

## **Agenda**



- Setting up a cluster of VMs
- Message Passing Interface
- Blocking / Non-blocking messages
- Data partitioning
- Lab2: Parallel and distributed data processing



# Lab 2: Use cluster to distribute work and do some parallel data processing



- Now that you have a group of nodes managed as a cluster
  - Create a distributed data processing capability.
  - Master node takes a dataset and distributes / partitions the data on multiple active slaves that have enough resources.
  - Slaves process the data and send results back to Master. Master merges / reduces the data and reports back to the user.
  - Use blocking vs non-blocking messages in OpenMPI.
  - Status of the work at slaves is reported back to the master as work progresses and finishes.
  - Master collects the performance data (CPU, mem, IO) and stores it for analysis.
  - Any errors are also reported to the master.
  - Use compression to transfer data. Quantify communication cost bytes sent over the network, latency of transfer.
  - Use OpenMPI for communication between nodes, zlib etc. compression libs
  - Handle a node crash Master should get work done on remaining nodes. What are the options should it restart the job or only shift part of the work to another node? How is a slave crash detected? Use the heartbeat mechanism.



## Setting up a cluster of VMs (MPI Cluster)



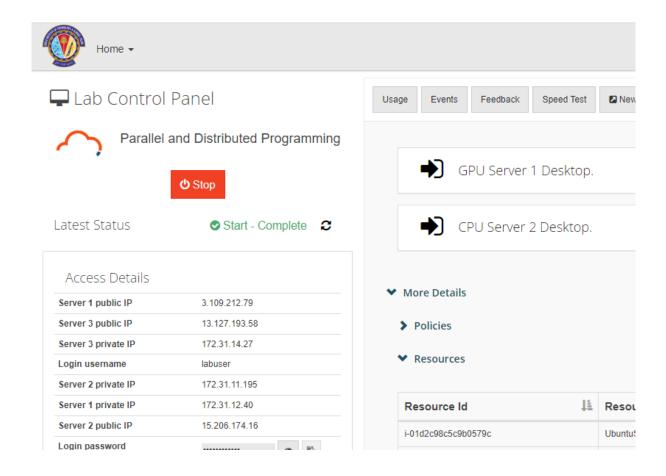
- Step 1: Configure your hosts file
- hosts file is used by the device operating system to map hostnames to IP addresses.
- Example of a host file in master.

```
$ sudo nano /etc/hosts
#MPI CLUSTERS
172.20.36.120 manager
172.20.36.153 worker1
172.20.36.143 worker2
172.20.36.116 worker3
```

- For worker (slave) node
- Example of a host file for worker2

```
#MPI CLUSTER SETUP
172.20.36.120 manager
172.20.36.143 worker2
```

\$ mpirun -hostfile /etc/hosts ./mpi\_hello





## **MPI (Message Passing Interface)**



- A standard message passing specification for the vendors to implement
- Context: distributed memory parallel computers
  - Each processor has its own memory and cannot access the memory of other processors
  - Any data to be shared must be explicitly transmitted from one to another
- Most message passing programs use the single program multiple data (SPMD) model
  - Each processor executes the same set of instructions
  - Parallelization is achieved by letting each processor operation a different piece of data
  - MIMD (Multiple Instructions Multiple Data)



## **SPMD** example



```
main(int argc, char **argv){
  if(process is assigned Master role){
    /* Assign work and coordinate workers and collect results */
    MasterRoutine(/*arguments*/);
   } else { /* it is worker process */
    /* interact with master and other workers. Do the work and send
  results to the master*/
    WorkerRoutine(/*arguments*/);
```



## Why MPI?



- Small
  - Many programs can be written with only 6 basic functions
- Large
  - MPI's extensive functionality from many functions
- Scalable
  - Point-to-point communication
- Flexible
  - Don't need to rewrite parallel programs across platforms

MPI\_INIT : Initiate an MPI computation.
MPI\_FINALIZE : Terminate a computation.

MPI\_COMM\_SIZE : Determine number of processes. MPI\_COMM\_RANK : Determine my process identifier.

MPI\_SEND : Send a message.
MPI\_RECV : Receive a message.



#### What we need to know...







### **Basic functions**



Function	Description
int MPI_Init(int *argc, char **argv)	Initialize MPI
int MPI_Finalize()	Exit MPI
int MPI_Comm_size(MPI_Comm comm, int *size)	Determine number of processes within a comm
int MPI_Comm_rank(MPI_Comm comm, int *rank)	Determine process rank within a comm
int MPI_Send(void *buf, int count, MPI_Datatype datatype, int dest, int tag, MPI_Comm comm)	Send a message
int MPI_Recv(void *buf, int count, MPI_Datatype datatype, int src, int tag, MPI_Comm comm, MPI_Status *status)	Receive a message



#### Communicator



- An identifier associated with a group of processes
  - Each process has a unique rank within a specific communicator from 0 to (nprocesses-1)
  - Always required when initiating a communication by calling an MPI function
- Default: MPI\_COMM\_WORLD
  - Contains all processes
- Several communicators can co-exist
  - A process can belong to different communicators at the same time



#### **Hello World**



```
#include "mpi.h"
int main( int argc, char *argv[] ) {
   int nproc, rank;
   MPI_Init (&argc,&argv); /* Initialize MPI */
   MPI_Comm_size(MPI_COMM_WORLD,&nproc); /* Get Comm Size*/
   MPI_Comm_rank(MPI_COMM_WORLD,&rank); /* Get rank */
   printf("Hello World from process %d\n", rank);
   MPI_Finalize(); /* Finalize */
   return 0;
```



## How to compile...



- Need to tell the compiler where to find the MPI include files and how to link to the MPI libraries.
- Fortunately, most MPI implementations come with scripts that take care of these issues:
  - mpicc mpi\_code.c –o a.out
- Two widely used (and free) MPI implementations
  - MPICH (<a href="http://www-unix.mcs.anl.gov/mpi/mpich">http://www-unix.mcs.anl.gov/mpi/mpich</a>)
  - OPENMPI (<a href="http://www.openmpi.org">http://www.openmpi.org</a>)



## **Blocking Message Passing**



- The call waits until the data transfer is done
  - The sending process waits until all data are transferred to the system buffer
  - The receiving process waits until all data are transferred from the system buffer to the receive buffer
  - Buffers can be freely reused

## **Blocking Message Send**



- buf Specifies the starting address of the buffer.
- count Indicates the number of buffer elements
- dtype Denotes the datatype of the buffer elements
- dest Specifies the rank of the destination process in the group associated with the communicator comm
- tag Denotes the message label
- comm Designates the communication context that identifies a group of processes



## **Blocking Message Send**



Standard (MPI_Send)	The sending process returns when the system can buffer the message or when the message is received and the buffer is ready for reuse.
Buffered (MPI_Bsend)	The sending process returns when the message is buffered in an application-supplied buffer.
Synchronous (MPI_Ssend)	The sending process returns only if a matching receive is posted and the receiving process has started to receive the message.
Ready (MPI_Rsend)	The message is sent as soon as possible.



## **Blocking Message Receive**



MPI\_Recv (void \*buf, int count, MPI\_Datatype dtype, int source, int tag, MPI\_Comm comm, MPI\_Status \*status);

- buf Specifies the starting address of the buffer.
- count Indicates the number of buffer elements
- dtype Denotes the datatype of the buffer elements
- source Specifies the rank of the source process in the group associated with the communicator comm
- tag Denotes the message label
- comm
   Designates the communication context that identifies a group of processes
- status Returns information about the received message



### **Example** (from http://mpi.deino.net/mpi\_functions/index.htm)



```
if (rank == 0) {
  for (i=0; i<10; i++) buffer[i] = i;
  MPI_Send(buffer, 10, MPI_INT, 1, 123, MPI_COMM_WORLD);
} else if (rank == 1) {
  for (i=0; i<10; i++) buffer[i] = -1;
  MPI_Recv(buffer, 10, MPI_INT, 0, 123, MPI_COMM_WORLD, &status);
  for (i=0; i<10; i++)
    if (buffer[i] != i)
     printf("Error: buffer[%d] = %d but is expected to be %d\n", i, buffer[i], i);
```



## Non-blocking Message Passing



- Returns immediately after the data transferred is initiated
- Allows to overlap computation with communication
- Need to be careful though
  - When send and receive buffers are updated before the transfer is over, the result will be wrong



## Non-blocking Message Passing



MPI\_Isend (void \*buf, int count, MPI\_Datatype dtype, int dest, int tag, MPI\_Comm comm, MPI\_Request \*req);

MPI\_Recv (void \*buf, int count, MPI\_Datatype dtype, int source, int tag, MPI\_Comm comm, MPI\_Request \*req);

MPI\_Wait(MPI\_Request \*req, MPI\_Status \*status);

• req Specifies the request used by a completion routine when called by the application to complete the send operation.

