

IF3230

Sistem Paralel dan Terdistribusi

paralel programming model: distributed memory

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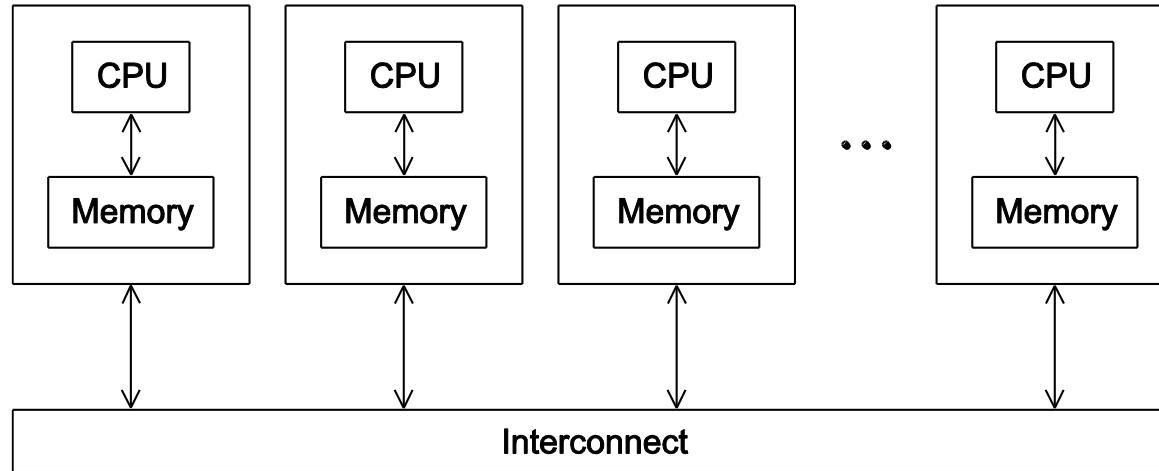
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Roadmap

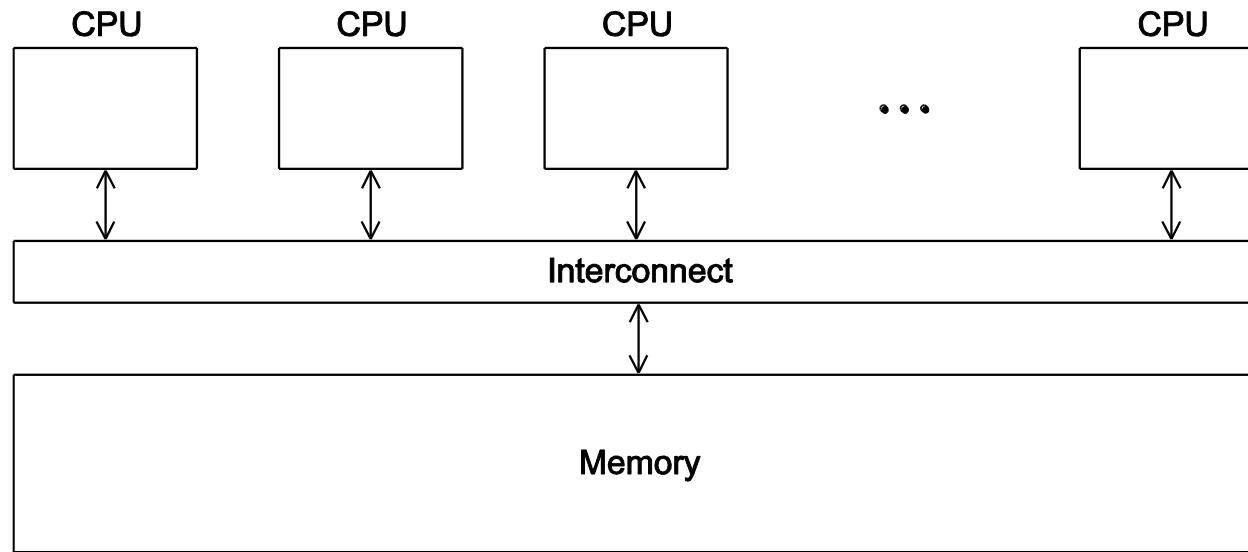
- ▶ Writing your first MPI program.
- ▶ Using the common MPI functions.
- ▶ The Trapezoidal Rule in MPI.
- ▶ Collective communication.
- ▶ MPI derived datatypes.
- ▶ Performance evaluation of MPI programs.
- ▶ Parallel sorting.
- ▶ Safety in MPI programs.



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 $0, 1, 2, \dots p-1$



Our first MPI program

```
-----  
1 #include <stdio.h>  
2 #include <string.h> /* For strlen */  
3 #include <mpi.h> /* For MPI functions, etc */  
4  
5 const int MAX_STRING = 100;  
6  
7 int main(void) {  
8     char greeting[MAX_STRING];  
9     int comm_sz; /* Number of processes */  
10    int my_rank; /* My process rank */  
11  
12    MPI_Init(NULL, NULL);  
13    MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);  
14    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);  
19        MPI_Send(greeting, strlen(greeting)+1, MPI_CHAR, 0, 0,  
20                  MPI_COMM_WORLD);  
21    } else {  
22        printf("Greetings from process %d of %d!\n", my_rank, comm_sz);  
23        for (int q = 1; q < comm_sz; q++) {  
24            MPI_Recv(greeting, MAX_STRING, MPI_CHAR, q,  
25                      0, MPI_COMM_WORLD, MPI_STATUS_IGNORE);  
26            printf("%s\n", greeting);  
27        }  
28    }  
29  
30    MPI_Finalize();  
31    return 0;  
32 } /* main */
```



Compilation

wrapper script to compile

`mpicc -g -Wall -o mpi_hello mpi_hello.c`

source file

produce debugging information

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

turns on all warnings

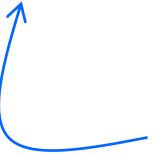


Execution

`mpirun -n <number of processes> <executable>`

`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



```
int MPI_Comm_size(  
    MPI_Comm    comm          /* in  */,  
    int*        comm_sz_p     /* out */);
```

number of processes in the communicator

```
int MPI_Comm_rank(  
    MPI_Comm    comm          /* in  */,  
    int*        my_rank_p     /* out */);
```

my rank

(the process making this call)



