Parallel Programming COMP5112

Distributed Memory Programming with MPI (3)

Slides adapted from the lecture notes by Peter Pacheco

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.

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 efficiently.
- Similarly, a function that receives data can distribute the items into their correct destinations in memory when they're received.

Derived datatypes

- A derived datatype 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
                                            /* in
                   count
     int
                                                   */,
                                           /* in
                   array_of_blocklengths[]
                   array_of_displacements[]
                                           /* in
                                                   */,
     MPI Aint
                                           /* in */
     MPI Datatype array of types[]
     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)

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.

```
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. The POSIX library function *gettimeofday* returns time in microseconds elapsed from some point in the past.
- Pacheco book example code provides a GET_TIME macro records the number of seconds since some time in the past.

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



Elapsed serial time

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

MPI_Barrier

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



MPI_Barrier

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

Run-times of serial and parallel matrixvector multiplication

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

(Seconds)

Speedup

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

Efficiency

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

Speedups of Parallel Matrix-Vector Multiplication

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

Efficiencies of Parallel Matrix-Vector Multiplication

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

Scalability

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

Scalability

 Programs that can maintain a constant efficiency without increasing the problem size are sometimes said to be strongly scalable.

 Programs that can maintain a constant efficiency if the problem size increases at the same rate as the number of processes are sometimes said to be weakly scalable.

A PARALLEL SORTING ALGORITHM

Sorting

- n keys and p = # 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.

Serial bubble sort

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

Odd-even transposition sort

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

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

Odd phases, compare swaps:

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

Example

Start: 5, 9, 4, 3

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

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

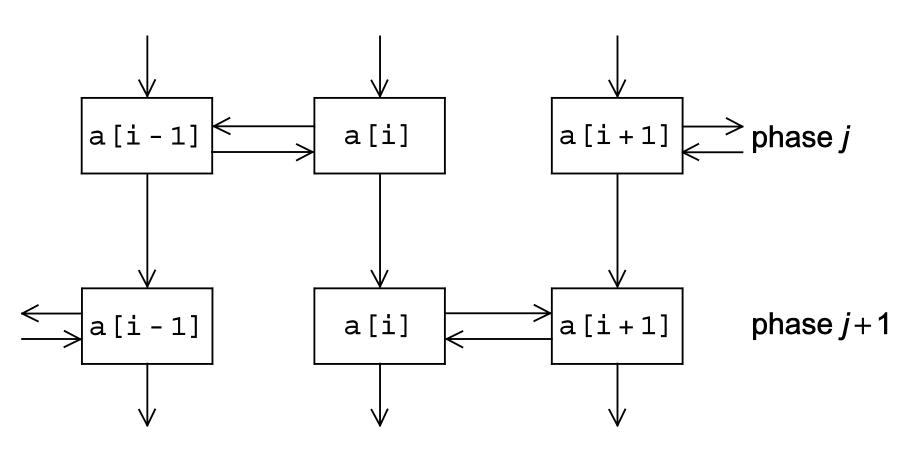
Even phase: compare-swap (5,3) and (9,4) getting the list 3, 5, 4, 9

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

Serial odd-even transposition sort

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

Communications among tasks in oddeven 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;
  else
                            /* Even rank */
     partner = my_rank + 1;
                       /* Odd phase */
else
   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.

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.

Restructuring communication

```
\label{eq:mpi_send} \begin{split} \text{MPI\_Send(msg, size, MPI\_INT, (my\_rank+1) \% comm\_sz, 0, comm);} \\ \text{MPI\_Recv(new\_msg, size, MPI\_INT, (my\_rank+comm\_sz-1) \% comm\_sz,} \\ 0, comm, MPI\_STATUS\_IGNORE. \end{split}
```



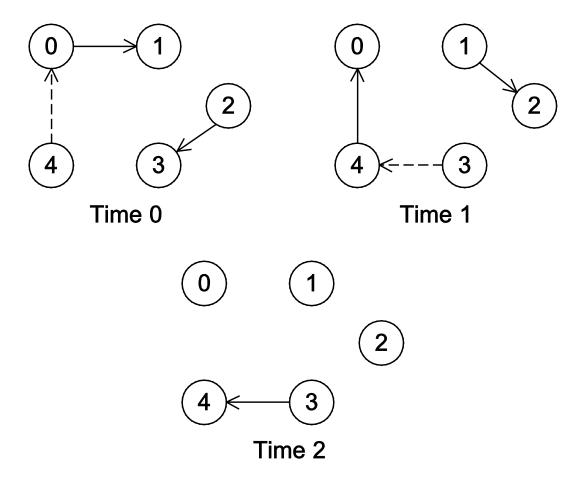
MPI_Sendrecv

- An alternative to scheduling the communications by ourselves.
- Carries out a blocking send and a receive in a single call.
- The destination 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 */,
                                /* in */,
     int
                  dest
     int
                                /* in */,
                  send_tag
     void*
                  recv_buf_p /* out */,
                  recv_buf_size /*in */.
     int
                  recv_buf_type /*in */,
     MPI_Datatype
     int
                                /* in */,
                  source
     int
                                /* in */,
                  recv_tag
                  communicator /*in */,
     MPI Comm
                                /* in */);
     MPI Status*
                  status p
```

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) {
     if (my_keys[m_i] \le 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

	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 singleprogram 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.