Blocking	MPI_Send	MPI_Bsend	MPI_Ssend	MPI_Rsend	MPI_Recv
Non-blocking	MPI_Isend	MPI_lbsend	MPI_Issend	MPI_Irsend	MPI_Irecv



## Non-blocking Message Passing



```
right = (rank + 1) % nproc;
left = rank - 1;
if (left < 0) left = nproc - 1;
MPI_Irecv(buffer, 10, MPI_INT, left, 123, MPI_COMM_WORLD, &request);
MPI_Isend(buffer2, 10, MPI_INT, right, 123, MPI_COMM_WORLD,
  &request2);
MPI_Wait(&request, &status);
MPI_Wait(&request2, &status);
```



#### How to execute MPI codes?



- The implementation supplies scripts to launch the MPI parallel calculation
  - mpirun –np #proc a.out
  - mpiexec –n #proc a.out
- A copy of the same program runs on each processor core within its own process (private address space)
- Communication
  - through the network interconnect in distributed memory
  - through the shared memory



#### **Collective communications**



- A single call handles the communication between all the processes in a communicator
- There are 3 types of collective communications
  - Data movement (e.g. MPI\_Bcast)
  - Reduction (e.g. MPI\_Reduce)
  - Synchronization (e.g. MPI\_Barrier)



#### **Broadcast**



- int MPI\_Bcast(void \*buffer, int count, MPI\_Datatype datatype, int root, MPI\_Comm comm);
  - One process (root) sends data to all the other processes in the same communicator

Must be called by all the processes with the same arguments

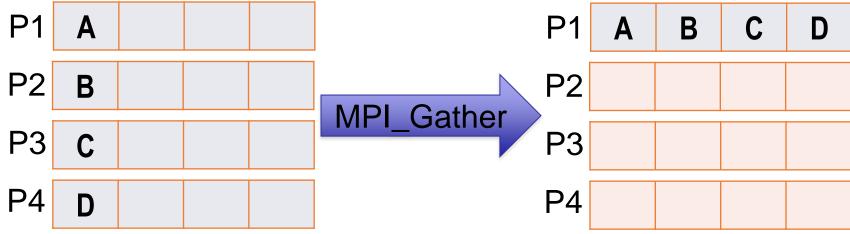
P1	Α	В	С	D	P1	Α	В	С	D
P2					MPI_Bcast	Α	В	С	D
P3					P3	Α	В	С	D
P4					P4	Α	В	С	D



#### **Gather**

- int MPI\_Gather(void \*sendbuf, int sendcnt, MPI\_Datatype
   sendtype, void \*recvbuf, int recvcnt, MPI\_Datatype recvtype, int
   root, MPI\_Comm comm)
  - One process (root) collects data to all the other processes in the same communicator

Must be called by all the processes with the same arguments

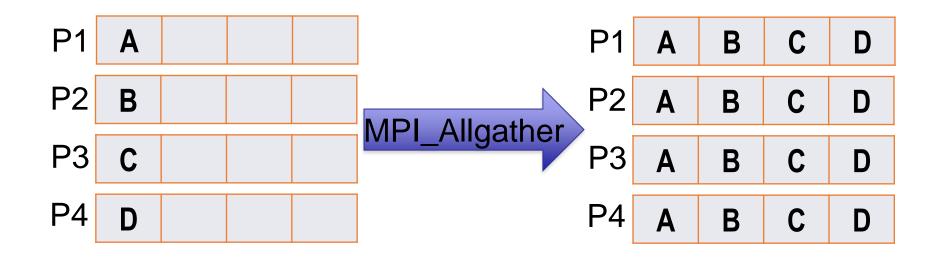




#### **Gather to All**



- int MPI\_Allgather(void \*sendbuf, int sendcnt, MPI\_Datatype sendtype, void \*recvbuf, int recvcnt, MPI\_Datatype recvtype, MPI\_Comm comm)
  - All the processes collects data to all the other processes in the same communicator
  - Must be called by all the processes with the same arguments

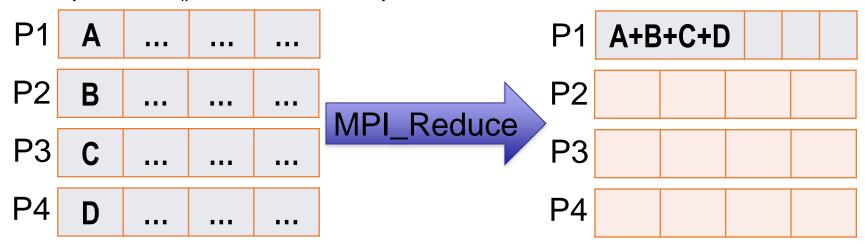




#### Reduction



- int MPI\_Reduce(void \*sendbuf, void \*recvbuf, int count, MPI\_Datatype datatype, MPI\_Op op, int root, MPI\_Comm comm)
  - One process (root) collects data to all the other processes in the same communicator, and performs an operation on the data
  - MPI\_SUM, MPI\_MIN, MPI\_MAX, MPI\_PROD, logical AND, OR, XOR, and a few more
  - MPI\_Op\_create(): User defined operator

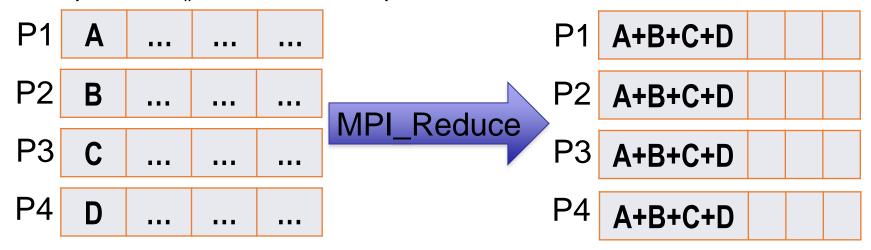




#### **Reduction to All**



- int MPI\_Allreduce(void \*sendbuf, void \*recvbuf, int count, MPI\_Datatype datatype, MPI\_Op op, MPI\_Comm comm)
  - All the processes collect data to all the other processes in the same communicator, and perform an operation on the data
  - MPI\_SUM, MPI\_MIN, MPI\_MAX, MPI\_PROD, logical AND, OR, XOR, and a few more
  - MPI\_Op\_create(): User defined operator





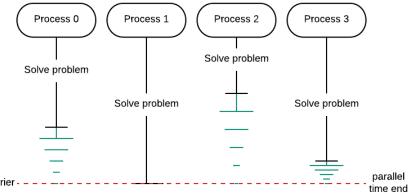
## **Synchronization**



#### **MPI** Barrier

- Blocks until all processes in the communicator have reached this routine.
- A barrier is used when you want all the processes to complete a portion of code before continuing.
- int MPI Barrier(MPI Comm comm)

```
int solveProblem(int id, int numProcs)
{ sleep( ((double)id+1) / numProcs); return 42; }
int main(int argc, char** argv) {
int id = -1, numProcesses = -1; double startTime = 0.0, totalTime = 0.0;
int answer = 0.0; MPI_Init(&argc, &argv); MPI_Comm_rank(MPI_COMM_WORLD, &id);
MPI Comm size(MPI COMM WORLD, &numProcesses);
MPI Barrier(MPI COMM WORLD);
if (id == MASTER) { startTime = MPI Wtime(); }
answer = solveProblem(id, numProcesses);
MPI_Barrier(MPI_COMM_WORLD);
if ( id == MASTER )
{ totalTime = MPI Wtime() - startTime;
printf("\nThe answer is %d; computing it took %f secs.\n\n", answer, totalTime);
MPI_Finalize();
return 0;
```



Reference: http://selkie.macalester.edu/csinparallel/modules/Patternlets/build/html/MessagePassing/Barrier\_Tags.html



#### **Data distributions**



$$\mathbf{x} + \mathbf{y} = (x_0, x_1, \dots, x_{n-1}) + (y_0, y_1, \dots, y_{n-1})$$

$$= (x_0 + y_0, x_1 + y_1, \dots, x_{n-1} + y_{n-1})$$

$$= (z_0, z_1, \dots, z_{n-1})$$

$$= \mathbf{z}$$

Compute a vector sum.

