

Looking at our table of parallel efficiencies (Table 3.7), we see that our matrix-vector multiplication program definitely doesn't have the same scalability as program A: in almost every case when  $p$  is increased, the efficiency decreases. On the other hand, the program is somewhat like program B: if  $p \geq 2$  and we increase both  $p$  and  $n$  by a factor of 2, the parallel efficiency, for the most part, actually increases. Furthermore, the only exceptions occur when we increase  $p$  from 2 to 4, and when computer scientists discuss scalability, they're usually interested in large values of  $p$ . When  $p$  is increased from 4 to 8 or from 8 to 16, our efficiency always increases when we increase  $n$  by a factor of 2.

Recall that 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**. Program A is strongly scalable, and program B is weakly scalable. Furthermore, our matrix-vector multiplication program is also apparently weakly scalable.

### 3.7 A PARALLEL SORTING ALGORITHM

What do we mean by a parallel sorting algorithm in a distributed-memory environment? What would its "input" be and what would its "output" be? The answers depend on where the keys are stored. We can start or finish with the keys distributed among the processes or assigned to a single process. In this section we'll look at an algorithm that starts and finishes with the keys distributed among the processes. In Programming Assignment 3.8 we'll look at an algorithm that finishes with the keys assigned to a single process.

If we have a total of  $n$  keys and  $p = \text{comm\_sz}$  processes, our algorithm will start and finish with  $n/p$  keys assigned to each process. (As usual, we'll assume  $n$  is evenly divisible by  $p$ .) At the start, there are no restrictions on which keys are assigned to which processes. However, when the algorithm terminates,

- the keys assigned to each process should be sorted in (say) increasing order, and
- 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  $r$ .

So if we lined up the keys according to process rank—keys from process 0 first, then keys from process 1, and so on—then the keys would be sorted in increasing order. For the sake of explicitness, we'll assume our keys are ordinary **ints**.

#### 3.7.1 Some simple serial sorting algorithms

Before starting, let's look at a couple of simple serial sorting algorithms. Perhaps the best known serial sorting algorithm is bubble sort (see Program 3.14). The array `a` stores the unsorted keys when the function is called, and the sorted keys when the function returns. The number of keys in `a` is `n`. The algorithm proceeds by comparing

```

1 void Bubble_sort(
2     int a[] /* in/out */,
3     int n   /* in   */) {
4     int list_length, i, temp;
5
6     for (list_length = n; list_length >= 2; list_length--)
7         for (i = 0; i < list_length-1; i++)
8             if (a[i] > a[i+1]) {
9                 temp = a[i];
10                a[i] = a[i+1];
11                a[i+1] = temp;
12            }
13
14 } /* Bubble_sort */

```

**Program 3.14:** Serial bubble sort

the elements of the list *a* pairwise: *a*[0] is compared to *a*[1], *a*[1] is compared to *a*[2], and so on. Whenever a pair is out of order, the entries are swapped, so in the first pass through the outer loop, when *list\_length* = *n*, the largest value in the list will be moved into *a*[*n*-1]. The next pass will ignore this last element and it will move the next-to-the-largest element into *a*[*n*-2]. Thus, as *list\_length* decreases, successively more elements get assigned to their final positions in the sorted list.

There isn't much point in trying to parallelize this algorithm because of the inherently sequential ordering of the comparisons. To see this, suppose that *a*[*i*-1] = 9, *a*[*i*] = 5, and *a*[*i*+1] = 7. The algorithm will first compare 9 and 5 and swap them, it will then compare 9 and 7 and swap them, and we'll have the sequence 5, 7, 9. If we try to do the comparisons out of order, that is, if we compare the 5 and 7 first and then compare the 9 and 5, we'll wind up with the sequence 5, 9, 7. Therefore, the order in which the "compare-swaps" take place is essential to the correctness of the algorithm.

A variant of bubble sort known as **odd-even transposition sort** has considerably more opportunities for parallelism. The key idea is to "decouple" the compare-swaps. The algorithm consists of a sequence of *phases*, of two different types. During *even* phases, compare-swaps are executed on the pairs

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

and during *odd* phases, compare-swaps are executed on the pairs

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

Here's a small 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.



