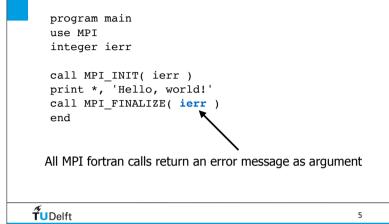


# A Minimal MPI Program (Fortran 90)



# Exiting the MPI Environment

```
• MPI_FINALIZE ( )
```

Cleans up all MPI state. Once this routine has been called, no MPI routine ( even **MPI INIT** ) may be called

### Syntax

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```
int MPI_Finalize ( );
```

```
MPI_FINALIZE ( IERROR )
INTEGER IERROR
```

**MUST** call MPI\_FINALIZE when you exit from an MPI program.

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# Starting the MPI Environment

• MPI\_INIT ( )

Initializes MPI environment. This function must be called and must be the first MPI function called in a program (exception: **MPI\_INITIALIZED**)

Syntax
 int MPI\_Init ( int \*argc, char \*\*\*argv )

MPI\_INIT ( IERROR ) INTEGER IERROR

NOTE: Both C and Fortran return error codes for all calls.

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# C and Fortran Language Considerations

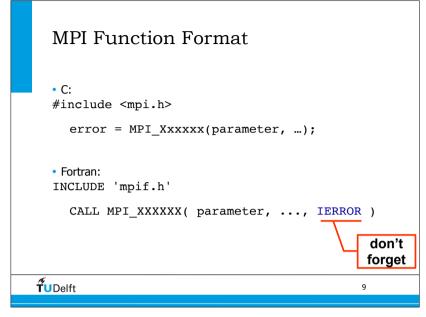
- Bindings
  - C
    - All MPI names have an MPI\_ prefix
    - Defined constants are in all capital letters
    - Defined types and functions have one capital letter after the prefix; the remaining letters are lowercase

6

8

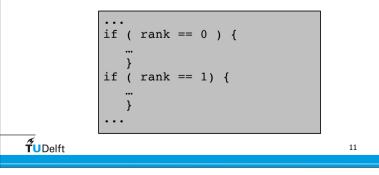
- Fortran
  - All MPI names have an **MPI\_** prefix
  - No capitalization rules apply
  - last argument is an returned error value

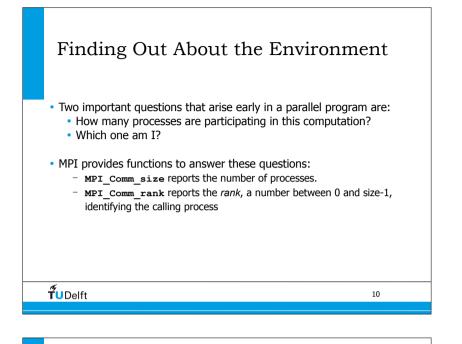
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## MPI Rank

- MPI runtime assigns each process a rank, which can be used as an ID of the processes
  - ranks start from 0 and extent to N-1
- Processes can perform different tasks and handle different data based on their rank

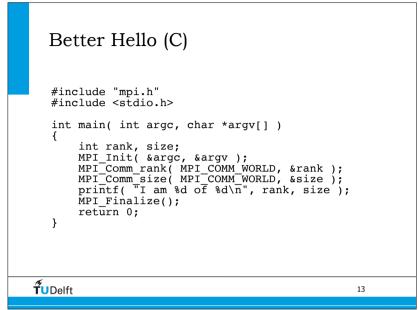




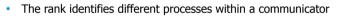
# Exercise: Hello World

- README.txt
  - Try to answer the questions in the README
  - How is the program compiled?
  - How do you run the parallel program?
- There is a C and Fortran version of the exercise.
- Use the job.slurm script to submit to the workload manager slurm.
- Try to run directly without submitting to slurm (but don't tell me).

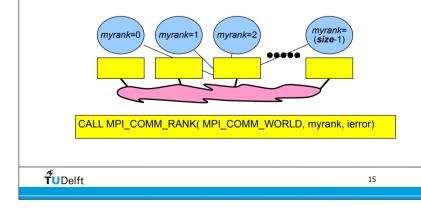
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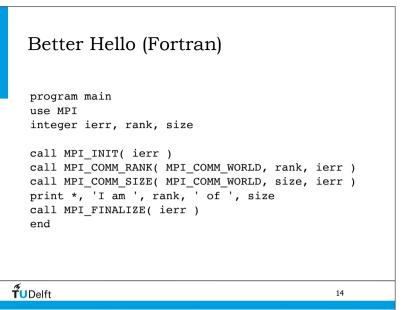


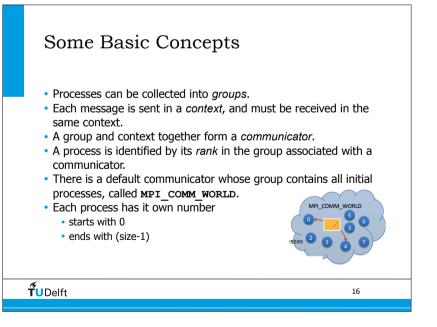
# Rank

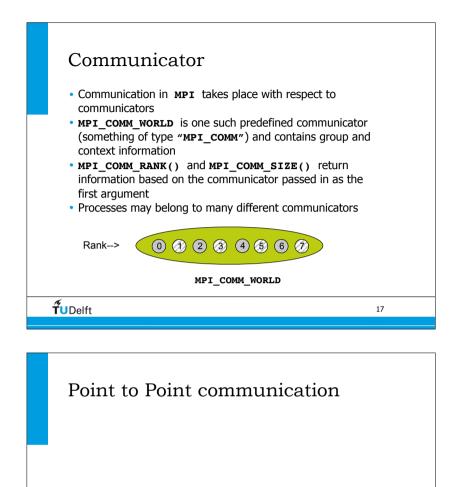


• The rank is the basis for any work and data distribution.



