

Cluster computing/Scripting/MPI

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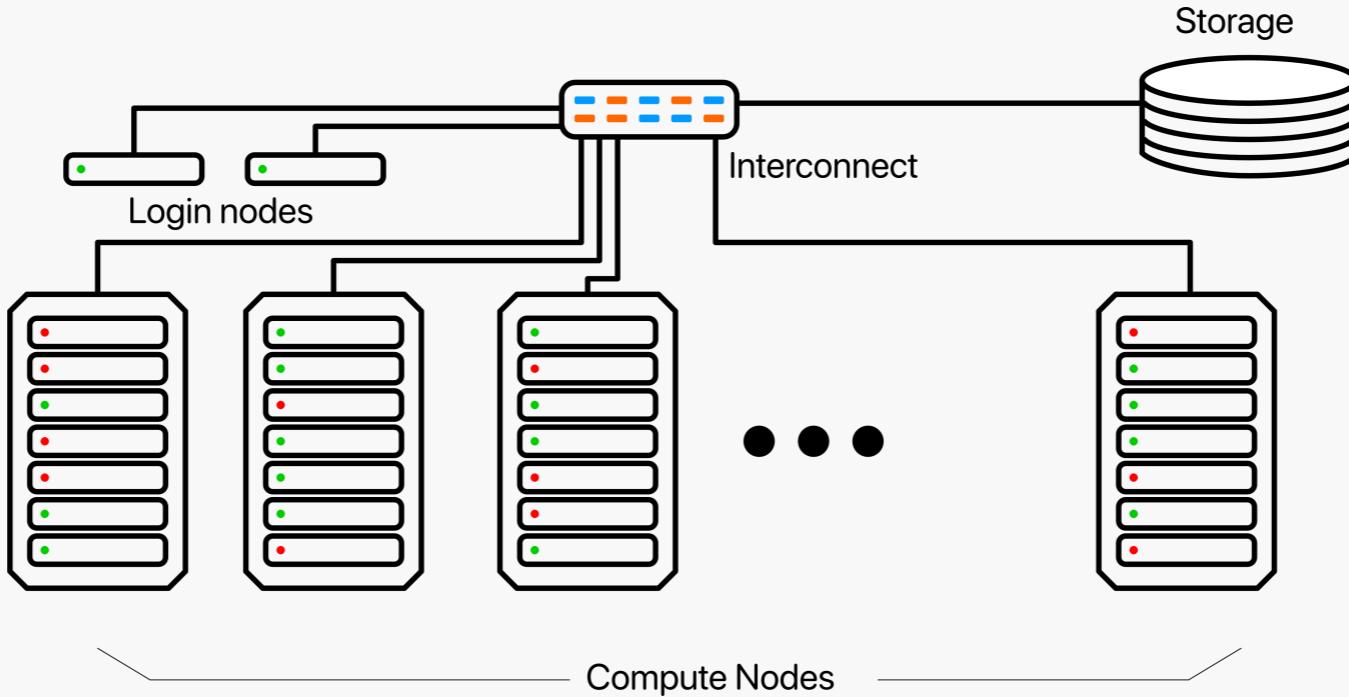
NCC Beginner Training Event, 14th - 16th November 2022

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Cluster Computing



Specific configuration of our cluster "Cyclone"

- 33 nodes in total
 - Hostnames: `cn{01, ..., 17}, gpu{01, ..., 16}`
- We will be using CPU nodes: `cn ??`
 - 2×20-core Intel Xeon (Gold) "Cascade Lake"
 - 192 GBytes RAM
- Common storage for our course: `/onyx/data/edu17/`

Cluster Computing

- Log in to a *login node* or *frontend node*. Login node in our case has hostname `front01`
- To run programs on *compute nodes*, a *job scheduler* is available
- Distinguish between *interactive* and *batch* jobs

Cluster Computing

- Log in to a *login node* or *frontend node*. Login node in our case has hostname `front01`
- To run programs on *compute nodes*, a *job scheduler* is available
- Distinguish between *interactive* and *batch* jobs

SLURM job scheduler

- See currently running and waiting jobs: `squeue`
- Ask for an interactive job: `salloc`
- Submit a batch job: `sbatch`
- Run an executable: `srun`

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

- Type `hostname`. This tells you the name of the node you are currently logged into:

```
[ikoutsou@front01 ~]$ hostname  
front01
```

this is the login node.

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

- Type `hostname`. This tells you the name of the node you are currently logged into:

```
[ikoutsou@front01 ~]$ hostname  
front01
```

this is the login node.

- Ask for one node (`-N 1`):

```
[ikoutsou@front01 ~]$ salloc -N 1 -p cpu --reservation=NCC -A edu17  
salloc: Granted job allocation 197848  
salloc: Waiting for resource configuration  
salloc: Nodes cn17 are ready for job  
[ikoutsou@cn17 ~]$
```

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

- Type `hostname`. This tells you the name of the node you are currently logged into:

```
[ikoutsou@front01 ~]$ hostname  
front01
```

this is the login node.

- Ask for one node (`-N 1`):

```
[ikoutsou@front01 ~]$ salloc -N 1 -p cpu --reservation=NCC -A edu17  
salloc: Granted job allocation 197848  
salloc: Waiting for resource configuration  
salloc: Nodes cn17 are ready for job  
[ikoutsou@cn17 ~]$
```

- `-A edu17`: charge project with name `edu17`

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

- Type `hostname`. This tells you the name of the node you are currently logged into:

```
[ikoutsou@front01 ~]$ hostname  
front01
```

this is the login node.

- Ask for one node (`-N 1`):

```
[ikoutsou@front01 ~]$ salloc -N 1 -p cpu --reservation=NCC -A edu17  
salloc: Granted job allocation 197848  
salloc: Waiting for resource configuration  
salloc: Nodes cn17 are ready for job  
[ikoutsou@cn17 ~]$
```

- `-A edu17`: charge project with name `edu17`
- `--reservation=NCC`: use reservation `NCC`

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

- Type `hostname`. This tells you the name of the node you are currently logged into:

```
[ikoutsou@front01 ~]$ hostname  
front01
```

this is the login node.

- Ask for one node (`-N 1`):

```
[ikoutsou@front01 ~]$ salloc -N 1 -p cpu --reservation=NCC -A edu17  
salloc: Granted job allocation 197848  
salloc: Waiting for resource configuration  
salloc: Nodes cn17 are ready for job  
[ikoutsou@cn17 ~]$
```

- `-A edu17`: charge project with name `edu17`
- `--reservation=NCC`: use reservation `NCC`
- `-p=cpu`: requests a node from the `cpu` partition

Cluster Computing Introductory Example

- Log in:

```
[localhost ~]$ ssh <username>@cyclone.hpcf.cyi.ac.cy
```

- Type `hostname`. This tells you the name of the node you are currently logged into:

```
[ikoutsou@front01 ~]$ hostname  
front01
```

this is the login node.

- Ask for one node (`-N 1`):

```
[ikoutsou@front01 ~]$ salloc -N 1 -p cpu --reservation=NCC -A edu17  
salloc: Granted job allocation 197848  
salloc: Waiting for resource configuration  
salloc: Nodes cn17 are ready for job  
[ikoutsou@cn17 ~]$
```

- `-A edu17`: charge project with name `edu17`
- `--reservation=NCC`: use reservation `NCC`
- `-p=cpu`: requests a node from the `cpu` partition

- Type `hostname` again:

```
[ikoutsou@cn17 ~]$ hostname  
cn17
```

`cn17` is a compute node.

