

# **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.





#### **Last session**

Mineran Hum 1

all



## **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.



# Safety in MPI programs, Deadlock?



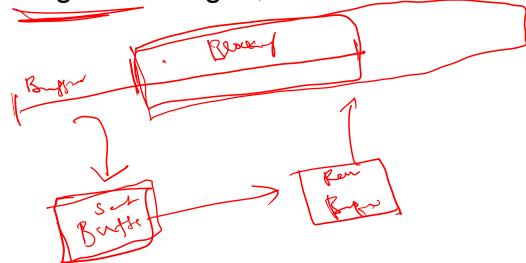
- The MPI standard allows MPI\_Send to behave in two different ways:
  - it can simply copy the message into an MPI managed buffer and return,
  - or it can block until the matching call to MPI\_Recv starts.

# 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.



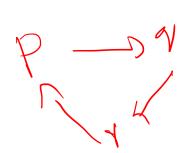




# Safety in MPI programs



- If the MPI\_Send is executed by each process block, 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 <u>buffering</u> is said to be <u>unsafe</u>.
- 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 <u>synchronous</u> 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

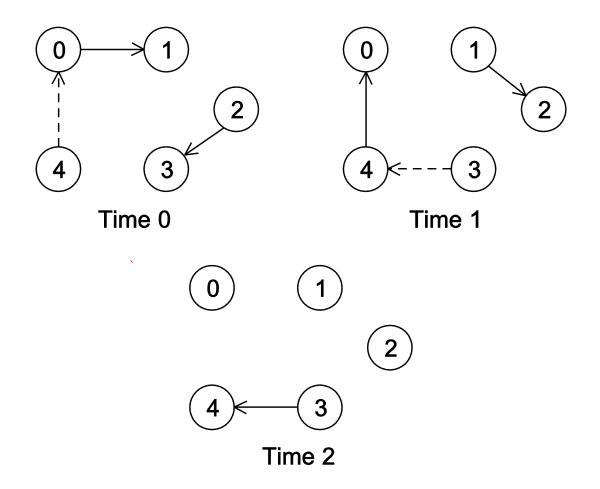


```
int MPI_Sendrecv(
    void*
         send_buf_p /*in */,
         send_buf_size /*in */,
    int
    MPI_Datatype send_buf_type /*in */,
               dest /*in */,
    int
               send_tag /*in */,
    int
             recv_buf_p /* out */,
    void*
         recv_buf_size /*in */,
    int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

MRZ