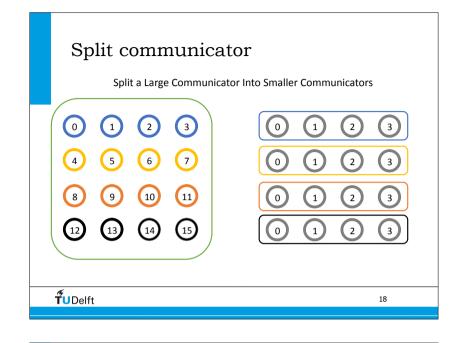


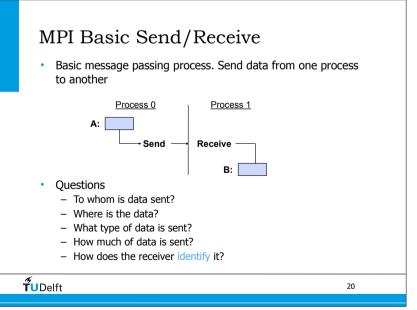


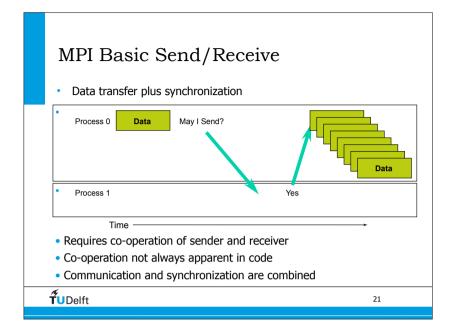


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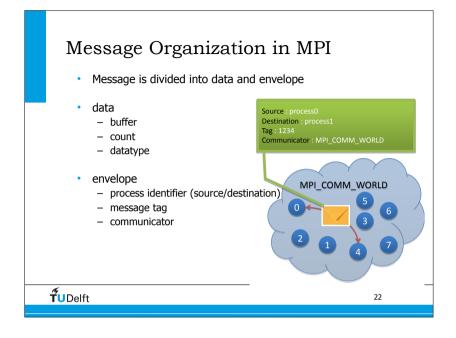


# MPI Basic Send/Receive

- Thus the basic (blocking) send has become: MPI\_Send ( buf, count, datatype, dest, tag, comm )
  - Blocking means the function does not return until it is safe to reuse the data in buffer. The message may not have been received by the target process.
- And the receive has become: MPI\_Recv( buf, count, datatype, source, tag, comm, status )
  - The source, tag, and the count of the message actually received can be retrieved from status

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C Datatypes	5
MPI datatype	
	C datatype
MPI_CHAR	signed char
MPI_SHORT	signed short int
MPI_INT	signed int
MPI_LONG	signed long int
MPI_UNSIGNED_CHAR	unsigned char
MPI_UNSIGNED_SHORT	unsigned short int
MPI_UNSIGNED_LONG	unsigned long_int
MPI_UNSIGNED	unsigned int
MPI FLOAT	float
 MPI_DOUBLE	double
MPI_LONG_DOUBLE	long double
MPI_BYTE	
MPI_PACKED	
MIT_INCAD	

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# **MPI** Fortran Datatypes

MPI FORTRAN	FORTRAN datatypes
MPI_INTEGER	INTEGER
MPI_REAL	REAL
MPI_REAL8	REAL*8
MPI_DOUBLE_PRECISION	DOUBLE PRECISION
MPI_COMPLEX	COMPLEX
MPI_LOGICAL	LOGICAL
MPI_CHARACTER	CHARACTER
MPI_BYTE	
MPI_PACKED	

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# Communication Envelope

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Envelope information is returned from • MPI RECV in status. From: source rank tag • C: status.MPI SOURCE To: status.MPI TAG destination rank <u>count</u> via MPI Get count() item-1 • Fortran: item-2 status(MPI\_SOURCE) status(MPI\_TAG) item-3 "count" item-4 elements <u>count</u> via MPI\_GET\_COUNT() item-n **T**UDelft 27

# Process Naming and Message Tags Naming a process - destination is specified by ( rank, group ) - Processes are named according to their rank in the group - Groups are defined by their distinct "communicator" - MPI ANY SOURCE wildcard rank permitted in a receive Tags are integer variables or constants used to uniquely identify individual messages • Tags allow programmers to deal with the arrival of messages in an orderly manner • MPI tags are guaranteed to range from 0 to 32767 by MPI-1 - Vendors are free to increase the range in their implementations • MPI ANY TAG can be used as a wildcard value **T**UDelft 26

Retrieving Further Information
<ul> <li>Status is a data structure allocated in the user's program.</li> <li>In C: <ul> <li>int recvd_tag, recvd_from, recvd_count;</li> <li>MPI_Status status;</li> <li>MPI_Recv(, MPI_ANY_SOURCE, MPI_ANY_TAG,, &amp;status)</li> <li>recvd_tag = status.MPI_TAG;</li> <li>recvd_from = status.MPI_SOURCE;</li> <li>MPI_Get_count( &amp;status, datatype, &amp;recvd_count );</li> </ul> </li> <li>In Fortran:</li> </ul>
<pre>integer recvd_tag, recvd_from, recvd_count integer status(MPI_STATUS_SIZE) call MPI_RECV(, MPI_ANY_SOURCE, MPI_ANY_TAG, status, ierr) tag_recvd = status(MPI_TAG) recvd_from = status(MPI_SOURCE) call MPI_GET_COUNT(status, datatype, recvd_count, ierr)</pre>
TUDelft 28

# Requirements for Point-to-Point Communications

For a communication to succeed:

- Sender must specify a valid destination rank.
