

Programming with MPI

Parallel Programming

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Programming with MPI

Review

Introduction

History

Groups and Communicators

Point-To-Point Communication

Collective Communication

Derived Datatypes

Summary

- Which aspect is not part of the POSIX Threads standard?
 1. Thread management
 2. Mutexes
 3. Semaphores
 4. Condition variables
 5. Synchronization

- Which thread-to-task mapping does Linux use?
 1. 1:1 mapping (each thread is mapped to a kernel task)
 2. n:1 mapping (all threads are mapped to one kernel task)
 3. m:n mapping (multiple threads are mapped to multiple kernel tasks)

- What happens if a thread is not joined after termination?
 1. The process can crash
 2. Zombie threads remain
 3. Stack memory can overflow

- When will a thread be canceled by `pthread_cancel`?
 1. Before the next function call
 2. After the next function call
 3. After a timeout of 100 ms
 4. After an I/O operation such as `printf`
 5. When a cancellation point function is called

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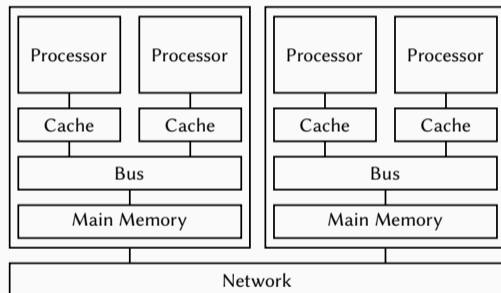
Point-To-Point Communication

Collective Communication

Derived Datatypes

Summary

- Shared memory systems have limited scalability
 - Two to four processors with a few dozen cores
- Complex problems require more nodes
 - Distributed memory can be scaled arbitrarily
- Nodes are connected via a network
 - Determines scalability and performance
- Different network technologies and topologies
 - Major competitors: Ethernet and InfiniBand



- OpenMP is a convenient and high-level programming concept
 - It is limited to shared memory systems
- Parallel applications across multiple nodes require message passing
 - Message Passing Interface (MPI) provides necessary functionality
- MPI supports basic and complex operations
 - Sending, receiving, reduction etc.
 - Process groups and synchronization
 - Point-to-point, collective or one-sided communication
- MPI also offers parallel I/O
 - Concurrent access to shared files

- MPI is a standard by the MPI Forum
 - Over 40 participating organizations
 - First standardized and vendor-independent API
 - MPI is not a library but a specification of one
- There are multiple implementations of the standard
 - MPICH, MVAPICH, OpenMPI, Intel MPI etc.
 - Vendors often provide their own implementations

- MPI implementations are not necessarily binary-compatible
 - They have the same API but different ABIs
 - Compiling an application works with any implementation
- Running compiled application requires original implementation
 - Different implementations might have different constants etc.
 - Way to start processes on different nodes might differ
- Some implementations promise ABI compatibility
 - MPICH ABI Compatibility Initiative for MPICH, Intel MPI, Cray MPT, MVAPICH2, Parastation MPI and RIKEN MPI [MPICH Collaborators, 2023]

- Parallel applications now run as independent processes
 - Processes can only access their own data, no shared memory
 - No risk of overwriting other processes' data accidentally
 - Results have to be communicated between processes
- Application code is typically still contained in one file
 - MPI allows us to write a generic version of the application
 - We can determine our rank and the number of processes

- MPI applications often use SPMD
 - All tasks execute same application but at different points
 - Tasks use different data (domain decomposition)
 - Additional logic to execute only parts of the application
- Decomposition is critical for achievable performance
 - Rows might be faster than columns depending on memory layout
 - Size of sub-domains determines load of each task
- Distribution also determines communication schema
 - Communication might have to be performed at boundaries

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0	1	2	3
4	5	6	7
8	9	10	11
12	13	14	15

