



# PARALLEL PROGRAMMING



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# The MPI\_BARRIER

Blocks until all processes have reached this routine

```
INCLUDE 'mpif.h'
```

```
MPI_BARRIER(COMM, IERROR)
```

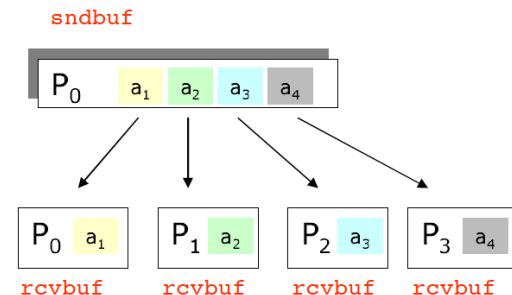
```
INTEGER    COMM, IERROR
```

# Collective Communications

- At the bottom line are based on point 2 point (you could build your own)
- The MPI include a series of subroutines to handle different patterns of communications: **1 to N**, **N to 1** and **N to N**
- Collective Communications imply a *synchronization point* among processes

# MPI\_Scatter

One-to-all communication: different data sent from root process to all others in the communicator.



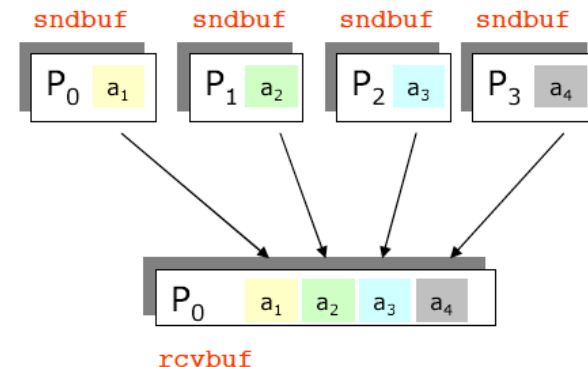
**MPI\_SCATTER( sndbuf, sendcount, sendtype, rcvbuf, rcvcount, recvtype, root, comm)**

- [ IN sndbuf] starting address of send buffer (choice)
- [ IN sendcount] number of elements sent to each process (integer, significant only at **root**)
- [ IN sendtype] data type of send buffer elements (significant only at **root**) (handle)
- [ OUT rcvbuf] address of receive buffer (choice)
- [ IN rcvcount] number of elements in receive buffer (integer)
- [ IN recvtype] data type of rcv buffer elements (handle)
- [ IN root] rank of receiving process (integer)
- [ IN comm] communicator (handle)

# MPI\_Gather

One-to-all communication: different data collected by the root process, from all others processes in the communicator.

It is the opposite of Scatter



**MPI\_GATHER( sendbuf, sendcount, sendtype, rcvbuf, rcvcount, recvttype, root, comm)**

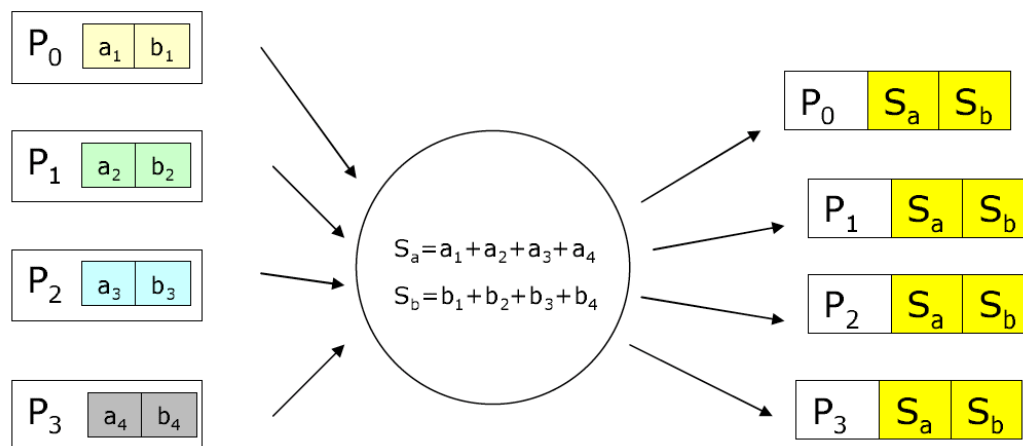
- [ IN sendbuf] starting address of send buffer (choice)
- [ IN sendcount] number of elements in send buffer (integer)
- [ IN sendtype] data type of send buffer elements (handle)
- [ OUT rcvbuf] address of receive buffer (choice, significant only at **root**)
- [ IN rcvcount] number of elements for any single receive (integer, significant only at **root**)
- [ IN recvttype] data type of recv buffer elements (significant only at **root**) (handle)
- [ IN root] rank of receiving process (integer)
- [ IN comm] communicator (handle)

