

1 MPI\_WTIME()

2  
3 double MPI\_Wtime(void)

4 DOUBLE PRECISION MPI\_WTIME()

5  
6 {double MPI::Wtime() (*binding deprecated, see Section 15.2*) }

7  
8 MPI\_WTIME returns a floating-point number of seconds, representing elapsed wall-clock time since some time in the past.

9  
10 The “time in the past” is guaranteed not to change during the life of the process. The user is responsible for converting large numbers of seconds to other units if they are preferred.

11  
12 This function is portable (it returns seconds, not “ticks”), it allows high-resolution, and carries no unnecessary baggage. One would use it like this:

13  
14  
15 {  
16     double starttime, endtime;  
17     starttime = MPI\_Wtime();  
18     .... stuff to be timed ...  
19     endtime = MPI\_Wtime();  
20     printf("That took %f seconds\n",endtime-starttime);  
21 }  
22

23  
24 The times returned are local to the node that called them. There is no requirement that different nodes return “the same time.” (But see also the discussion of MPI\_WTIME\_IS\_GLOBAL).

25  
26  
27  
28 MPI\_WTICK()

29 double MPI\_Wtick(void)

30 DOUBLE PRECISION MPI\_WTICK()

31  
32 {double MPI::Wtick() (*binding deprecated, see Section 15.2*) }

33  
34 MPI\_WTICK returns the resolution of MPI\_WTIME in seconds. That is, it returns, as a double precision value, the number of seconds between successive clock ticks. For example, if the clock is implemented by the hardware as a counter that is incremented every millisecond, the value returned by MPI\_WTICK should be  $10^{-3}$ .

## 35 36 37 38 39 8.7 Startup

40  
41 One goal of MPI is to achieve *source code portability*. By this we mean that a program written using MPI and complying with the relevant language standards is portable as written, and must not require any source code changes when moved from one system to another. This explicitly does *not* say anything about how an MPI program is started or launched from the command line, nor what the user must do to set up the environment in which an MPI program will run. However, an implementation may require some setup to be performed

before other MPI routines may be called. To provide for this, MPI includes an initialization routine `MPI_INIT`.

`MPI_INIT()`

```
int MPI_Init(int *argc, char ***argv)
```

```
MPI_INIT(IERROR)
```

```
INTEGER IERROR
```

```
{void MPI::Init(int& argc, char**& argv) (binding deprecated, see Section 15.2) }
```

```
{void MPI::Init() (binding deprecated, see Section 15.2) }
```

[ All MPI programs must contain exactly one call to an MPI initialization routine: `MPI_INIT` or `MPI_INIT_THREAD`. ] Each MPI process must call an MPI initialization routine, `MPI_INIT` or `MPI_INIT_THREAD`, exactly once. Subsequent calls to any initialization routine are erroneous. The only MPI functions that may be invoked before the MPI initialization routines are called are `MPI_GET_VERSION`, `MPI_INITIALIZED`, and `MPI_FINALIZED`.

The version for ISO C accepts the `argc` and `argv` that are provided by the arguments to `main` or `NULL`:

```
int main(int argc, char **argv)
```

```
{
```

```
    MPI_Init(&argc, &argv);
```

```
    /* parse arguments */
```

```
    /* main program */
```

```
    MPI_Finalize();    /* see below */
```

```
}
```

The Fortran version takes only `IERROR`.

Conforming implementations of MPI are required to allow applications to pass `NULL` for both the `argc` e `argv` arguments of `main` in C. [ and C++. In C++, there is an alternative binding for `MPI::Init` that does not have these arguments at all.

*Rationale.* In some applications, libraries may be making the call to `MPI_Init`, and may not have access to `argc` and `argv` from `main`. It is anticipated that applications requiring special information about the environment or information supplied by `mpixec` can get that information from environment variables. (*End of rationale.*)

]

After MPI is initialized, the application can access information about the execution environment by querying the predefined info object `MPI_INFO_ENV`. The following keys are predefined for this object, corresponding to the arguments of `MPI_COMM_SPAWN` or of `mpixec`:

`command` name of program executed

`argv` (space separated) arguments to command  
`maxprocs` Maximum number of MPI processes to start.  
`soft` Allowed values for number of processors  
`host` Hostname.  
`arch` Architecture name.  
`wdir` Working directory of the MPI process  
`file` Value is the name of a file in which additional information is specified.  
`thread_level` Requested level of thread support (if requested before the program started execution)

The info object `MPI_INFO_ENV` need not contain a (key,value) pair for each of these predefined keys; the set of (key,value) pairs provided is implementation-dependent. Implementations may provide additional, implementation specific, (key,value) pairs.

