rank = 0,..,n-1).

MPI\_COMM\_WORLD is a handle defined in mpi.h





#### Rank and size

To know its rank, a process has to call:

MPI\_Comm\_rank(MPI\_Comm comm, int \*rank)

To know the size of its communicator, a process has to call:

MPI\_Comm\_size(MPI\_Comm comm, int \*size)

## Initializing and exiting MPI

The first MPI routine to be called is:

int MPI\_Init(int \*argc, char \*\*\*argv)

which initializes the default communicator (MPI\_COMM\_WORLD).

To exit from the MPI environment:

int MPI\_Finalize()



### Example: Hello, World!

Write a program that prints the rank of the process and the size of its communicator.

```
#include <stdio.h>
#include "mpi.h"

#define MASTER 0

int main (int argc, char **argv)
{
  int numtasks, taskid, len;
  char hostname[MPI_MAX_PROCESSOR_NAME];

MPI_Init(&argc, &argv);
  MPI_Comm_size(MPI_COMM_WORLD, &numtasks);
  MPI_Comm_rank(MPI_COMM_WORLD, &taskid);
  MPI_Get_processor_name(hostname, &len);
  printf ("Hello from task %d on %s!\n", taskid, hostname);

if (taskid == MASTER)
  printf("MASTER: Number of MPI tasks is: %d\n", numtasks);

MPI_Finalize();
}
```

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## Example: Hello, World!

#### Lesson learned:

- Compiling produces a unique executable
- Each command is executed by each process independently
- The runtime system (e.g., Open MPI) controls how the executable is replicated over computing nodes, how processes are created, how standard output/error are managed.



#### Point-to-Point communication

#### where:

- buf is the pointer to the message to be sent (application buffer)
- count is the number of elements in the message
- datatype specifies the type of the elements in the message
- dest is the rank of the recipient
- tag is a non-negative integer whose purpose is left to the user
- comm is the communicator



#### where:

- buf is the pointer to the array where the received message must be stored
- count is the number of elements in the message
- datatype specifies the type of the elements in the message
- source is the rank of the sender
- only messages with the specified tag are considered for reception
- comm is the communicator
- status contains info about the envelope of the message to be received



A communication is successful iff:

- sender specifies a valid receiver's rank
- receiver specifies a valid source's rank
- sender and receiver are in the same communicator
- tags match
- datatypes correspond
- the receiver has a sufficiently large buffer



### Wildcarding:

It is possible to leave the sender unspecified (in the MPI\_Recv), by means of MPI\_ANY\_SOURCE.

It is also possible to leave the tag unspecified, by means of MPI\_ANY\_TAG.

There is no wildcard for the communicator.



The envelope of a message is given by the MPI\_RECV struct, that can be retrieved from the status variable.

Other useful information are:

- Source: status.MPI\_SOURCE

- Tag: status.MPI TAG

Why retrieving source and tag from the status? In case of receive with wildcard, it is the only way to know the sender and the type of the received message.



To get the number of elements with a specific datatype in the received message, the following function must be called on status:

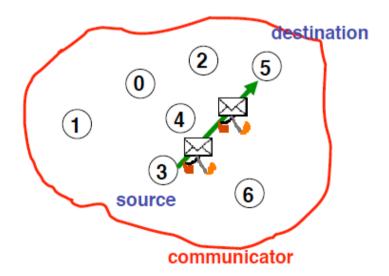
int MPI\_Get\_count(MPI\_Status status, MPI\_Datatype datatype, int \*count)

Such a number is reported in count, after the function has been executed.



### **Message order preservation:**

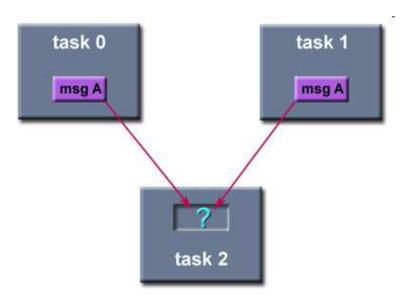
messages from the same source, in the same communicator, with the same tag, to the same receiver are delivered in the same order they have been sent.





The programmer must avoid starvation!

Process 0 and process 1 send the same message to process 2, that has set just one MPI\_Recv. Thus, only one MPI\_Send will succeed.





# Example: send/receive an integer

```
#include <stdio.h>
#include "mpi.h"
int main(int argc, char **argv)
     MPI Status status;
     int rank, size:
     /* data to communicate */
     int data int;
     /* initialize MPI */
     MPI Init(&argc, &argv);
     MPI Comm rank(MPI COMM WORLD, &rank);
     MPI Comm size(MPI COMM WORLD, &size);
     if (rank == 0) {
           data int = 10:
           MPI Send(&data_int, 1, MPI_INT, 1, 666, MPI_COMM_WORLD);
     } else if (rank == 1) {
           MPI_Recv(&data_int, 1, MPI_INT, 0, 666, MPI_COMM_WORLD, &status);
           printf("Process 1 receives %d from process 0.\n", data int);
     MPI Finalize();
     return 0;
```



# Example: send/receive an array of floats

```
#include <stdio.h>
#include "mpi.h"
#define MSIZE 10
int main(int argc, char **argv)
      MPI Status status;
      int rank, size;
      int i, j;
      /*data to communicate */
      float a[MSIZE];
      MPI Init(&argc, &argv);
      MPI Comm rank(MPI COMM WORLD, &rank);
      if (rank == 0) {
             for (i = 0; i < MSIZE; i++)
                    a[i] = (float) i;
             MPI Send(a, MSIZE, MPI FLOAT, 1, 666, MPI COMM WORLD);
      } else if (rank == 1) {
             MPI Recv(a, MSIZE, MPI FLOAT, 0, 666, MPI COMM WORLD, &status);
             printf("Process 1 receives the following array from process 0.\n");
             for (i = 0; i < MSIZE; i++)
                   printf("%6.2f\n", a[i]);
      MPI Finalize();
      return 0;
```