(times are in milliseconds)



## 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 <u>buffer size</u>, 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**

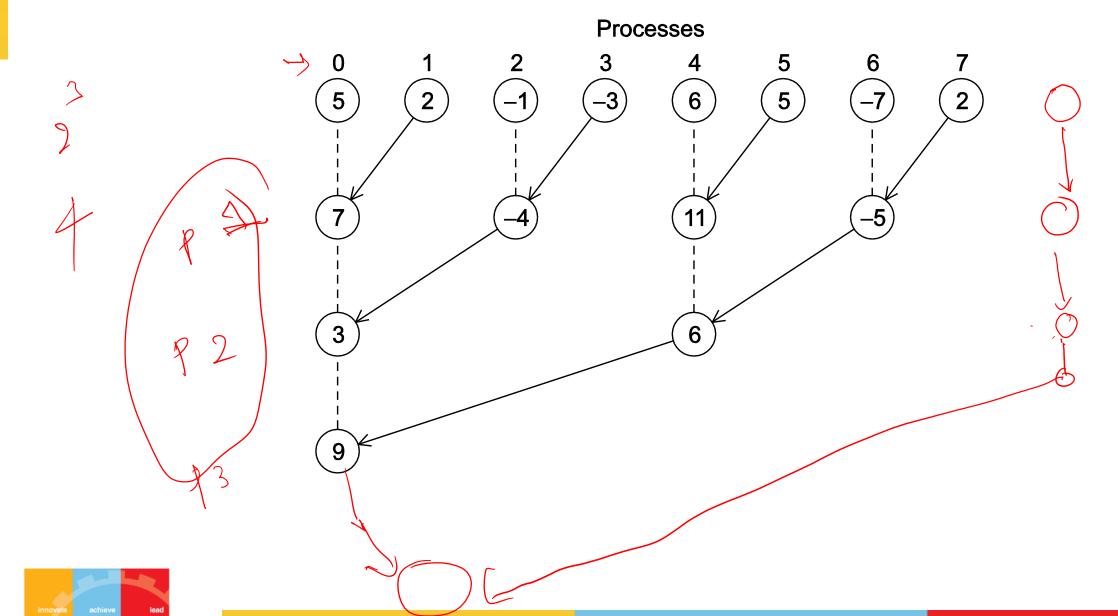


- 1. In the first phase:
  - (a) Process 1 sends to 0, 3 sends to 2, 5 sends to 4, and 7 sends to 6.
  - (b) Processes 0, 2, 4, and 6 add in the received values.
  - (c) Processes 2 and 6 send their new values to processes 0 and 4, respectively.
  - (d) Processes 0 and 4 add the received values into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



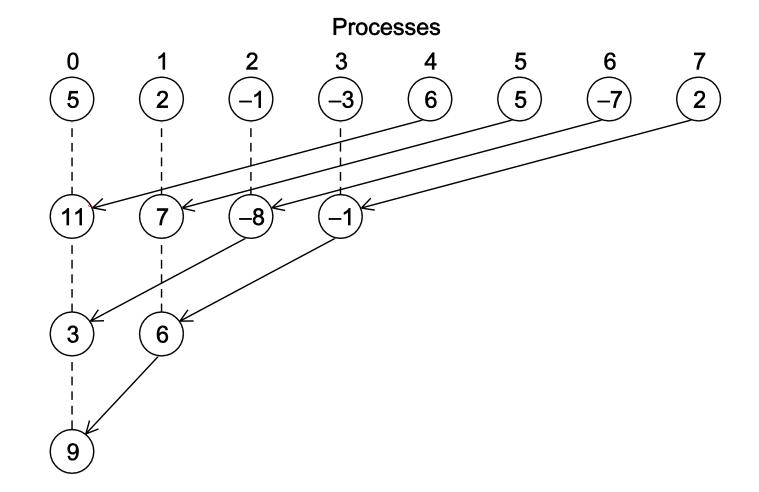
# A tree-structured global sum





# 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

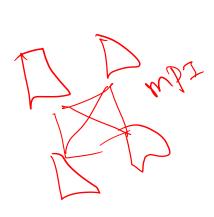


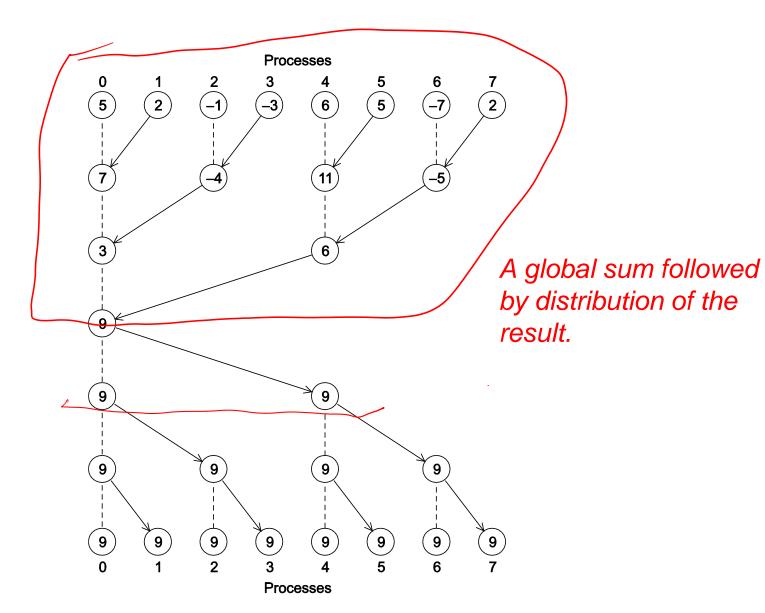
 Useful in a situation in which all of the processes need the result of a global sum in order to complete some larger computation.









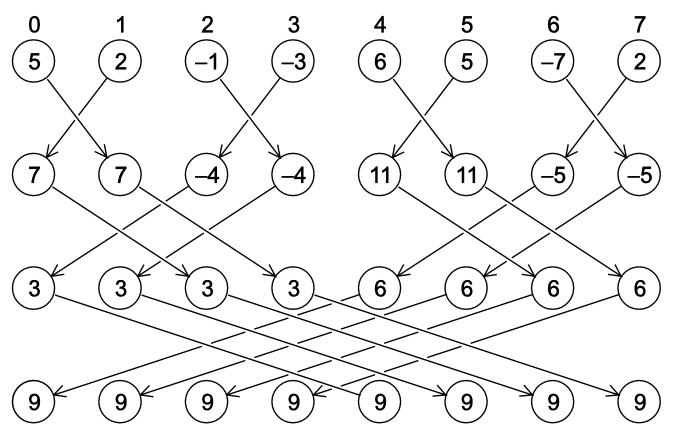








#### **Processes**



A butterfly-structured global sum.



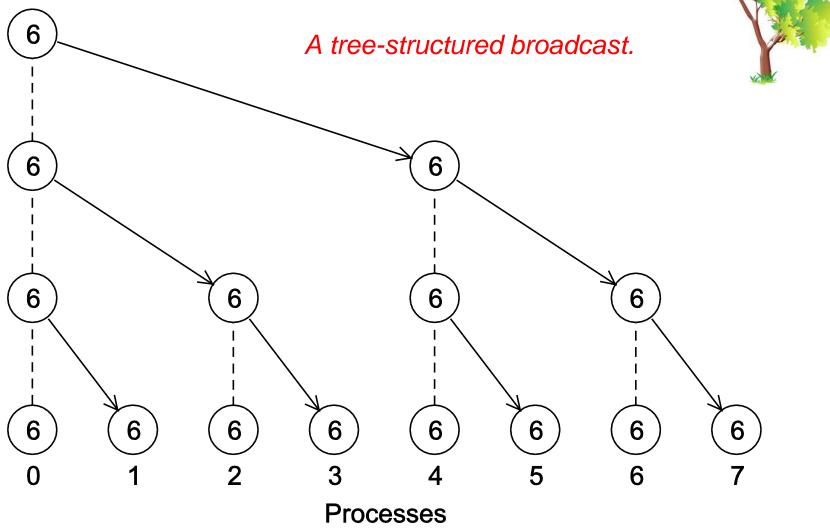
#### **Broadcast**



 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**



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

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

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

$$= \mathbf{z}$$

2+4

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 */,
               send_count /*in */,
    int
               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

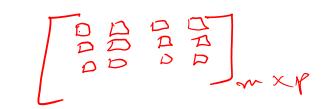
$$y_i = a_{i0}x_0 + a_{i1}x_1 + a_{i2}x_2 + \cdots + a_{i,n-1}x_{n-1}$$

i-th component of y

Dot product of the ith row of A with x.



# **Matrix-vector multiplication**



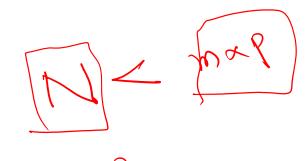


<i>a</i> <sub>00</sub>	<i>a</i> <sub>01</sub>		$a_{0,n-1}$
$a_{10}$	$a_{11}$		$a_{1,n-1}$
:	:		:
$a_{i0}$	$a_{i1}$	•••	$a_{i,n-1}$
<i>a</i> <sub>i0</sub> :	<i>a</i> <sub>i1</sub> :	•••	$a_{i,n-1}$ :

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

1 R 1

 $m \times k$   $m \times k$ 







### 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 */
    v[i] = 0.0;

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

Serial pseudo-code



### C style arrays



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



#### 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

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



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



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



#### 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.