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- Application has to be made available on multiple nodes
 - This is normally achieved by using a common file system on all nodes
 - For instance, an NFS file system can be mounted everywhere
- Processes have to be started on participating nodes
 - Many implementations include support for spawning processes via SSH
 - The batch scheduler can also take care of it, requires coordination
- Processes have to locate each other and coordinate
 - Similar to previous point, implementation often takes care of both
 - If the scheduler is involved, it has to pass information to the implementation
 - Process Management Interface (PMI) is typically used to connect components

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- MPI is the current de-facto standard in HPC
 - Previously, Parallel Virtual Machine (PVM) was widely used
- MPI is developed by the MPI Forum, started in 1992
 - MPI-1.0 in 1995: Basic features, communication only
 - MPI-2.0 in 1997: Additional features, including I/O
 - MPI-3.0 in 2012: Better support for one-sided communication
 - MPI-4.0 in 2021: Large-count routines, persistent collectives
- Standard is important for portability across different systems
 - MPI also offers high performance and convenience

- MPI standard defines an API for C and Fortran
 - C++ used to be available but has been deprecated
 - Bindings are also available for Python, Java etc.
- Abstraction to support efficient communication and I/O
 - Functions have to be high-level enough to be able to apply optimizations
- Standard allows thread-safe implementations but does not require them
 - MPI implementations are typically thread-unsafe by default
 - Thread-safety does have a performance impact due to locking etc.

- MPI defines syntax and semantics
 - Syntax determines arguments, semantics how a function behaves
- Example: Function for sending data
 - Standard includes description of behavior and rationale
 - “The send call [...] is blocking: it does not return until the message data and envelope have been safely stored away so that the sender is free to modify the send buffer.”
[Message Passing Interface Forum, 2015]
 - Abstract: `MPI_SEND(buf, count, datatype, dest, tag, comm)`
 - Arguments are annotated as IN/OUT/INOUT and described
 - C: `int MPI_Send(const void* buf, ...)`
 - Return value via normal method
 - Fortran: `MPI_Send(buf, ..., ierror)`
 - Return value via extra argument (`ierror`)

- Non-blocking
 - Call returns before operation has been completed
 - User might not be allowed to reuse specified resources (for example, buffers)
- Blocking
 - User is allowed to reuse resources
- Local
 - Completion of a call depends only on the local process
- Non-local
 - Completion of a call might depend on remote processes
 - Communication might be required to happen before completion
- Collective
 - All processes in a communicator have to be involved in a call

- 1992: “Standards for Message Passing in a Distributed Memory Environment”
 - Working group established and prepares draft for MPI-1
 - Group consists of 175 people from 40 organizations
- 1994: MPI-1.0 is released
 - MPI-1.1 in 1995, MPI-1.2 in 1997 and MPI-1.3 in 2008
 - Point-to-point and collective communication
 - Groups, communicators and topologies
 - Environment checks
 - Profiling interface

- 1998: MPI-2.0 is released
 - MPI-2.1 in 2008 and MPI-2.2 in 2009
 - One-sided communication
 - Dynamic process management
 - Parallel I/O
- 2012: MPI-3.0 is released
 - MPI-3.1 in 2015
 - Improved one-sided communication
 - Non-blocking collectives

- 2021: MPI-4.0 is released
 - Large-count versions of many routines
 - Persistent collectives
 - Partitioned communication

- MPI implementations consist of headers and libraries
 - Main header (`mpi.h`) has to be included
 - Applications have to be linked to MPI libraries
- MPI provides own compilers for convenience
 - `mpicc` for C and `mpifort` for Fortran
 - These are usually compiler wrappers around the underlying compiler
- Compiler wrappers take care of linking etc.
 - Compiler flags can usually be extracted if linking should be done manually

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Summary

- MPI needs to be initialized and finalized
 - Has to be done manually
 - Do as little as possible before and after
- MPI_Init expects application's arguments
 - MPI might parse certain arguments
 - It is possible to pass NULL to ignore

```
1  int main(void) {
2      MPI_Init(NULL, NULL);
3      hello();
4      MPI_Finalize();
5
6      return 0;
7  }
```