- Receiver must specify a valid source rank.
- The communicator must be the same.
- Tags must match.
- Message datatypes must match.
- Receiver's buffer must be large enough.

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	<pre>if (myid &gt; 0){</pre>					
		MPI-task	0	P	2	3
		-1-	MPI_send	MPI_send	MPI_send	
		-2-		MPI_recv	MPI_recv	MPI_recv
		-3-		MPI_Get_c	MPI_Get_c	MPI_Get_c
TUDelft 31					31	

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# Exercise: SendRecv (1)

Write a simple program where every processor sends data to the next one. You may use as a starting point the basic.c or basic.f90. The program should work as follows:

- Let ntasks be the number of the tasks.
- Every task with a rank less than ntasks-1 sends a message to task myid+1. For example, task 0 sends a message to task 1.
- The message content is an integer array of 100 elements.
- The message tag is the receiver's id number.
- The sender prints out the number of elements it sends and the tag number.

• All tasks with rank  $\geq 1$  receive messages. You should check the MPI\_SOURCE and MPI\_TAG fields of the status variable (in Fortran you should check the corresponding array elements). Also check the number of elements that the process received using MPI\_Get\_count.

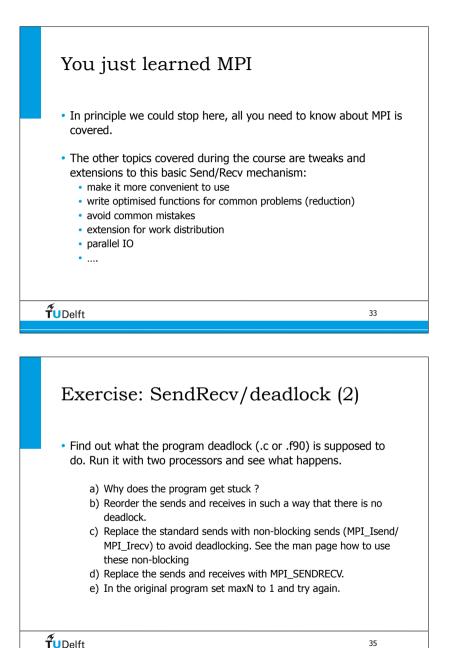
• Each receiver prints out the number of elements it received, the message tag, and the rank.

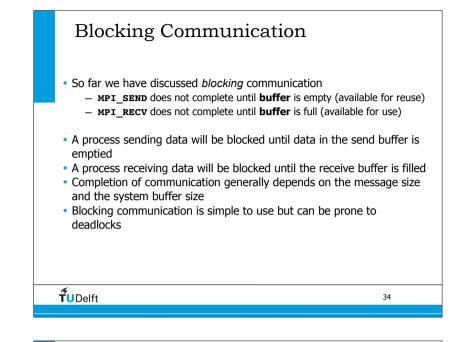
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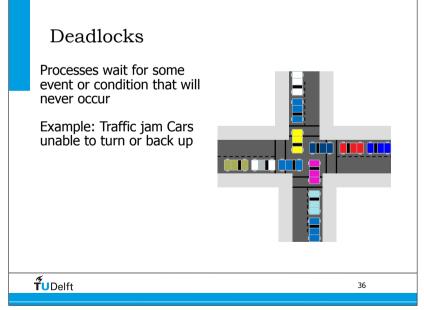
Write the program using MPI\_Send and MPI\_Recv

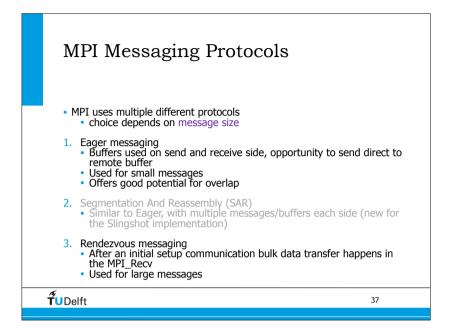
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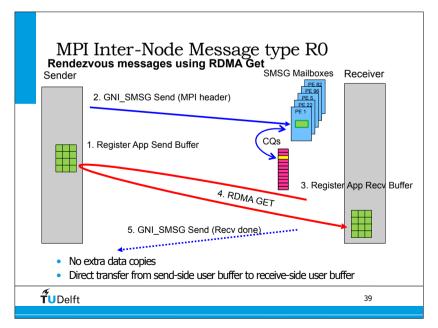
### Is MPI Large or Small? • Is MPI large (300+ functions) or small (6 functions)? - MPI's extensive functionality requires many functions - Number of functions not necessarily a measure of complexity - Many programs can be written with just 6 basic functions MPI INIT MPI COMM SIZE MPI SEND MPI FINALIZE MPI COMM RANK MPI RECV • MPI is just right A small number of concepts - Large number of functions provides flexibility, robustness, efficiency, modularity, and convenience - One need not master all parts of MPI to use it **T**UDelft 32

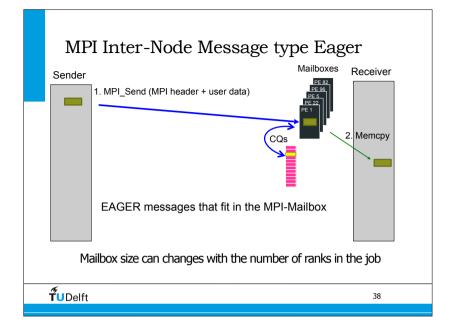