### Serial implementation of vector addition



```
void Vector_sum(double x[], double y[], double z[], int n) {
  int i;

for (i = 0; i < n; i++)
    z[i] = x[i] + y[i];
} /* Vector_sum */</pre>
```

## Different partitions of a 12-component vector among 3 processes



	Components											
									Block-cyclic			
Process	Block				Cyclic			Blocksize = 2				
0	0	1	2	3	0	3	6	9	0	1	6	7
1	4	5	6	7	1	4	7	10	2	3	8	9
2	8	9	10	11	2	5	8	11	4	5	10	11



## **Partitioning options**



- Block partitioning
  - Assign blocks of consecutive components to each process.
- Cyclic partitioning
  - Assign components in a round robin fashion.
- Block-cyclic partitioning
  - Use a cyclic distribution of blocks of components.



## Parallel implementation of vector addition



```
void Parallel_vector_sum(
    double local_x[] /* in */,
    double local_y[] /* in */,
    double local_z[] /* out */,
    int local_n /* in */) {
    int local_i;

    for (local_i = 0; local_i < local_n; local_i++)
        local_z[local_i] = local_x[local_i] + local_y[local_i];
} /* Parallel_vector_sum */</pre>
```



#### **Scatter**



 MPI\_Scatter can be used in a function that reads in an entire vector on process 0 but only sends the needed components to each of the other processes.



## Reading and distributing a vector



```
void Read vector(
     double
              local a[] /* out */,
             local n /*in */,
     int
                   /* in */
     int
              n
     char vec name [] /* in */,
     int my_rank /* in */,
     MPI_Comm comm /*in */) {
  double* a = NULL;
  int i;
  if (my_rank == 0) {
     a = malloc(n*sizeof(double));
     printf("Enter the vector %s\n", vec name);
     for (i = 0; i < n; i++)
        scanf("%lf", &a[i]);
     MPI Scatter(a, local n, MPI DOUBLE, local a, local n, MPI DOUBLE,
           0, comm);
     free(a):
  } else {
     MPI_Scatter(a, local_n, MPI_DOUBLE, local_a, local_n, MPI_DOUBLE,
           0. \text{comm}:
  /* Read_vector */
```

#### **Gather**



 Collect all of the components of the vector onto process 0, and then process 0 can process all of the components.

```
int MPI Gather(
    void*
               send_buf_p /* in */,
    int
               send_count /* in */,
               send_type /*in */,
    MPI_Datatype
    void*
               recv_buf_p /* out */,
    int recv_count /* in */,
               recv_type /*in */,
    MPI_Datatype
               dest\_proc /* in */,
    int
               comm /* in */);
    MPI_Comm
```



### Print a distributed vector (1)



```
void Print_vector(
    double local_b[] /* in */,
    int
          local_n /* in */,
          n /* in */,
    int
    char title[] /* in */,
    int
           my_rank /* in */,
            comm /* in */) {
    MPI_Comm
  double* b = NULL;
  int i;
```



### Print a distributed vector (2)



```
if (my_rank == 0) 
  b = malloc(n*sizeof(double));
  MPI_Gather(local_b, local_n, MPI_DOUBLE, b, local_n, MPI_DOUBLE,
         0, comm);
   printf("%s\n", title);
   for (i = 0; i < n; i++)
     printf("%f ", b[i]);
   printf("\n");
   free(b);
} else {
   MPI Gather(local b, local n, MPI DOUBLE, b, local n, MPI DOUBLE,
         0, comm);
/* Print_vector */
```

### **Allgather**



- Concatenates the contents of each process' send\_buf\_p and stores this in each process' recv\_buf\_p.
- As usual, recv\_count is the amount of data being received from each process.



### **Matrix-vector multiplication**



$$A = (a_{ij})$$
 is an  $m \times n$  matrix

 $\mathbf{x}$  is a vector with n components

y = Ax is a vector with m components

$$y_i = a_{i0}x_0 + a_{i1}x_1 + a_{i2}x_2 + \cdots + a_{i,n-1}x_{n-1}$$

i-th component of y

Dot product of the ith row of A with x.





- The MPI standard allows MPI\_Send to behave in two different ways:
  - it can simply copy the message into an MPI managed buffer and return,
  - or it can block until the matching call to MPI\_Recv starts.



- Many implementations of MPI set a threshold at which the system switches from buffering to blocking.
- Relatively small messages will be buffered by MPI\_Send.
- Larger messages, will cause it to block.





- If the MPI\_Send executed by each process blocks, no process will be able to start executing a call to MPI\_Recv, and the program will hang or deadlock.
- Each process is blocked waiting for an event that will never happen.

(see pseudo-code)





A program that relies on MPI provided buffering is said to be unsafe.

 Such a program may run without problems for various sets of input, but it may hang or crash with other sets.

### MPI\_Ssend



- An alternative to MPI\_Send defined by the MPI standard.
- The extra "s" stands for synchronous and MPI\_Ssend is guaranteed to block until the matching receive starts.



### Restructuring communication







### MPI\_Sendrecv



- An alternative to scheduling the communications ourselves.
- Carries out a blocking send and a receive in a single call.
- The dest and the source can be the same or different.
- Especially useful because MPI schedules the communications so that the program won't hang or crash.



### MPI\_Sendrecv

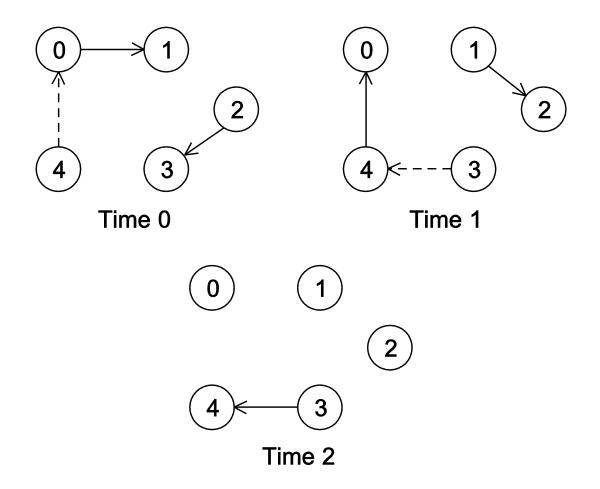


```
int MPI_Sendrecv(
    void*
         send_buf_p /*in */,
         send_buf_size /*in */,
    int
    MPI_Datatype send_buf_type /*in */,
               dest /*in */,
    int
               send_tag /*in */,
    int
             recv_buf_p /* out */,
    void*
         recv_buf_size /*in */,
    int
    MPI_Datatype recv_buf_type /*in */,
    int
               source /*in */,
    int
               recv tag /*in */,
    MPI_Comm communicator /*in */,
    MPI_Status* status_p /*in */);
```