- MPI uses communicators
 - Basically a group of processes
- We can determine our rank
 - Same as OpenMP's thread ID
- We can query the communicator's size
 - This is the total amount of processes

```
1 void hello(void) {
2     int rank;
3     int size;
4
5     MPI_Comm_rank(MPI_COMM_WORLD,
6                   &rank);
7     MPI_Comm_size(MPI_COMM_WORLD,
8                   &size);
9
10    printf("Hello from %d/%d.\n",
11           rank, size);
12 }
```

- We can start the application directly
 - It will usually start with one process
- `mpiexec` allows spawning more processes
 - Optional and specified by the standard
 - There is also often `mpirun`
- The `-n` argument is standardized
 - Implementations provide additional ones

```
$ ./hello
Hello from 0/1.

$ mpiexec -n 1 ./hello
Hello from 0/1.

$ mpiexec -n 4 ./hello
Hello from 0/4.
Hello from 3/4.
Hello from 1/4.
Hello from 2/4.
```

- MPI_Init only allows serial processes
 - That is, no threads are allowed to run
- Thread-safety requires locks
 - MPI is tuned for high performance
 - Locking overhead should be avoided
- MPI_Init_thread allows requesting a thread-safety level
 - Implementations may not support all

```
1  int main(void) {
2      int thread_level;
3
4      MPI_Init_thread(NULL, NULL,
5                      MPI_THREAD_MULTIPLE,
6                      &thread_level);
7
8      printf("thread_level=%d\n",
9             thread_level);
10
11     MPI_Finalize();
12     return 0;
13 }
```

- MPI_THREAD_SINGLE
 - Only one thread will run
- MPI_THREAD_FUNNELED
 - Process can be multi-threaded but only the main thread will make MPI calls
- MPI_THREAD_SERIALIZED
 - All threads can make MPI calls but not at the same time
- MPI_THREAD_MULTIPLE
 - Threads can make MPI calls in parallel

```
1  int main(void) {
2      int thread_level;
3
4      MPI_Init_thread(NULL, NULL,
5                     MPI_THREAD_MULTIPLE,
6                     &thread_level);
7
8      printf("thread_level=%d\n",
9             thread_level);
10
11     MPI_Finalize();
12     return 0;
13 }
```


- MPI_THREAD_SINGLE
 - Only one thread will run
- MPI_THREAD_FUNNELED
 - Process can be multi-threaded but only the main thread will make MPI calls
- MPI_THREAD_SERIALIZED
 - All threads can make MPI calls but not at the same time
- MPI_THREAD_MULTIPLE
 - Threads can make MPI calls in parallel

```
$ ./init_thread
thread_level=3

$ mpiexec -n 4 ./init_thread
thread_level=3
thread_level=3
thread_level=3
thread_level=3
```

- `MPI_Get_processor_name`
 - Returns an implementation-defined processor name
 - This typically returns the hostname of the current node
- `MPI_Initialized`
 - Checks whether MPI has been initialized
 - Useful if libraries want to check for MPI support
- `MPI_Wtime`
 - Returns wall-clock time for time measurements
- `MPI_Wtick`
 - Returns resolution of `MPI_Wtime`

- Communicators allow separating different sets of processes
 - Groups contain processes
 - Communicators are based on groups
- All processes are available by default (MPI_COMM_WORLD)
 - Ranks are numbered from 0 to n-1
- Communicators can be used to define independent contexts
 - For instance, MPI-aware library should not interfere with application
- Some operations should only be performed by the local process
 - If they require a communicator, MPI_COMM_SELF can be used