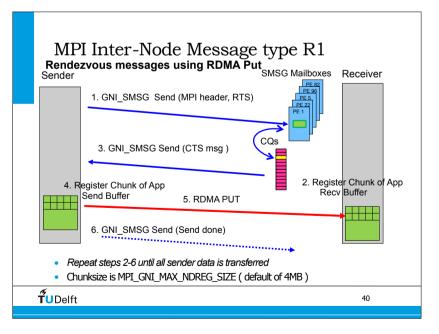


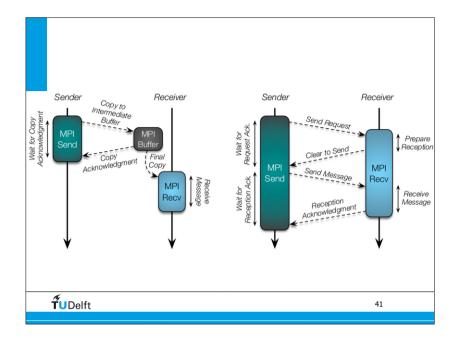


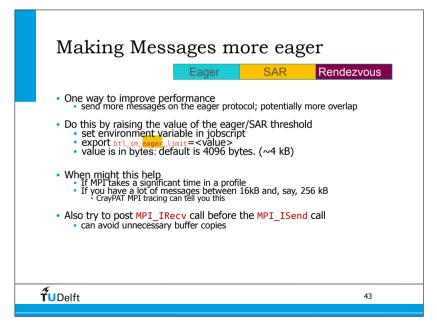


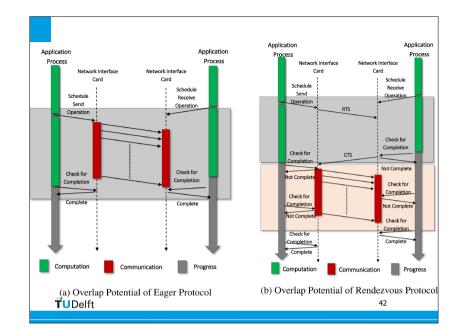












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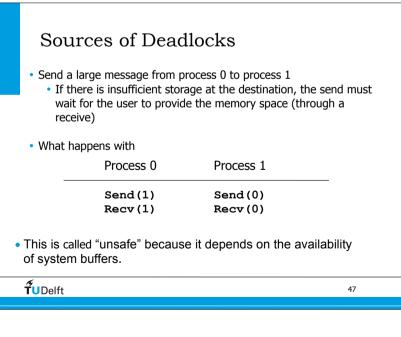
# Tuning MPI message protocols

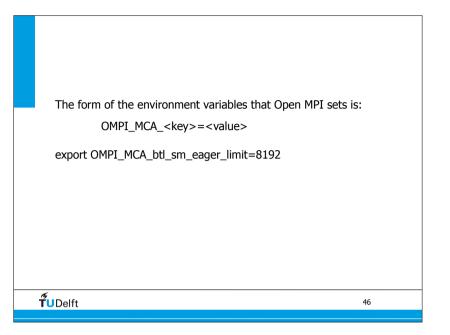
- btl\_sm\_eager\_limit: If message data plus header information fits within this limit, the message is sent "eagerly" that is, a sender attempts to write its entire message to shared buffers without waiting for a receiver to be ready. Above this size, a sender will only write the first part of a message, then wait for the receiver to acknowledge its readiness before continuing. Eager sends can improve performance by decoupling senders from receivers.
  btl\_sm\_max\_send\_size: Large messages are sent in fragments of this size. Larger segments can lead to greater efficiencies, though they could perhaps also inhibit pipelining between sender and receiver
- receiver.
- btl\_sm\_free\_list\_num: This is the initial number of fragments on each (eager and max) free list. The free lists can grow in response to resource congestion, but you can increase this parameter to pre-reserve space for more fragments.

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### ompi info --param btl all

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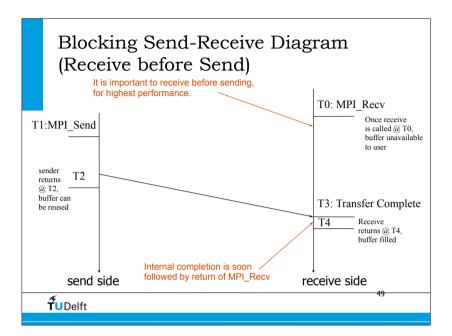




# Some Solutions to the "unsafe" Problem

• Order the operations more carefully:

	Process 0	Process 1	
	Send(1) Recv(1)	Recv(0) Send(0)	
• Use	e non-blocking operations: Process 0	Process 1	
	Isend(1)	Isend(0)	
	Irecv(1)	Irecv(0)	
	Waitall	Waitall	
Ť∪De	lft		48



# Non-Blocking Communications

- Separate communication into three phases:
- Initiate non-blocking communication
  - returns Immediately
  - routine name starting with MPI\_I...