```
\ S & X & X & X
```

```
#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	N
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	

7

(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**

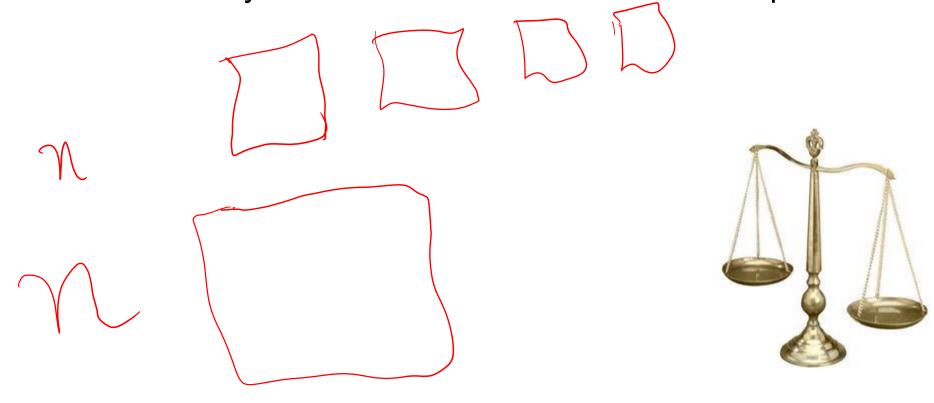


	Order of Matrix					
comm_sz	1024	2048	4096	8192	16,384	
1	1.00	1.00	1.00	1.00	1.00	
2	0.89	0.94	0.97	0.96	0.98	
4	0.51	0.78	0.89	0.96	0.98	
8	0.30	0.61	0.82	0.94	0.98	
16	0.15	0.39	0.68	0.89	0.97	

# **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 <u>number of processes</u> 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 >= 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

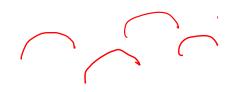
Odd phase: compare-swap (5,4) getting the list 3, 4, 5, 9



## Serial odd-even transposition sort



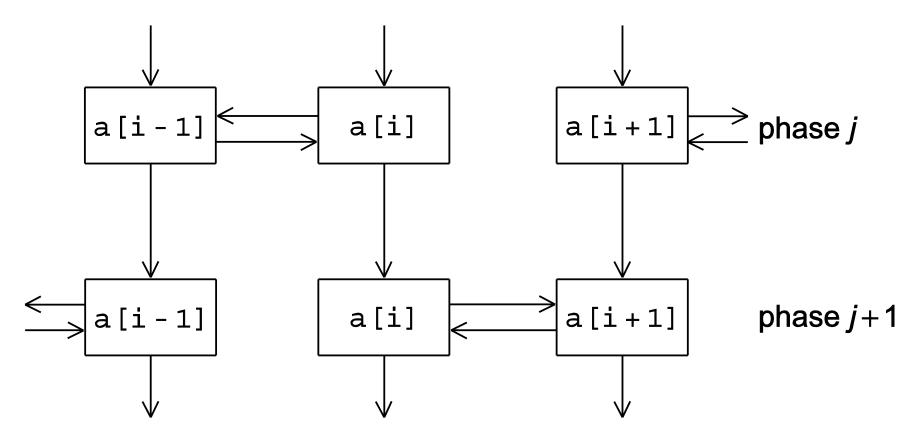
```
void Odd_even_sort(
     int a[] /* in/out */,
     int n /* in */) {
  int phase, i, temp;
  for (phase = 0; phase < n; phase++)
     if (phase % 2 == 0) { /* Even phase */
        for (i = 1; i < n; i += 2)
           if (a[i-1] > a[i]) {
              temp = a[i];
              a[i] = a[i-1];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

#### 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	

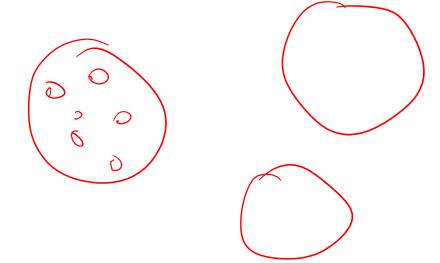
(times are in milliseconds)



# **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.



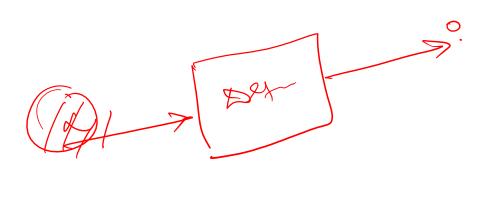


# **Concluding Remarks (2)**



- 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.