# Example: send/receive part of an array of floats

```
#include <stdio.h>
#include "mpi.h"
#define VSIZE 50
#define BORDER 12
int main (int argc, char *argv[]) {
      MPI Status status;
      int indx, rank, size, nprocs;
      int start send buf = BORDER;
      int start recv buf = VSIZE - BORDER;
      int length = 10:
      int vector[VSIZE];
      MPI Init(&argc, &argv);
      MPI Comm rank(MPI COMM WORLD, &rank);
      MPI_Comm_size(MPI_COMM_WORLD, &size);
      /* all processes initialize vector */
      for (indx=0; indx<VSIZE; indx++) vector[indx] = indx;
      /* send length integers starting from the "start send buf"-th position of vector */
      if (rank == 0) {
             MPI Send(&vector[start send buf], length, MPI INT, 1, 666, MPI COMM WORLD);
      /* receive length integers in the "start_recv_buf"-th position of vector */
      if (rank == 1) {
             MPI_Recv(&vector[start_recv_buf], length, MPI_INT, 0, 666, MPI_COMM_WORLD, &status);
      MPI Finalize();
      return 0:
```

# Measuring execution time

double MPI\_Wtime(void);

Standard timers (e.g., POSIX) are not adequate for MPI programs:

- insufficient accuracy
- not portable

The execution time of a task is measured by getting the current time before the execution starts, and after.

MPI\_Wtime() returns the local time, measured from a prefixed time.

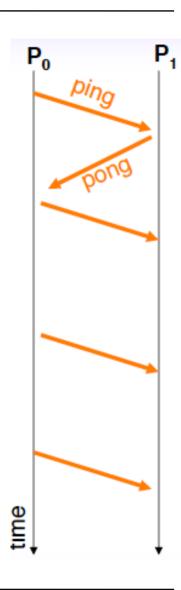
To know the resolution of the timer:

double MPI\_Wtick(void);



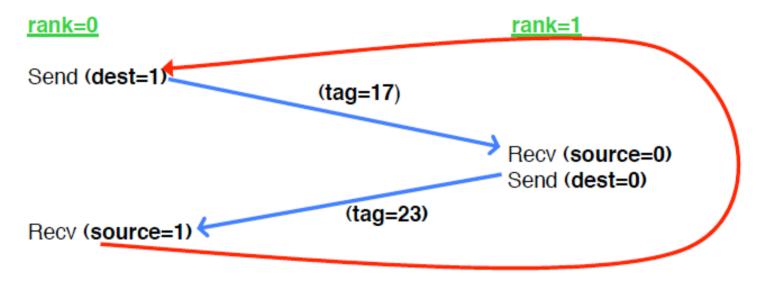
Write a program where two processes send to each other a message, according to the scheme in figure.

- process 0 sends the message to process 1 (PING)
- after receiving the message, process 1 sends it back to process 0 (PONG)
- etc. (repeat 50 times)



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# **Example: PING-PONG**





```
#include <stdio.h>
#include <mpi.h>
#define proc A 0
#define proc_B 1
#define ping 100
#define pong 101
#define number_of_messages 50
#define length_of_message 1
int main(int argc, char *argv[])
    int my rank;
    float buffer[length_of_message];
    int i;
    double start, finish, time;
    MPI Status status;
    MPI_Init(&argc, &argv);
    MPI Comm rank(MPI COMM WORLD, &my rank);
```

**MPI** 



```
if (my_rank == proc_A)
{
    start = MPI_Wtime();
    for (i = 1; i <= number_of_messages; i++)
    {
        MPI_Ssend(buffer, length_of_message, MPI_FLOAT, proc_B, ping, MPI_COMM_WORLD);
        MPI_Recv(buffer, length_of_message, MPI_FLOAT, proc_B, pong, MPI_COMM_WORLD, &status);
    }
    finish = MPI_Wtime();
    time = finish - start;
    printf("Time for one message: %f seconds.\n", (float)(time / (2 * number_of_messages)));
}</pre>
```





# Completed communications

A communication is **locally completed** on a process if the latter has completed its part of operations related to communication (the last one being the emptying of the exit buffer).

With respect to execution, local completion means that the process can execute the instruction that follows the SEND or RECV.

A communication is **globally completed** if all involved processes have completed their operations related to the communication.

A communication is globally completed if and only if it is locally completed on all involved processes.



# Completed communications

The completion phase of the SEND function depends on the message size:

- Buffered for small size
- Synchronous for large size

In the first case the process can exit from the SEND after the message has been copied in a system buffer.

In the second case the process can exit from the SEND only when a RECV function is being executed.

### Consider the following case:

```
if (myRank = 0)
SEND A to Process 1
RECV B from Process 1
else if (myRank = 1)
SEND B to Process 0
RECV A from Process 0
endif
```

There are two SEND that wait, and two RECV that can be executed only after the related SEND are completed.



# Blocking and non blocking communication modes

### Blocking:

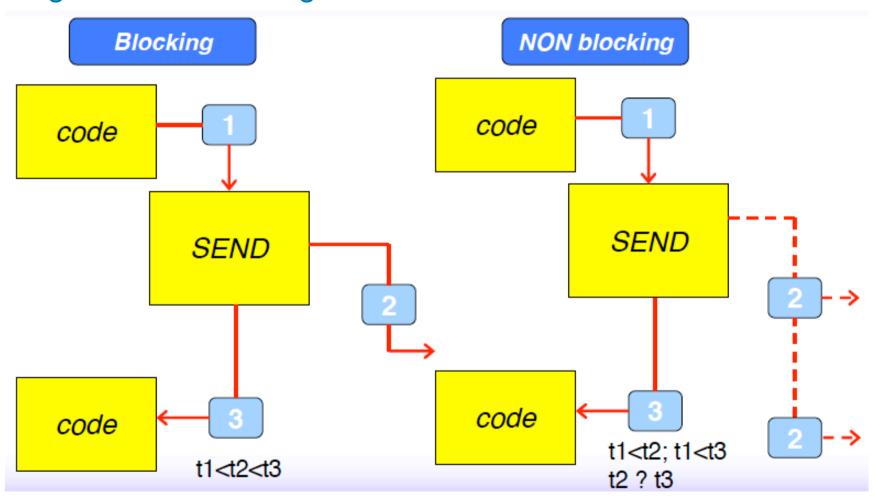
 control is returned to the process that has invoked the communication primitive only when the latter has completed