### Safe communication with five processes







#### Parallel odd-even transposition sort



```
void Merge_low(
     int my_keys[], /* in/out */
     int recv_keys[], /* in */
     int temp_keys[], /* scratch */
     int local_n /* = n/p, in */) {
  int mi, ri, ti;
  m i = r i = t i = 0;
   while (t_i < local_n) {</pre>
     if (my_keys[m_i] <= recv_keys[r_i]) {</pre>
        temp keys[t i] = my keys[m i];
        t i++; m i++;
     } else {
        temp_keys[t_i] = recv_keys[r_i];
        t_i++; r_i++;
   for (m_i = 0; m_i < local_n; m_i++)
     my_keys[m_i] = temp_keys[m_i];
  /* Merge_low */
```



## Run-times of parallel odd-even sort



	Number of Keys (in thousands)				
Processes	200	400	800	1600	3200
1	88	190	390	830	1800
2	43	91	190	410	860
4	22	46	96	200	430
8	12	24	51	110	220
16	7.5	14	29	60	130

(times are in milliseconds)



### **Thank You**



Extra slides follows

#### **Outline**

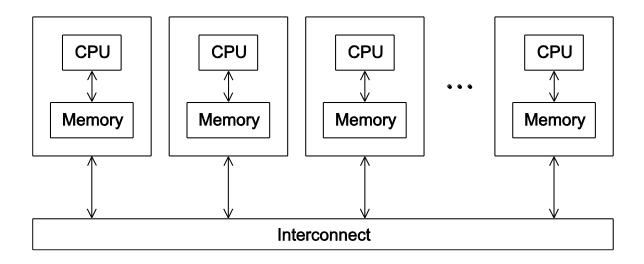


- Writing your first MPI program.
- Using the common MPI functions.
- Collective communication.
- MPI derived datatypes.
- Performance evaluation of MPI programs.
- Parallel sorting.
- Safety in MPI programs.



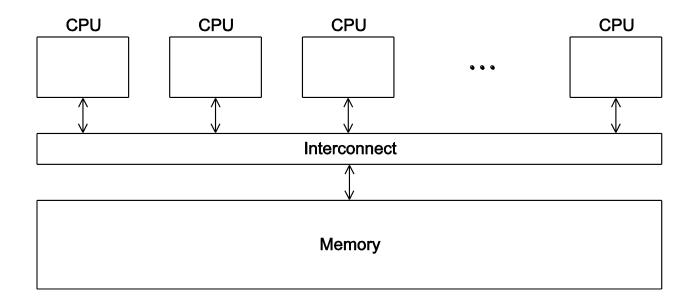
## A distributed memory system





# A shared memory system







#### **Hello World!**



```
#include <stdio.h>
int main(void) {
   printf("hello, world\n");
   return 0;
}
```



(a classic)



### **Identifying MPI processes**



- Common practice to identify processes by nonnegative integer ranks.
- p processes are numbered 0, 1, 2, .. p-1

### Our first MPI program

```
#include < stdio.h>
  #include <string.h> /* For strlen
  #include <mpi.h>
                     /* For MPI functions, etc */
   const int MAX_STRING = 100;
   int main(void) {
                 greeting[MAX_STRING];
      char
                comm_sz; /* Number of processes */
      int
                 my_rank; /* My process rank
      int
11
12
      MPI_Init(NULL, NULL);
13
      MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);
      MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);
15
16
      if (my_rank != 0) {
17
         sprintf(greeting, "Greetings from process %d of %d!",
18
               my_rank, comm_sz);
         MPI_Send(greeting, strlen(greeting)+1, MPI_CHAR, 0, 0,
20
               MPI_COMM_WORLD);
        else {
22
         printf("Greetings from process %d of %d!\n", my_rank, comm_sz);
23
         for (int q = 1; q < comm_sz; q++) {
            MPI_Recv(greeting, MAX_STRING, MPI_CHAR, q,
               0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
26
            printf("%s\n", greeting);
27
28
29
30
      MPI_Finalize();
      return 0;
      /* main */
32
```





## Compilation





source file

mpicc -g -Wall -o mpi\_hello mpi\_hello.c



produce debugging information

create this executable file name (as opposed to default a.out)

turns on all warnings



#### **Execution**



mpiexec -n <number of processes> <executable>

mpiexec -n 1 ./mpi\_hello

run with 1 process

mpiexec -n 4 ./mpi\_hello

run with 4 processes



#### **Execution**



```
mpiexec -n 1 ./mpi_hello
```

Greetings from process 0 of 1!

mpiexec -n 4 ./mpi\_hello

Greetings from process 0 of 4!

Greetings from process 1 of 4!

Greetings from process 2 of 4!

Greetings from process 3 of 4!



### **MPI Programs**



- Written in C.
  - Has main.
  - Uses stdio.h, string.h, etc.
- Need to add mpi.h header file.
- Identifiers defined by MPI start with "MPI\_".
- First letter following underscore is uppercase.
  - For function names and MPI-defined types.
  - Helps to avoid confusion.



### **MPI Components**



- MPI\_Init
  - Tells MPI to do all the necessary setup.

```
int MPI_Init(
   int* argc_p /* in/out */,
   char*** argv_p /* in/out */);
```

- MPI\_Finalize
  - Tells MPI we're done, so clean up anything allocated for this program.

```
int MPI_Finalize(void);
```



#### **Basic Outline**



```
#include <mpi.h>
int main(int argc, char* argv[]) {
   /* No MPI calls before this */
   MPI_Init(&argc, &argv);
   MPI_Finalize();
   /* No MPI calls after this */
   return 0;
```

#### Communicators



- A collection of processes that can send messages to each other.
- MPI\_Init defines a communicator that consists of all the processes created when the program is started.
- Called MPI\_COMM\_WORLD.



#### Communicators





number of processes in the communicator



#### **SPMD**



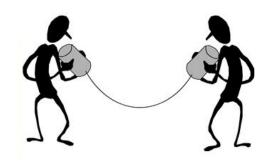
- Single-Program Multiple-Data
- We compile <u>one</u> program.
- Process 0 does something different.
  - Receives messages and prints them while the other processes do the work.
- The if-else construct makes our program SPMD.



#### Communication



```
int MPI_Send(
```





# **Data types**

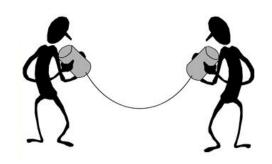


MPI datatype	C datatype
MPI_CHAR	signed char
MPI_SHORT	signed short int
MPI_INT	signed int
MPI_LONG	signed long int
MPI_LONG_LONG	signed long long int
MPI_UNSIGNED_CHAR	unsigned char
MPI_UNSIGNED_SHORT	unsigned short int
MPI_UNSIGNED	unsigned int
MPI_UNSIGNED_LONG	unsigned long int
MPI_FLOAT	float
MPI_DOUBLE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



#### Communication







### Message matching



```
MPI_Send(send_buf_p, send_buf_sz, send_type, dest, send_tag
         send_comm);
                 MPI_Send
                 src = q
                                          MPI_Recv
\texttt{MPI\_Recv(recv\_buf\_p, recv\_buf\_sz, recv\_type, src,}
          recv_comm, &status);
```



### Receiving messages



- A receiver can get a message without knowing:
  - the amount of data in the message,
  - the sender of the message,
  - or the tag of the message.







### status\_p argument



MPI\_Status\*

**MPI\_Status\*** status;

status.MPI\_SOURCE status.MPI\_TAG

MPI\_SOURCE
MPI\_TAG
MPI\_ERROR



# How much data am I receiving?







#### Issues with send and receive



- Exact behavior is determined by the MPI implementation.
- MPI\_Send may behave differently with regard to buffer size, cutoffs and blocking.
- MPI\_Recv always blocks until a matching message is received.
- Know your implementation; don't make assumptions!