- Chicken and egg problem
 - Creating new communicator requires an existing communicator
 - `MPI_COMM_WORLD` can be used
- Processes can have multiple ranks
 - Rank only valid in a communicator
 - Processes can belong to multiple groups and communicators

```
1 void comm(void) {
2     MPI_Comm new_comm;
3     MPI_Group new_group;
4     MPI_Group world_group;
5
6     MPI_Comm_group(MPI_COMM_WORLD,
7                   &world_group);
8     MPI_Group_incl(world_group,
9                   size, reverse_ranks,
10                  &new_group);
11     MPI_Comm_create(MPI_COMM_WORLD,
12                   new_group, &new_comm);
13
14     print_rank(new_comm);
15 }
```

- Chicken and egg problem
 - Creating new communicator requires an existing communicator
 - `MPI_COMM_WORLD` can be used
- Processes can have multiple ranks
 - Rank only valid in a communicator
 - Processes can belong to multiple groups and communicators

```
1 void print_rank(MPI_Comm comm) {  
2     int new_rank;  
3  
4     MPI_Comm_rank(comm, &new_rank);  
5     printf("rank=%d (world=%d)\n",  
6           new_rank, rank);  
7 }
```

```
$ mpiexec -n 4 ./comm  
rank=3 (world=0)  
rank=2 (world=1)  
rank=1 (world=2)  
rank=0 (world=3)
```

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Summary

- Message order is guaranteed
 - If a process sends two messages, the first one will be received first
 - If a process posts two receives, the first one will get the message
- Rules do not apply when multi-threaded
 - If two threads send one message each, their order is undefined
 - Would require coordinating threads, that is, introduce overhead
- There are no fairness guarantees
 - A message might never be received because of other matching messages

- Point-to-point between two processes
- Sending
 - Buffer: Data to send
 - Count: Number of elements
 - Datatype: Type of elements
 - Destination: Target rank
 - Tag: Distinguish messages
 - Communicator: Process mapping

```
1 void mysend(void) {
2     char str[100];
3     snprintf(str, 100,
4         "Hello from %d\n", rank);
5
6     MPI_Send(str, 100, MPI_CHAR,
7         (rank + 1) % size,
8         0, MPI_COMM_WORLD);
9     MPI_Recv(str, 100, MPI_CHAR,
10        (size + rank - 1) % size,
11        0, MPI_COMM_WORLD,
12        MPI_STATUS_IGNORE);
13
14    printf("%d: %s", rank, str);
15 }
```


- Point-to-point between two processes
- Receiving
 - Buffer: Where to receive data
 - Count: Number of elements
 - Datatype: Type of elements
 - Source: Source rank
 - Tag: Distinguish messages
 - Communicator: Process mapping
 - Status: Query information

```
1 void mysend(void) {
2     char str[100];
3     snprintf(str, 100,
4         "Hello from %d\n", rank);
5
6     MPI_Send(str, 100, MPI_CHAR,
7         (rank + 1) % size,
8         0, MPI_COMM_WORLD);
9     MPI_Recv(str, 100, MPI_CHAR,
10        (size + rank - 1) % size,
11        0, MPI_COMM_WORLD,
12        MPI_STATUS_IGNORE);
13
14    printf("%d: %s", rank, str);
15 }
```

- Point-to-point between two processes
- Ring communication
 - Send to next process
 - Receive from previous process
 - Output order might be mixed

```
$ mpiexec -n 4 ./send
1: Hello from 0
0: Hello from 3
3: Hello from 2
2: Hello from 1
```

- Might not be clear from which process to receive
 - Functions require specifying a source rank and tag
- Wildcards allow matching any source or any tag
 - `MPI_ANY_SOURCE` instead of actual source rank
 - `MPI_ANY_TAG` instead of actual source tag
- We still might be interested to know which rank and tag a message came from
 - Can be queried via `MPI_Status`'s `MPI_SOURCE` and `MPI_TAG` members
- `MPI_Get_count` returns the number of received elements