This example required four phases to sort a four-element list. In general, it may require fewer phases, but the following theorem guarantees that we can sort a list of  $n$  elements in at most  $n$  phases:

**Theorem.** Suppose  $A$  is a list with  $n$  keys, and  $A$  is the input to the odd-even transposition sort algorithm. Then, after  $n$  phases  $A$  will be sorted.

Program 3.15 shows code for a serial odd-even transposition sort function.

```

1 void Odd_even_sort(
2     int a[] /* in/out */,
3     int n /* in */) {
4     int phase, i, temp;
5
6     for (phase = 0; phase < n; phase++)
7         if (phase % 2 == 0) { /* Even phase */
8             for (i = 1; i < n; i += 2)
9                 if (a[i-1] > a[i]) {
10                     temp = a[i];
11                     a[i] = a[i-1];
12                     a[i-1] = temp;
13                 }
14             } else { /* Odd phase */
15                 for (i = 1; i < n-1; i += 2)
16                     if (a[i] > a[i+1]) {
17                         temp = a[i];
18                         a[i] = a[i+1];
19                         a[i+1] = temp;
20                     }
21             }
22 } /* Odd_even_sort */

```

Program 3.15: Serial odd-even transposition sort

### 3.7.2 Parallel odd-even transposition sort

It should be clear that odd-even transposition sort has considerably more opportunities for parallelism than bubble sort, because all of the compare-swaps in a single phase can happen simultaneously. Let's try to exploit this.

There are a number of possible ways to apply Foster's methodology. Here's one:

- **Tasks:** Determine the value of  $a[i]$  at the end of phase  $j$ .
- **Communications:** The task that's determining the value of  $a[i]$  needs to communicate with either the task determining the value of  $a[i-1]$  or  $a[i+1]$ . Also the value of  $a[i]$  at the end of phase  $j$  needs to be available for determining the value of  $a[i]$  at the end of phase  $j+1$ .

This is illustrated in Figure 3.12, where we've labeled the tasks determining the value of  $a[i]$  with  $a[i]$ .

$a[1]$  is compared to  
re swapped, so in the  
argest value in the list  
it element and it will  
st.length decreases,  
in the sorted list.  
because of the inher-  
those that  $a[i-1] = 9$ ,  
d 5 and swap them, it  
equence 5, 7, 9. If we  
5 and 7 first and then  
herefore, the order in  
ness of the algorithm.  
ort has considerably  
'the compare-swaps.  
it types. During even

5, 9, 3, 4.

3, 5, 4, 9.

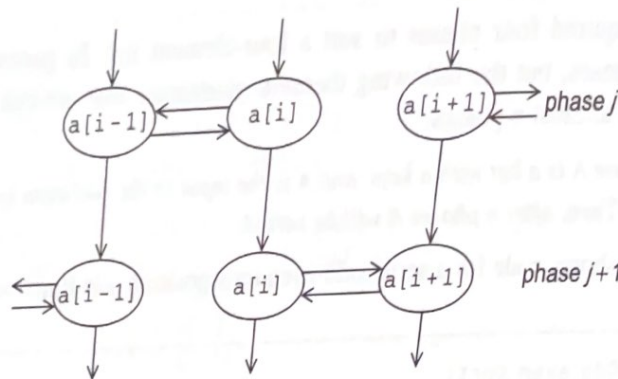


FIGURE 3.12

Communications among tasks in an odd-even sort. Tasks determining  $a[i]$  are labeled with  $a[i]$ .

Now recall that when our sorting algorithm starts and finishes execution, each process is assigned  $n/p$  keys. In this case our aggregation and mapping are at least partially specified by the description of the problem. Let's look at two cases.

When  $n = p$ , Figure 3.12 makes it fairly clear how the algorithm should proceed. Depending on the phase, process  $i$  can send its current value,  $a[i]$ , either to process  $i - 1$  or process  $i + 1$ . At the same time, it should receive the value stored on process  $i - 1$  or process  $i + 1$ , respectively, and then decide which of the two values it should store as  $a[i]$  for the next phase.