- Do some work

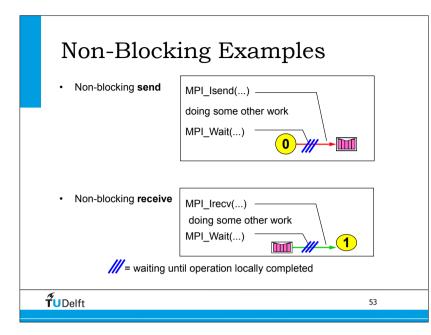
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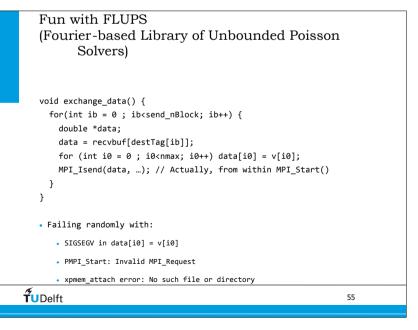
- "latency hiding"
- Wait for non-blocking communication to complete

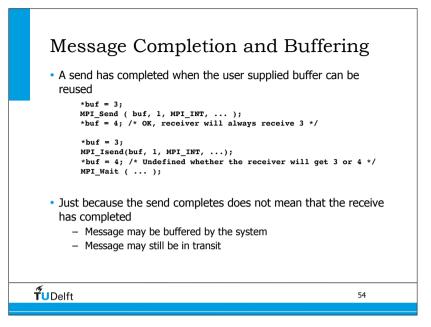
51

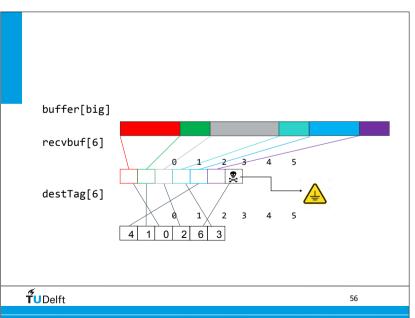
### Non-Blocking Send-Receive Diagram **High Performance Implementations** Offer Low Overhead for Non-blocking Calls Once receive is called @ T0, buffer unavailable to user T0: MPI Irecv T2: MPI Isend T1: Returns returns @ T3 sender buffer unavailable T4: MPI Wait called sender Т5 completes @ T5 buffer available T6 after MPI\_Wait T7: transfer T6: MPI Wait = =T8 T9: Wait returns finishes MPI Wait, returns @ T8 here, receive buffer filled Internal completion is soon followed by return of MPI Wait receive side send side **T**UDelft

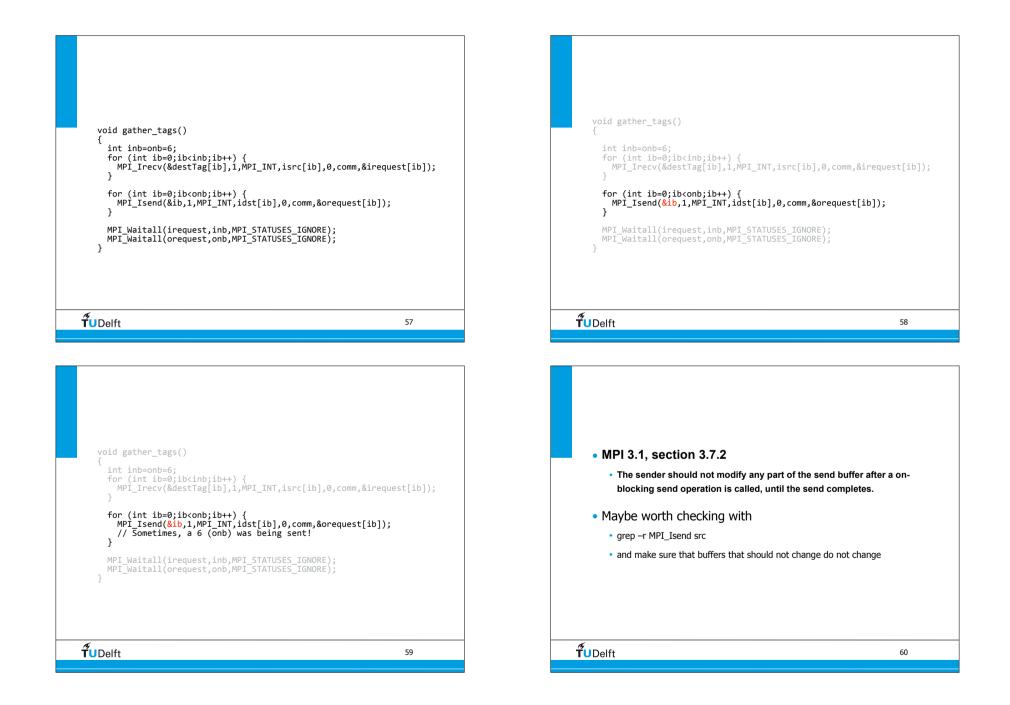
Non-Blocking Communicat • Non-blocking (asynchronous) operations return "request handles" that can be waited on and qu MPI ISEND( start, count, datatype,	(immediately) ieried			
comm, request )	,,			
MPI_IRECV( start, count, datatype,	, src, tag,			
comm, request )				
<ul> <li>MPI_WAIT( request, status )</li> <li>Non-blocking operations allow overlapping comp communication.</li> </ul>	outation and			
• Anywhere you use MPI_send or MPI_Recv, you can use the pair				
Of MPI_Isend/MPI_Wait Of MPI_Irecv/MPI_Wait				
<ul> <li>Combinations of blocking and non-blocking sends/receives can be used to synchronize execution instead of barriers</li> </ul>				
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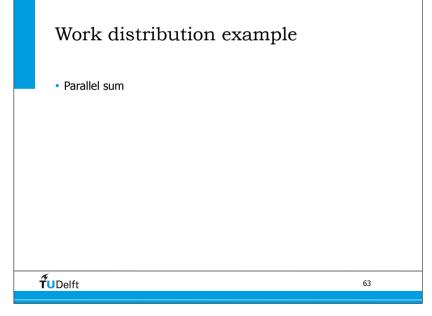


# Multiple Completion's

- It is often desirable to wait on multiple requests