In case where the MPI processes were started with `MPI_COMM_SPAWN_MULTIPLE` or, equivalently, with a startup mechanism that supports multiple process specifications, then the values stored in the info object `MPI_INFO_KEY` at a process are those values that affect the local MPI process.

**Example 8.3** If MPI is started with a call to

```
mpirun -n 5 -arch sun ocean : -n 10 -arch rs6000 atmos
```

Then the first 5 processes will have in their `MPI_INFO_ENV` object the pairs (command, ocean), (maxprocs,5), and (arch, sun). The next 10 processes will have in `MPI_INFO_KEY` (command, atmos), (maxprocs,10), and (arch, rs6000)

*Advice to users.* The values passed in `MPI_INFO_KEY` are the values of the arguments passed to the mechanism that started the MPI execution – not the actual value provided. Thus, the value associated with `maxprocs` is the number of MPI processes requested; it can be larger than the actual number of processes obtained, if the `soft` option was used. (*End of advice to users.*)

*Advice to implementors.* Good quality implementations will provide a (key,value) pair for each parameter that can be passed to the command that starts an MPI program. (*End of advice to implementors.*)

`MPI_FINALIZE()`

```
int MPI_Finalize(void)
```

```
MPI_FINALIZE(IERROR)
```

```
INTEGER IERROR
```

```
{void MPI::Finalize() (binding deprecated, see Section 15.2) }
```

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This routine cleans up all MPI state. [ Each process must call `MPI_FINALIZE` before it exits. Unless there has been a call to `MPI_ABORT`, before each process exits process must ensure that all pending nonblocking communications are (locally) complete before calling `MPI_FINALIZE`. Further, at the instant at which the last process calls `MPI_FINALIZE`, all pending sends must be matched by a receive, and all pending receives must be matched by a send.

For example, the following program is correct

] Each MPI process that terminates normally (i.e., not due to a call to `MPI_ABORT` or an unrecoverable error) must call `MPI_FINALIZE` before it exits.

Before MPI is finalized at an MPI process, the process must locally complete all MPI calls. When the last process calls `MPI_FINALIZE`, all non-local MPI calls at each process must be matched by MPI calls at the other processes that are needed to complete the relevant operation: For example, for each send, the matching receive has occurred, and for each receive, a matching send has occurred; each collective operation has been invoked at all involved processes, etc.

The call to `MPI_FINALIZE` does not free objects created by MPI commands – i.e., objects that the user can free using MPI calls.

`MPI_FINALIZE` is collective over all connected processes. If no processes were spawned, accepted or connected then this means over `MPI_COMM_WORLD`; otherwise it is collective over the union of all processes that have been and continue to be connected, as explained in Section 10.5.4 on page 362.

The following examples illustrates these rules

**Example 8.4** The following code is correct

Process 0	Process 1
-----	-----
<code>MPI_Init();</code>	<code>MPI_Init();</code>
<code>MPI_Send(dest=1);</code>	<code>MPI_Recv(src=0);</code>
<code>MPI_Finalize();</code>	<code>MPI_Finalize();</code>

**Example 8.5** Without a matching receive, the program is erroneous

Process 0	Process 1
-----	-----
<code>MPI_Init();</code>	<code>MPI_Init();</code>
<code>MPI_Send (dest=1);</code>	
<code>MPI_Finalize();</code>	<code>MPI_Finalize();</code>

[ deleted in April Since `MPI_FINALIZE` is a collective call, a correct MPI program will naturally ensure that all participants in pending collective operations have made the call before calling `MPI_FINALIZE`.

A successful return from a blocking communication operation or from `MPI_WAIT` or `MPI_TEST` tells the user that the buffer can be reused and means that the communication is completed by the user, but does not guarantee that the local process has no more work to do. A successful return from `MPI_REQUEST_FREE` with a request handle generated by an `MPI_ISEND` nullifies the handle but provides no assurance of operation completion. The `MPI_ISEND` is complete only when it is known by some means that a matching receive has

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completed. `MPI_FINALIZE` guarantees that all local actions required by communications the user has completed will, in fact, occur before it returns.