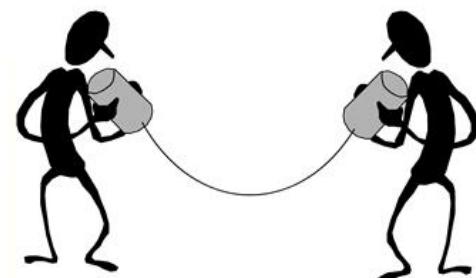
SPMD

- ▶ Single-Program Multiple-Data
- ▶ We compile one 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(  
    void*           msg_buf_p /* in */,  
    int             msg_size   /* in */,  
    MPI_Datatype   msg_type   /* in */,  
    int             dest       /* in */,  
    int             tag        /* in */,  
    MPI_Comm       communicator /* in */);
```



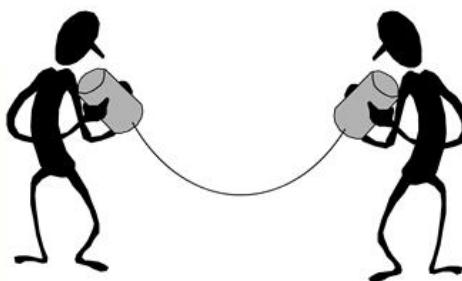
Data types

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	

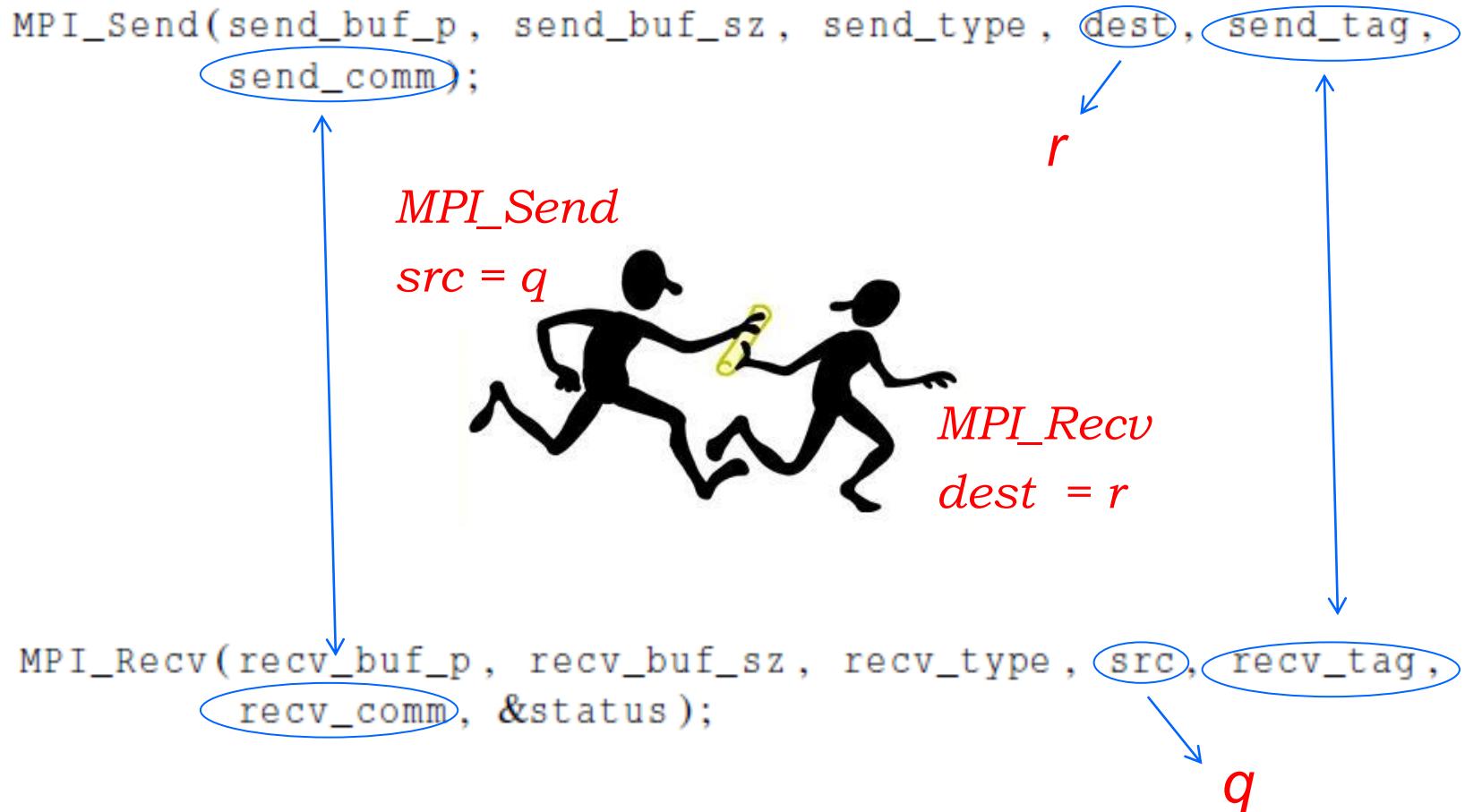


Communication

```
int MPI_Recv(  
    void*           msg_buf_p    /* out */,  
    int             buf_size     /* in */,  
    MPI_Datatype   buf_type    /* in */,  
    int             source       /* in */,  
    int             tag          /* in */,  
    MPI_Comm        communicator /* in */,  
    MPI_Status*    status_p     /* out */);
```

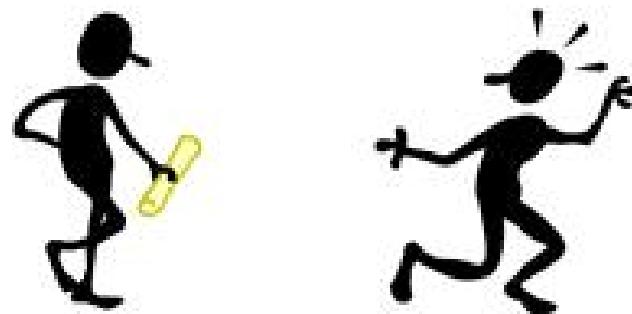


Message matching



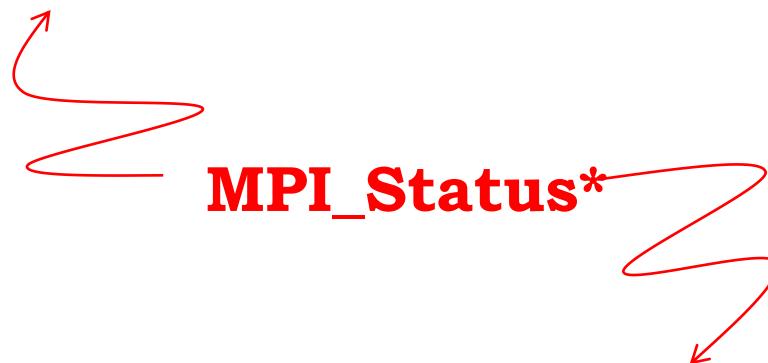
Receiving messages

- ▶ 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_Recv(recv_buf_p, recv_buf_sz, recv_type, src, recv_tag,  
recv_comm, &status);
```



```
MPI_Status* status;  
status.MPI_SOURCE  
status.MPI_TAG
```

MPI_SOURCE
MPI_TAG
MPI_ERROR



How much data am I receiving?

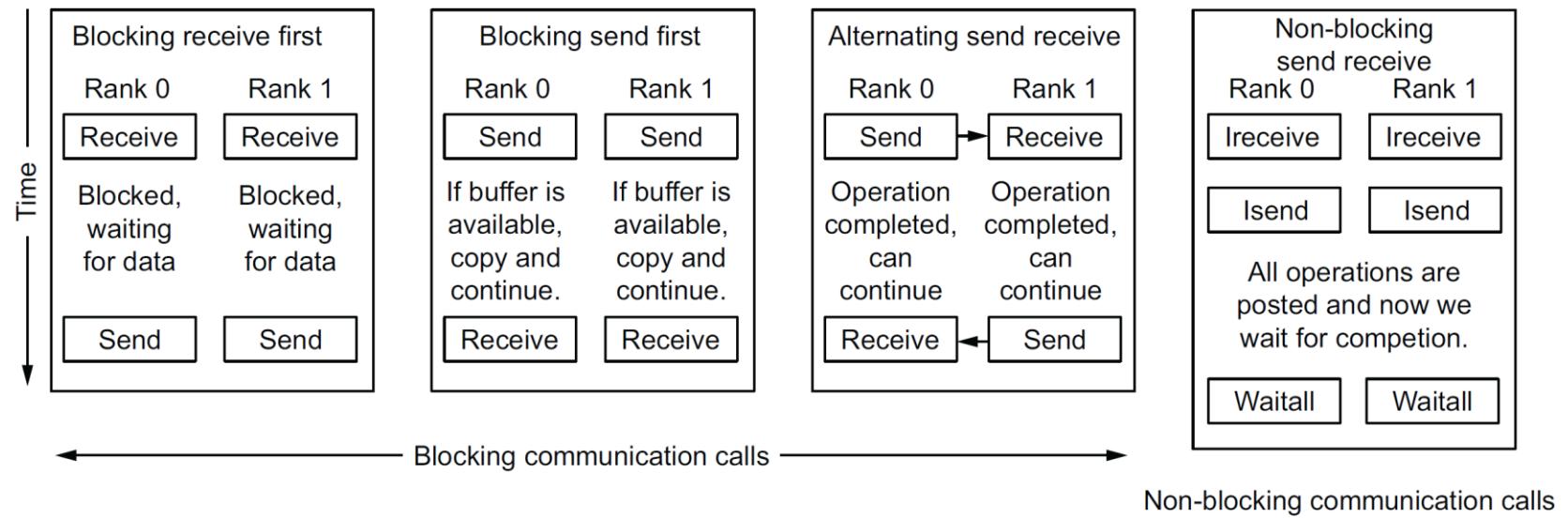
```
int MPI_Get_count(  
    MPI_Status* status_p /* in */,  
    MPI_Datatype type      /* in */,  
    int*           count_p /* out */);
```

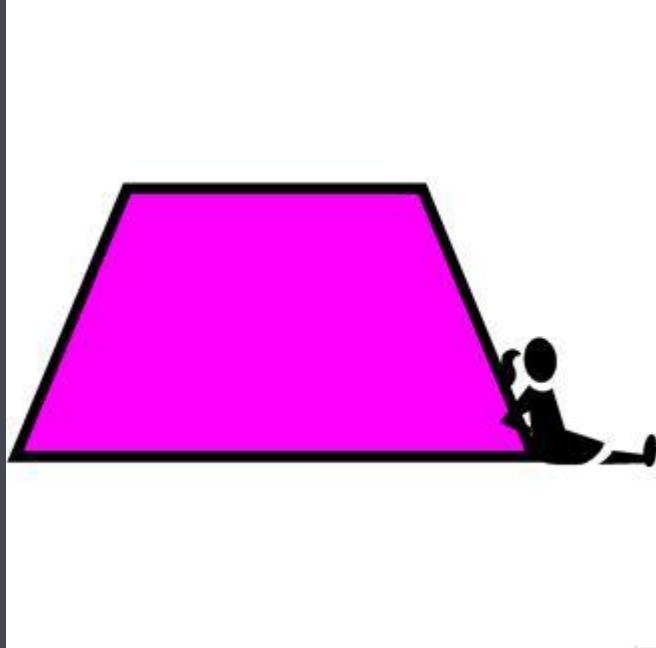


Issues with send and receive

- ▶ 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!