Cluster Computing Introductory Example

- Release the node (type `exit` or hit `ctrl-d`):

```
[ikoutsou@cn17 ~]$ exit
salloc: Relinquishing job allocation 197848
[ikoutsou@front01 ~]$
```

we're back on `front01`

Cluster Computing Introductory Example

- Release the node (type `exit` or hit `ctrl-d`):

```
[ikoutsou@cn17 ~]$ exit
salloc: Relinquishing job allocation 197848
[ikoutsou@front01 ~]$
```

we're back on `front01`

Please **do not** hold nodes unnecessarily; when you have nodes `salloced` you may be blocking other users from using those nodes.

Cluster Computing Introductory Example

- Release the node (type `exit` or hit `ctrl-d`):

```
[ikoutsou@cn17 ~]$ exit
salloc: Relinquishing job allocation 197848
[ikoutsou@front01 ~]$
```

we're back on `front01`

Please **do not** hold nodes unnecessarily; when you have nodes `salloced` you may be blocking other users from using those nodes.

- Use `srun` instead of `salloc`:

```
[ikoutsou@front01 ~]$ srun -N 1 -p cpu -A edu17 --reservation=NCC hostname
cn15
```

Cluster Computing Introductory Example

- Release the node (type `exit` or hit `ctrl-d`):

```
[ikoutsou@cn17 ~]$ exit
salloc: Relinquishing job allocation 197848
[ikoutsou@front01 ~]$
```

we're back on `front01`

Please **do not** hold nodes unnecessarily; when you have nodes `salloced` you may be blocking other users from using those nodes.

- Use `srun` instead of `salloc`:

```
[ikoutsou@front01 ~]$ srun -N 1 -p cpu -A edu17 --reservation=NCC hostname
cn15
```

- Allocates a node, runs the specified command (in this case `hostname`), and then exits the node, releasing the allocation

Cluster Computing Introductory Example

- Release the node (type `exit` or hit `ctrl-d`):

```
[ikoutsou@cn17 ~]$ exit
salloc: Relinquishing job allocation 197848
[ikoutsou@front01 ~]$
```

we're back on `front01`

Please **do not** hold nodes unnecessarily; when you have nodes `salloced` you may be blocking other users from using those nodes.

- Use `srun` instead of `salloc`:

```
[ikoutsou@front01 ~]$ srun -N 1 -p cpu -A edu17 --reservation=NCC hostname
cn15
```

- Allocates a node, runs the specified command (in this case `hostname`), and then exits the node, releasing the allocation
- The output, `cn15`, reveals that we were allocated node `cn15` for this specific — very short — job

Cluster Computing Introductory Example

- Run multiple instances of `hostname` in parallel:

```
[ikoutsou@front01 ~]$ srun -N 1 -n 2 -p cpu -A edu17 --reservation=NCC hostname  
cn15  
cn15
```

Cluster Computing Introductory Example

- Run multiple instances of `hostname` in parallel:

```
[ikoutsou@front01 ~]$ srun -N 1 -n 2 -p cpu -A edu17 --reservation=NCC hostname  
cn15  
cn15
```

- `-N 1`: use one node
- `-n 2`: use two processes

Cluster Computing Introductory Example

- Run multiple instances of `hostname` in parallel:

```
[ikoutsou@front01 ~]$ srun -N 1 -n 2 -p cpu -A edu17 --reservation=NCC hostname  
cn15  
cn15
```

- `-N 1`: use one node
- `-n 2`: use two processes

- Run on more than one node:

```
[ikoutsou@front01 ~]$ srun -N 2 -n 2 -p cpu -A edu17 --reservation=NCC hostname  
cn02  
cn01
```

runs one instance of `hostname` on each node

Cluster Computing Introductory Example

- Run multiple instances of `hostname` in parallel:

```
[ikoutsou@front01 ~]$ srun -N 1 -n 2 -p cpu -A edu17 --reservation=NCC hostname  
cn15  
cn15
```

- `-N 1`: use one node
- `-n 2`: use two processes

- Run on more than one node:

```
[ikoutsou@front01 ~]$ srun -N 2 -n 2 -p cpu -A edu17 --reservation=NCC hostname  
cn02  
cn01
```

runs one instance of `hostname` on each node

- Try:

```
[ikoutsou@front01 ~]$ srun -N 2 -n 1 -p cpu -A edu17 --reservation=NCC hostname
```

Cluster Computing Introductory Example

- Run multiple instances of `hostname` in parallel:

```
[ikoutsou@front01 ~]$ srun -N 1 -n 2 -p cpu -A edu17 --reservation=NCC hostname
cn15
cn15
```

- N 1: use one node
 - n 2: use two processes

- Run on more than one node:

```
[ikoutsou@front01 ~]$ srun -N 2 -n 2 -p cpu -A edu17 --reservation=NCC hostname
cn02
cn01
```

runs one instance of `hostname` on each node

- Try:

```
[ikoutsou@front01 ~]$ srun -N 2 -n 1 -p cpu -A edu17 --reservation=NCC hostname
```

```
srun: Warning: can't run 1 processes on 2 nodes, setting nnodes to 1
cn17
```

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

- `pwd` will tell you where you are in the file system:

```
[ikoutsou@front01 NCC-training]$ pwd  
/nvme/h/ikoutsou/NCC-training
```

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

- `pwd` will tell you where you are in the file system:

```
[ikoutsou@front01 NCC-training]$ pwd  
/nvme/h/ikoutsou/NCC-training
```

- `/` is referred to as the *root directory*

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

- `pwd` will tell you where you are in the file system:

```
[ikoutsou@front01 NCC-training]$ pwd  
/nvme/h/ikoutsou/NCC-training
```

- `/` is referred to as the *root directory*
- `.` is an alias for the *current directory*

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

- `pwd` will tell you where you are in the file system:

```
[ikoutsou@front01 NCC-training]$ pwd  
/nvme/h/ikoutsou/NCC-training
```

- `/` is referred to as the *root directory*
- `.` is an alias for the *current directory*
- `..` is an alias for the directory one level above

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

- `pwd` will tell you where you are in the file system:

```
[ikoutsou@front01 NCC-training]$ pwd  
/nvme/h/ikoutsou/NCC-training
```

- `/` is referred to as the *root directory*
- `.` is an alias for the *current directory*
- `..` is an alias for the directory one level above
- `~` is an alias for your *home directory*

Cluster Computing Introductory Example

- Make a directory. List it, see that it is there:

```
[ikoutsou@front01 ~]$ mkdir NCC-training  
[ikoutsou@front01 ~]$ ls  
NCC-training  
[ikoutsou@front01 ~]$
```

- Change into it:

```
[ikoutsou@front01 ~]$ cd NCC-training/  
[ikoutsou@front01 NCC-training]$
```

- `pwd` will tell you where you are in the file system:

```
[ikoutsou@front01 NCC-training]$ pwd  
/nvme/h/ikoutsou/NCC-training
```

- `/` is referred to as the *root directory*
- `.` is an alias for the *current directory*
- `..` is an alias for the directory one level above
- `~` is an alias for your *home directory*
- E.g.:

```
[ikoutsou@front01 NCC-training]$ cd .. / .. /  
[ikoutsou@front01 h]$ pwd  
/nvme/h
```

Cluster Computing Introductory Example

- `cd` without any additional arguments takes you home (equivalent to `cd ~`)

```
[ikoutsou@front01 h]$ cd  
[ikoutsou@front01 ~]$ pwd  
/nvme/h/ikoutsou
```