# Input



- Most MPI implementations only allow process 0 in MPI\_COMM\_WORLD access to stdin.
- Process 0 must read the data (scanf) and send to the other processes.

```
MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);
MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);
Get_data(my_rank, comm_sz, &a, &b, &n);
h = (b-a)/n;
. . .
```



# Function for reading user input



```
void Get_input(
     int
              my rank /*in */.
          comm_sz /*in */,
     int
     double* a_p /* out */,
     double* b_p /* out */,
             n_p /* out */) {
     int*
  int dest;
  if (my rank == 0) {
     printf("Enter a, b, and n\n");
     scanf("%lf %lf %d", a p, b p, n p);
     for (dest = 1; dest < comm_sz; dest++) {</pre>
        MPI_Send(a_p, 1, MPI_DOUBLE, dest, 0, MPI_COMM_WORLD);
        MPI_Send(b_p, 1, MPI_DOUBLE, dest, 0, MPI_COMM_WORLD);
        MPI Send(n p, 1, MPI INT, dest, 0, MPI COMM WORLD);
   else { /* my\_rank != 0 */}
     MPI_Recv(a_p, 1, MPI_DOUBLE, 0, 0, MPI_COMM_WORLD,
           MPI STATUS_IGNORE);
     MPI_Recv(b_p, 1, MPI_DOUBLE, 0, 0, MPI_COMM_WORLD,
           MPI_STATUS_IGNORE);
     MPI_Recv(n_p, 1, MPI_INT, 0, 0, MPI_COMM_WORLD,
           MPI STATUS IGNORE);
  /* Get_input */
```

### **Collective communication**







#### **Tree-structured communication**

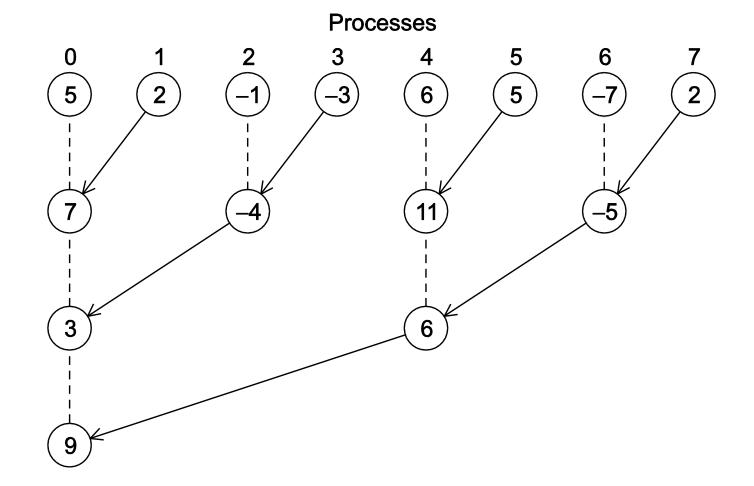


- 1. In the first phase:
  - (a) Process 1 sends to 0, 3 sends to 2, 5 sends to 4, and 7 sends to 6.
  - (b) Processes 0, 2, 4, and 6 add in the received values.
  - (c) Processes 2 and 6 send their new values to processes 0 and 4, respectively.
  - (d) Processes 0 and 4 add the received values into their new values.
- 2. (a) Process 4 sends its newest value to process 0.
  - (b) Process 0 adds the received value to its newest value.



# A tree-structured global sum

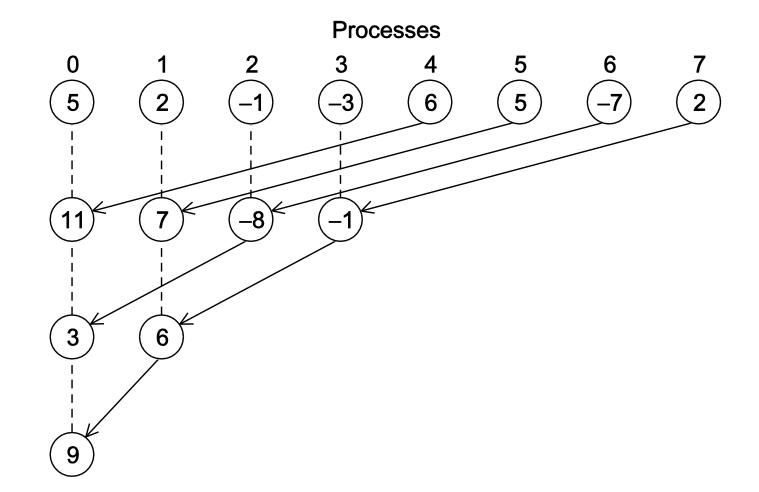






# An alternative tree-structured global sum







#### MPI\_Reduce





# Predefined reduction operators in MPI



Operation Value	Meaning
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 of maximum
MPI_MINLOC	Minimum and location of minimum





- All the processes in the communicator must call the same collective function.
- For example, a program that attempts to match a call to MPI\_Reduce on one process with a call to MPI\_Recv on another process is erroneous, and, in all likelihood, the program will hang or crash.





- The arguments passed by each process to an MPI collective communication must be "compatible."
- For example, if one process passes in 0 as the dest\_process and another passes in 1, then the outcome of a call to MPI\_Reduce is erroneous, and, once again, the program is likely to hang or crash.





- The output\_data\_p argument is only used on dest\_process.
- However, all of the processes still need to pass in an actual argument corresponding to output\_data\_p, even if it's just NULL.



- Point-to-point communications are matched on the basis of tags and communicators.
- Collective communications don't use tags.
- They're matched solely on the basis of the communicator and the order in which they're called.



# Example (1)



Time	Process 0	Process 1	Process 2
0	a = 1; c = 2	a = 1; c = 2	a = 1; c = 2
1	MPI_Reduce(&a, &b,)	MPI_Reduce(&c, &d,)	MPI_Reduce(&a, &b,)
2	MPI_Reduce(&c, &d,)	MPI_Reduce(&a, &b,)	MPI_Reduce(&c, &d,)

### Multiple calls to MPI\_Reduce



# Example (2)



- Suppose that each process calls MPI\_Reduce with operator MPI\_SUM, and destination process 0.
- At first glance, it might seem that after the two calls to MPI\_Reduce, the
  value of b will be 3, and the value of d will be 6.



# Example (3)



- However, the names of the memory locations are irrelevant to the matching of the calls to MPI\_Reduce.
- The order of the calls will determine the matching so the value stored in b will be 1+2+1=4, and the value stored in d will be 2+1+2=5.

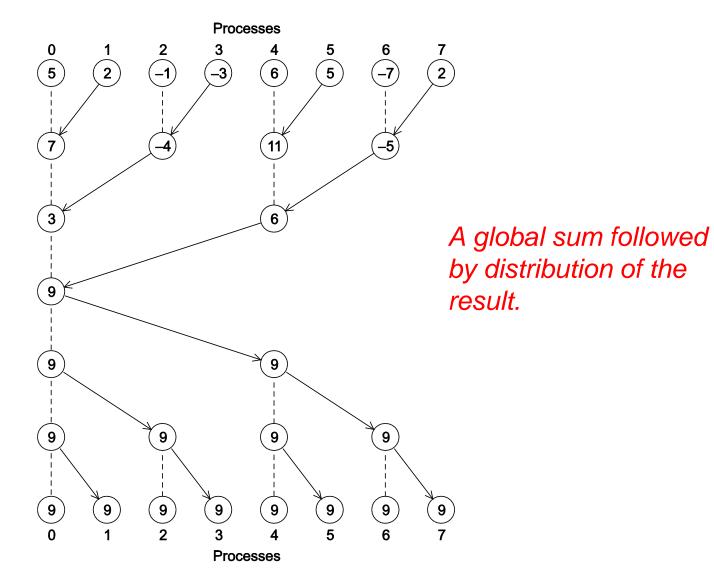
### MPI\_Allreduce



 Useful in a situation in which all of the processes need the result of a global sum in order to complete some larger computation.





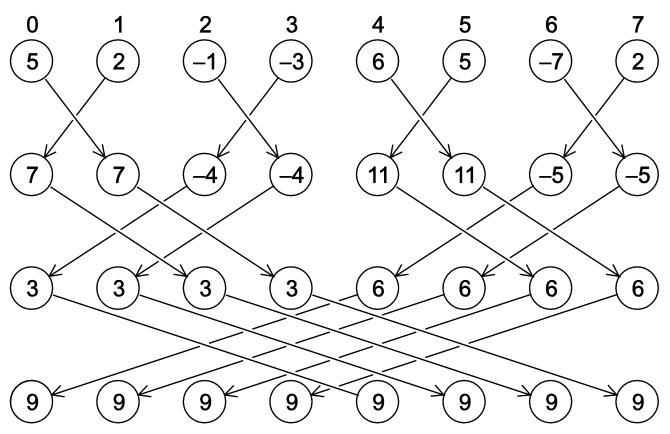








#### **Processes**



A butterfly-structured global sum.