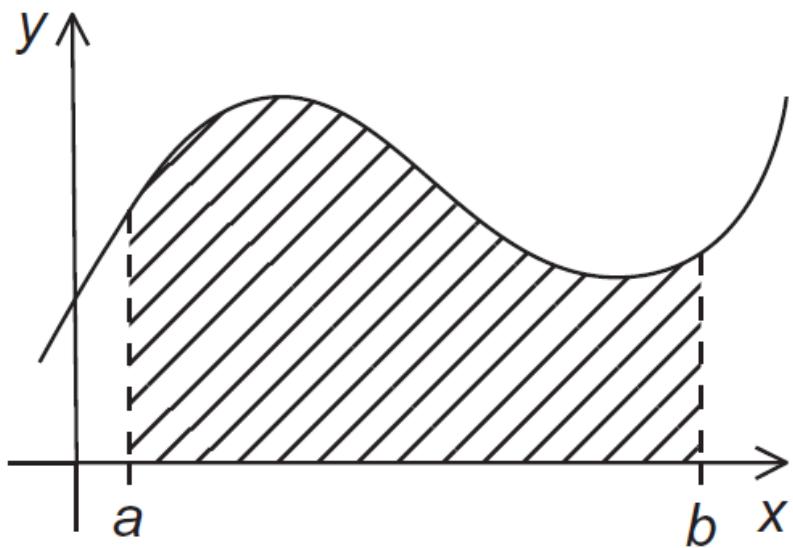




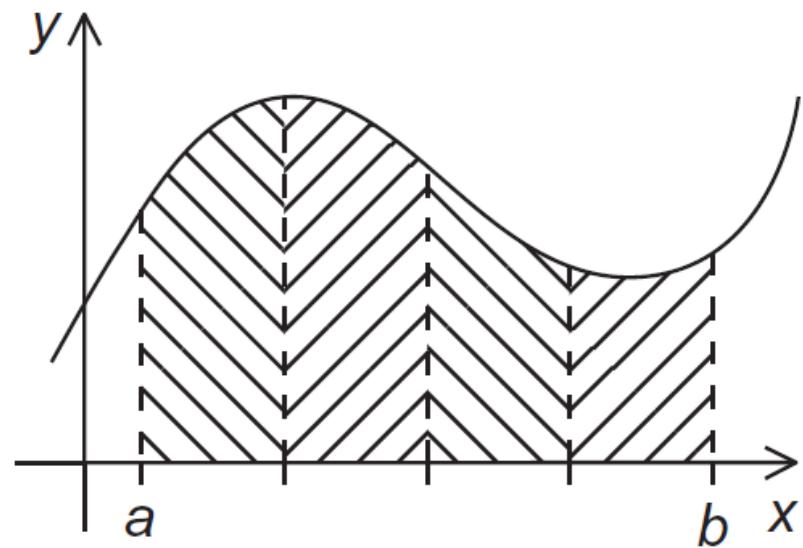


Trapezoidal rule in mpi

The Trapezoidal Rule



(a)



(b)



The Trapezoidal Rule

$$\text{Area of one trapezoid} = \frac{h}{2}[f(x_i) + f(x_{i+1})]$$

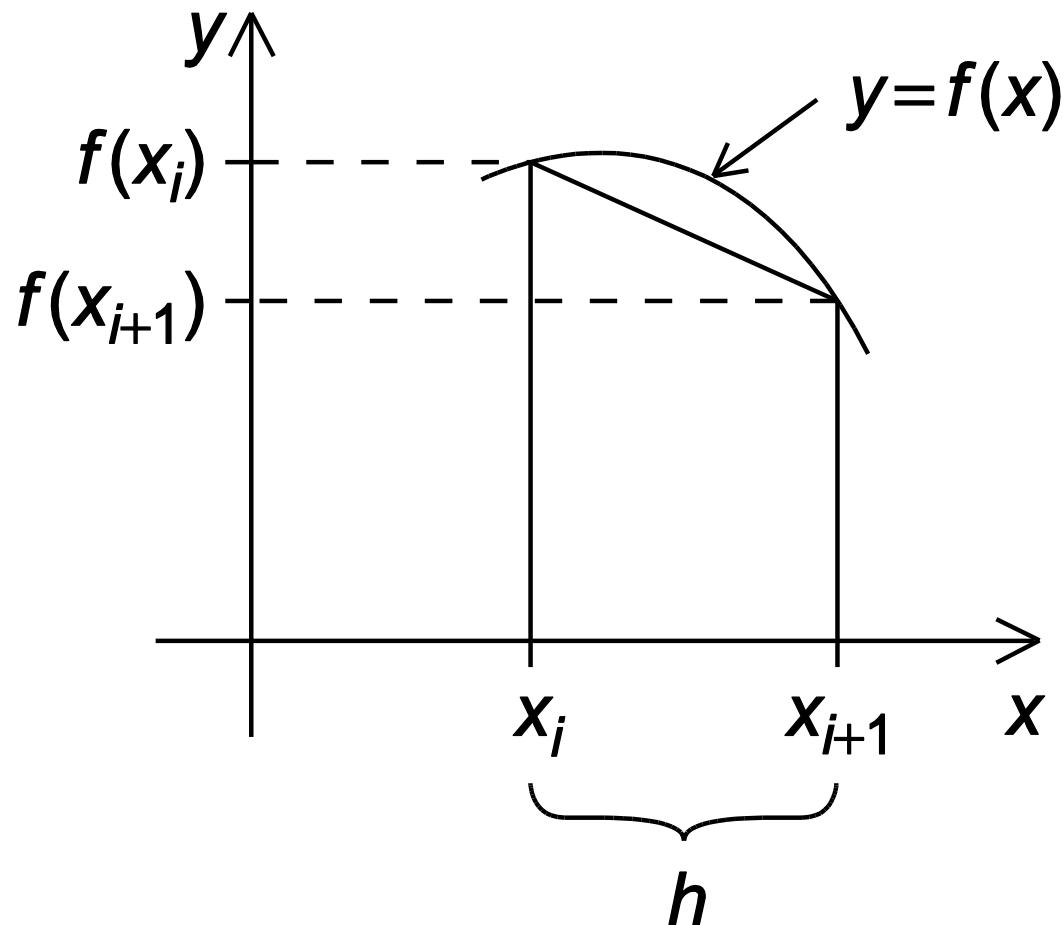
$$h = \frac{b - a}{n}$$

$$x_0 = a, x_1 = a + h, x_2 = a + 2h, \dots, x_{n-1} = a + (n-1)h, x_n = b$$

$$\text{Sum of trapezoid areas} = h[f(x_0)/2 + f(x_1) + f(x_2) + \cdots + f(x_{n-1}) + f(x_n)/2]$$



One trapezoid



Pseudo-code for a serial program

```
/* Input: a, b, n */
h = (b-a)/n;
approx = (f(a) + f(b))/2.0;
for (i = 0; i <= n-1; i++) {
    x_i = a + i*h;
    approx += f(x_i);
}
approx = h*approx;
```



Parallelizing the Trapezoidal Rule

1. Partition problem solution into tasks.
2. Identify communication channels between tasks.
3. Aggregate tasks into composite tasks.
4. Map composite tasks to cores.