Cluster Computing Introductory Example

- `cd` without any additional arguments takes you home (equivalent to `cd ~`)

```
[ikoutsou@front01 ~]$ cd  
[ikoutsou@front01 ~]$ pwd  
/nvme/h/ikoutsou
```

- Make a subdirectory under `NCC-training` for our first C program

```
[ikoutsou@front01 ~]$ cd NCC-training  
[ikoutsou@front01 NCC-training]$ mkdir 01  
[ikoutsou@front01 NCC-training]$ cd 01  
[ikoutsou@front01 01]$
```

Cluster Computing Introductory Example

- `cd` without any additional arguments takes you home (equivalent to `cd ~`)

```
[ikoutsou@front01 ~]$ cd  
[ikoutsou@front01 ~]$ pwd  
/nvme/h/ikoutsou
```

- Make a subdirectory under `NCC-training` for our first C program

```
[ikoutsou@front01 ~]$ cd NCC-training  
[ikoutsou@front01 NCC-training]$ mkdir 01  
[ikoutsou@front01 NCC-training]$ cd 01  
[ikoutsou@front01 01]$
```

- Use `emacs` or `vim` (or any other text editor you're comfortable with) to type out our first program

Cluster Computing Introductory Example

- `cd` without any additional arguments takes you home (equivalent to `cd ~`)

```
[ikoutsou@front01 ~]$ cd  
[ikoutsou@front01 ~]$ pwd  
/nvme/h/ikoutsou
```

- Make a subdirectory under `NCC-training` for our first C program

```
[ikoutsou@front01 ~]$ cd NCC-training  
[ikoutsou@front01 NCC-training]$ mkdir 01  
[ikoutsou@front01 NCC-training]$ cd 01  
[ikoutsou@front01 01]$
```

- Use `emacs` or `vim` (or any other text editor you're comfortable with) to type out our first program
- E.g.:

```
[ikoutsou@front01 01]$ emacs -nw prog-01.c
```

Cluster Computing Introductory Example

- Type out the following program:

```
#include <unistd.h>
#include <stdio.h>
#include <sys/types.h>

int
main(int argc, char *argv[])
{
    char hname[256];
    pid_t p;
    gethostname(hname, 256);
    p = getpid();
    printf(" Hostname: %s, pid: %lu\n", hname, p);
    return 0;
}
```

Cluster Computing Introductory Example

- Type out the following program:

```
#include <unistd.h>           // ← provides definitions for gethostname() and getpid()
#include <stdio.h>            // ← provides definitions for printf()
#include <sys/types.h>         // ← defines the pid_t type
// main()
//   → argv[] is an array of strings which holds all command line arguments
//   → argc holds the number of elements of argv
int
main(int argc, char *argv[])
{
    char hname[256];          // ← declare hname[] as an array of 256 characters (a string of length 256)
    pid_t p;                  // ← declare p as a pid_t type, in this case, an unsigned long integer
    gethostname(hname, 256);   // ← call gethostname(), return hostname in hname which is 256 characters long
    p = getpid();              // ← call getpid(), return value in p
    printf(" Hostname: %s, pid: %lu\n", hname, p); // print statement
    return 0; // ← return a value of 0 to the operating system. By convention 0 means success.
}
```

Cluster Computing Introductory Example

After saving and exiting, you are back at the command line

Cluster Computing Introductory Example

After saving and exiting, you are back at the command line

- `ls` should show the file `prog-01.c`, which you just typed in and saved:

```
[ikoutsou@front01 01]$ ls  
prog-01.c
```

Cluster Computing Introductory Example

After saving and exiting, you are back at the command line

- `ls` should show the file `prog-01.c`, which you just typed in and saved:

```
[ikoutsou@front01 01]$ ls  
prog-01.c
```

- It's time to *compile* it into an executable. Use the `gompi` module:

```
[ikoutsou@front01 01]$ module load gOMPI  
[ikoutsou@front01 01]$ gcc prog-01.c -o p01
```

Cluster Computing Introductory Example

After saving and exiting, you are back at the command line

- `ls` should show the file `prog-01.c`, which you just typed in and saved:

```
[ikoutsou@front01 01]$ ls  
prog-01.c
```

- It's time to *compile* it into an executable. Use the `gomp` module:

```
[ikoutsou@front01 01]$ module load gomp  
[ikoutsou@front01 01]$ gcc prog-01.c -o p01
```

- `-o p01` means "name the resulting executable `p01`". If you don't specify `-o` the executable name defaults to `a.out`

Cluster Computing Introductory Example

After saving and exiting, you are back at the command line

- `ls` should show the file `prog-01.c`, which you just typed in and saved:

```
[ikoutsou@front01 01]$ ls  
prog-01.c
```

- It's time to *compile* it into an executable. Use the `gompi` module:

```
[ikoutsou@front01 01]$ module load gOMPI  
[ikoutsou@front01 01]$ gcc prog-01.c -o p01
```

- `-o p01` means "name the resulting executable `p01`". If you don't specify `-o` the executable name defaults to `a.out`
- Type `ls` to make sure it has been created. Then run it on the frontend node:

```
[ikoutsou@front01 01]$ ls  
p01  prog-01.c  
[ikoutsou@front01 01]$ ./p01  
Hostname: front01, pid: 14848
```

Cluster Computing Introductory Example

After saving and exiting, you are back at the command line

- `ls` should show the file `prog-01.c`, which you just typed in and saved:

```
[ikoutsou@front01 01]$ ls  
prog-01.c
```

- It's time to *compile* it into an executable. Use the `gomp` module:

```
[ikoutsou@front01 01]$ module load gomp  
[ikoutsou@front01 01]$ gcc prog-01.c -o p01
```

- `-o p01` means "name the resulting executable `p01`". If you don't specify `-o` the executable name defaults to `a.out`
- Type `ls` to make sure it has been created. Then run it on the frontend node:

```
[ikoutsou@front01 01]$ ls  
p01  prog-01.c  
[ikoutsou@front01 01]$ ./p01  
Hostname: front01, pid: 14848
```

Note: commands like `gcc` or `ls` are globally accessible because their locations are included in your shell environment's search path. for `p01` though, which we just created, you need to explicitly give its path, in this case via `./` which means "current directory".

Cluster Computing Introductory Example

- Run your new program `p01` using `srun` on two nodes with two processes each

Cluster Computing Introductory Example

- Run your new program `p01` using `srun` on two nodes with two processes each

```
[ikoutsou@front01 01]$ srun -N 2 -n 4 -p cpu --reservation=NCC -A edu17 ./p01
Hostname: cn02, pid: 11037
Hostname: cn01, pid: 8207
Hostname: cn01, pid: 8209
Hostname: cn01, pid: 8208
[ikoutsou@front01 01]$
```

Cluster Computing Introductory Example

- Run your new program `p01` using `srun` on two nodes with two processes each

```
[ikoutsou@front01 01]$ srun -N 2 -n 4 -p cpu --reservation=NCC -A edu17 ./p01
Hostname: cn02, pid: 11037
Hostname: cn01, pid: 8207
Hostname: cn01, pid: 8209
Hostname: cn01, pid: 8208
[ikoutsou@front01 01]$
```

- Go nuts 😊 :

```
[ikoutsou@front01 01]$ srun -N 2 -n 48 -p cpu --reservation=NCC -A edu17 ./p01
Hostname: cn02, pid: 11294
Hostname: cn02, pid: 11300
...
Hostname: cn01, pid: 8482
Hostname: cn01, pid: 8474
[ikoutsou@front01 01]$
```