# Reduction

The reduction operation allow to:

- Collect data from each process
- Reduce the data to a single value
- Store the result on the root processes
- Store the result on all processes
- **Overlap of communication and computing**

MPI op	Function
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 and location
MPI_MINLOC	Minimum and location



# Calling MPI\_REDUCE

**MPI\_REDUCE(in, out, count, type, op, receiver, comm, err)**

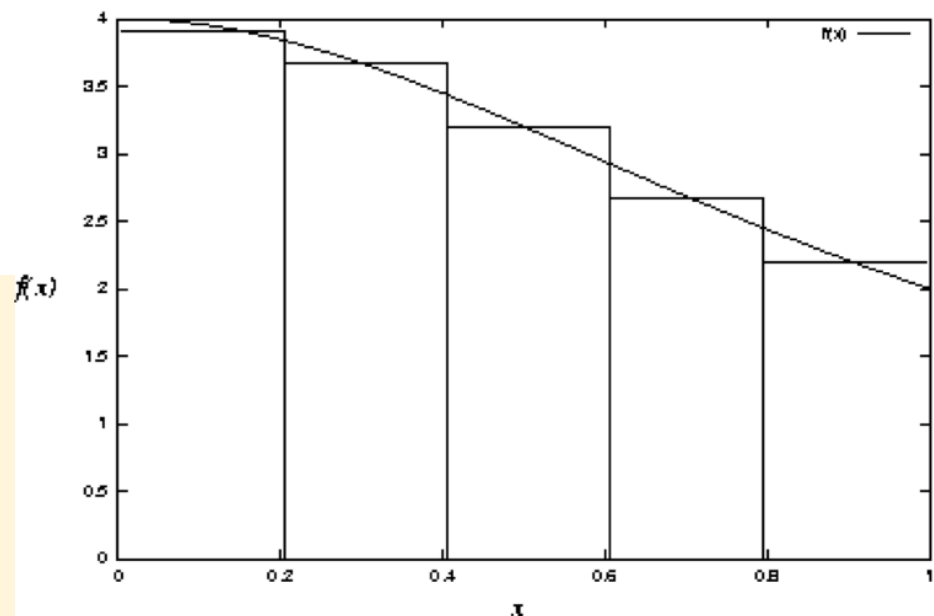
in:	data to be sent (from all)
out:	storage for reduced data (on receiver)
count:	number of data items to be reduced
type:	type (=size) of data items
op:	reduction operation, e.g. <b>MPI_SUM</b>
receiver:	rank of sending processor of data
communicator:	group identifier, <b>MPI_COMM_WORLD</b>
err:	error status or <b>MPI_SUCCESS</b>

# Approximate PI Using MPI collectives

$$\int_0^1 \frac{1}{1+x^2} dx = \arctan(x) \Big|_0^1 = \arctan(1) - \arctan(0) = \arctan(1) = \frac{\pi}{4}$$

$$\pi = 4 \int_0^1 \frac{1}{1+x^2} dx$$

Integrate, i.e determine area under function numerically using slices of  $h * f(x)$  at midpoints





```
#include <stdio.h>

int main(){
    long n , i ;
    double w,x,sum,pi,f,a;

    n = 100000000;
    w = 1.0/n;
    sum = 0.0;

    for ( i = 1 ; i <= n ; i++ ) {
        x = w * (i - 0.5);
        sum = sum + (4.0 / (1.0 + x * x ) );
    }

    pi = w * sum ;
    printf("Value of pi: %.16g\n", pi);

    return 0;
}
```

# Assignment

- 1) Implement the PI approximation in parallel using the Message Passing paradigm

# STANDARD BLOCKING SEND - RECV

**MPI\_SEND(buf, count, type, dest, tag, comm, ierr)**

**MPI\_RECV(buf, count, type, dest, tag, comm, status, ierr)**

**Buf** array of MPI type **type**.

**Count** (INTEGER) number of element of **buf** to be sent/recv

**Type** (INTEGER) MPI type of **buf**

**Dest** (INTEGER) rank of the destination process

**Tag** (INTEGER) number identifying the message

**Comm** (INTEGER) communicator of the sender and receiver

\* **Status** (INTEGER) array of size **MPI\_STATUS\_SIZE** containing communication status information