### Nonblocking:

- control is immediately returned to the process that has invoked the communication primitive
- a check on the actual completion of the communication must be performed afterwards
- meanwhile, the process can perform other operations

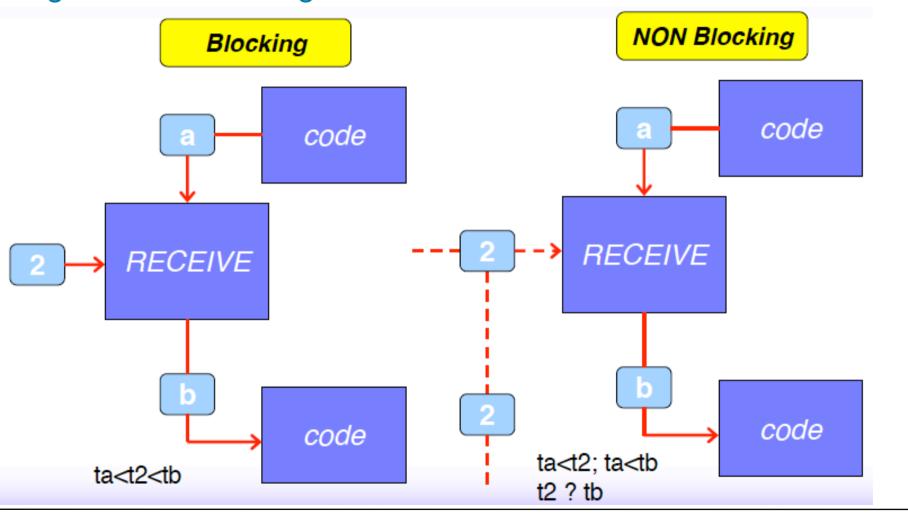


# Blocking and non blocking communication modes





# Blocking and non blocking communication modes





# Communication completion criteria

### Synchronous send

→ completed when the application buffer can be reused and the message reception has started

#### Buffered send

→ completed when the message has been completely copied to the transfer buffer

#### Standard send

→ completed when the application buffer can be reused

#### Reiceve

→ completed when the message has arrived

the I stands for "immediate"



### Communication functions

### Blocking:

- Synchronous send → MPI\_Ssend
- Buffered send → MPI\_Bsend
- Standard send → MPI\_Send
- Ready send → MPI\_Rsend
- Receive → MPI\_Recv

### Nonblocking:

- Synchronous non blocking send → MPI\_Issend
- Buffered non blocking send → MPI\_Ibsend
- Standard non blocking send → MPI\_Isend
- Ready non blocking send → MPI\_Irsend
- Non blocking receive → MPI\_Irecv

To check for completion → MPI\_Test, MPI\_Wait



# Synchronous send: MPI\_Ssend

Sends the message and stays blocked until the application buffer of the sender is ready for being used for another operation **and** the receiver has started to receive the message.

Arguments are the same of the MPI\_Send

**Pros:** This send is more secure than the standard one, because the network is not overloaded with pending messages. Sender and receiver are always synchronized (the resulting parallel program is more deterministic).

Cons: Risk of deadlocks. Risk of idle time.



# Buffered send: MPI\_Bsend

It is immediately completed as soon as the process has copied the message on a transfer buffer that must be explicitly managed by the program:

### MPI\_Buffer\_attach(void \*buf, int size)

allocates a memory area as transfer buffer

### MPI\_Buffer\_detach(void \*buf, int \*size)

deallocates the memory area of the transfer buffer

Sender and receiver are not synchronized.

- buf is the memory address of the buffer
- size is an int (INTEGER) that defines the size IN BYTES of the buffer

Note: the transfer buffer is a user-managed alternative to the system buffer, which is MPI-managed.



# Standard send: MPI\_Send

Base blocking send operation: completed when the message has been sent and the application buffer can be reused.

The message can stay in the communication network for a while.

### Ready send: MPI Rsend

A send that uses the ready communication mode may be started only if the matching receive is already posted. Otherwise, the operation is erroneous and its outcome is undefined.

It is better to avoid it, unless you are 200% sure that the related receive has been set.



# Send-receive: MPI\_Sendrecv

Executes a blocking send and receive operation. Both send and receive use the same communicator, but possibly different tags. The send buffer and receive buffers must be disjoint, and may have different lengths and datatypes.

int MPI\_Sendrecv(void \*sbuf, int scount, MPI\_Datatype s\_dtype, int dest, int stag, void \*dbuf, int dcount, MPI\_Datatype d\_type, int src, int dtag, MPI\_Comm comm, MPI\_Status \*status)

The first half of the list of arguments is related to the send, the second half to the receive.

- [IN] **dest** is the *rank* of the *receiver* within the communicator **comm**
- [IN] stag is the identifier of the sent message
- [IN] **src** is the *rank* of the *sender* within the communicator **comm**
- [IN] dtag is the identifier of the received message



# Nonblocking communications

A nonblocking communication is tipically made by three phases:

- 1. the beginning of the send/receive of the message;
- 2. the execution of an activity that does not imply the access to the data involved in the communication;
- 3. the wait for the completion of the communication;

#### Pros:

Performance: a nonblocking communication allows one to:

- overlap communication phases with computing phases;
- reduce the effects of the communication latency.