- What happens if we send 100,000 bytes?
 1. The same as with 100
 2. Application deadlocks
 3. Crash due to stack overflow
 4. MPI warns about too many elements

```
1 void mysend(void) {
2     char str[100];
3     snprintf(str, 100,
4         "Hello from %d\n", rank);
5
6     MPI_Send(str, 100, MPI_CHAR,
7         (rank + 1) % size,
8         0, MPI_COMM_WORLD);
9     MPI_Recv(str, 100, MPI_CHAR,
10        (size + rank - 1) % size,
11        0, MPI_COMM_WORLD,
12        MPI_STATUS_IGNORE);
13
14    printf("%d: %s", rank, str);
15 }
```

- MPI_Send is the default blocking send function
 - Standard allows using a buffer but does not mandate it
 - “The send call [...] uses the standard communication mode. In this mode, **it is up to MPI to decide whether outgoing messages will be buffered.** [...] In such a case, the send call may complete before a matching receive is invoked. On the other hand, [...] MPI may choose not to buffer outgoing messages, for performance reasons. In this case, the send call will not complete until a matching receive has been posted, and the data has been moved to the receiver. [...] The standard mode send is non-local: successful completion of the send operation may depend on the occurrence of a matching receive.” [Message Passing Interface Forum, 2015]
- Buffering is typically only used for small messages
 - Larger messages make the send operation synchronous

- There are a number of different send/receive variants
 - Synchronous send (MPI_Ssend)
 - Blocks until the destination process has started to receive the message
 - Behaves like MPI_Send for large messages
 - Blocking and non-blocking (MPI_Send and MPI_Isend)
 - Blocking behavior specifies when calls return and buffers can be reused
 - Non-blocking allows overlapping communication with computation
 - Buffered (MPI_Bsend)
 - Data is explicitly buffered, buffers have to be provided manually
 - Behaves like MPI_Send for small messages
 - Ready send (MPI_Rsend)
 - Requires matching receive operation to be started already, otherwise undefined
 - Combined blocking send and receive (MPI_Sendrecv)
 - Avoids deadlocks due to blocking sends waiting for receives to be posted

- Non-blocking send does not deadlock
 - I stands for immediate
- MPI_Wait blocks until completion
 - Functions to wait for multiple requests (all, any or some)
 - It is an error not to wait or access the buffer before the send has finished
- Alternatively, MPI_Test or MPI_Probe

```
1 void mysend(char* str, char* buf) {
2     MPI_Request req;
3
4     MPI_Isend(str, 100000, MPI_CHAR,
5              (rank + 1) % size,
6              0, MPI_COMM_WORLD, &req);
7     MPI_Recv(buf, 100000, MPI_CHAR,
8              (size + rank - 1) % size,
9              0, MPI_COMM_WORLD,
10             MPI_STATUS_IGNORE);
11     MPI_Wait(&req,
12             MPI_STATUS_IGNORE);
13
14     printf("%d: %s", rank, buf);
15 }
```

- Non-blocking send does not deadlock
 - I stands for immediate
- `MPI_Wait` blocks until completion
 - Functions to wait for multiple requests (all, any or some)
 - It is an error not to wait or access the buffer before the send has finished
- Alternatively, `MPI_Test` or `MPI_Probe`

```
$ mpiexec -n 4 ./isend
2: Hello from 1.
0: Hello from 3.
1: Hello from 0.
3: Hello from 2.
```


- Combined blocking send and receive
 - Still blocking but avoids deadlock
- Abstraction to achieve typical use case
 - For example, send to and receive from neighboring processes
 - Implementation can handle this specific use case efficiently and correctly

```
1 void mysend(void) {
2     char str[100000];
3     char buf[100000];
4     snprintf(str, 100000,
5             "Hello from %d.\n", rank);
6
7     MPI_Sendrecv(str, 100000,
8                 MPI_CHAR, (rank + 1) % size,
9                 0, buf, 100000, MPI_CHAR,
10                (size + rank - 1) % size,
11                0, MPI_COMM_WORLD,
12                MPI_STATUS_IGNORE);
13
14     printf("%d: %s", rank, buf);
15 }
```