However, it's unlikely that we'll actually want to apply the algorithm when  $n = p$ , since we're unlikely to have more than a few hundred or a few thousand processors at our disposal, and sorting a few thousand values is usually a fairly trivial matter for a single processor. Furthermore, even if we do have access to thousands or even millions of processors, the added cost of sending and receiving a message for each compare-exchange will slow the program down so much that it will be useless. Remember that the cost of communication is usually much greater than the cost of "local" computation—for example, a compare-swap.

How should this be modified when each process is storing  $n/p > 1$  elements? (Recall that we're assuming that  $n$  is evenly divisible by  $p$ .) Let's look at an example. Suppose we have  $p = 4$  processes and  $n = 16$  keys assigned, as shown in Table 3.8. In the first place, we can apply a fast serial sorting algorithm to the keys assigned to each process. For example, we can use the C library function `qsort` on each process to sort the local keys. Now if we had one element per process, 0 and 1 would exchange elements, and 2 and 3 would exchange. So let's try this: Let's have 0 and 1 exchange *all* their elements and 2 and 3 exchange all of theirs. Then it would seem natural for 0 to keep the four smaller elements and 1 to keep the larger. Similarly, 2 should keep the smaller and 3 the larger. This gives us the situation shown in the third row of the table. Once again, looking at the one element per process case, in phase 1, processes 1 and 2 exchange their elements and processes 0 and 3 are idle. If process 1 keeps the smaller and 2 the larger elements, we get the distribution shown in the fourth row. Continuing this process for two more phases results in a sorted list. That is, each process' keys are stored in increasing order, and if  $q < r$ ,

However,  
algorithm i  
First, hc  
a process is  
my\_rank-1 a  
calculations a  
if my\_rank =  
But when eith  
We can use the  
is idle:

```
if (phase
    if (my_r
        partn
    else
        partn
```



Table 3.8 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

then the keys assigned to process  $q$  are less than or equal to the keys assigned to process  $r$ .

In fact, our example illustrates the worst-case performance of this algorithm:

**Theorem.** *If parallel odd-even transposition sort is run with  $p$  processes, then after  $p$  phases, the input list will be sorted.*

The parallel algorithm is clear to a human computer:

```
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;
    }
}
```

However, there are some details that we need to clear up before we can convert the algorithm into an MPI program.

First, how do we compute the partner rank? And what is the partner rank when a process is idle? If the phase is even, then odd-ranked partners exchange with  $my\_rank-1$  and even-ranked partners exchange with  $my\_rank+1$ . In odd phases, the calculations are reversed. However, these calculations can return some invalid ranks: if  $my\_rank = 0$  or  $my\_rank = comm\_sz-1$ , the partner rank can be  $-1$  or  $comm\_sz$ . But when either  $partner = -1$  or  $partner = comm\_sz$ , the process should be idle. We can use the rank computed by `Compute_partner` to determine whether a process is idle:

```
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;

```

MPI\_PROC\_NULL is a constant defined by MPI. When it's used as the source or destination rank in a point-to-point communication, no communication will take place and the call to the communication will simply return.

### 3.7.3 Safety in MPI programs

If a process is not idle, we might try to implement the communication with a call to MPI\_Send and a call to MPI\_Recv:

```

MPI_Send(my_keys, n/comm_sz, MPI_INT, partner, 0, comm);
MPI_Recv(temp_keys, n/comm_sz, MPI_INT, partner, 0, comm,
        MPI_STATUS_IGNORE);

```

This, however, might result in the programs' hanging or crashing. Recall that 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. Furthermore, many implementations of MPI set a threshold at which the system switches from buffering to blocking. That is, messages that are relatively small will be buffered by MPI\_Send, but for larger messages, it will 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**, that is, each process is blocked waiting for an event that will never happen.

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. If we use MPI\_Send and MPI\_Recv in this way, our program will be unsafe, and it's likely that for small values of  $n$  the program will run without problems, while for larger values of  $n$ , it's likely that it will hang or crash.

There are a couple of questions that arise here:

1. In general, how can we tell if a program is safe?
2. How can we modify the communication in the parallel odd-even sort program so that it is safe?