Non blocking communications avoid deadlock situations.

#### Cons:

Programming nonblocking message passing operations is more complicate.



# Synchronous nonblocking send

MPI\_Issend(buf, count, datatype, dest, tag, comm, &request\_handle)



MPI\_Wait(&request\_handle, &status)

- buf cannot be used between Issend and Wait
- Issend immediately followed by Wait is equivalent to Ssend
- status is not used in the Issend, but in the Wait
- request\_handle allows the Wait to refer to a precise send
- wait and test allow to know when the receiver has received the message

NOTE: "I" stands for "immediate" (because the function returns the control to the process immediately)



# Standard nonblocking send

MPI\_Isend(void \*buf, int count, MPI\_Datatype datatype,int dest,int tag, MPI\_Comm comm, MPI\_Request \*request)

- Initialize the send operation
- Identifies a memory area that must be used as a buffer; the execution goes on without waiting that the message is being copied by the application buffer
- a request handle is returned to manage the state of the pending message
- a program cannot modify the application buffer until a subsequent call to MPI\_Wait or MPI\_Test indicate that the send has completed
- may be related to a blocking receive
- improves the overlapping degree between computing and communication



# Nonblocking receive

# int MPI\_Irecv(void buf, int count, MPI\_Datatype datatype, int source, int tag, MPI\_Comm comm, MPI\_Request request)

- Initializes the receive operation
- Identifies a memory area as receive buffer
- The function returns immediately, not waiting for the message to be fully copied in the local application buffer
- buf cannot be used while the request is pending (use MPI\_Wait or MPI\_Test to know when the request has been fulfilled)
- to check for the status of the receive request, use the request handle that is returned (the blocking MPI\_Recv has the status variable, instead)



# Completion test

When using point-to-point nonblocking communications, it is fundamental to check for the completion of the communication, before

- using the receive buffer
- using the send buffer

MPI provides two types of check operations:

- WAIT allows to stop the execution of the process until the communication completes
- *TEST* returns to the calling process a value that is TRUE if the communication has completed, FALSE otherwise

MPI\_Wait(MPI\_request request, MPI\_Status status)

MPI\_Test(MPI\_Request request, int flag, MPI\_Status status)

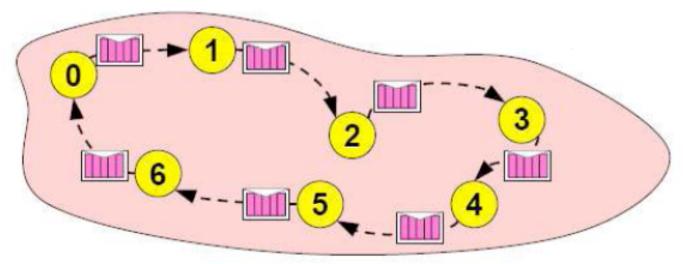


### **Deadlocks**

Consider the following code:

```
MPI_Ssend(..., right_rank, ...)
MPI_Recv(..., left_rank, ...)
```

The Ssend does not complete until the corresponding Recv has been started, but since all processes are blocked on SSend, no one starts the Recv.





### **Deadlocks**

How to avoid deadlocks:

### 1) change the order of the calls

- dangerous if dependencies involve more than 2 processes
- require clear knowledge of the communication pattern (it must be deterministic)
- does not add complexity to the program

### 2) use non blocking operations

- adds complexity to the program
- may improve efficiency

### 3) use buffered modes

- buffer must be carefully managed
- does not add too much complexity to the program
- adds determinism to the program



# Multiple non blocking communications

It is possible that many non blocking communications are posted at the same time, for which MPI provides operations for simultaneously checking their completion:

a) Wait or test for completing **one** message in a list: MPI\_Waitany / MPI\_Testany

b) Wait or test for completing **all** messages: MPI\_Waitall / MPI\_Testall

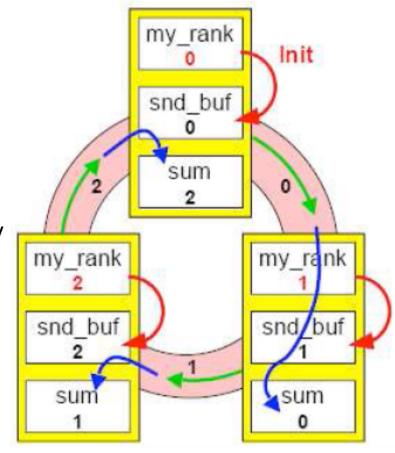
c) Wait or test for completin **some** messages: MPI\_Waitsome / MPI\_Testsome



# Example - propagate message on a ring

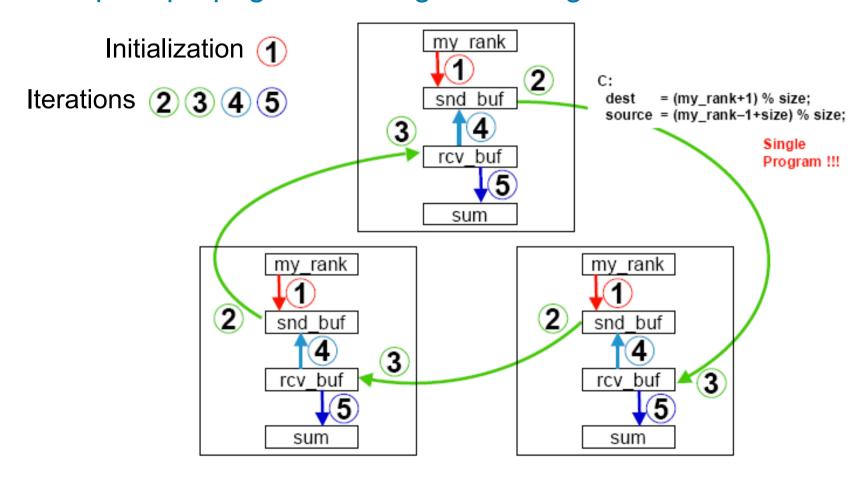
Consider a set of N processes organized as a ring.