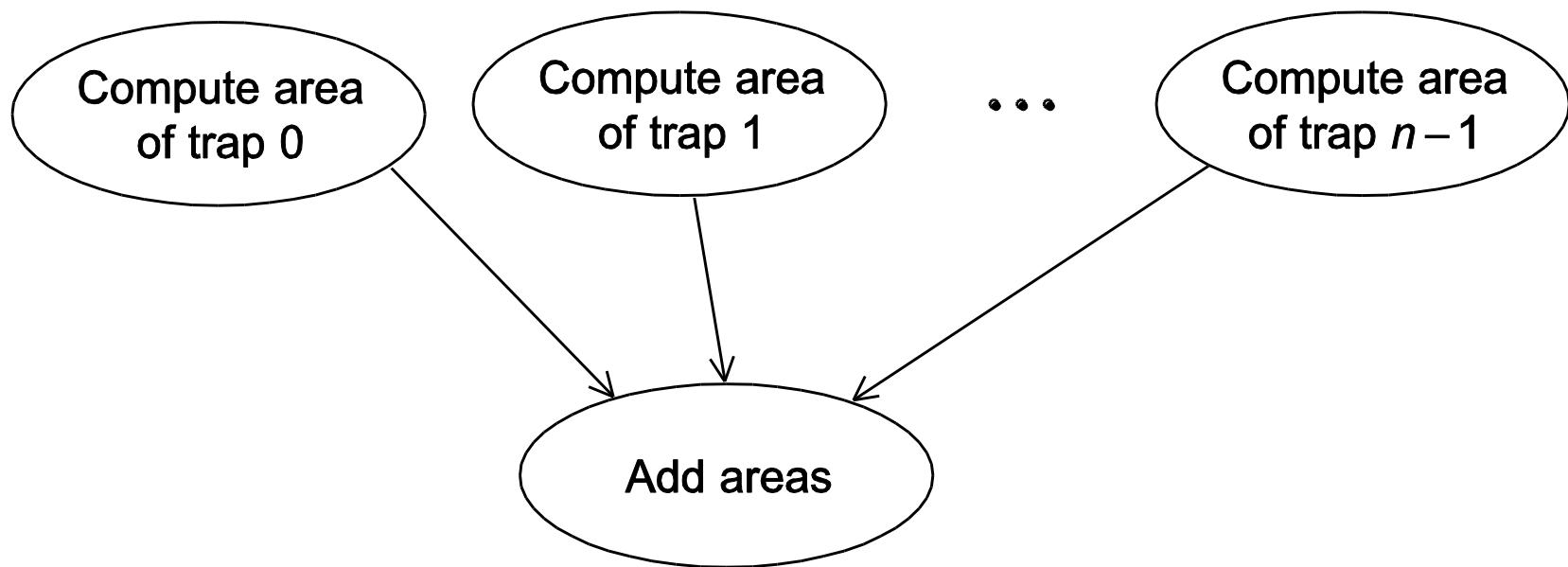


Parallel pseudo-code

```
1      Get a, b, n;
2      h = (b-a)/n;
3      local_n = n/comm_sz;
4      local_a = a + my_rank*local_n*h;
5      local_b = local_a + local_n*h;
6      local_integral = Trap(local_a, local_b, local_n, h);
7      if (my_rank != 0)
8          Send local_integral to process 0;
9      else /* my_rank == 0 */
10         total_integral = local_integral;
11         for (proc = 1; proc < comm_sz; proc++) {
12             Receive local_integral from proc;
13             total_integral += local_integral;
14         }
15     }
16     if (my_rank == 0)
17         print result;
```



Tasks and communications for Trapezoidal Rule



First version (1)

```
1 int main(void) {
2     int my_rank, comm_sz, n = 1024, local_n;
3     double a = 0.0, b = 3.0, h, local_a, local_b;
4     double local_int, total_int;
5     int source;
6
7     MPI_Init(NULL, NULL);
8     MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);
9     MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);
10
11    h = (b-a)/n;           /* h is the same for all processes */
12    local_n = n/comm_sz;   /* So is the number of trapezoids */
13
14    local_a = a + my_rank*local_n*h;
15    local_b = local_a + local_n*h;
16    local_int = Trap(local_a, local_b, local_n, h);
17
18    if (my_rank != 0) {
19        MPI_Send(&local_int, 1, MPI_DOUBLE, 0, 0,
20                  MPI_COMM_WORLD);
```



First version (2)

```
21 } else {
22     total_int = local_int;
23     for (source = 1; source < comm_sz; source++) {
24         MPI_Recv(&local_int, 1, MPI_DOUBLE, source, 0,
25                  MPI_COMM_WORLD, MPI_STATUS_IGNORE);
26         total_int += local_int;
27     }
28 }
29
30 if (my_rank == 0) {
31     printf("With n = %d trapezoids, our estimate\n", n);
32     printf("of the integral from %f to %f = %.15e\n",
33            a, b, total_int);
34 }
35 MPI_Finalize();
36 return 0;
37 } /* main */
```



First version (3)

```
1 double Trap(
2     double left_endpt /* in */,
3     double right_endpt /* in */,
4     int trap_count /* in */,
5     double base_len /* in */) {
6     double estimate, x;
7     int i;
8
9     estimate = (f(left_endpt) + f(right_endpt))/2.0;
10    for (i = 1; i <= trap_count -1; i++) {
11        x = left_endpt + i*base_len;
12        estimate += f(x);
13    }
14    estimate = estimate*base_len;
15
16    return estimate;
17 } /* Trap */
```



Dealing with I/O

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

int main(void) {
    int my_rank, comm_sz;

    MPI_Init(NULL, NULL);
    MPI_Comm_size(MPI_COMM_WORLD, &comm_sz);
    MPI_Comm_rank(MPI_COMM_WORLD, &my_rank);

    printf("Proc %d of %d > Does anyone have a toothpick?\n",
           my_rank, comm_sz);

    MPI_Finalize();
    return 0;
} /* main */
```

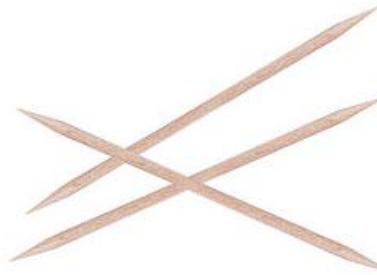
Each process just prints a message.



Running with 6 processes

```
Proc 0 of 6 > Does anyone have a toothpick?  
Proc 1 of 6 > Does anyone have a toothpick?  
Proc 2 of 6 > Does anyone have a toothpick?  
Proc 4 of 6 > Does anyone have a toothpick?  
Proc 3 of 6 > Does anyone have a toothpick?  
Proc 5 of 6 > Does anyone have a toothpick?
```

unpredictable output



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;
    . . .
```