**ierr** (INTEGER) error code

*\* used only for receive operations*

# NON-BLOCKING SEND - RECV

**MPI\_ISEND(buf, count, type, dest, tag, comm, request, ierr)**

**MPI\_Irecv(buf, count, type, dest, tag, comm, request, ierr)**

**Buf** array of MPI type **type**.

**Count** (INTEGER) number of element of **buf** to be sent/recv

**Type** (INTEGER) MPI type of **buf**

**Dest** (INTEGER) rank of the destination process

**Tag** (INTEGER) number identifying the message

**Comm** (INTEGER) communicator of the sender and receiver

**Request** (INTEGER) request handler, used for checking the communication status

**ierr** (INTEGER) error code

# No-Blocking Checkpoint

## **MPI\_WAIT(request, status, ierr)**

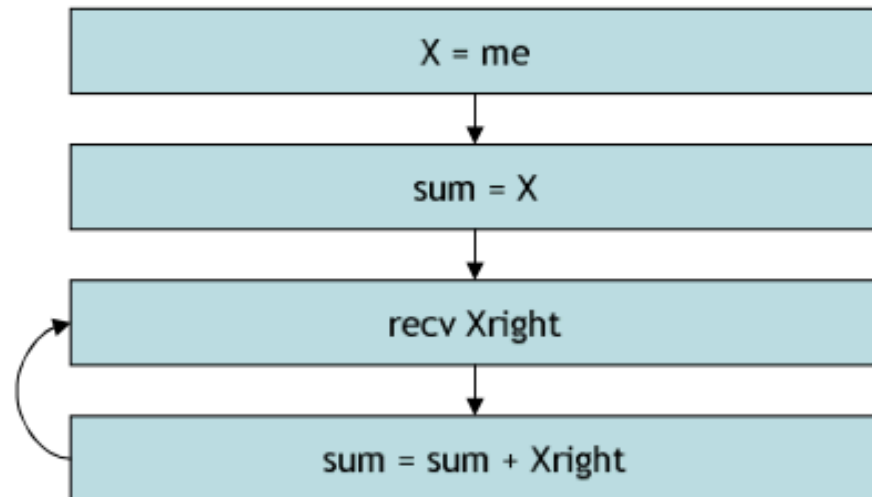
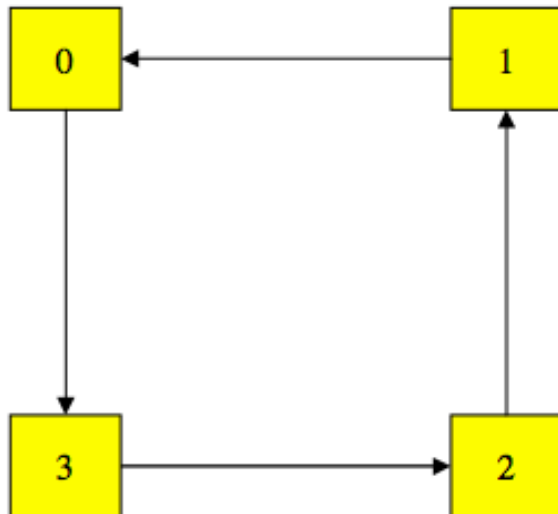
**Request** (INTEGER) request handler, used for checking the communication status

**Status** (INTEGER) array of size **MPI\_STATUS\_SIZE** containing communication status information

**ierr** (INTEGER) error code

Wait until the communication handled by the object request is terminated. For test only use MPI\_TEST, for checkpoint of many communication use MPI\_WAITALL





Who/when does the SEND?

What does it send?

How many times?

sum = 6 (on 4 processors)

# Exercise

Implement the proposed exercise, first exchanging one single element (mype) among processes as illustrated in class as well as on the previous slide. Try to optimize the code for sending in the ring a large set of data and overlapping the computation ( $\Sigma$ ) and the communication (send-recv). In case of a dataset larger than one element the local sum is considered a vector sum (element by element).



# External MPI Resources

Here are some links to tutorials and literature:

CI-Tutor at NCSA: <http://www.citutor.org/>

MPI reference and mini tutorial at LLNL:

<http://computing.llnl.gov/tutorials/mpi/>

Designing and Building // Programs, by Ian Foster:

<http://www.mcs.anl.gov/~itf/dbpp/>

MPI standards: <http://www.mpi-forum.org/>

OpenMPI: <http://www.open-mpi.org>

MPICH: <http://www.mcs.anl.gov/research/projects/mpich2>