- Initially, every process saves its MPI\_COMM\_WORLD rank in a variable called snd\_buf.
- Then, every process repeats N times:
  - send the snd\_buf value to the neighbor on the right;
  - receive a value, add it to the sum variable, then copy it to snd\_buf
- MPI\_Issend (nonblocking) is used
- to avoid deadlocks;
- to check the correctness, because the syncronous send would cause a deadlock.





#### Example - propagate message on a ring





# Example - propagate message on a ring

```
#include <stdio.h>
#include <mpi.h>
#define to_right 201
int main (int argc, char *argv[])
    int my rank, size;
    int snd buf, recv buf;
    int right, left;
    int sum, i;
    MPI Status status;
    MPI Request request;
    MPI_Init(&argc, &argv);
    MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);
    MPI_Comm_size(MPI_COMM_WORLD, &size);
    right = (my rank+1) % size;
    left = (my rank-1+size) % size;
```

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# Example - propagate message on a ring

```
/* ... this SPMD-style neighbor computation with modulo has the same meaning as: */
/* right = my rank + 1; */
/* if (right == size) right = 0; */
/* left = my rank - 1; */
/* if (left == -1) left = size-1;*/
sum = 0:
snd buf = my rank;
for( i = 0; i < size; i++)
     MPI Issend(&snd buf, 1, MPI INT, right, to right, MPI COMM WORLD, &request);
     MPI_Recv(&recv_buf, 1, MPI_INT, left, to_right, MPI_COMM_WORLD, &status);
     MPI_Wait(&request, &status);
     sum += recv buf;
     snd buf = recv buf;
printf ("PE%i:\tSum = %i\n", my rank, sum);
MPI Finalize();
```

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Derived datatypes are used for **polymorphic messages**, i.e., messages that are made of compound items comprising various generic datatypes, e.g., integers, floats, and characters (all in a single data block).

Derived datatypes are also used to transfer data from non-contiguous buffers, e.g., a portion of a matrix.

To this effect MPI lets programmers specify mixed and non-contiguous communication buffers, so that objects of various shapes and sizes can be transferred directly without copying.

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Problem: to specify non-contiguous data of a single type, or contiguous data of mixed types, or non-contiguous data of mixed types.

A few possible solutions are that you could:

- make multiple MPI calls to send and receive each data element
- use MPI\_Pack to combine different datatypes into contiguous memory for sending and MPI\_Unpack to unpack the data back into non-contiguous memory after being received (datatype MPI\_PACKED)
- use MPI\_BYTE to get around the datatype-matching rules. MPI\_BYTE can be used to match any byte of storage (on a byte-addressable machine), irrespective of the datatype of the variable that contains this byte.

Generally, however, these solutions are slow, clumsy, and wasteful of memory. Using MPI\_BYTE or MPI\_PACKED might also result in a program that isn't portable to a heterogeneous system of machines. The idea of MPI derived datatypes is to provide a portable and efficient way of communicating non-contiguous or mixed types in a message. MPI derived datatypes provide a simpler, cleaner, more elegant and efficient way to handle this type of data.

Derived datatypes are created at run-time before being used in a communication, by means of the appropriate constructor, which populates a datatype descriptor (whose type is MPI\_Datatype):

- int MPI\_Type\_contiguous(.., MPI\_Datatype \*newtype)
- int MPI\_Type\_vector(.., MPI\_Datatype \*newtype)
- int MPI\_Type\_struct(.., MPI\_Datatype \*newtype)

A created datatype must be committed, using int MPI\_Type\_commit(MPI\_Datatype \*newtype)

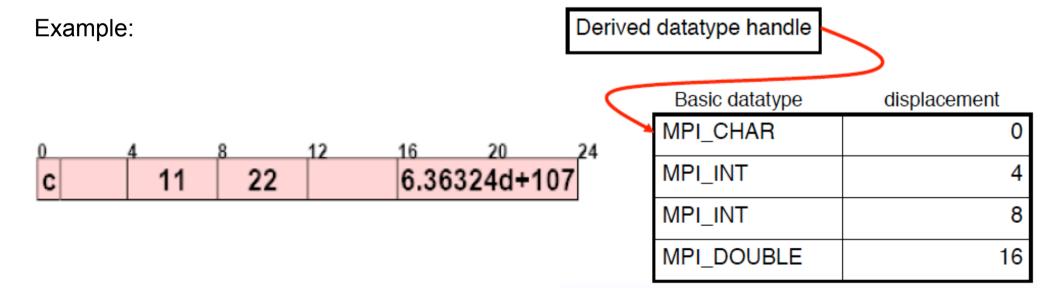
After being used, derived datatypes mus be deallocated, using int MPI\_Type\_free(MPI\_Datatype \*newtype)

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# Derived datatypes

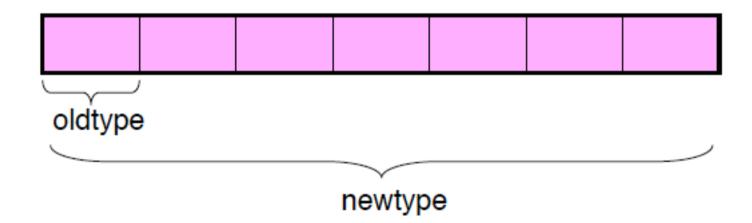
Type map of a derived datatype:

basic datatype 0	displacement of datatype 0
basic datatype 1	displacement of datatype 1
basic datatype n-1	displacement of datatype n-1