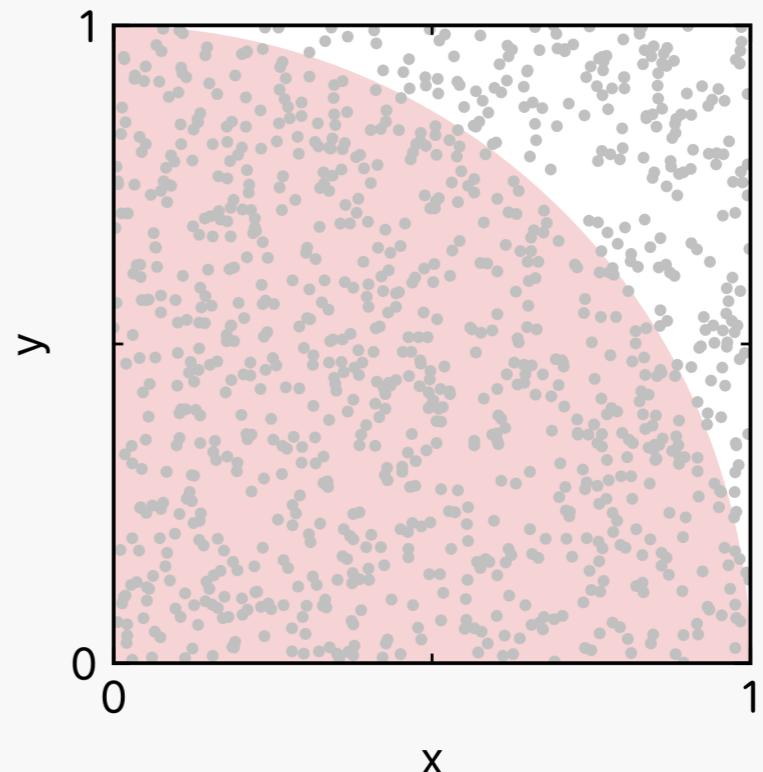
(You will see 48 lines; I suppressed some above)

Cluster Computing Introductory Example

- Time for step 2 — writing a simple program to compute π in parallel

Cluster Computing Introductory Example

- Time for step 2 — writing a simple program to compute π in parallel
- Calculation of π via simple Monte Carlo:
 - Define a unit square. Set $n_{\text{hit}} = 0$
 - Randomly pick points (x, y) within the unit square
 - If $x^2 + y^2 < 1$, $n_{\text{hit}} += 1$
 - Repeat N times
- The ratio n_{hit}/N approaches the area of a circle quadrant $\Rightarrow \frac{\pi}{4}$



Cluster Computing Introductory Example

- First, make a new directory under `~/NCC-training/`.

```
[ikoutsou@front01 01]$ cd ../  
[ikoutsou@front01 NCC-training]$ mkdir 02  
[ikoutsou@front01 01]$ cd 02/  
[ikoutsou@front01 02]$
```

Cluster Computing Introductory Example

- First, make a new directory under `~/NCC-training/`.

```
[ikoutsou@front01 01]$ cd ../  
[ikoutsou@front01 NCC-training]$ mkdir 02  
[ikoutsou@front01 01]$ cd 02/  
[ikoutsou@front01 02]$
```

- Copy the program `pi.c` from `/onyx/data/edu17/02/pi.c`

```
[ikoutsou@front01 02]$ cp /onyx/data/edu17/02/pi.c .
```

- Inspect `pi.c`, e.g.:

```
[ikoutsou@front01 02]$ emacs -nw pi.c
```

Cluster Computing Introductory Example

- First, make a new directory under `~/NCC-training/`.

```
[ikoutsou@front01 01]$ cd ../  
[ikoutsou@front01 NCC-training]$ mkdir 02  
[ikoutsou@front01 01]$ cd 02/  
[ikoutsou@front01 02]$
```

- Copy the program `pi.c` from `/onyx/data/edu17/02/pi.c`

```
[ikoutsou@front01 02]$ cp /onyx/data/edu17/02/pi.c .
```

- Inspect `pi.c`, e.g.:

```
[ikoutsou@front01 02]$ emacs -nw pi.c
```

Cluster Computing Introductory Example

```
#include <unistd.h>
#include <stdio.h>
#include <stdlib.h>
#include <sys/types.h>

int
main(int argc, char *argv[])
{
    unsigned long int N = 10000;
    unsigned long int nhit = 0;
    for(int i=0; i<N; i++) {
        double x = drand48();
        double y = drand48();
        if((x*x + y*y) < 1)
            nhit++;
    }
    double pi = 4.0 * (double)nhit/(double)N;
    printf(" N = %16d    pi = %lf\n", N, pi);
    return 0;
}
```

Cluster Computing Introductory Example

```
#include <unistd.h>
#include <stdio.h>
#include <stdlib.h>
#include <sys/types.h>

int
main(int argc, char *argv[])
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    return 0;
}
```

- Compile and run on frontend:

```
[ikoutsou@front01 02]$ gcc pi.c -o pi
[ikoutsou@front01 02]$ ./pi
N =          10000    pi = 3.136400
```

Cluster Computing Introductory Example

- Now run, e.g. on 4 processes:

Cluster Computing Introductory Example

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```
[ikoutsov@front01 02]$ srun -N 1 -n 4 -p cpu -A edu17 --reservation=NCC ./pi
N =      10000    pi = 3.148800
```

Cluster Computing Introductory Example

- Now run, e.g. on 4 processes:

```
[ikoutsov@front01 02]$ srun -N 1 -n 4 -p cpu -A edu17 --reservation=NCC ./pi  
N =      10000    pi = 3.148800  
N =      10000    pi = 3.148800  
N =      10000    pi = 3.148800  
N =      10000    pi = 3.148800
```

We get exactly the same result four times 😊

Cluster Computing Introductory Example

- Now run, e.g. on 4 processes:

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We get exactly the same result four times 😱

We need to seed the random number generator differently for each process

Cluster Computing Introductory Example

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We get exactly the same result four times 😱

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- Use the process id (`pid`) from the previous example, to seed the random number generator

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main(int argc, char *argv[])
{
    unsigned long int N = 10000;
    unsigned long int nhit = 0;

    pid_t p = getpid(); // ← Add this
    srand48(p); // ← Add this

    for(int i=0; i<N; i++) {
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        double y = drand48();
        if((x*x + y*y) < 1)
            nhit++;
    }
    double pi = 4.0 * (double)nhit/(double)N;
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    return 0;
}
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Cluster Computing Introductory Example

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- `srand48()` sets the random number generator **seed**

Cluster Computing Introductory Example

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    double pi = 4.0 * (double)nhit/(double)N;
    printf(" N = %16d    pi = %lf\n", N, pi);
    return 0;
}
```

- `srand48()` sets the random number generator **seed**
- Need a unique seed for each instance of the program ⇒ use process id.