To answer the first question, we can use an alternative to MPI\_Send defined by the MPI standard. It's called MPI\_Ssend. The extra "s" stands for *synchronous* and MPI\_Ssend is guaranteed to block until the matching receive starts. So, we can check whether a program is safe by replacing the calls to MPI\_Send with calls to MPI\_Ssend. If the program doesn't hang or crash when it's run with appropriate input and comm\_sz, then the original program was safe. The arguments to MPI\_Ssend are the same as the arguments to MPI\_Send:

```

int MPI_Ssend(
    void*      ms
    int        ms
    MPI_Datatype ms
    int        de
    int        ta
    MPI_Comm   co

```

The answer to the second question is: The most common cause of an error is first sending to each other and then receiving. Another example is a process with rank  $q + 1$ , except

```

MPI_Send(msg, size, MPI_INT, partner, 0, comm);
MPI_Recv(new_msg, size, MPI_INT, partner, 0, comm, MPI_STATUS_IGNORE);

```

In both settings, we need to restructure the communication as follows:

```

if (my_rank % 2 == 0)
    MPI_Send(msg, size, MPI_INT, partner, 0, comm);
    MPI_Recv(new_msg, size, MPI_INT, partner, 0, comm, MPI_STATUS_IGNORE);
} else {
    MPI_Recv(new_msg, size, MPI_INT, partner, 0, comm, MPI_STATUS_IGNORE);
    MPI_Send(msg, size, MPI_INT, partner, 0, comm);
}

```

It's fairly clear that this will work: processes 0 and 2 will first send, then receive from 0 and 2, respectively; processes 1 and 3 will first receive from 0 and 2, then send to 0 and 2.

However, it may not be clear that this works for a sequence of events greater than 1. Suppose, for example, that the dashed arrows show a correct sequence of events.

MPI provides an alternative to MPI\_Send that can call the function MPI\_Ssendrecv:

```

int MPI_Ssendrecv(
    void*      ms
    int        ms
    MPI_Datatype ms
    int        de
    int        ta
    MPI_Comm   co

```



```

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 */);

```

The answer to the second question is that the communication must be restructured. The most common cause of an unsafe program is multiple processes simultaneously first sending to each other and then receiving. Our exchanges with partners is one example. Another example is a "ring pass," in which each process  $q$  sends to the process with rank  $q + 1$ , except that process  $\text{comm\_sz} - 1$  sends to 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);

```

In both settings, we need to restructure the communications so that some of the processes receive before sending. For example, the preceding communications could be restructured as follows:

```

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);
}

```

It's fairly clear that this will work if  $\text{comm\_sz}$  is even. If, say,  $\text{comm\_sz} = 4$ , then processes 0 and 2 will first send to 1 and 3, respectively, while processes 1 and 3 will receive from 0 and 2, respectively. The roles are reversed for the next send-receive pairs: processes 1 and 3 will send to 2 and 0, respectively, while 2 and 0 will receive from 1 and 3.

However, it may not be clear that this scheme is also safe if  $\text{comm\_sz}$  is odd (and greater than 1). Suppose, for example, that  $\text{comm\_sz} = 5$ . Then, Figure 3.13 shows a possible sequence of events. The solid arrows show a completed communication, and the dashed arrows show a communication waiting to complete.

MPI provides an alternative to scheduling the communications ourselves—we can call the function `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    /* in */,
    int        recv_buf_size /* in */,
    MPI_Datatype recv_buf_type /* in */,
    int        src          /* in */,
    int        recv_tag     /* in */,
    MPI_Comm   communicator /* in */);

```

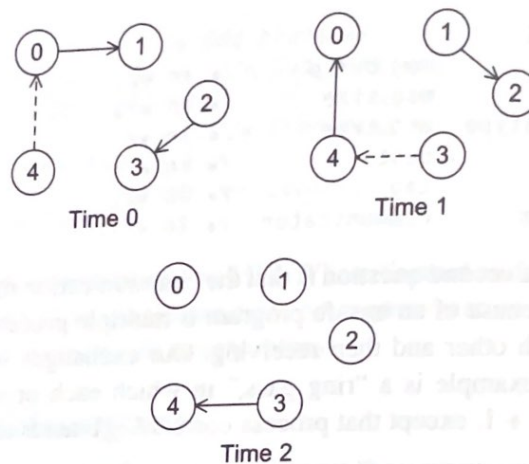


FIGURE 3.13

Safe communication with five processes

```

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  */);
  