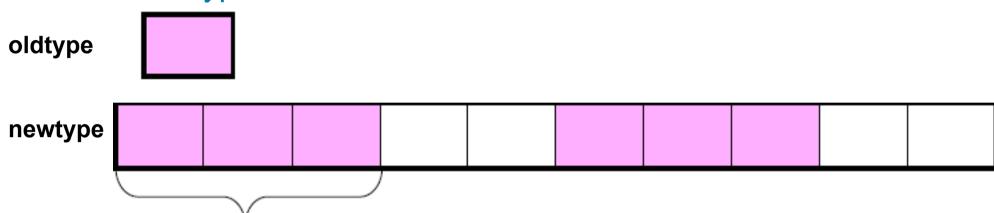
The most simple derived datatype is a set of count elements of the same type:



int MPI\_Type\_contiguous(int count, MPI\_Datatype \*oldtype, MPI\_Datatype \*newtype)

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# **Derived datatypes**



blocklength = 3 elements per block

stride = 5 (step between two blocks of elements)

count = 2 (blocks)

int MPI\_Type\_vector(int count, int blocklength, int stride, MPI\_Datatype \*oldtype, MPI\_Datatype \*newtype)

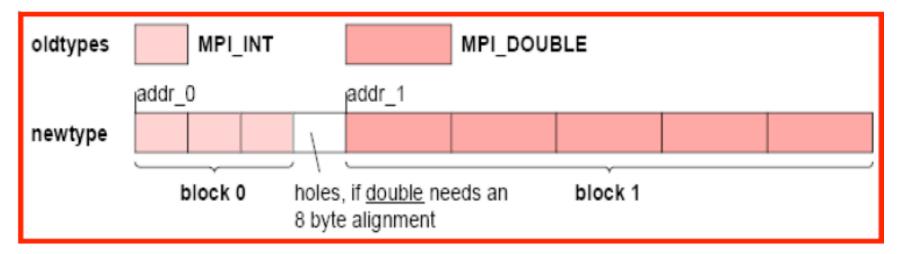


Type struct:

```
count: number of blocks (integer) - also number of entries in arrays array_of_types, array_of_displacements and array_of_blocklengths array_of_blocklengths: number of elements in each block (array) array_of_displacements: byte displacement of each block (array) array_of_types: type of elements in each block (array of handles to datatype objects) newtype: new datatype (handle)
```



#### Type struct example:



```
count = 2

array_of_blocklengths = (3, 5 )

array_of_displacements = (0, addr_1 - addr_0)

array_of_types = (MPI_INT, MPI_DOUBLE )
```



struct buff\_layout array of types[0]=MPI INT; Data layout and buffer: array of blocklengths[0]=3; { int i\_val[3]; array of displacements[0]=0; array\_of\_types[1]=MPI\_DOUBLE; double d\_val[5]; array of blocklengths[1]=5; buffer: array\_of\_displacements[1]=...; MPI Type struct(2, array of blocklengths, array of displacements, array of types, &buff datatype); Compiler MPI Type commit(&buff datatype); MPI\_Send(&buffer, 1, buff\_datatype, ...) &buffer = the start the datatype handle address of the data describes the data layout int double



Memory layout of the struct datatype in the previous example:

```
Fixed memory layout:
```

```
struct buff_layout
{
    int i_val[3];
    double d_val[5];
}
```

Arbitrary memory layout:

- Each array is allocated independently
- Each buffer is a copy of a 3-int-array and of a 5-double-array
- For each buffer it is necessary a specific datatype handle



To compute a displacement:

array\_of\_displacements[i] := address(block\_i) - address(block\_0)

where address (of a location in memory) is obtained by means of

int MPI\_Get\_address(void \*location, MPI\_Aint \*address)

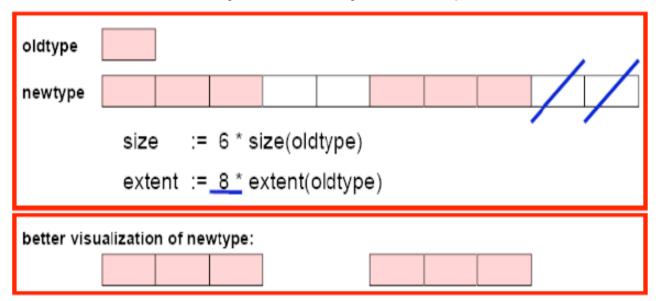


Size and extent of a datatype:

**Size** = number of bytes that must be transferred **Extent** = interval between the first and the last byte

For base types: Size = Extent = number of bytes used by the compiler

For derived datatypes:





Size and extent of a datatype:

int MPI\_Type\_size(MPI\_Datatype datatype, int \*size)

int MPI\_Type\_get\_extent (MPI\_Datatype datatype, MPI\_Aint \*ab, MPI\_Aint \*extent)



# Derived datatypes - Example of contiguous data

- create a datatype that represents the row of a matrix
- send a different row to each process

```
#include <stdio.h>
#include "mpi.h"
#define SIZF 4
int main(argc,argv)
int argc;
char *argv[]; {
     int numtasks, rank, source=0, dest, tag=1, i;
     float a[SIZE][SIZE] = {1.0, 2.0, 3.0, 4.0, // matrix
     5.0, 6.0, 7.0, 8.0,
     9.0, 10.0, 11.0, 12.0,
     13.0, 14.0, 15.0, 16.0};
     float b[SIZE];
                    // array
     MPI Status stat;
     MPI Datatype rowtype:
     MPI Init(&argc,&argv);
```



# Derived datatypes - Example of contiguous data

```
MPI Comm rank(MPI COMM WORLD, &rank);
MPI_Comm_size(MPI_COMM_WORLD, &numtasks);
MPI Type contiguous(SIZE, MPI_FLOAT, &rowtype);
MPI Type commit(&rowtype);
if (numtasks == SIZE) {
if (rank == 0) { // master
  for (i=0; i<numtasks; i++)
   MPI Send(&a[i][0], 1, rowtype, i, tag,
  MPI COMM WORLD):
 MPI Recv(b, SIZE, MPI FLOAT, source, tag, MPI COMM WORLD, &stat); // source is 0 (master)
 printf("rank = %d, b = %3.1f %3.1f %3.1f %3.1f\n", rank,b[0],b[1],b[2],b[3]);
else
    printf("Must specify %d processors. Terminating.\n",SIZE);
MPI_Type_free(&rowtype);
MPI Finalize();
```

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#### Collective communications

MPI provides function that implement communication patterns involving more processes, thus avoiding the programmer to implement them by combining point-to-point communications (which is ennoying and error-prone).