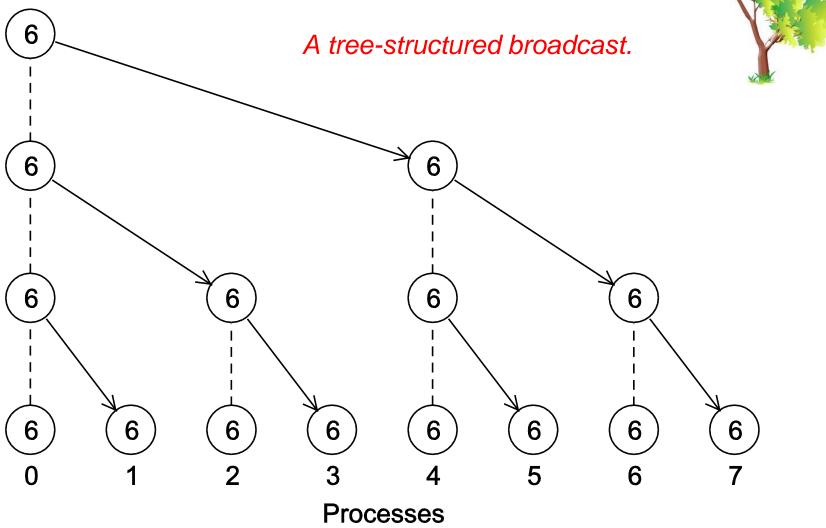
#### **Broadcast**



 Data belonging to a single process is sent to all of the processes in the communicator.









# A version of Get\_input that uses MPI\_Bcast



```
void Get input(
     int my_rank /* in */,
     int comm_sz /* in */,
     double * a_p /* out */,
     double* b_p /* out */,
     int * n_p /* out */) {
  if (my_rank == 0) 
     printf("Enter a, b, and n\n");
     scanf("%lf %lf %d", a_p, b_p, n_p);
  MPI_Bcast(a_p, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
  MPI_Bcast(b_p, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
  MPI_Bcast(n_p, 1, MPI_INT, 0, MPI_COMM_WORLD);
  /* Get_input */
```



#### **Data distributions**



$$\mathbf{x} + \mathbf{y} = (x_0, x_1, \dots, x_{n-1}) + (y_0, y_1, \dots, y_{n-1})$$

$$= (x_0 + y_0, x_1 + y_1, \dots, x_{n-1} + y_{n-1})$$

$$= (z_0, z_1, \dots, z_{n-1})$$

$$= \mathbf{z}$$

Compute a vector sum.

# Serial implementation of vector addition



```
void Vector_sum(double x[], double y[], double z[], int n) {
  int i;

for (i = 0; i < n; i++)
    z[i] = x[i] + y[i];
} /* Vector_sum */</pre>
```

# Different partitions of a 12-component vector among 3 processes



	Components											
								Block-cyclic				
Process	Block				Cyclic			Blocksize = 2				
0	0	1	2	3	0	3	6	9	0	1	6	7
1	4	5	6	7	1	4	7	10	2	3	8	9
2	8	9	10	11	2	5	8	11	4	5	10	11



# **Partitioning options**



- Block partitioning
  - Assign blocks of consecutive components to each process.
- Cyclic partitioning
  - Assign components in a round robin fashion.
- Block-cyclic partitioning
  - Use a cyclic distribution of blocks of components.



# Parallel implementation of vector addition



```
void Parallel_vector_sum(
    double local_x[] /* in */,
    double local_y[] /* in */,
    double local_z[] /* out */,
    int local_n /* in */) {
    int local_i;

    for (local_i = 0; local_i < local_n; local_i++)
        local_z[local_i] = local_x[local_i] + local_y[local_i];
} /* Parallel_vector_sum */</pre>
```



#### **Scatter**



 MPI\_Scatter can be used in a function that reads in an entire vector on process 0 but only sends the needed components to each of the other processes.



# Reading and distributing a vector



```
void Read vector(
     double
              local a[] /* out */,
             local n /*in */,
     int
                   /* in */
     int
              n
     char vec name [] /* in */,
     int my_rank /* in */,
     MPI_Comm comm /*in */) {
  double* a = NULL;
  int i;
  if (my_rank == 0) {
     a = malloc(n*sizeof(double));
     printf("Enter the vector %s\n", vec name);
     for (i = 0; i < n; i++)
        scanf("%lf", &a[i]);
     MPI Scatter(a, local n, MPI DOUBLE, local a, local n, MPI DOUBLE,
           0, comm);
     free(a):
  } else {
     MPI_Scatter(a, local_n, MPI_DOUBLE, local_a, local_n, MPI_DOUBLE,
           0. \text{comm}):
  /* Read_vector */
```

#### **Gather**



 Collect all of the components of the vector onto process 0, and then process 0 can process all of the components.

```
int MPI Gather(
    void*
               send_buf_p /* in */,
    int
               send_count /* in */,
               send_type /*in */,
    MPI_Datatype
    void*
               recv_buf_p /* out */,
    int recv_count /* in */,
               recv_type /* in */,
    MPI_Datatype
               dest\_proc /* in */,
    int
               comm /* in */);
    MPI_Comm
```



# Print a distributed vector (1)



```
void Print_vector(
    double local_b[] /* in */,
    int
          local_n /* in */,
          n /* in */,
    int
    char title[] /* in */,
    int
           my_rank /* in */,
            comm /* in */) {
    MPI_Comm
  double* b = NULL;
  int i;
```



# Print a distributed vector (2)



```
if (my_rank == 0) 
  b = malloc(n*sizeof(double));
   MPI_Gather(local_b, local_n, MPI_DOUBLE, b, local_n, MPI_DOUBLE,
         0, comm);
   printf("%s\n", title);
   for (i = 0; i < n; i++)
     printf("%f ", b[i]);
   printf("\n");
   free(b);
} else {
   MPI Gather(local b, local n, MPI DOUBLE, b, local n, MPI DOUBLE,
         0, comm);
/* Print_vector */
```

# **Allgather**



- Concatenates the contents of each process' send\_buf\_p and stores this in each process' recv\_buf\_p.
- As usual, recv\_count is the amount of data being received from each process.



# **Matrix-vector multiplication**



$$A = (a_{ij})$$
 is an  $m \times n$  matrix

 $\mathbf{x}$  is a vector with n components

y = Ax is a vector with m components

$$y_i = a_{i0}x_0 + a_{i1}x_1 + a_{i2}x_2 + \cdots + a_{i,n-1}x_{n-1}$$

i-th component of y

Dot product of the ith row of A with x.



# **Matrix-vector multiplication**



<i>a</i> <sub>00</sub>	<i>a</i> <sub>01</sub>	• • •	$a_{0,n-1}$
$a_{10}$	$a_{11}$	• • •	$a_{1,n-1}$
:	:		:
$a_{i0}$	$a_{i1}$		$a_{i,n-1}$
:	:		:

$x_0$	
$x_1$	
:	=
$x_{n-1}$	

		У0
<i>x</i> <sub>0</sub>		У1
<i>x</i> <sub>1</sub>		÷ :
:	=	$y_i = a_{i0}x_0 + a_{i1}x_1 + \cdots + a_{i,n-1}x_{n-1}$
n-1		:.
		<i>y</i> <sub>m</sub> −1

#### Multiply a matrix by a vector



```
/* For each row of A */
for (i = 0; i < m; i++) {
    /* Form dot product of ith row with x */
    y[i] = 0.0;

for (j = 0; j < n; j++)
    y[i] += A[i][j]*x[j];
}</pre>
```

Serial pseudo-code



#### C style arrays



0 1 2 3 4 5 6 7 8 9 10 11



#### Serial matrix-vector multiplication



```
void Mat_vect_mult(
     double A[] /* in */,
     double x[] /* in */,
     double y[] /* out */,
     int m /*in */,
     int n /* in */) {
  int i, j;
  for (i = 0; i < m; i++) {
     y[i] = 0.0;
     for (j = 0; j < n; j++)
       y[i] += A[i*n+j]*x[j];
  /* Mat_vect_mult */
```



# An MPI matrix-vector multiplication function (1)



```
void Mat_vect_mult(
    double local_A[] /* in */,
    double local_x[] /* in */,
    double local_y[] /* out */,
    int local_m /* in */,
    int n /*in */,
    int local_n /* in */,
    MPI_Comm comm /*in */) {
  double * x;
  int local_i, j;
  int local_ok = 1;
```



# An MPI matrix-vector multiplication function (2)



# Mpi derived datatypes





# **Derived datatypes**



- Used to represent any collection of data items in memory by storing both the types of the items and their relative locations in memory.
- The idea is that if a function that sends data knows this information about a collection of data items, it can collect the items from memory before they are sent.
- Similarly, a function that receives data can distribute the items into their correct destinations in memory when they're received.