`MPI_FINALIZE` guarantees nothing about pending communications that have not been completed (completion is assured only by `MPI_WAIT`, `MPI_TEST`, or `MPI_REQUEST_FREE` combined with some other verification of completion). ]

**Example 8.6** This program is correct HEADER SKIP ENDHEADER

```
rank 0                                rank 1
=====
...
MPI_Isend();                          MPI_Recv();
MPI_Request_free();                   MPI_Barrier();
MPI_Barrier();                       MPI_Finalize();
MPI_Finalize();                      exit();
exit();
```

**Example 8.7** This program is erroneous and its behavior is undefined: HEADER SKIP ENDHEADER

```
rank 0                                rank 1
=====
...
MPI_Isend();                          MPI_Recv();
MPI_Request_free();                   MPI_Finalize();
MPI_Finalize();                      exit();
exit();
```

**Example 8.8** This program is correct: The send operation on process 0 is locally complete when `MPI_Finalize` is called: the local buffer can be reused and no further MPI calls are required on the sender side.

Process 0	Process 1
-----	-----
<code>MPI_Init();</code>	<code>MPI_Init();</code>
<code>MPI_Isend();</code>	<code>MPI_Recv();</code>
<code>MPI_Request_free();</code>	<code>MPI_Barrier();</code>
<code>MPI_Barrier();</code>	<code>MPI_Finalize();</code>
<code>MPI_Finalize();</code>	

**Example 8.9** This program is erroneous: The send operation on process 0 is not locally complete when `MPI_Finalize` is called

```

Process 0                                Proces 1
-----                                -----
MPI_Init();                             MPI_Init();
MPI_Isend();                             MPI_Recv();
MPI_Request_free();                     MPI_Finalize();
MPI_Finalize();                         exit();
exit();

```

[ If no `MPI_BUFFER_DETACH` occurs between an `MPI_BSEND` (or other buffered send) and `MPI_FINALIZE`, the `MPI_FINALIZE` implicitly supplies the `MPI_BUFFER_DETACH`.

**Example 8.10** This program is correct, and after the `MPI_Finalize`, it is as if the buffer had been detached. HEADER SKIP ENDHEADER

```

rank 0                                rank 1
=====
...
buffer = malloc(1000000);              MPI_Recv();
MPI_Buffer_attach();                   MPI_Finalize();
MPI_Bsend();                           exit();
MPI_Finalize();
free(buffer);
exit();

]

```

**Example 8.11** This program is correct. The attached buffer is a resource allocated by the user, not by MPI; it is available to the user after MPI is finalized.

```

Process 0                                Process 1
-----                                -----
MPI_Init(0;                             MPI_Init();
buffer = malloc(1000000);               MPI_Recv();
MPI_Buffer_attach();                   MPI_Finalize();
MPI_Bsend();                           exit();
MPI_Finalize();
free(buffer);
exit();

[

```

**Example 8.12** In this example, `MPI_lprobe()` must return a `FALSE` flag. `MPI_Test_cancelled()` must return a `TRUE` flag, independent of the relative order of execution of `MPI_Cancel()` in process 0 and `MPI_Finalize()` in process 1.

The `MPI_lprobe()` call is there to make sure the implementation knows that the “tag1” message exists at the destination, without being able to claim that the user knows about it.

HEADER SKIP ENDHEADER

```

1
2 rank 0 rank 1
3 =====
4 MPI_Init(); MPI_Init();
5 MPI_Isend(tag1);
6 MPI_Barrier(); MPI_Barrier();
7 MPI_Barrier(); MPI_Iprobe(tag2);
8 MPI_Barrier(); MPI_Barrier();
9 MPI_Finalize();
10 exit();
11 MPI_Cancel();
12 MPI_Wait();
13 MPI_Test_cancelled();
14 MPI_Finalize();
15 exit();
16