Cluster Computing Introductory Example

- Compile again and run:

```
[ikoutsou@front01 02]$ gcc pi.c -o pi
[ikoutsou@front01 02]$ srun -N 1 -n 4 -p cpu -A edu17 --reservation=NCC ./pi
N =          10000    pi = 3.150800
N =          10000    pi = 3.143200
N =          10000    pi = 3.151200
N =          10000    pi = 3.152800
```

Cluster Computing Introductory Example

- Compile again and run:

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[ikoutsou@front01 02]$ srun -N 1 -n 4 -p cpu -A edu17 --reservation=NCC ./pi
N =          10000    pi = 3.150800
N =          10000    pi = 3.143200
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- Now we would like to average over these four values to obtain a better estimate of π

Cluster Computing Introductory Example

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```

- Now we would like to average over these four values to obtain a better estimate of π
- First redirect the output to a file, e.g.:

```
[ikoutsou@front01 02]$ gcc pi.c -o pi
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```

Cluster Computing Introductory Example

- Compile again and run:

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[ikoutsou@front01 02]$ srun -N 1 -n 4 -p cpu -A edu17 --reservation=NCC ./pi
N =          10000  pi = 3.150800
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- Now we would like to average over these four values to obtain a better estimate of π
- First redirect the output to a file, e.g.:

```
[ikoutsou@front01 02]$ gcc pi.c -o pi
[ikoutsou@front01 02]$ srun -N 1 -n 4 -p cpu -A edu17 --reservation=NCC ./pi > pi-out.txt
```

- Now `pi-out.txt` contains the four lines of output

```
[ikoutsou@front01 02]$ ls
pi  pi.c  pi-out.txt
[ikoutsou@front01 02]$ more pi-out.txt
N =          10000  pi = 3.113600
N =          10000  pi = 3.128400
N =          10000  pi = 3.156800
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[ikoutsou@front01 02]$
```

Cluster Computing Introductory Example

- The program `awk` allows us to add over columns of a file.

Cluster Computing Introductory Example

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- E.g.:

```
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3.136800
```

Cluster Computing Introductory Example

- The program `awk` allows us to add over columns of a file.
- E.g.:

```
[ikoutsou@front01 02]$ cat pi-out.txt | awk '{pi_sum+=$6}; END {printf "%8.6f\n", pi_sum/NR}'  
3.136800
```

- Use more processes

```
[ikoutsou@front01 02]$ srun -N 2 -n 80 -p cpu -A edu17 --reservation=NCC ./pi > pi-out.txt  
[ikoutsou@front01 02]$ cat pi-out.txt | awk '{pi_sum+=$6}; END {printf "%8.6f\n", pi_sum/NR}'  
3.142435
```

Cluster Computing Introductory Example

- The program `awk` allows us to add over columns of a file.
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- `$6` is the sixth column in the file, the value for π per process
- `pi_sum` is our summation variable
- `NR` is an AWK internal variable, the number of rows

Cluster Computing Introductory Example

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 - `pi_sum` is our summation variable
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- Let's wrap this up in a script "for posterity"
 - In fact, we'll write a *Slurm batch script*

Cluster Computing Introductory Example

- Copy from `/onyx/data/edu17/02/pi.sh`

```
#!/bin/bash
#SBATCH -J pi
#SBATCH -o pi.txt
#SBATCH -e pi.err
#SBATCH -p cpu
#SBATCH -A edu17
#SBATCH --reservation=NCC
#SBATCH -t 00:02:00
#SBATCH -n 80
#SBATCH -N 2

### Add these two lines
srun ./pi > pi-out.txt
cat pi-out.txt | awk '{sum+=$6}; END {printf "%8.6f\n", sum/NR}'
```

Cluster Computing Introductory Example

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- All Slurm options — that you so far used after `srun` — are now included in the lines starting with `#SBATCH`

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```

- All Slurm options — that you so far used after `srun` — are now included in the lines starting with `#SBATCH`
- Thus `srun` is now run without options
- Additional options include:
 - `-J`: sets the job name
 - `-o` and `-e`: set the files where the output and error should be redirected
 - `-t`: sets a time limit. The job will be killed if it exceeds this time (here 2 minutes)

Cluster Computing Introductory Example

- Submit the job

```
[ikoutsou@front01 02]$ sbatch pi.sh
Submitted batch job 198021
```

Cluster Computing Introductory Example

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Submitted batch job 198021
```

- Query its status. Filter only your jobs:

```
[ikoutsou@front01 02]$ squeue -u $(whoami)
JOBID PARTITION      NAME      USER ST      TIME   NODES NODELIST(REASON)
198021      cpu        pi ikoutsou PD      0:00       2 (Reservation)
```

- Status: PD, R, CG: "Pending", "Running", "Completing"

Cluster Computing Introductory Example

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```

- Status: PD, R, CG: "Pending", "Running", "Completing"
- After the program completes:
 - File `pi.err` contains any errors (hopefully empty)
 - File `pi.txt` contains the output of `awk` — what would be printed to the screen had you used `srun` like before
 - File `pi-out.txt` should also contain new values from the `srun` that was run during the script

MPI — Outline

- Overview of the Message Passing Interface (MPI)
- Basics of MPI
 - Distributed memory paradigm (as compared to shared memory)
 - Start-up and initialization
- Synchronization
- Collectives
- Point-to-point communication

The Message Passing Interface

- MPI: An Application Programmer Interface (API)
 - A *library specification*; determines functions, their names and arguments, and their functionality
- A *de facto* standard for programming *distributed memory* systems
- Current specification is version 4 (MPI-4.0), released June 9, 2021
 - For most systems you can reliably assume MPI-3.1 is in place
- Several free (open) or vendor-provided implementations, e.g.:
 - Mvapich
 - OpenMPI
 - IntelMPI

The Message Passing Interface

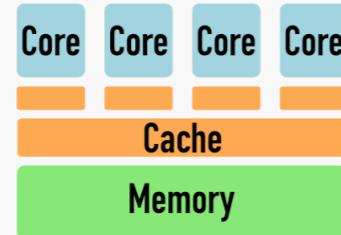
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Distributed memory programming

- Each process has its own memory domain
- MPI functions facilitate:
 - Obtaining environment information about the running process, e.g., process id, number of processes, etc.
 - Achieving *communication* between processes, e.g. synchronization, copying of data, etc.

Shared vs Distributed memory paradigm

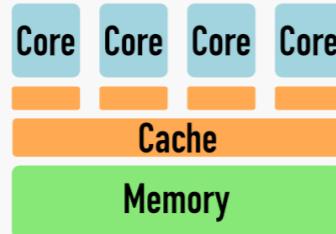
Shared memory



- Multiple processes share common memory (common *memory address space*)
- E.g. multi-core CPU, multi-socket node, GPU threads, etc.
- Programming models: OpenMP, pthreads, MPI, CUDA (sort of)

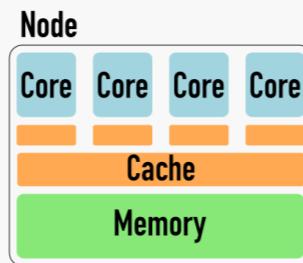
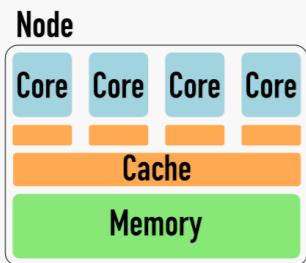
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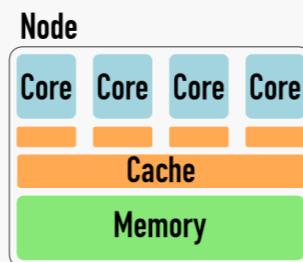
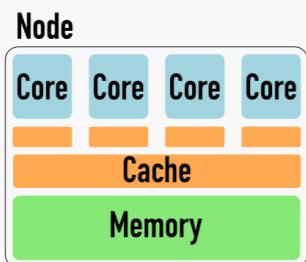