- An example is a worker/manager program, where the manager waits for one or more workers to send it a message
  - MPI\_WAITALL( count, array\_of\_requests, array\_of\_statuses )

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- MPI\_WAITANY( count, array\_of\_requests, index, status )
- MPI\_WAITSOME( incount, array\_of\_requests, outcount, array\_of\_indices, array\_of\_statuses )
- There are corresponding versions of **TEST** for each of these
- **T**UDelft



# Send Mode ( MPI\_Send, MPI\_Isend ) Standard mode ( MPI\_Send, MPI\_Isend ) The standard MPI send, the send will not complete until the send buffer is empty Synchronous mode ( MPI\_Ssend, MPI\_Issend ) The send does not complete until after a matching receive has been posted. Buffered mode ( MPI\_Bsend, MPI\_Ibsend ) User supplied buffer space is used for system buffering The send will complete as soon as the send buffer is copied to the system buffer

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 Porallel Sum

 Memory

 Porallel Principality

 Porallel algorithm

 - Parallel algorithm

 - Scatter

 - Half of the array is sent to process 0

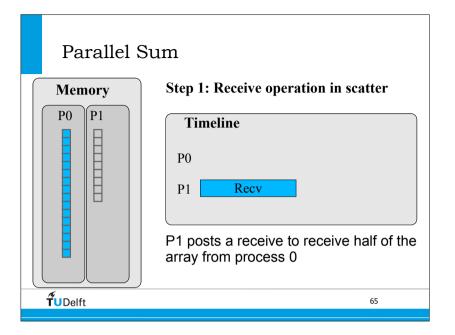
 - Compute

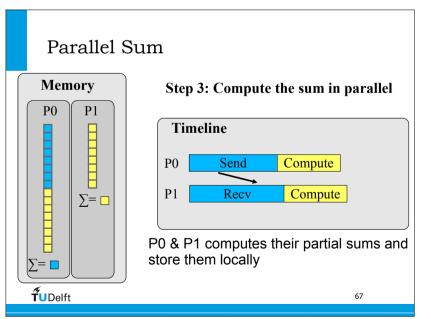
 - Po & P1 sum independently their segment

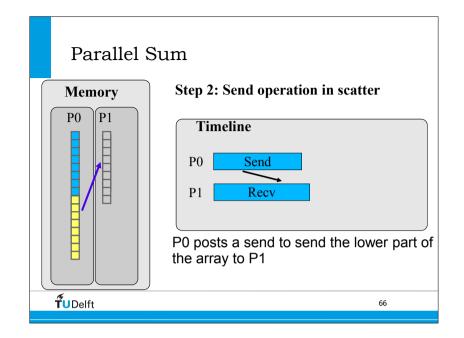
 - Reduction

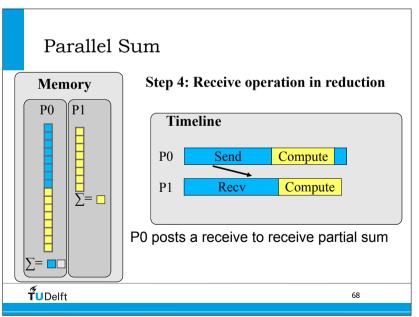
 - Potial sum on P1 sent to P0

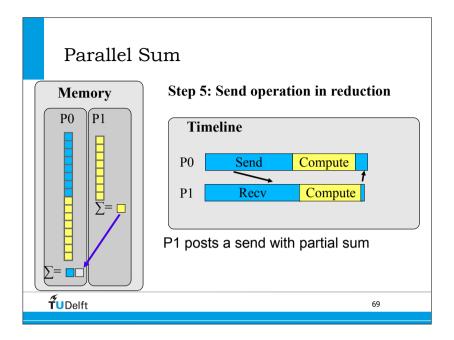
 - Po sums the partial sums

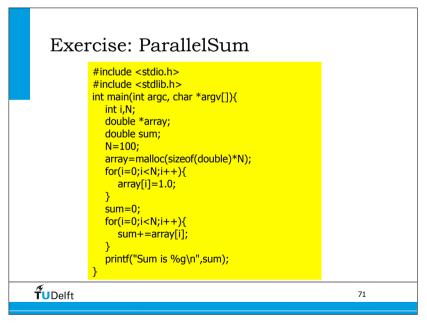


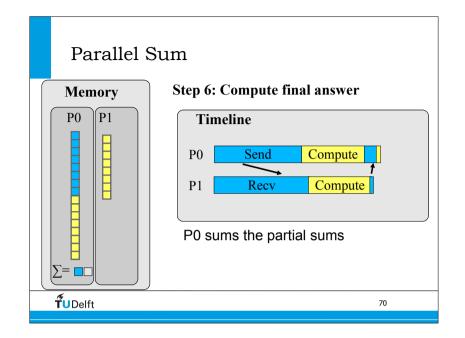


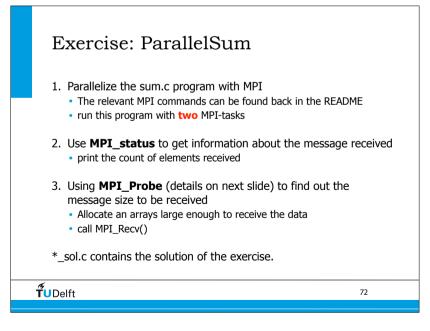


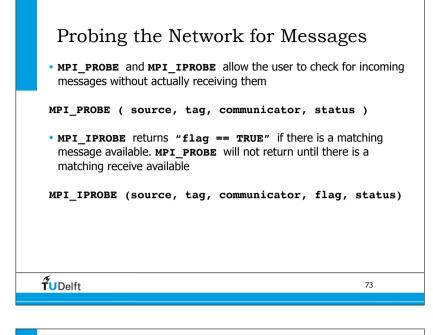












# Exercise: PingPong/Basic

Ping-pong is a standard test in which two processes repeatedly pass a message back and forth.