```

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**Example 8.13** This program is correct. The cancel operation must succeed, since the send cannot complete normally.

```

23 Process 0 Process 1
24 -----
25 MPI_Issend(); MPI_Finalize();
26 MPI_Cancel();
27 MPI_Wait();
28 MPI_Finalize();
29

```

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*Advice to implementors.* An implementation may need to delay the return from MPI\_FINALIZE until all potential future message cancellations have been processed. One possible solution is to place a barrier inside MPI\_FINALIZE (*End of advice to implementors.*)

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*Advice to implementors.*

Even though a process has completed all the communications it initiated, such communication may not yet be completed from the viewpoint of the underlying MPI system. E.g., a blocking send may have returned, even though the data is still buffered at the sender. The MPI implementation must ensure that a process has completed any involvement in MPI communication before MPI\_FINALIZE returns. Thus, if a process exits after the call to MPI\_FINALIZE, this will not cause an ongoing communication to fail.

The MPI implementation should also complete freeing all objects marked for deletion by MPI calls that freed them.

An implementation may need to delay the return from `MPI_FINALIZE` on a process even if all communications related to MPI calls by that process have completed; the process may still receive cancel requests for messages it has completed receiving. One possible solution is to place a barrier inside `MPI_FINALIZE`.

*(End of advice to implementors.)*

*Advice to users.* If a process continues execution after the call to `MPI_FINALIZE` then it is recommended that the user explicitly free all the objects allocated by MPI calls before the call to `MPI_FINALIZE`. *(End of advice to users.)*

Once `MPI_FINALIZE` returns, no MPI routine (not even `MPI_INIT`) may be called, except for `MPI_GET_VERSION`, `MPI_INITIALIZED`, and `MPI_FINALIZED`.

[ Each process must complete any pending communication it initiated before it calls `MPI_FINALIZE`. If the call returns, each process may continue local computations, or exit, without participating in further MPI communication with other processes. ]

[ `MPI_FINALIZE` is collective over all connected processes. If no processes were spawned, accepted or connected then this means over `MPI_COMM_WORLD`; otherwise it is collective over the union of all processes that have been and continue to be connected, as explained in Section 10.5.4 on page 362. ]

*Advice to implementors.* Even though a process has completed all the communication it initiated, such communication may not yet be completed from the viewpoint of the underlying MPI system. E.g., a blocking send may have completed, even though the data is still buffered at the sender. The MPI implementation must ensure that a process has completed any involvement in MPI communication before `MPI_FINALIZE` returns. Thus, if a process exits after the call to `MPI_FINALIZE`, this will not cause an ongoing communication to fail. *(End of advice to implementors.)*

Although it is not required that all processes return from `MPI_FINALIZE`, it is required that at least process 0 in `MPI_COMM_WORLD` return, so that users can know that the MPI portion of the computation is over. In addition, in a POSIX environment, they may desire to supply an exit code for each process that returns from `MPI_FINALIZE`.

**Example 8.14** The following illustrates the use of requiring that at least one process return and that it be known that process 0 is one of the processes that return. One wants code like the following to work no matter how many processes return.

```
...
MPI_Comm_rank(MPI_COMM_WORLD, &myrank);
...
MPI_Finalize();
if (myrank == 0) {
    resultfile = fopen("outfile","w");
    dump_results(resultfile);
    fclose(resultfile);
}
exit(0);
```



The level(s) of thread support that can be provided by `MPI_INIT_THREAD` will depend on the implementation, and may depend on information provided by the user before the program started to execute (e.g., with arguments to `mpiexec`). If possible, the call will return `provided = required`. Failing this, the call will return the least supported level such that `provided > required` (thus providing a stronger level of support than required by the user). Finally, if the user requirement cannot be satisfied, then the call will return in `provided` the highest supported level.

A **thread compliant** MPI implementation will be able to return `provided = MPI_THREAD_MULTIPLE`. Such an implementation may always return `provided = MPI_THREAD_MULTIPLE`, irrespective of the value of `required`. [ At the other extreme, an MPI library that is not thread compliant may always return `provided = MPI_THREAD_SINGLE`, irrespective of the value of `required`. ]

An MPI library that is not thread compliant must always return `provided=MPI_THREAD_SINGLE`, even if `MPI_INIT_THREAD` is called on a multithreaded process. The library should also return correct values for the MPI calls that can be executed before initialization, even if multiple threads have been spawned.