- Combined blocking send and receive
 - Still blocking but avoids deadlock
- Abstraction to achieve typical use case
 - For example, send to and receive from neighboring processes
 - Implementation can handle this specific use case efficiently and correctly

```
$ mpiexec -n 4 ./sendrecv
0: Hello from 3.
3: Hello from 2.
1: Hello from 0.
2: Hello from 1.
```

	Send	Receive	Testing
Blocking	MPI_Send MPI_Ssend MPI_Rsend	MPI_Recv	MPI_Probe MPI_Wait
	MPI_Sendrecv		
Non-blocking	MPI_Isend MPI_Issend	MPI_Irecv	MPI_Iprobe MPI_Test

- Most functions are available as blocking and non-blocking versions
 - There are also non-blocking synchronous functions
- Blocking is easier to use, non-blocking is more efficient

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- Point-to-point communication happens between two ranks
 - Collective communication happens between all ranks
- Which ranks are involved depends on communicator
 - By default, we only have `MPI_COMM_WORLD` and `MPI_COMM_SELF`
- MPI contains a wide range of collective communication functions
 - Broadcast
 - Barrier
 - Distributing or collecting data
- One collective call is often more efficient than many point-to-point calls
 - InfiniBand hardware typically has support for efficient collectives

- 1:1 communication
 - Traditional point-to-point communication such as send and receive
- 1:n communication
 - Collective communication such as broadcast
- n:1 communication
 - Collective communication such as reduction
- n:n communication
 - Collective communication such as reduction to all

P0	A	B	C
P1			
P2			

Broadcast



P0	A	B	C
P1	A	B	C
P2	A	B	C

P0	A	B	C
P1			
P2			

P0	A	B	C
P1			
P2			

Broadcast



P0	A	B	C
P1	A	B	C
P2	A	B	C

Scatter



P0	A	B	C
P1	B		
P2	C		

P0	A	B	C
P1			
P2			

Broadcast

→

P0	A	B	C
P1	A	B	C
P2	A	B	C

P0	A	B	C
P1			
P2			

Scatter

→

P0	A	B	C
P1	B		
P2	C		

P0	A		
P1	B		
P2	C		

Gather

→

P0	A	B	C
P1	B		
P2	C		

- Reducing
 - Send buffer: Data to reduce
 - Receive buffer: Root needs separate buffer
 - Count: Number of elements
 - Datatype: Type of elements
 - Operation: Reduction to perform
 - Root: Rank to reduce at
 - Communicator: Process mapping
- Reduction operations known from OpenMP
 - Apply a given function to multiple buffers, reducing it to one buffer
- Ordering is arbitrary, might influence result

```
1 void reduce(void) {  
2     int buf = 42;  
3  
4     MPI_Reduce(&rank, &buf, 1,  
5               MPI_INT, MPI_MAX,  
6               0, MPI_COMM_WORLD);  
7  
8     printf("%d: %d\n", rank, buf);  
9 }
```

- Reducing
 - Send buffer: Data to reduce
 - Receive buffer: Root needs separate buffer
 - Count: Number of elements
 - Datatype: Type of elements
 - Operation: Reduction to perform
 - Root: Rank to reduce at
 - Communicator: Process mapping
- Reduction operations known from OpenMP
 - Apply a given function to multiple buffers, reducing it to one buffer
- Ordering is arbitrary, might influence result

```
$ mpiexec -n 4 ./reduce
0: 3
1: 42
2: 42
3: 42
```

- Reducing to all
 - Send buffer: Data to reduce
 - Receive buffer: Needs separate buffer
 - Count: Number of elements
 - Datatype: Type of elements
 - Operation: Reduction to perform
 - Communicator: Process mapping
- No root rank specified anymore
 - Reduced buffer is available for all ranks

```
1 void reduce(void) {
2     int buf = 42;
3
4     MPI_Allreduce(&rank, &buf, 1,
5                 MPI_INT, MPI_MAX,
6                 MPI_COMM_WORLD);
7
8     printf("%d: %d\n", rank, buf);
9 }
```