Three classes: all-to-one, one-to-all and all-to-all.

#### Collective operations:

- all processes must communicate, i.e., call the collective routine with same arguments
- blocking and nonblocking (MPI 3.0)
- no tags
- receive buffers must fit exactly the message size



#### **Barrier synchronization**:

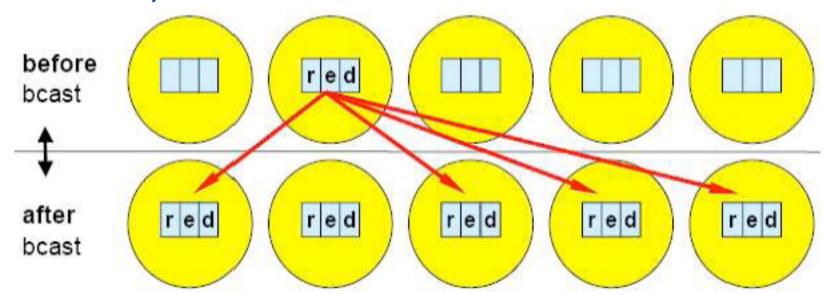
int MPI\_Barrier(MPI\_Comm comm)

Usually it is not necessary, because processes already synchronize with communication (a process cannot advance without all needed data).



#### **Broadcast**:

int MPI\_Bcast(void \*buffer, int count, MPI\_Datatype datatype, int root, MPI\_Comm comm)





#### **Broadcast:**

- replicates the content of the root's buffer to all other processes within the communicator
- does not imply synchronization, necessarily

#### Parameters:

- buffer (in/out) address of the send/receive buffer
- count (in) number of elements forming the data to be sent
- datatype (in) type of the data to be sent
- root (in) rank of the sender
- comm (in) name of the comunicator

Broadcast belongs to the one-to-all class.



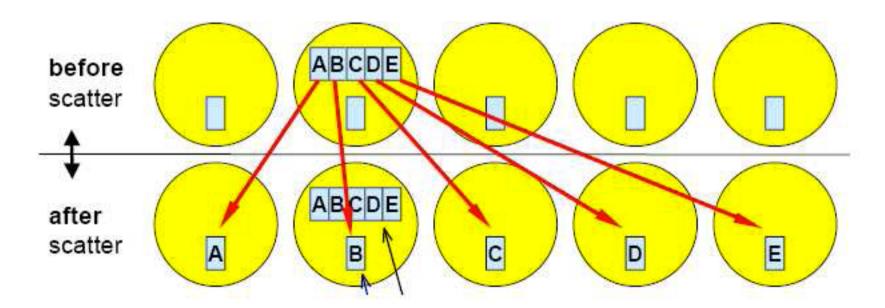
```
#include <mpi.h>
#include <stdio.h>
int main(int argc, char** argv) {
    int rank;
    int buf:
    const int root = 0:
    MPI Init(&argc, &argv);
    MPI Comm rank(MPI COMM WORLD, &rank);
    if (rank == root) {
      buf = 777;
    printf("[%d]: Before Bcast, buf is %d\n", rank, buf);
    /* everyone calls bcast, data is taken from root and ends up in everyone's buf */
    MPI Bcast(&buf, 1, MPI INT, root, MPI COMM WORLD);
    printf("[%d]: After Bcast, buf is %d\n", rank, buf);
    MPI Finalize();
    return 0;
```

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#### Scatter:

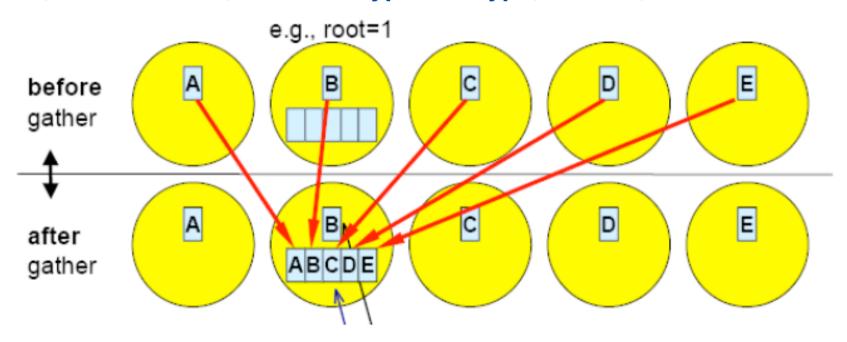
int MPI\_Scatter(void \*sendbuf, int sendcount, MPI\_Datatype sendtype, void \*recvbuf, int recvcount, MPI\_Datatype recvtype, int root, MPI\_Comm comm)





#### Gather:

int MPI\_Gather(void \*sendbuf, int sendcount, MPI\_Datatype sendtype, void \*recvbuf, int recvcount, MPI\_Datatype recvtype, int root, MPI\_Comm comm)





#### Gather:

- Every process (included the root one) sends the content of its send buffer to the root process
- The root process receives the data and order them based on the rank of the sender

Gather belongs to the "all-to-one" class.

There is also MPI\_All\_gather, which is equivalent to a MPI\_Gather operation followed by a broadcast executed by the root process.

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#### Reduction:

used to perform computations that involve distributed data within a group of processes.