*Rationale.* Such code is erroneous, but the error cannot be detected until `MPI_INIT_THREAD` is called. The requirements in the previous paragraph ensure that the error can be properly detected. (*End of rationale.*)

A call to `MPI_INIT` has the same effect as a call to `MPI_INIT_THREAD` with a `required = MPI_THREAD_SINGLE`.

Vendors may provide (implementation dependent) means to specify the level(s) of thread support available when the MPI program is started, e.g., with arguments to `mpiexec`. This will affect the outcome of calls to `MPI_INIT` and `MPI_INIT_THREAD`. Suppose, for example, that an MPI program has been started so that only `MPI_THREAD_MULTIPLE` is available. Then `MPI_INIT_THREAD` will return `provided = MPI_THREAD_MULTIPLE`, irrespective of the value of `required`; a call to `MPI_INIT` will also initialize the MPI thread support level to `MPI_THREAD_MULTIPLE`. Suppose, on the other hand, that an MPI program has been started so that all four levels of thread support are available. Then, a call to `MPI_INIT_THREAD` will return `provided = required`; on the other hand, a call to `MPI_INIT` will initialize the MPI thread support level to `MPI_THREAD_SINGLE`.

*Rationale.* Various optimizations are possible when MPI code is executed single-threaded, or is executed on multiple threads, but not concurrently: mutual exclusion code may be omitted. Furthermore, if only one thread executes, then the MPI library can use library functions that are not thread safe, without risking conflicts with user threads. Also, the model of one communication thread, multiple computation threads fits many applications well, e.g., if the process code is a sequential Fortran/C/C++ program with MPI calls that has been parallelized by a compiler for execution on an SMP node, in a cluster of SMPs, then the process computation is multi-threaded, but MPI calls will likely execute on a single thread.

The design accommodates a static specification of the thread support level, for environments that require static binding of libraries, and for compatibility for current multi-threaded MPI codes. (*End of rationale.*)

*Advice to implementors.* If `provided` is not `MPI_THREAD_SINGLE` then the MPI library should not invoke C/ C++/Fortran library calls that are not thread safe, e.g., in an

environment where `malloc` is not thread safe, then `malloc` should not be used by the MPI library.

Some implementors may want to use different MPI libraries for different levels of thread support. They can do so using dynamic linking and selecting which library will be linked when `MPI_INIT_THREAD` is invoked. If this is not possible, then optimizations for lower levels of thread support will occur only when the level of thread support required is specified at link time. (*End of advice to implementors.*)

The following function can be used to query the current level of thread support.

`MPI_QUERY_THREAD(provided)`

OUT      provided      provided level of thread support (integer)

`int MPI_Query_thread(int *provided)`

`MPI_QUERY_THREAD(PROVIDED, IERROR)`

INTEGER PROVIDED, IERROR

{`int MPI::Query_thread()` (*binding deprecated, see Section 15.2*) }

The call returns in `provided` the current level of thread support. This support, which will be the value returned in `provided` by `MPI_INIT_THREAD`, if MPI was initialized by a call to `MPI_INIT_THREAD`.

`MPI_IS_THREAD_MAIN(flag)`

OUT      flag      true if calling thread is main thread, false otherwise (logical)

`int MPI_Is_thread_main(int *flag)`

`MPI_IS_THREAD_MAIN(FLAG, IERROR)`

LOGICAL FLAG

INTEGER IERROR

{`bool MPI::Is_thread_main()` (*binding deprecated, see Section 15.2*) }

This function can be called by a thread to find out whether determine if it is the main thread (the thread that called `MPI_INIT` or `MPI_INIT_THREAD`).

All routines listed in this section must be supported by all MPI implementations.

*Rationale.* MPI libraries are required to provide these calls even if they do not support threads, so that portable code that contains invocations to these functions [be able to] can link correctly. `MPI_INIT` continues to be supported so as to provide compatibility with current MPI codes. (*End of rationale.*)

*Advice to users.* It is possible to spawn threads before MPI is initialized, but no MPI call other than [ `MPI_INITIALIZED` ] `MPI_GET_VERSION`, `MPI_INITIALIZED`, or `MPI_FINALIZED` should be executed by these threads, until `MPI_INIT_THREAD` is