Write a program that sends a 'float' array of fixed length, say, ten times back (ping) and forth (pong) to obtain an average time for one ping-pong.

Time the ping-pongs with MPI\_WTIME() calls.

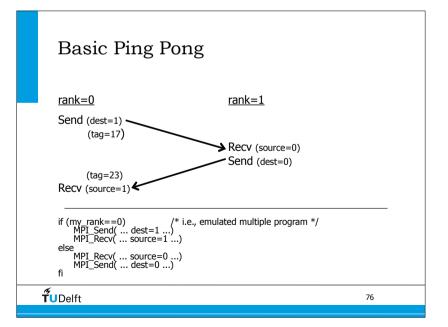
You may use pingpong.c or pingpong.f90 as a starting point for this exercise.

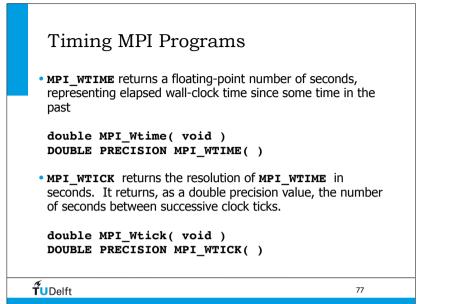
Investigate how the bandwidth varies with the size of the message.

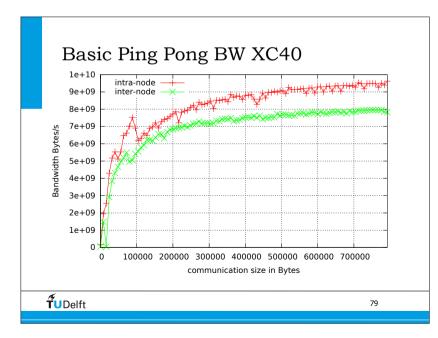
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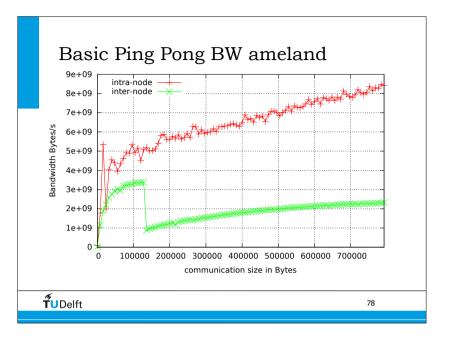
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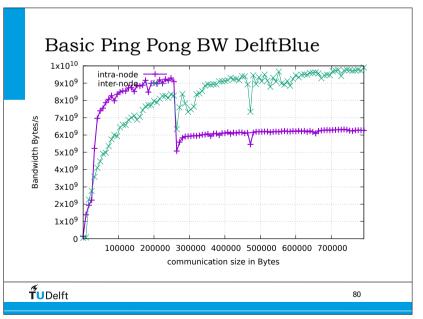
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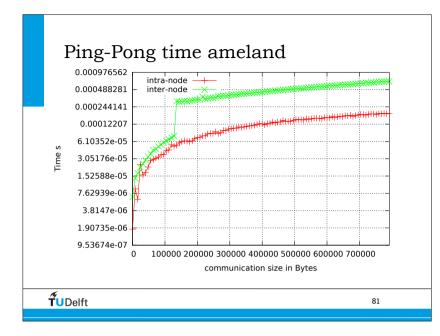


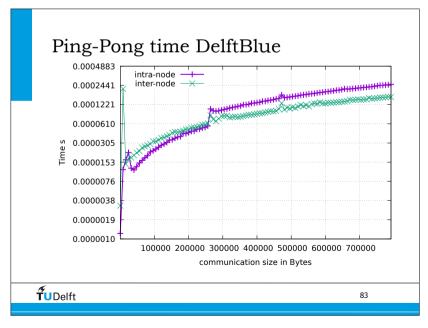


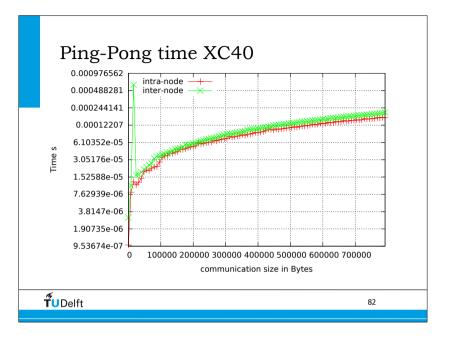


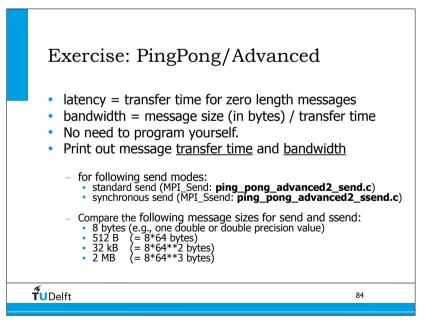


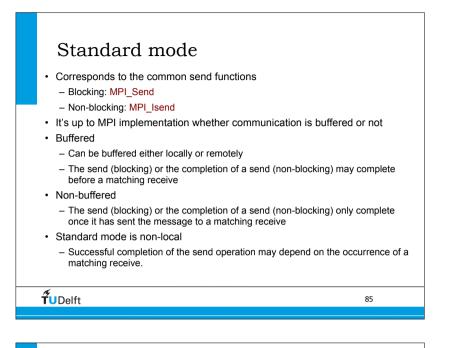












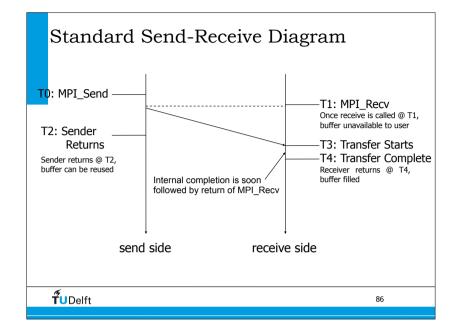
# Synchronous mode

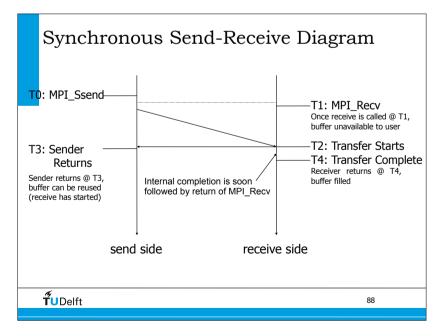
### Blocking: MPI Ssend

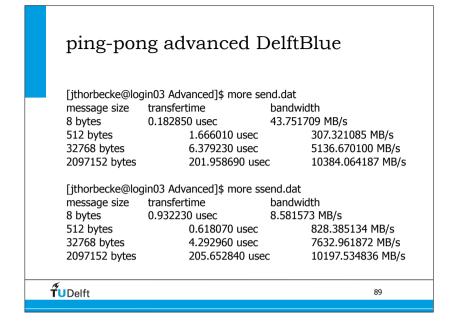
- Blocking send only returns once the corresponding receive has been posted
- Same parameters as for standard mode send MPI Send
- Uses
  - Debugging potential deadlocks in the program are found by using synchronous sends