#### Examples:

- global sum and product
- global max and min
- user-defined global operations

Suppose to have an **associative and commutative operator o** that, given two elements of the same datatype, produces a result of the same type:

do o d1 o d2 o d3 o ... o ds-2 o ds-1

 $d_i$  = data in process with rank **i** (single variable or vector)



#### Reduction:

int MPI\_Reduce(void \*sendbuf, void \*recvbuf, int count, MPI\_Datatype datatype, MPI\_Op op, int root, MPI\_Comm comm)

#### Example:

```
root = 0;
MPI_Reduce(&localvalue, &globalvalue, 1, MPI_INT, MPI_SUM, root, MPI_COMM_WORLD);
```

The sum of all localvalues (each one being an integer) is returned in globalvalue of process with rank 0.



#### Reduction:

Handle predefiniti per le operazioni	Funzione
MPI_MAX	Maximum
MPI_MIN	Minimum
MPI_SUM	Sum
MPI_PROD	Product
MPI_LAND	Logical AND
MPI_BAND	Bitwise AND
MPI_LOR	Logical OR
MPI_BOR	Bitwise OR
MPI_LXOR	Logical exclusive OR
MPI_BXOR	Bitwise exclusive OR
MPI_MAXLOC	Maximum & location
MPI_MINLOC	Minimum & location

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#### Reduction:

User-defined operators must execute:

```
for (i = 1 to len)
inoutvec[i] = inoutvec[i] OP invec[i]
```

where OP is an arbitrary operator respecting the following syntax:

void my\_operator(void \*invec, void \*inoutvec, int \*len, MPI\_Datatype \*datatype)



#### Reduction:

To register an user-defined operation:

int MPI\_Op\_create(MPI\_User\_function \*func, int commute, MPI\_Op \*op)

to delete the operation:

int MPI\_Op\_Free(op)



#### Reduction:

Example: 1-norm

$$N_1(x) = \sum_{j=0}^{p-1} |\chi_j|$$

where x is a vector of size p.

Step 1: Implement your own reduction operator.

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#### Reduction:

Example: 1-norm

Step 2: Register the function onenorm with MPI by:

MPI\_Op\_create((MPI\_user\_function \*)onenorm, commute, &myop);

#### where

- onenorm is the function to be registered.
- **commute** is an input int (scalar) variable which must be set to **1** if the function satisfies the commutative property. Otherwise, it must be set to **0**.
- myop is an int (output) to be used in subsequent reduction operation call in place of onenorm.

Step 3: Call MPI reduction operation function MPI\_Reduce using the registered myop.



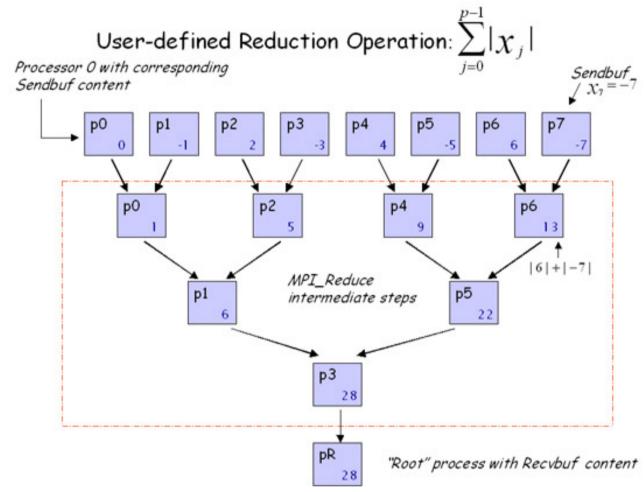
```
#include <mpi.h>
Reduction:
                         #include <stdio>
                         #include <math.h>
Example: 1-norm
                         void main(int argc, char* argv[]) {
                           int root=0, p, myid;
                           float sendbuf, recvbuf;
                           MPI Op myop;
                           int commute = 0;
                           MPI Init(&argc, &argv);
                           MPI Comm size(MPI COMM WORLD, &p);
                           MPI Comm rank(MPI COMM WORLD, &myid);
                           MPI Op create((MPI user function *)onenorm, commute, &myop);
                           sendbuf = myid*((int)pow((double)-1,myid));
                           MPI Barrier(MPI COMM WORLD);
                           MPI Reduce(&sendbuf, &recvbuf, 1, MPI FLOAT, myop,
                                     root, MPI COMM WORLD);
                           if (myid == root)
                             printf("The operation yields %f\n", recvbuf);
                           MPI Finalize();
```

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Reduction:

Example: 1-norm





As in the nonblocking point-to-point case, all calls are local and return immediately, irrespective of the status of other processes. The call initiates the operation, which indicates that the system may start to copy data out of the send buffer and into the receive buffer.

Once initiated, all associated send buffers and buffers associated with input arguments (such as arrays of counts, displacements, or datatypes in the vector versions of the collectives) should not be modified, and all associated receive buffers should not be accessed, until the collective operation completes.

The call returns a request handle, which must be passed to a completion call.



int MPI\_Ibarrier(MPI\_Comm comm, MPI\_request \*request)

int MPI\_Ibcast(void\* buffer, int count, MPI\_Datatype datatype, int root, MPI\_Comm comm, MPI\_Request \*request)

int MPI\_Iscatter(const void\* sendbuf, int sendcount, MPI\_Datatype sendtype, void\* recvbuf, int recvcount, MPI\_Datatype recvtype, int root, MPI\_Comm comm, MPI\_Request \*request)

int MPI\_Igather(const void\* sendbuf, int sendcount, MPI\_Datatype sendtype, void\* recvbuf, int recvcount, MPI\_Datatype recvtype, int root, MPI\_Comm comm, MPI\_Request \*request)



#### References

- <a href="http://www.mpi-forum.org">http://www.mpi-forum.org</a>

- http://www.open-mpi.org