```

This function carries out a blocking send and a receive in a single call. The dest and the source can be the same or different. What makes it especially useful is that the MPI implementation schedules the communications so that the program won't hang or crash. The complex code we used earlier—the code that checks whether the process rank is odd or even—can be replaced with a single call to MPI\_Sendrecv. If it happens that the send and the receive buffers should be the same, MPI provides the alternative:

```

int MPI_Sendrecv_replace(
void*      buf_p          /* in/out */,
int        buf_size       /* in  */,
MPI_Datatype buf_type     /* in  */,
int        dest           /* in  */,
int        send_tag       /* in  */,
int        source         /* in  */,
int        recv_tag       /* in  */,
MPI_Comm   communicator   /* in  */,
MPI_Status* status_p      /* in  */);
  
```

### 3.7.4 Final details of parallel odd-even sort

Recall that we had developed the following parallel odd-even transposition sort algorithm:



2

```

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;
    }
}

```

In light of our discussion of safety in MPI, it probably makes sense to implement the send and the receive with a single call to `MPI_Sendrecv`:

```

MPI_Sendrecv(my_keys, n/comm_sz, MPI_INT, partner, 0,
             recv_keys, n/comm_sz, MPI_INT, partner, 0, comm,
             MPI_Status_ignore);

```

It only remains to identify which keys we keep. Suppose for the moment that we want to keep the smaller keys. Then we want to keep the smallest  $n/p$  keys in a collection of  $2n/p$  keys. An obvious approach to doing this is to sort (using a serial sorting algorithm) the list of  $2n/p$  keys and keep the first half of the list. However, sorting is a relatively expensive operation, and we can exploit the fact that we already have two sorted lists of  $n/p$  keys to reduce the cost by *merging* the two lists into a single list. In fact, we can do even better, because we don't need a fully general merge: once we've found the smallest  $n/p$  keys, we can quit. See Program 3.16.

To get the largest  $n/p$  keys, we simply reverse the order of the merge, that is, start with `local_n-1` and work backwards through the arrays. A final improvement avoids copying the arrays and simply swaps pointers (see Exercise 3.28).

Run-times for the version of parallel odd-even sort with the "final improvement" are shown in Table 3.9. Note that if parallel odd-even sort is run on a single processor, it will use whatever serial sorting algorithm we use to sort the local keys, so the times for a single process use serial quicksort, not serial odd-even sort, which would be *much* slower. We'll take a closer look at these times in Exercise 3.27.

**Table 3.9** Run-Times of Parallel Odd-Even Sort (times are in milliseconds)

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

a single call. The destination is especially useful is that the program won't check whether the data is sent to `MPI_Sendrecv`. If the same, MPI provides the

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 */

```

**Program 3.16:** The Merge\_low function in parallel odd-even transposition sort

### 3.8 SUMMARY

MPI, or the Message-Passing Interface, is a library of functions that can be called from C, C++, or Fortran programs. Many systems use `mpicc` to compile MPI programs and `mpiexec` to run them. C MPI programs should include the `mpi.h` header file to get function prototypes and macros defined by MPI.

`MPI_Init` does the setup needed to run MPI. It should be called before other MPI functions are called. When your program doesn't use `argc` and `argv`, `NULL` can be passed for both arguments.

In MPI a **communicator** is a collection of processes that can send messages to each other. After an MPI program is started, MPI always creates a communicator consisting of all the processes. It's called `MPI_COMM_WORLD`.

Many parallel programs use the **single program, multiple data**, or **SPMD**, approach, whereby running a single program obtains the effect of running multiple different programs by including branches on data such as the process rank. When you're done using MPI, you should call `MPI_Finalize`.

To send a message from one MPI process to another, you can use `MPI_Send`. To receive a message, you can use `MPI_Recv`. The arguments to `MPI_Send` describe the contents of the message and its destination. The arguments to `MPI_Recv` describe the storage that the message can be received into, and where the message should be received from. `MPI_Recv` is **blocking**, that is, a call to `MPI_Recv` won't return until the