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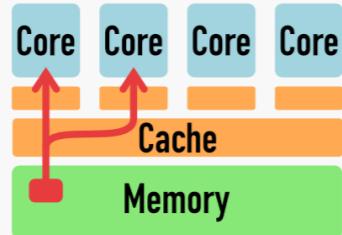


- Processes have distinct memory domains (different *memory address space*)
- E.g. multiple nodes within a cluster, multiple GPUs within a node
- Programming models: MPI



Shared vs Distributed memory paradigm

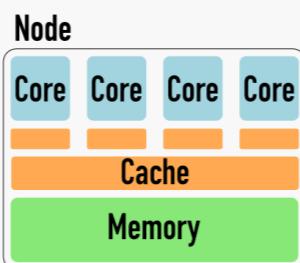
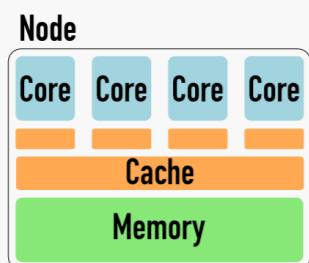
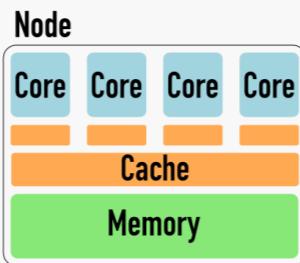
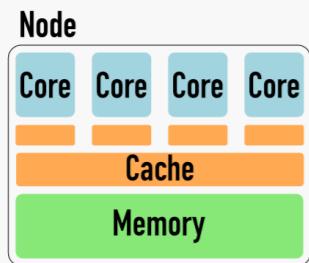
Shared memory



Data shared via memory

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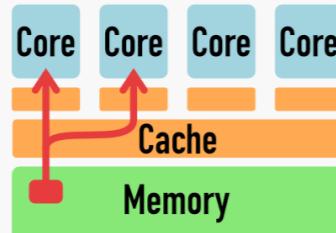
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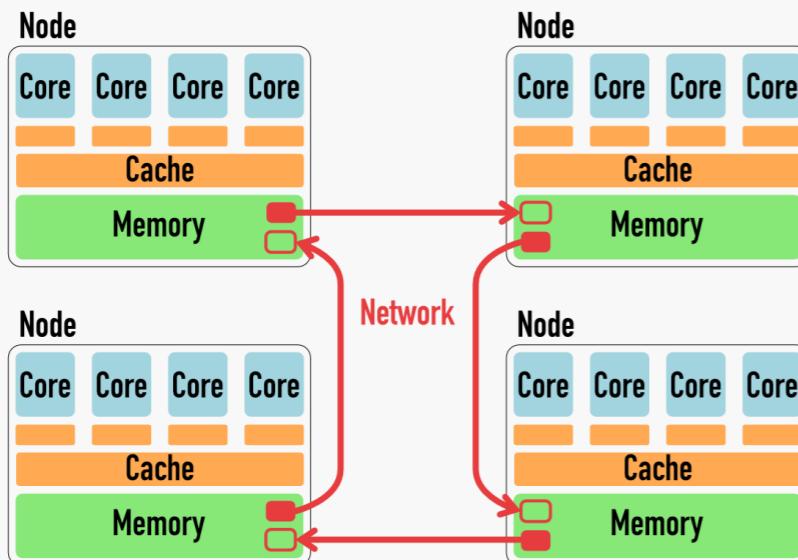
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Distributed memory



Data shared via explicit communication over a network

- Processes have distinct memory domains (different *memory address space*)
- E.g. multiple nodes within a cluster, multiple GPUs within a node
- Programming models: MPI

Running a program in parallel

- Trivially, in Linux it is simple to run a program in parallel

```
ssh node01 ./my_program &
ssh node02 ./my_program &
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```

`my_program` will run on each node identically

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- Depending on the system, instead of `mpirun` you may be required `mpiexec` or `srun` which take similar (but not identical) arguments

Compiling an MPI program

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gcc -o my_mpi_program my_mpi_program.c -I/opt/mpi/include -L/opt/mpi/lib -lmpi
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```

- Thankfully, knowing the locations of the MPI library and include files is never needed in practice; implementations come with wrappers that set the appropriate include paths and linker options:

```
mpicc -o my_mpi_program my_mpi_program.c
```

Initialization

- MPI functions begin with the `MPI_` prefix in C

Initialization

- MPI functions begin with the `MPI_` prefix in C
- Call `MPI_Init()` first, before any other MPI call:

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

where `argc` and `argv` are the typical names used for the command line variables passed to `main()`

Initialization

- MPI functions begin with the `MPI_` prefix in C
- Call `MPI_Init()` first, before any other MPI call:

```
MPI_Init(&argc, &argv);
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where `argc` and `argv` are the typical names used for the command line variables passed to `main()`

- Before the end of the program, call `MPI_Finalize()`, otherwise the MPI runtime may assume your program finished in error

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```
#include <mpi.h>

int
main(int argc, char *argv[])
{
    MPI_Init(&argc, &argv);
    /*
     ...
     ...
     ...
    */
    MPI_Finalize();
    return 0;
}
```

Initialization

- Two functions you will almost always call
 - `MPI_Comm_size()`: gives the number of parallel process running (n_{proc})
 - `MPI_Comm_rank()`: determines the *rank* of the process, i.e. a unique number between 0 and $n_{\text{proc}} - 1$ that identifies the calling process
- A complete example:

```
#include <stdio.h>
#include <mpi.h>

int
main(int argc, char *argv[])
{
    MPI_Init(&argc, &argv);
    int nproc, rank;
    MPI_Comm_size(MPI_COMM_WORLD, &nproc);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    printf(" This is rank = %d of nproc = %d\n", rank, nproc);
    MPI_Finalize();
    return 0;
}
```

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main(int argc, char *argv[])
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    MPI_Finalize();
    return 0;
}
```

- `MPI_COMM_WORLD` is an *MPI communicator*. This specific communicator is the default communicator, defined in `mpi.h`, and trivially specifies *all* processes
- A user can partition processes into subgroups by defining custom communicators, but this will not be covered here

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    printf(" This is rank = %d of nproc = %d\n", rank, nproc);
    MPI_Finalize();
    return 0;
}
```

- No assumptions can safely be made about the order in which the `printf()` statements occur, i.e. the order in which each process prints is practically random

Synchronization

- Compiling and running the previous program (assuming it is saved as `example.c`)