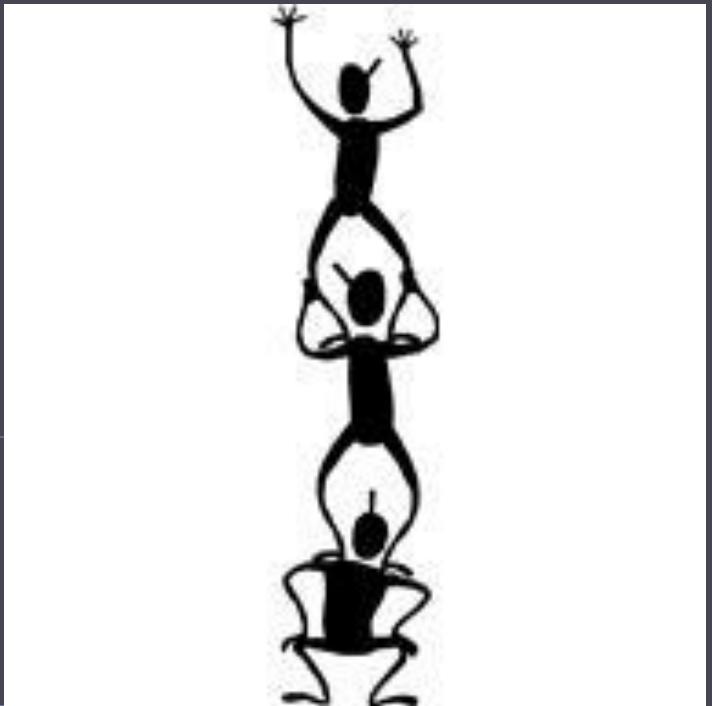


Function for reading user input

```
void Get_input(
    int      my_rank    /* in  */,
    int      comm_sz   /* in  */,
    double*  a_p        /* out */,
    double*  b_p        /* out */,
    int*     n_p        /* out */) {
    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++) {
            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

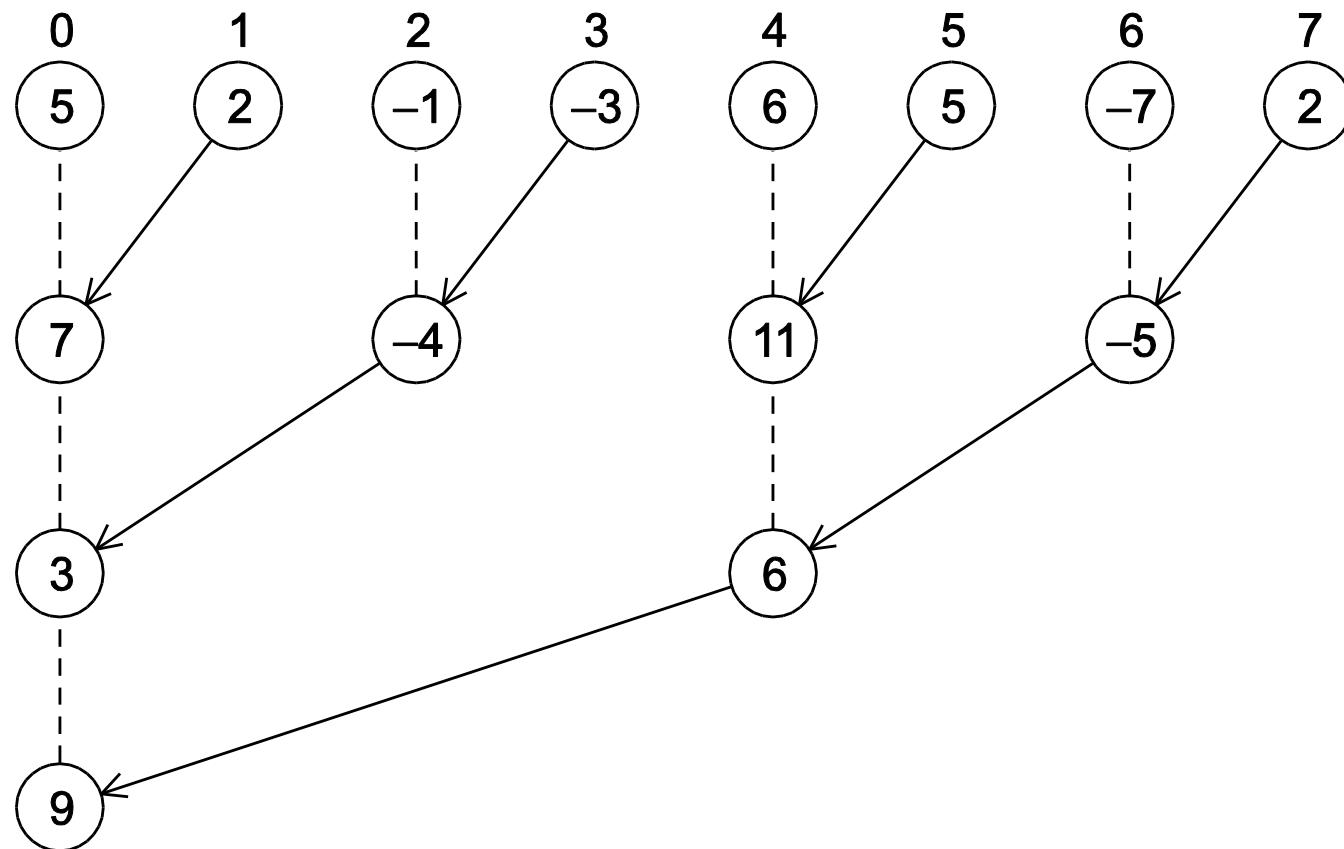
Tree-structured communication

- I. 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.

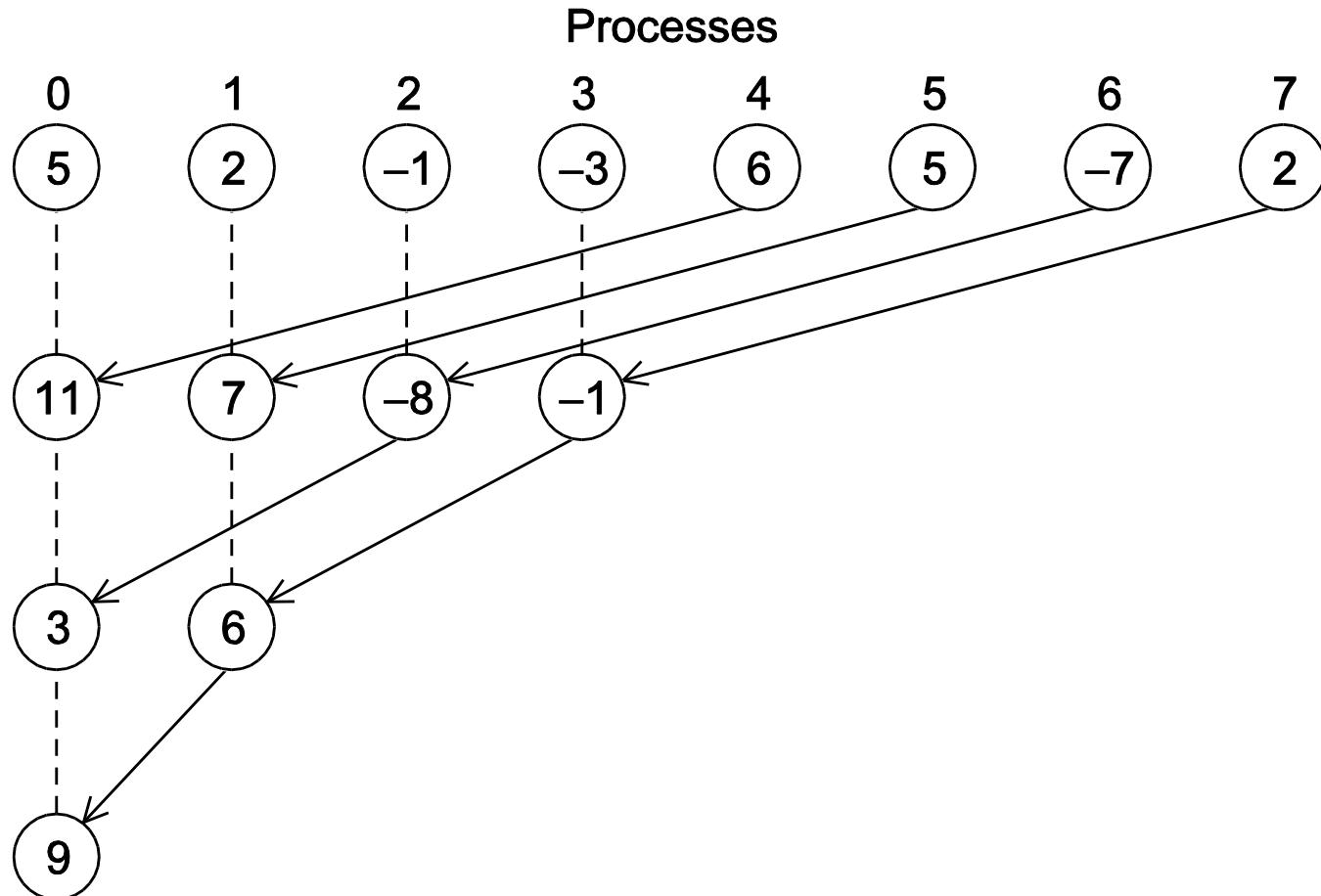


A tree-structured global sum

Processes



An alternative tree-structured global sum



MPI_Reduce

```
int MPI_Reduce(  
    void*           input_data_p /* in */,  
    void*           output_data_p /* out */,  
    int             count        /* in */,  
    MPI_Datatype   datatype    /* in */,  
    MPI_Op          operator     /* in */,  
    int             dest_process /* in */,  
    MPI_Comm        comm        /* in */);
```

```
MPI_Reduce(&local_int, &total_int, 1, MPI_DOUBLE, MPI_SUM, 0,  
MPI_COMM_WORLD);
```