## **Derived datatypes**



- Formally, consists of a sequence of basic MPI data types together with a displacement for each of the data types.
- Trapezoidal Rule example:

Variable	Address	
a	24	
b	40	
n	48	

 $\{(\texttt{MPI\_DOUBLE}, 0), (\texttt{MPI\_DOUBLE}, 16), (\texttt{MPI\_INT}, 24)\}$ 



#### **MPI\_Type create\_struct**



 Builds a derived datatype that consists of individual elements that have different basic types.



#### MPI\_Get\_address



- Returns the address of the memory location referenced by location\_p.
- The special type MPI\_Aint is an integer type that is big enough to store an address on the system.

```
int MPI_Get_address(
    void* location_p /* in */,
    MPI_Aint* address_p /* out */);
```



## MPI\_Type\_commit



 Allows the MPI implementation to optimize its internal representation of the datatype for use in communication functions.

```
int MPI_Type_commit(MPI_Datatype* new_mpi_t_p /* in/out */);
```



#### MPI\_Type\_free



 When we're finished with our new type, this frees any additional storage used.

```
int MPI_Type_free(MPI_Datatype* old_mpi_t_p /* in/out */);
```



# Get input function with a derived datatype (1)



# Get input function with a derived datatype (2)





# Get input function with a derived datatype (3)



```
void Get_input(int my_rank, int comm_sz, double* a_p, double* b_p,
     int* n_p) 
  MPI Datatype input mpi t;
  Build_mpi_type(a_p, b_p, n_p, &input_mpi_t);
   if (my rank == 0) 
     printf("Enter a, b, and n\n");
     scanf("%lf %lf %d", a_p, b_p, n_p);
  MPI_Bcast(a_p, 1, input_mpi_t, 0, MPI_COMM_WORLD);
  MPI_Type_free(&input_mpi_t);
  /* Get_input */
```

#### **Performance evaluation**





#### Elapsed parallel time



 Returns the number of seconds that have elapsed since some time in the past.



#### **Elapsed serial time**



- In this case, you don't need to link in the MPI libraries.
- Returns time in microseconds elapsed from some point in the past.

```
#include "timer.h"
. . .
double now;
. . .
GET_TIME(now);
```





#### **Elapsed serial time**



```
#include "timer.h"
. . .
double start, finish;
. . .
GET_TIME(start);
/* Code to be timed */
. . .
GET_TIME(finish);
printf("Elapsed time = %e seconds\n", finish-start);
```

#### **MPI\_Barrier**



 Ensures that no process will return from calling it until every process in the communicator has started calling it.





#### **MPI\_Barrier**



```
double local start, local finish, local elapsed, elapsed;
. . .
MPI_Barrier(comm);
local start = MPI Wtime();
/* Code to be timed */
. . .
local finish = MPI Wtime();
local_elapsed = local_finish - local_start;
MPI_Reduce(&local_elapsed, &elapsed, 1, MPI_DOUBLE,
  MPI MAX, 0, comm);
if (my rank == 0)
   printf("Elapsed time = %e seconds\n", elapsed);
```

# Run-times of serial and parallel matrix-vector multiplication



	Order of Matrix				
comm_sz	1024	2048	4096	8192	16,384
1	4.1	16.0	64.0	270	1100
2	2.3	8.5	33.0	140	560
4	2.0	5.1	18.0	70	280
8	1.7	3.3	9.8	36	140
16	1.7	2.6	5.9	19	71

(Seconds)



#### Speedup



$$S(n, p) = \frac{T_{\text{serial}}(n)}{T_{\text{parallel}}(n, p)}$$

#### **Efficiency**



$$E(n,p) = \frac{S(n,p)}{p} = \frac{T_{\text{serial}}(n)}{p \times T_{\text{parallel}}(n,p)}$$

# Speedups of Parallel Matrix-Vector Multiplication

	Order of Matrix				
comm_sz	1024	2048	4096	8192	16,384
1	1.0	1.0	1.0	1.0	1.0
2	1.8	1.9	1.9	1.9	2.0
4	2.1	3.1	3.6	3.9	3.9
8	2.4	4.8	6.5	7.5	7.9
16	2.4	6.2	10.8	14.2	15.5



## **Efficiencies of Parallel Matrix-Vector Multiplication**



	Order of Matrix				
comm_sz	1024	2048	4096	8192	16,384
1	1.00	1.00	1.00	1.00	1.00
2	0.89	0.94	0.97	0.96	0.98
4	0.51	0.78	0.89	0.96	0.98
8	0.30	0.61	0.82	0.94	0.98
16	0.15	0.39	0.68	0.89	0.97



# **Scalability**



• A program is scalable if the problem size can be increased at a rate so that the efficiency doesn't decrease as the number of processes increase.





# **Scalability**



- Programs that can maintain a constant efficiency without increasing the problem size are sometimes said to be strongly scalable.
- Programs that can maintain a constant efficiency if the problem size increases at the same rate as the number of processes are sometimes said to be weakly scalable.



# A parallel sorting algorithm





# **Sorting**



- n keys and p = comm sz processes.
- n/p keys assigned to each process.
- No restrictions on which keys are assigned to which processes.
- When the algorithm terminates:
  - The keys assigned to each process should be sorted in (say) increasing order.
  - If 0 ≤ q < r < p, then each key assigned to process q should be less than or equal to every key assigned to process r.</li>



#### Serial bubble sort



```
void Bubble_sort(
     int a[] /* in/out */,
     int n /* in */) {
  int list_length, i, temp;
  for (list_length = n; list_length \geq 2; list_length--)
     for (i = 0; i < list_length -1; i++)
        if (a[i] > a[i+1]) {
           temp = a[i];
           a[i] = a[i+1];
           a[i+1] = temp;
  /* Bubble_sort */
```