- Reducing to all
 - Send buffer: Data to reduce
 - Receive buffer: Needs separate buffer
 - Count: Number of elements
 - Datatype: Type of elements
 - Operation: Reduction to perform
 - Communicator: Process mapping
- No root rank specified anymore
 - Reduced buffer is available for all ranks

```
$ mpiexec -n 4 ./allreduce
0: 3
1: 3
2: 3
3: 3
```

- Why not use MPI_Reduce followed by MPI_Broadcast?
 1. More optimization potential
 2. Two collectives could deadlock
 3. Data could be broadcasted before reduction is finished

```
1 void reduce(void) {
2     int buf = 42;
3
4     MPI_Allreduce(&rank, &buf, 1,
5                 MPI_INT, MPI_MAX,
6                 MPI_COMM_WORLD);
7
8     printf("%d: %d\n", rank, buf);
9 }
```

- Barrier
 - Communicator: Process mapping
- Waits for all processes
 - Can cause significant overhead
 - Often not necessary due to implicit synchronization via messages

```
1 void barrier(void) {  
2     printf("%d: before barrier\n",  
3         rank);  
4  
5     MPI_Barrier(MPI_COMM_WORLD);  
6  
7     printf("%d: after barrier\n",  
8         rank);  
9 }
```

- Barrier
 - Communicator: Process mapping
- Waits for all processes
 - Can cause significant overhead
 - Often not necessary due to implicit synchronization via messages
- Does not work for everything
 - Terminal output might be buffered
 - Output has to be collected from nodes

```
$ mpiexec -n 4 ./barrier
1: before barrier
0: before barrier
2: before barrier
2: after barrier
3: before barrier
3: after barrier
0: after barrier
1: after barrier
```


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Summary

- MPI supports most basic data types out of the box
 - char, int, long, float, double etc.
- Applications often use their own data types
 - For instance, structures containing multiple values
- MPI allows handling these data types directly
 - Developers have to replicate the data types for MPI
 - MPI might be able to handle them more efficiently then
- Data types can then be specified like normal ones
 - Every function that accepts a data type also accepts derived ones

- Example: Diagonal of a 3×3 matrix
 - For instance, within a function doing parallel matrix calculations

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- MPI supports a vector data type
 - Count: Number of blocks
 - Block length: Elements per block
 - Stride: Elements between blocks
 - Old type: Old data type
 - New type: New data type

```
1  int MPI_Type_vector (
2      int count,
3      int blocklength,
4      int stride,
5      MPI_Datatype oldtype,
6      MPI_Datatype* newtype)
```

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```

```
1  MPI_Type_vector(3, 1, 4,  
2      MPI_DOUBLE, &newtype);  
3  MPI_Type_commit(&newtype);  
4  MPI_Send(matrix, 1, newtype,  
5      rank, 0, MPI_COMM_WORLD);
```

- Matrix is stored in row- or column-major order
 - 3×3 matrix has three diagonal elements
 - Each diagonal element is a double value
 - Diagonal elements are four values apart
- Can be generalized for arbitrary dimensions
 - Sender and receiver have to agree on data type
- There are many more data type constructors
 - Interactive tools can help create own derived data types [RookieHPC, 2022]

```
1 MPI_Type_vector(3, 1, 4,  
2 MPI_DOUBLE, &newtype);
```

1	2	3
4	5	6
7	8	9

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Summary

- MPI is a standard for parallel programming on distributed memory systems
 - It supports communication, synchronization, I/O and much more
- Groups of processes can be assigned to communicators
 - Allows separating different parts of an application or library
- Point-to-point communication allows sending messages between two processes
 - There are various versions of basic send and receive functions
- Collective communication involves all processes in a communicator
 - This includes actual communication as well as synchronization functionality
- Derived data types allow MPI to handle application-specific data types directly
 - Allows the MPI implementation to make access more convenient and efficient

References

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