```
double local_x[N], sum[N];  
.  
.  
.  
MPI_Reduce(local_x, sum, N, MPI_DOUBLE, MPI_SUM, 0,  
MPI_COMM_WORLD);
```



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



Collective vs. Point-to-Point Communications

- ▶ 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.



Collective vs. Point-to-Point Communications

- ▶ 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.



Collective vs. Point-to-Point Communications

- ▶ 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`.



Collective vs. Point-to-Point Communications

- ▶ 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 $2+1+2 = 5$.

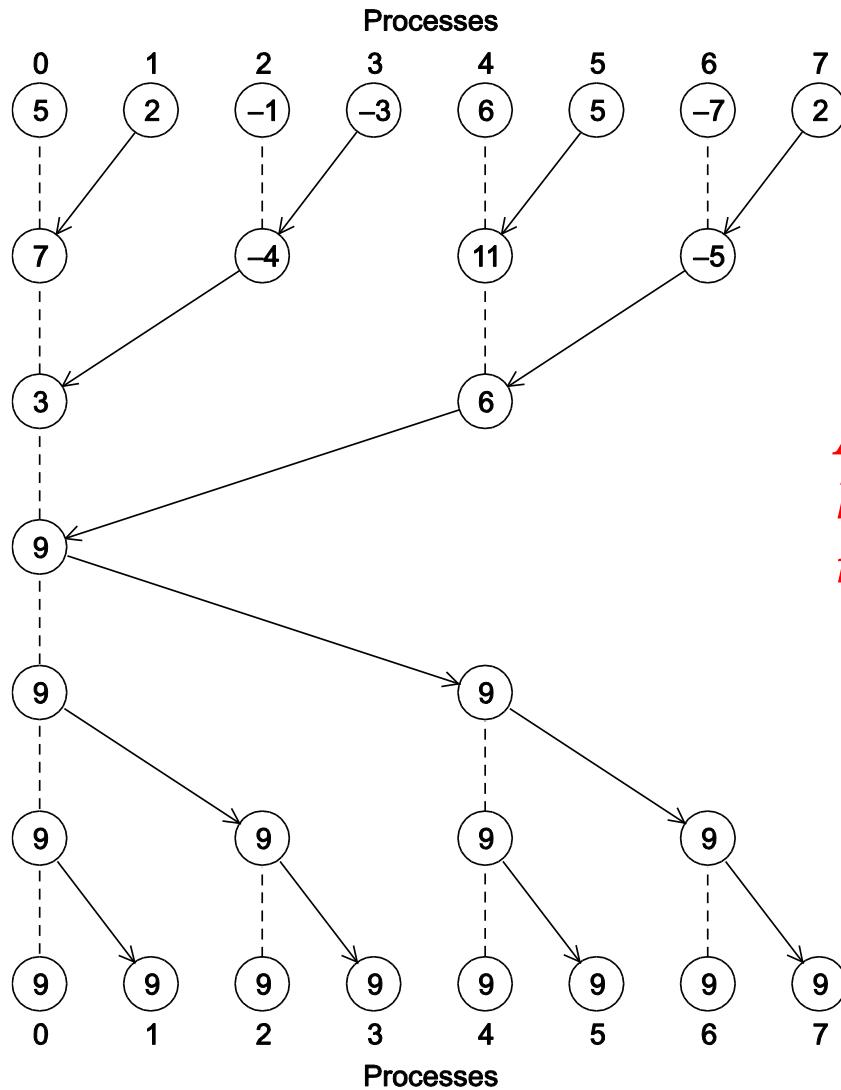


MPI_Allreduce

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

```
int MPI_Allreduce(  
    void *           input_data_p /* in */,  
    void *           output_data_p /* out */,  
    int              count        /* in */,  
    MPI_Datatype     datatype   /* in */,  
    MPI_Op            operator    /* in */,  
    MPI_Comm          comm        /* in */);
```



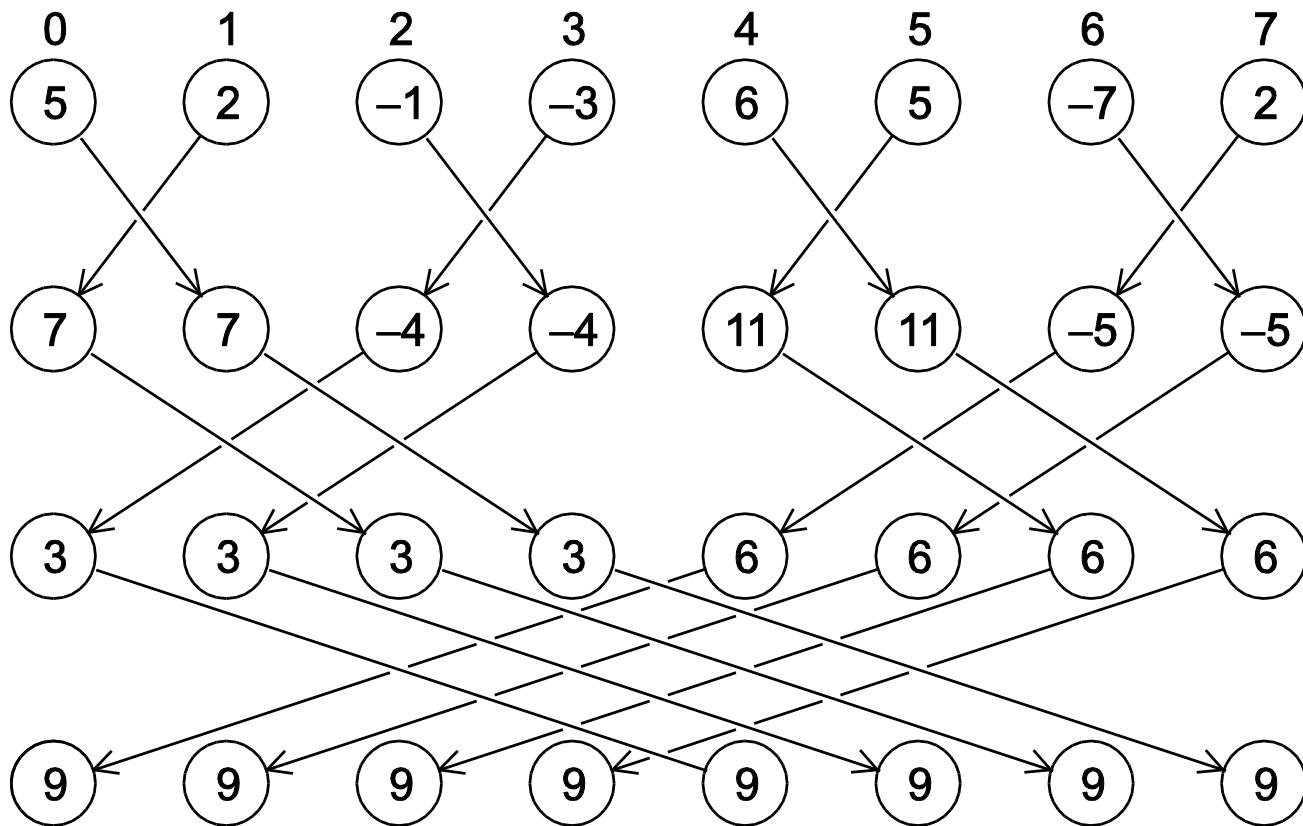


*A global sum followed
by distribution of the
result.*





Processes



A butterfly-structured global sum.

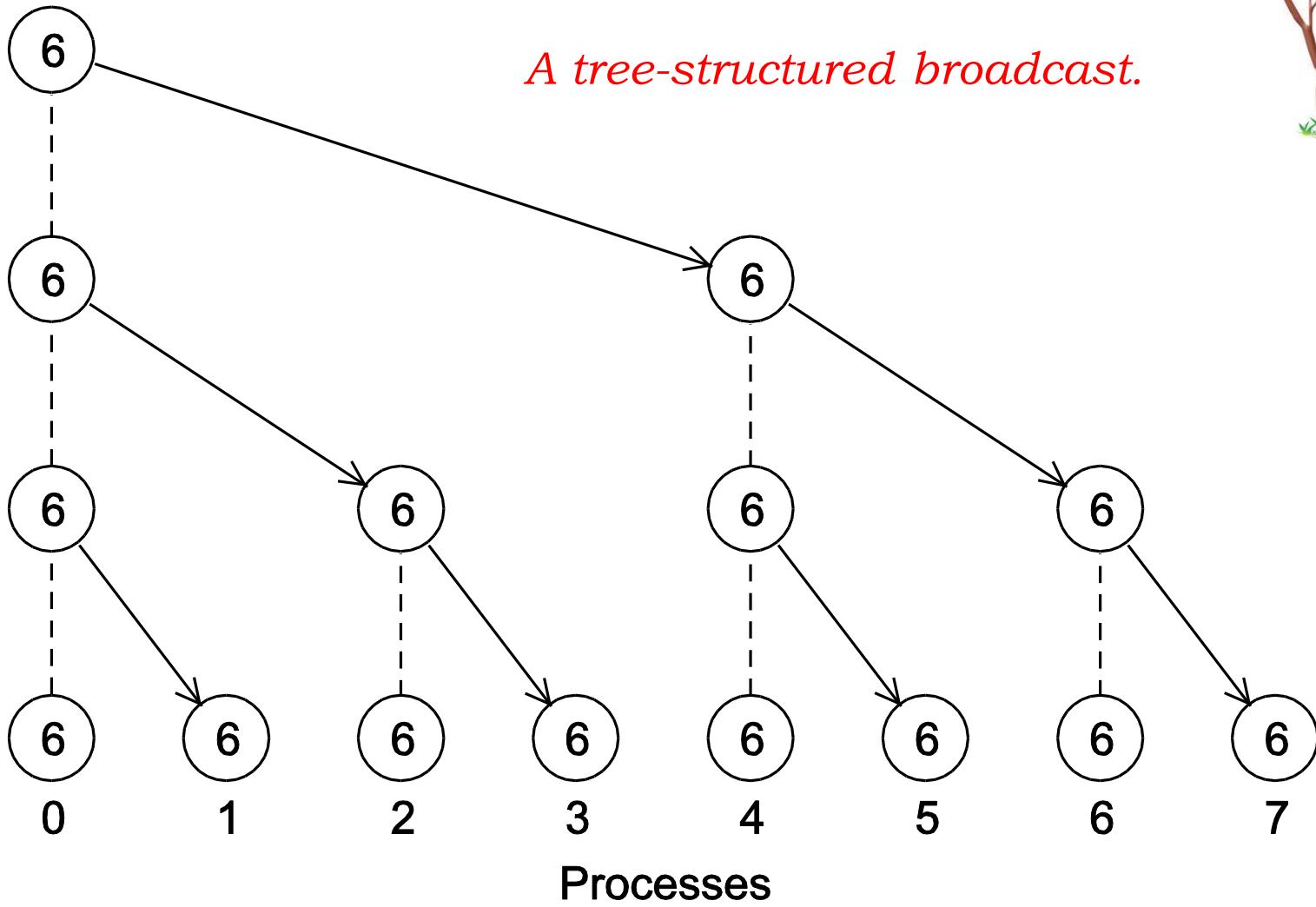


Broadcast

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