# **Odd-even transposition sort**



- A sequence of phases.
- Even phases, compare swaps:

$$(a[0], a[1]), (a[2], a[3]), (a[4], a[5]), \dots$$

Odd phases, compare swaps:

$$(a[1], a[2]), (a[3], a[4]), (a[5], a[6]), \dots$$



#### **Example**



Start: 5, 9, 4, 3

Even phase: compare-swap (5,9) and (4,3) getting the list 5, 9, 3, 4

Odd phase: compare-swap (9,3) getting the list 5, 3, 9, 4

Even phase: compare-swap (5,3) and (9,4) getting the list 3, 5, 4, 9

Odd phase: compare-swap (5,4) getting the list 3, 4, 5, 9



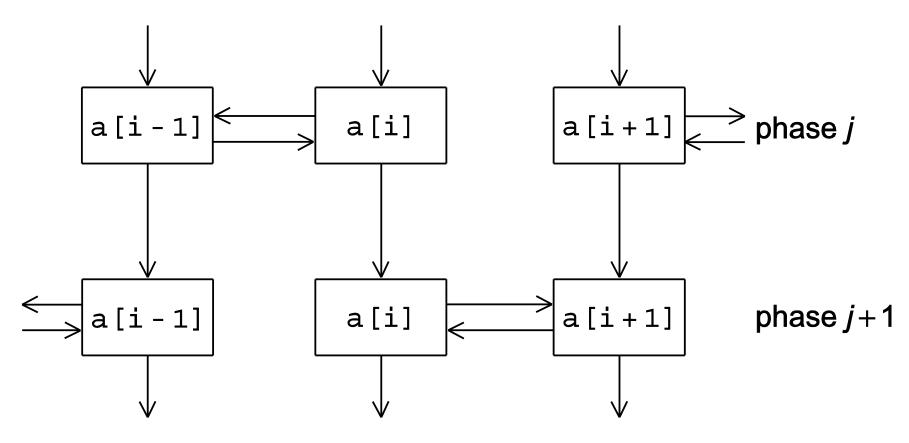
## Serial odd-even transposition sort



```
void Odd_even_sort(
     int a [] /* in/out */,
     int n /* in */) {
  int phase, i, temp;
  for (phase = 0; phase < n; phase++)
      if (phase % 2 == 0) { /* Even phase */
        for (i = 1; i < n; i += 2)
           if (a[i-1] > a[i]) {
              temp = a[i];
              a[i] = a[i-1];
              a[i-1] = temp;
       else { /* Odd phase */
        for (i = 1; i < n-1; i += 2)
           if (a[i] > a[i+1]) {
              temp = a[i];
              a[i] = a[i+1];
              a[i+1] = temp;
  /* Odd_even_sort */
```

# Communications among tasks in odd-even sort





Tasks determining a[i] are labeled with a[i].



## Parallel odd-even transposition sort



	Process						
Time	0	1	2	3			
Start	15, 11, 9, 16	3, 14, 8, 7	4, 6, 12, 10	5, 2, 13, 1			
After Local Sort	9, 11, 15, 16	3, 7, 8, 14	4, 6, 10, 12	1, 2, 5, 13			
After Phase 0	3, 7, 8, 9	11, 14, 15, 16	1, 2, 4, 5	6, 10, 12, 13			
After Phase 1	3, 7, 8, 9	1, 2, 4, 5	11, 14, 15, 16	6, 10, 12, 13			
After Phase 2	1, 2, 3, 4	5, 7, 8, 9	6, 10, 11, 12	13, 14, 15, 16			
After Phase 3	1, 2, 3, 4	5, 6, 7, 8	9, 10, 11, 12	13, 14, 15, 16			

#### Pseudo-code



```
Sort local keys;
for (phase = 0; phase < comm_sz; phase++) {
   partner = Compute_partner(phase, my_rank);
   if (I'm not idle) {
      Send my keys to partner;
      Receive keys from partner;
      if (my_rank < partner)</pre>
         Keep smaller keys;
      else
         Keep larger keys;
```

### Compute\_partner



```
if (phase % 2 == 0) /* Even phase */
  if (my_rank % 2 != 0) /* Odd rank */
     partner = my_rank - 1;
                            /* Even rank */
  else
     partner = my_rank + 1;
                       /* Odd phase */
else
  if (my_rank % 2 != 0) /* Odd rank */
     partner = my_rank + 1;
                            /* Even rank */
  else
     partner = my_rank - 1;
if (partner == -1 || partner == comm_sz)
  partner = MPI_PROC_NULL;
```



- The MPI standard allows MPI\_Send to behave in two different ways:
  - it can simply copy the message into an MPI managed buffer and return,
  - or it can block until the matching call to MPI\_Recv starts.



- Many implementations of MPI set a threshold at which the system switches from buffering to blocking.
- Relatively small messages will be buffered by MPI\_Send.
- Larger messages, will cause it to block.





- If the MPI\_Send executed by each process blocks, no process will be able to start executing a call to MPI\_Recv, and the program will hang or deadlock.
- Each process is blocked waiting for an event that will never happen.

(see pseudo-code)





A program that relies on MPI provided buffering is said to be unsafe.

 Such a program may run without problems for various sets of input, but it may hang or crash with other sets.

#### MPI\_Ssend



- An alternative to MPI\_Send defined by the MPI standard.
- The extra "s" stands for synchronous and MPI\_Ssend is guaranteed to block until the matching receive starts.



### Restructuring communication







#### MPI\_Sendrecv



- An alternative to scheduling the communications ourselves.
- Carries out a blocking send and a receive in a single call.
- The dest and the source can be the same or different.
- Especially useful because MPI schedules the communications so that the program won't hang or crash.



#### MPI\_Sendrecv

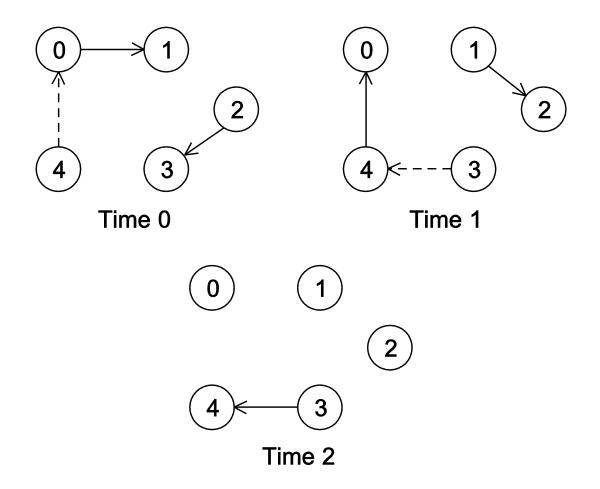


```
int MPI_Sendrecv(
    void*
         send_buf_p /*in */,
         send_buf_size /*in */,
    int
    MPI_Datatype send_buf_type /*in */,
               dest /*in */,
    int
               send_tag /*in */,
    int
             recv_buf_p /* out */,
    void*
         recv_buf_size /*in */,
    int
    MPI_Datatype recv_buf_type /* in */,
    int
               source /*in */,
    int
               recv tag /*in */,
    MPI_Comm communicator /*in */,
    MPI_Status* status_p /*in */);
```



### Safe communication with five processes







#### Parallel odd-even transposition sort



```
void Merge_low(
     int my_keys[], /* in/out */
     int recv_keys[], /* in */
     int temp_keys[], /* scratch */
     int local_n /* = n/p, in */) {
  int mi, ri, ti;
  m i = r i = t i = 0;
   while (t_i < local_n) {</pre>
     if (my_keys[m_i] <= recv_keys[r_i]) {</pre>
        temp keys[t i] = my keys[m i];
        t i++; m i++;
     } else {
        temp_keys[t_i] = recv_keys[r_i];
        t_i++; r_i++;
   for (m_i = 0; m_i < local_n; m_i++)
     my_keys[m_i] = temp_keys[m_i];
  /* Merge_low */
```



## Run-times of parallel odd-even sort



	Number of Keys (in thousands)						
Processes	200	400	800	1600	3200		
1	88	190	390	830	1800		
2	43	91	190	410	860		
4	22	46	96	200	430		
8	12	24	51	110	220		
16	7.5	14	29	60	130		

(times are in milliseconds)



# **Concluding Remarks (1)**



- MPI or the Message-Passing Interface is a library of functions that can be called from C, C++, or Fortran programs.
- A communicator is a collection of processes that can send messages to each other.
- Many parallel programs use the single-program multiple data or SPMD approach.



# **Concluding Remarks (2)**



- Most serial programs are deterministic: if we run the same program with the same input we'll get the same output.
- Parallel programs often don't possess this property.
- Collective communications involve all the processes in a communicator.



# **Concluding Remarks (3)**



- When we time parallel programs, we're usually interested in elapsed time or "wall clock time".
- Speedup is the ratio of the serial run-time to the parallel run-time.
- Efficiency is the speedup divided by the number of parallel processes.



## **Concluding Remarks (4)**



- If it's possible to increase the problem size (n) so that the efficiency doesn't decrease as p is increased, a parallel program is said to be scalable.
- An MPI program is unsafe if its correct behavior depends on the fact that MPI\_Send is buffering its input.