  - If many processes send messages to one process its unexpected message buffer can run out if it doesn't pre-post receives. By using MPI\_Ssend this can be avoided! Typical example is IO where single process writes data
- Non-blocking: MPI\_Issend
  - The completion (wait/test) of the send only returns once the corresponding receive has been posted
  - Same parameters as for standard mode send MPI\_Isend
  - Useful for debugging can be used to measure worst case scenario for how long the completion command has to wait

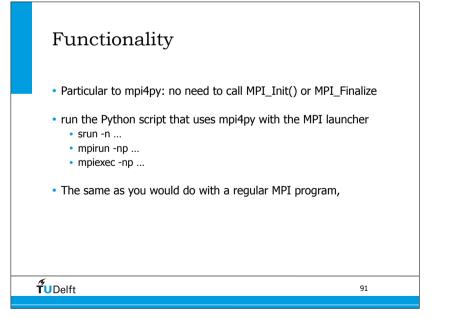
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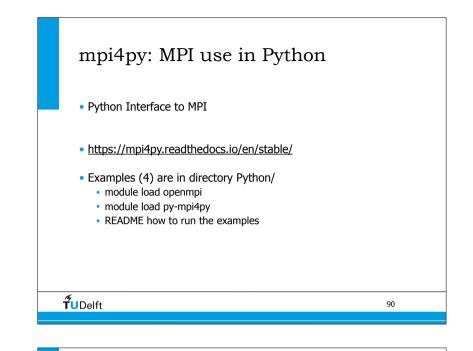
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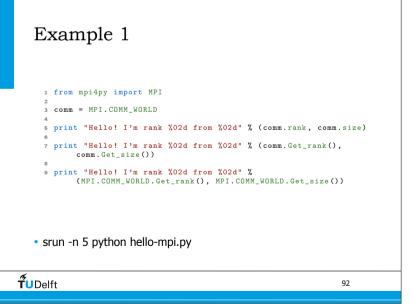












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### Example 2 1 from mpi4py import MPI 2 3 comm = MPI.COMM\_WORLD 4 assert comm.size == 2 5 6 if comm.rank == 0: sendmsg = 123 7 comm.send(sendmsg, dest=1, tag=11) 8 recvmsg = comm.recv(source=1, tag=22) 9 10 print "[%02d] Received message: %s" % (comm.rank, recvmsg) 11 else: recvmsg = comm.recv(source=0, tag=11) 12 print "[%02d] Received message: %d" % (comm.rank, recvmsg) 13 sendmsg = "Message from 1" 14 comm.send(sendmsg, dest=0, tag=22) 15 **T**UDelft 94