```
int MPI_Bcast(  
    void*           data_p          /* in/out */,  
    int             count           /* in */,  
    MPI_Datatype   datatype        /* in */,  
    int             source_proc    /* in */,  
    MPI_Comm       comm            /* in */);
```





A tree-structured broadcast.



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

$$\begin{aligned}\mathbf{x} + \mathbf{y} &= (x_0, x_1, \dots, x_{n-1}) + (y_0, y_1, \dots, y_{n-1}) \\ &= (x_0 + y_0, x_1 + y_1, \dots, x_{n-1} + y_{n-1}) \\ &= (z_0, z_1, \dots, z_{n-1}) \\ &= \mathbf{z}\end{aligned}$$

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 */
```



Different partitions of a 12-component vector among 3 processes

Process	Components												Block-cyclic Blocksize = 2						
	Block					Cyclic				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 */
```



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.

```
int MPI_Scatter(
    void*          send_buf_p /* in */,
    int            send_count /* in */,
    MPI_Datatype  send_type  /* in */,
    void*          recv_buf_p /* out */,
    int            recv_count /* in */,
    MPI_Datatype  recv_type  /* in */,
    int            src_proc   /* in */,
    MPI_Comm       comm       /* in */);
```



Reading and distributing a vector

```
void Read_vector(
    double    local_a[]    /* out */,
    int       local_n      /* in */,
    int       n            /* in */,
    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, 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 */,  
    MPI_Datatype   send_type  /* in */,  
    void*           recv_buf_p /* out */,  
    int             recv_count /* in */,  
    MPI_Datatype   recv_type  /* in */,  
    int             dest_proc  /* in */,  
    MPI_Comm        comm       /* in */);
```



Print a distributed vector (1)

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

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.

```
int MPI_Allgather(  
    void* send_buf_p /* in */,  
    int send_count /* in */,  
    MPI_Datatype send_type /* in */,  
    void* recv_buf_p /* out */,  
    int recv_count /* in */,  
    MPI_Datatype recv_type /* in */,  
    MPI_Comm comm /* in */);
```



Matrix-vector multiplication

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

\mathbf{x} is a vector with n components

$\mathbf{y} = A\mathbf{x}$ 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

a_{00}	a_{01}	\cdots	$a_{0,n-1}$
a_{10}	a_{11}	\cdots	$a_{1,n-1}$
\vdots	\vdots		\vdots
a_{i0}	a_{i1}	\cdots	$a_{i,n-1}$
\vdots	\vdots		\vdots
$a_{m-1,0}$	$a_{m-1,1}$	\cdots	$a_{m-1,n-1}$

$$\begin{matrix} x_0 \\ x_1 \\ \vdots \\ x_{n-1} \end{matrix} = \begin{matrix} y_0 \\ y_1 \\ \vdots \\ y_i = a_{i0}x_0 + a_{i1}x_1 + \cdots + a_{i,n-1}x_{n-1} \\ \vdots \\ y_{m-1} \end{matrix}$$



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];  
}
```

Serial pseudo-code



C style arrays

$$\begin{pmatrix} 0 & 1 & 2 & 3 \\ 4 & 5 & 6 & 7 \\ 8 & 9 & 10 & 11 \end{pmatrix}$$

stored as

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];
    }
}
```



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;
```

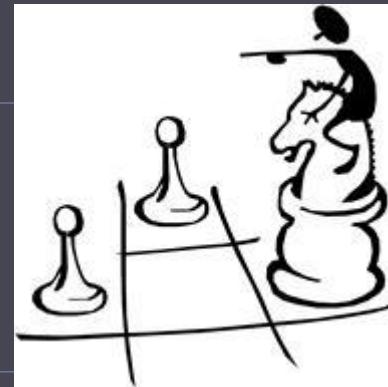


An MPI matrix-vector multiplication function (2)

```
x = malloc(n*sizeof(double));
MPI_Allgather(local_x, local_n, MPI_DOUBLE,
              x, local_n, MPI_DOUBLE, comm);

for (local_i = 0; local_i < local_m; local_i++) {
    local_y[local_i] = 0.0;
    for (j = 0; j < n; j++)
        local_y[local_i] += local_A[local_i*n+j]*x[j];
}
free(x);
} /* Mat_vect_mult */
```





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

{(MPI_DOUBLE,0),(MPI_DOUBLE,16),(MPI_INT,24)}



MPI_Type create_struct

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

```
int MPI_Type_create_struct(
    int             count          /* in     */,
    int             array_of_blocklengths[] /* in     */,
    MPI_Aint        array_of_displacements[] /* in     */,
    MPI_Datatype   array_of_types[]      /* in     */,
    MPI_Datatype*  new_type_p        /* out    */);
```



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)

```
void Build_mpi_type(
    double*          a_p           /* in */,
    double*          b_p           /* in */,
    int*             n_p           /* in */,
    MPI_Datatype*   input_mpi_t_p /* out */) {

    int array_of_blocklengths[3] = {1, 1, 1};
    MPI_Datatype array_of_types[3] = {MPI_DOUBLE, MPI_DOUBLE, MPI_INT};
    MPI_Aint a_addr, b_addr, n_addr;
    MPI_Aint array_of_displacements[3] = {0};
```



Get input function with a derived datatype (2)

```
MPI_Get_address(a_p, &a_addr);
MPI_Get_address(b_p, &b_addr);
MPI_Get_address(n_p, &n_addr);
array_of_displacements[1] = b_addr-a_addr;
array_of_displacements[2] = n_addr-a_addr;
MPI_Type_create_struct(3, array_of_blocklengths,
                      array_of_displacements, array_of_types,
                      input_mpi_t_p);
MPI_Type_commit(input_mpi_t_p);
} /* Build_mpi_type */
```



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 */*



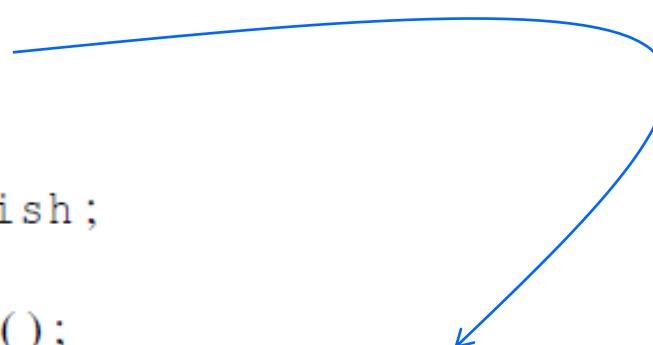


Performance evaluation

Elapsed parallel time

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

```
double MPI_Wtime( void );  
  
double start, finish;  
. . .  
start = MPI_Wtime();  
/* Code to be timed */  
. . .  
finish = MPI_Wtime();  
printf("Proc %d > Elapsed time = %e seconds\n"  
      my_rank, finish-start);
```



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.

```
int MPI_Barrier(MPI_Comm comm /* in */);
```



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

comm_sz	Order of Matrix				
	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

comm_sz	Order of Matrix				
	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

comm_sz	Order of Matrix				
	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 $p = \text{comm sz}$ 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 $0 \leq q < r < p$, then each key assigned to process q should be less than or equal to every key assigned to process r .



Serial bubble sort

```
void Bubble_sort(  
    int a[] /* in/out */,  
    int n      /* in      */) {  
    int list_length, i, temp;  
  
    for (list_length = n; list_length >= 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



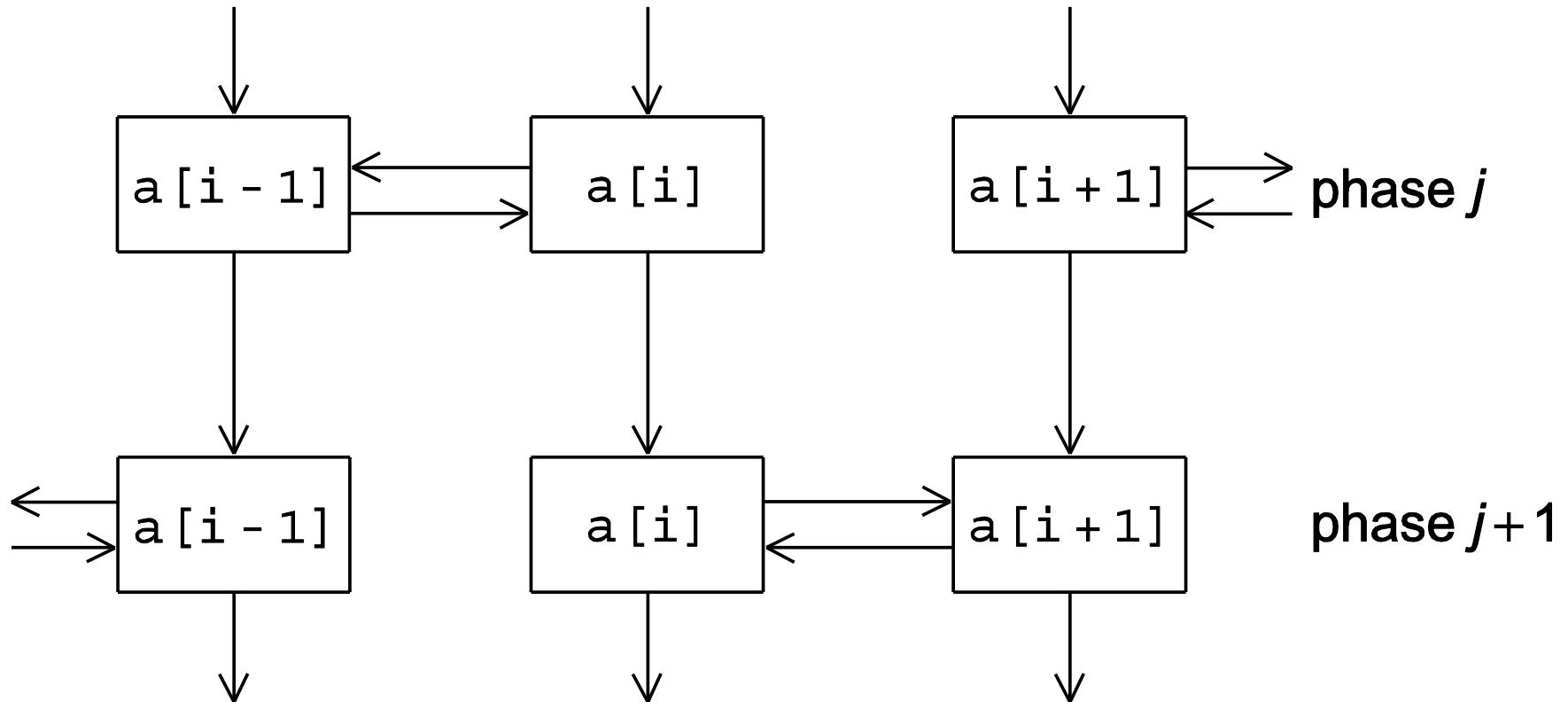
Serial odd-even transposition sort

```
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

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



Safety in MPI programs

- ▶ 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.



Safety in MPI programs

- ▶ 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.



Safety in MPI programs

- ▶ 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*)



Safety in MPI programs

- ▶ 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.

```
int MPI_Ssend(  
    void*           msg_buf_p      /* in */,  
    int             msg_size       /* in */,  
    MPI_Datatype   msg_type       /* in */,  
    int             dest           /* in */,  
    int             tag            /* in */,  
    MPI_Comm        communicator /* in */);
```



Restructuring communication

```
MPI_Send(msg, size, MPI_INT, (my_rank+1) % comm_sz, 0, comm);
MPI_Recv(new_msg, size, MPI_INT, (my_rank+comm_sz-1) % comm_sz,
         0, comm, MPI_STATUS_IGNORE.
```



```
if (my_rank % 2 == 0) {
    MPI_Send(msg, size, MPI_INT, (my_rank+1) % comm_sz, 0, comm);
    MPI_Recv(new_msg, size, MPI_INT, (my_rank+comm_sz-1) % comm_sz,
             0, comm, MPI_STATUS_IGNORE.
} else {
    MPI_Recv(new_msg, size, MPI_INT, (my_rank+comm_sz-1) % comm_sz,
             0, comm, MPI_STATUS_IGNORE.
    MPI_Send(msg, size, MPI_INT, (my_rank+1) % comm_sz, 0, comm);
}
```



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.

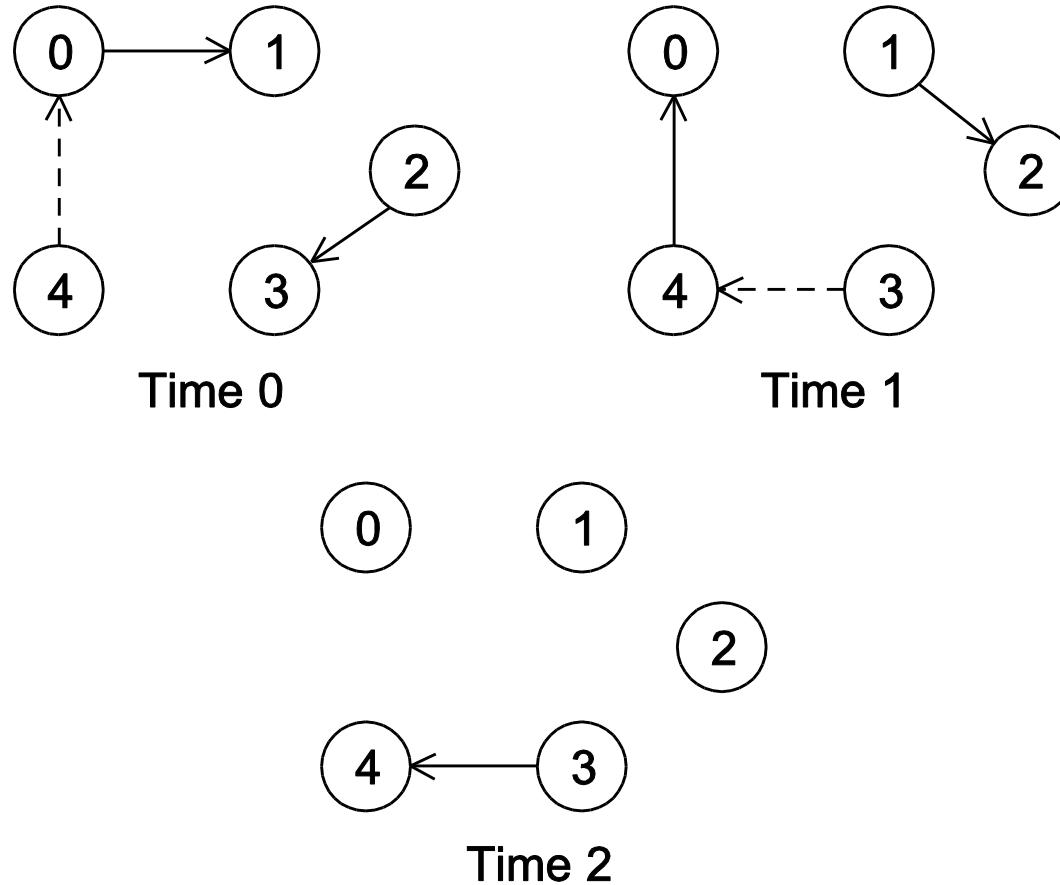


MPI_Sendrecv

```
int MPI_Sendrecv(
    void* send_buf_p          /* in */,
    int send_buf_size          /* in */,
    MPI_Datatype send_buf_type /* in */,
    int dest                  /* in */,
    int send_tag               /* in */,
    void* recv_buf_p           /* out */,
    int recv_buf_size          /* in */,
    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 m_i, r_i, t_i;

    m_i = r_i = t_i = 0;
    while (t_i < local_n) {
        if (my_keys[m_i] <= recv_keys[r_i]) {
            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

Processes	Number of Keys (in thousands)				
	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.