```
[user@front01 ~]$ mpicc -o example example.c
[user@front01 ~]$ mpirun -n 5 example
This is rank = 3 of nproc = 5
This is rank = 1 of nproc = 5
This is rank = 2 of nproc = 5
This is rank = 4 of nproc = 5
This is rank = 0 of nproc = 5
```

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This is rank = 3 of nproc = 5
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- Note that the order is random; Synchronization between processes can be achieved using the `MPI_Barrier()` function

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Synchronization

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This is rank = 1 of nproc = 5
This is rank = 2 of nproc = 5
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```

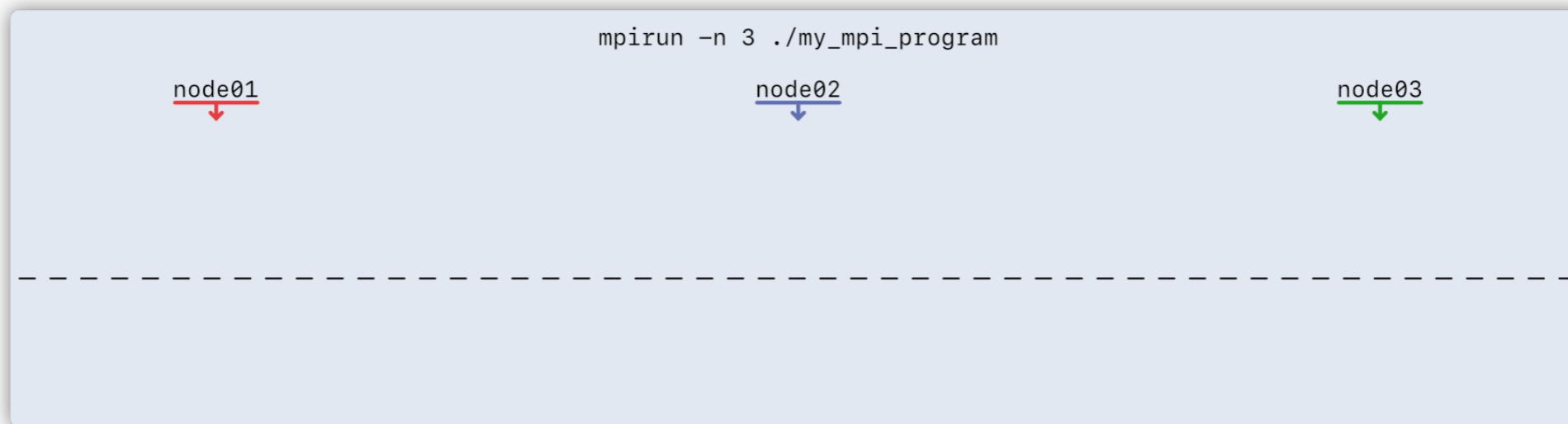
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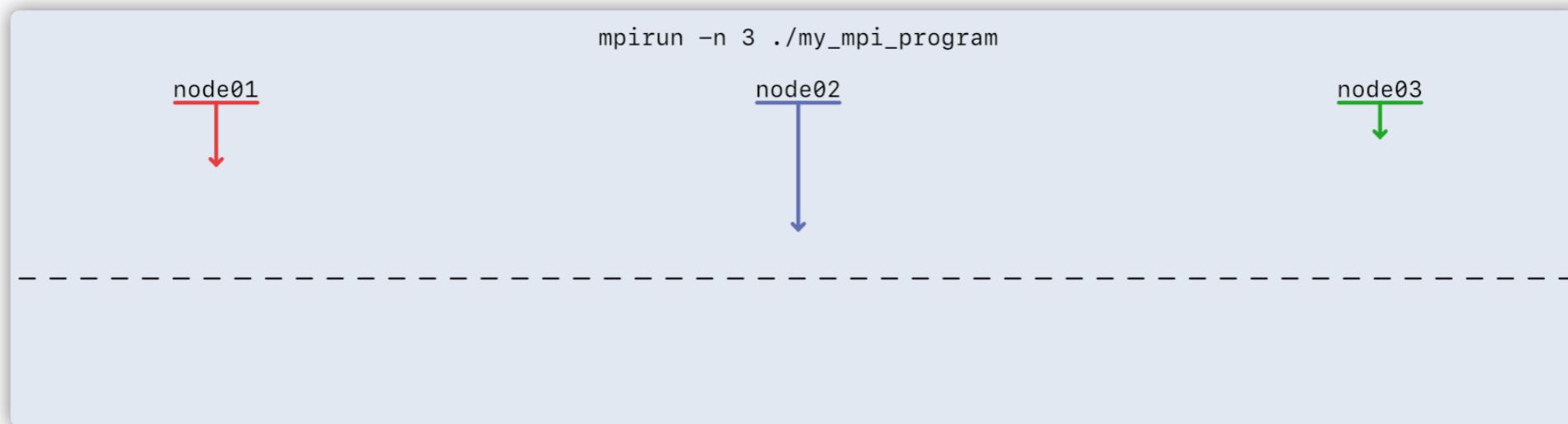


Synchronization

- Compiling and running the previous program (assuming it is saved as `example.c`)

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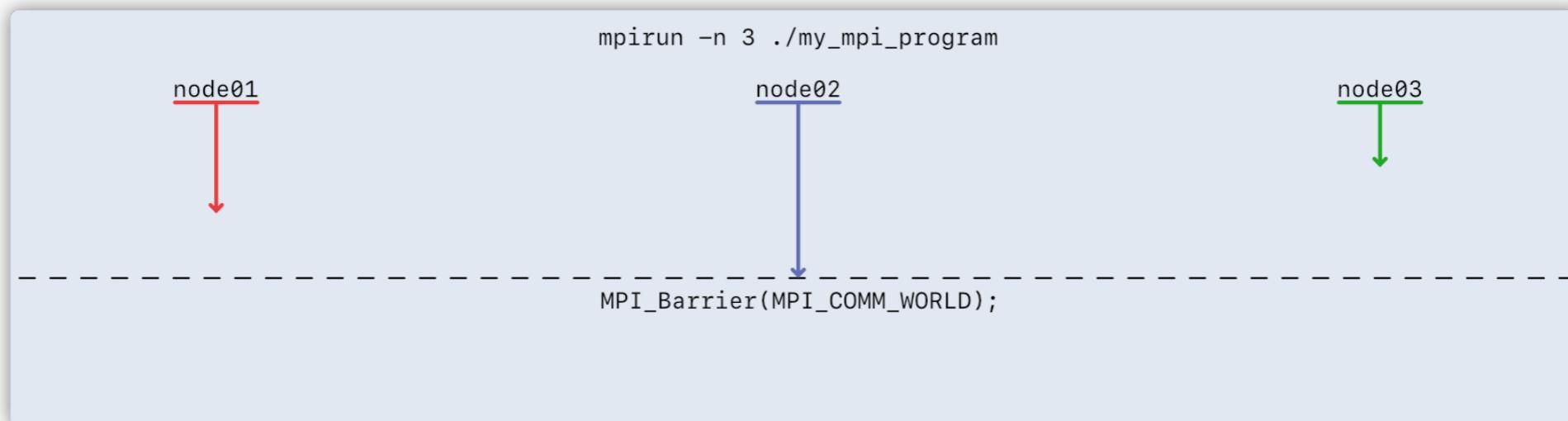


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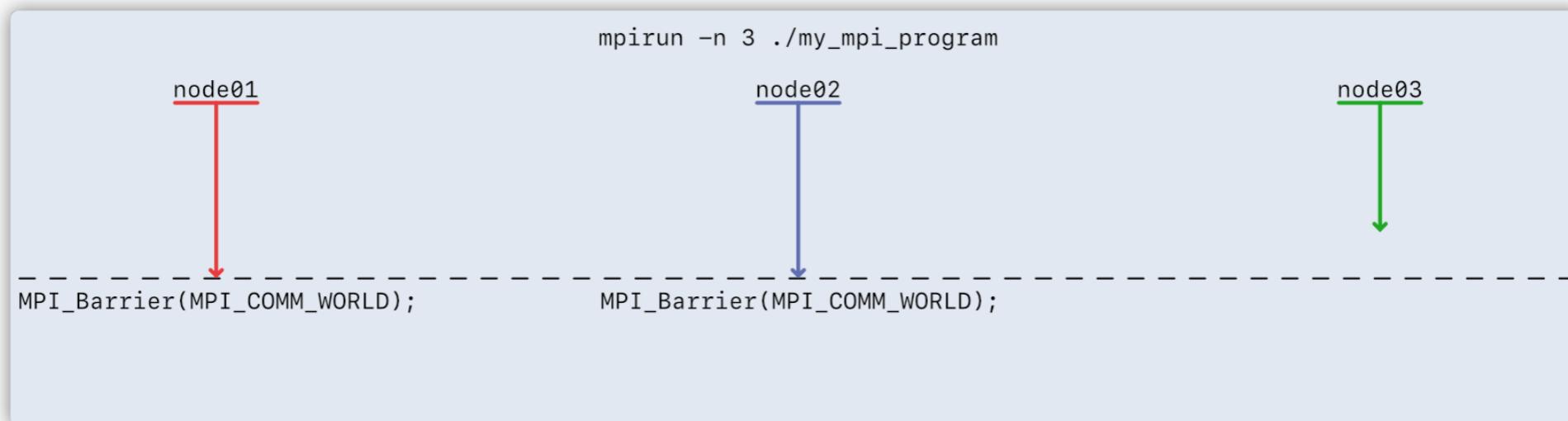


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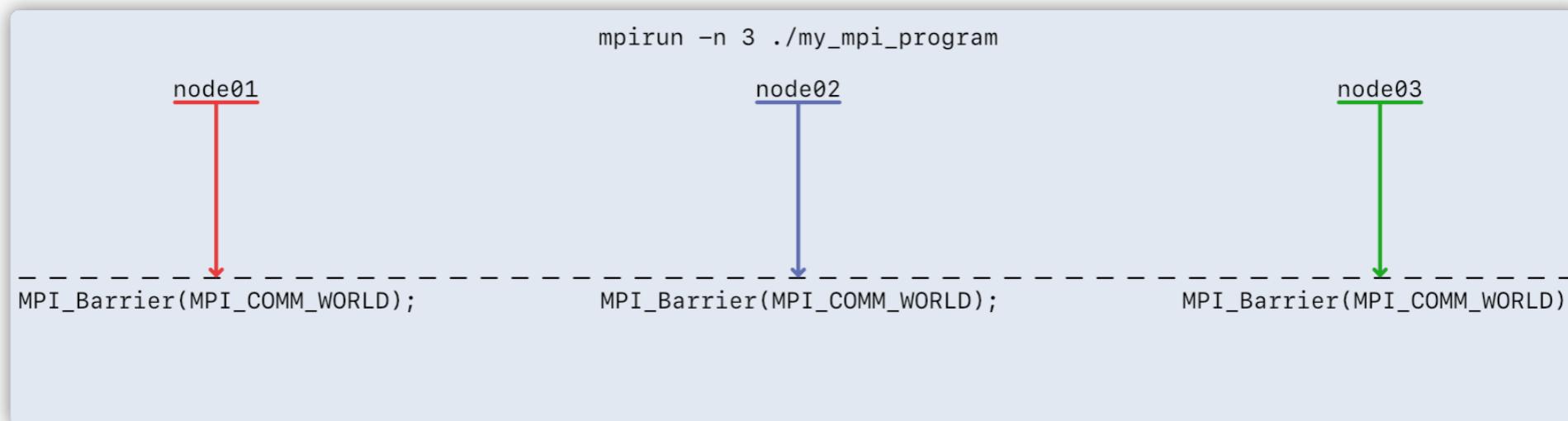


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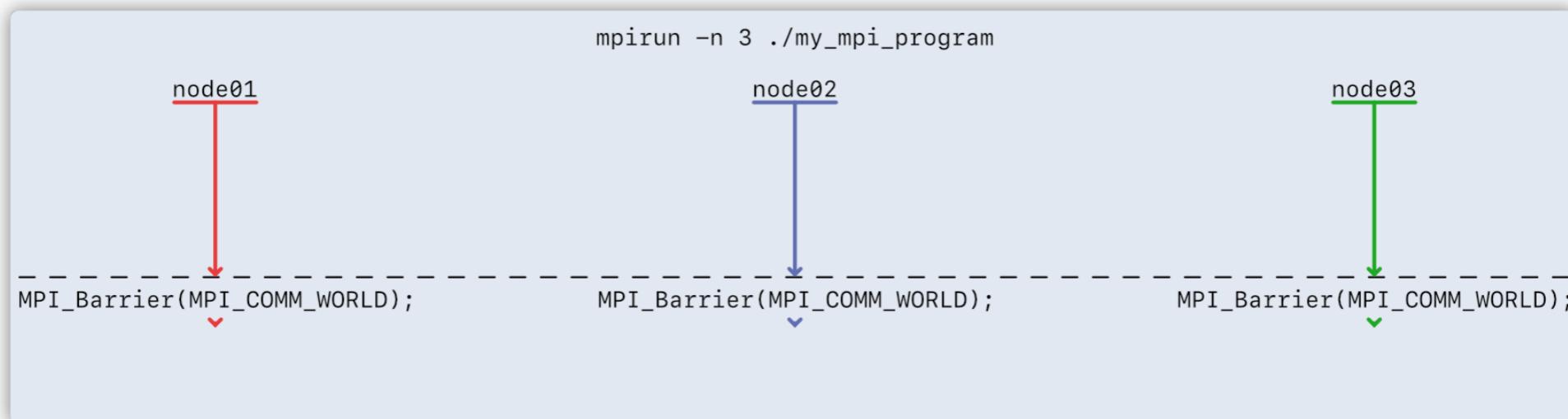


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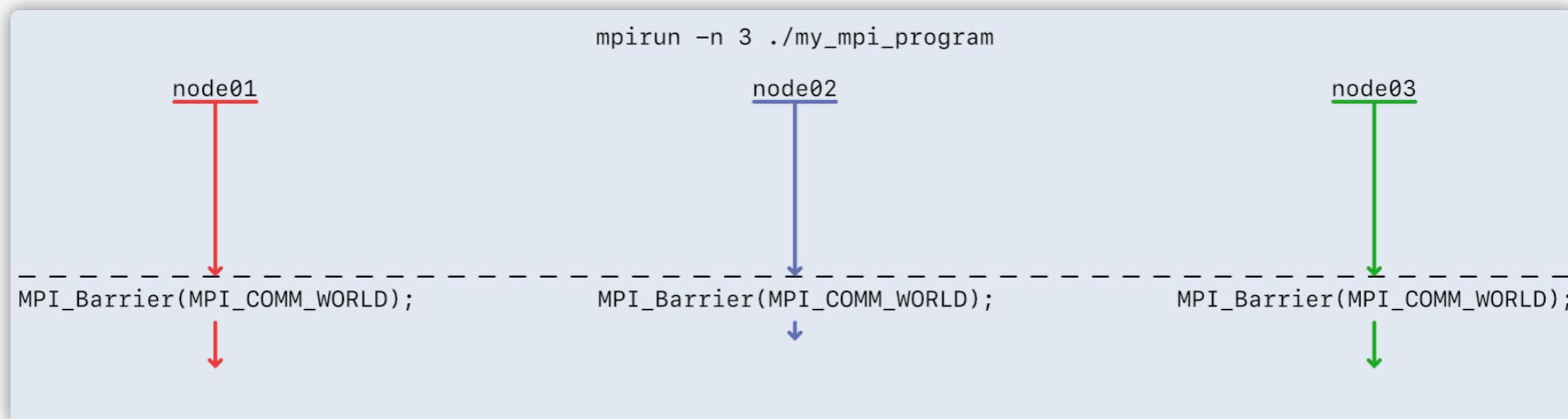


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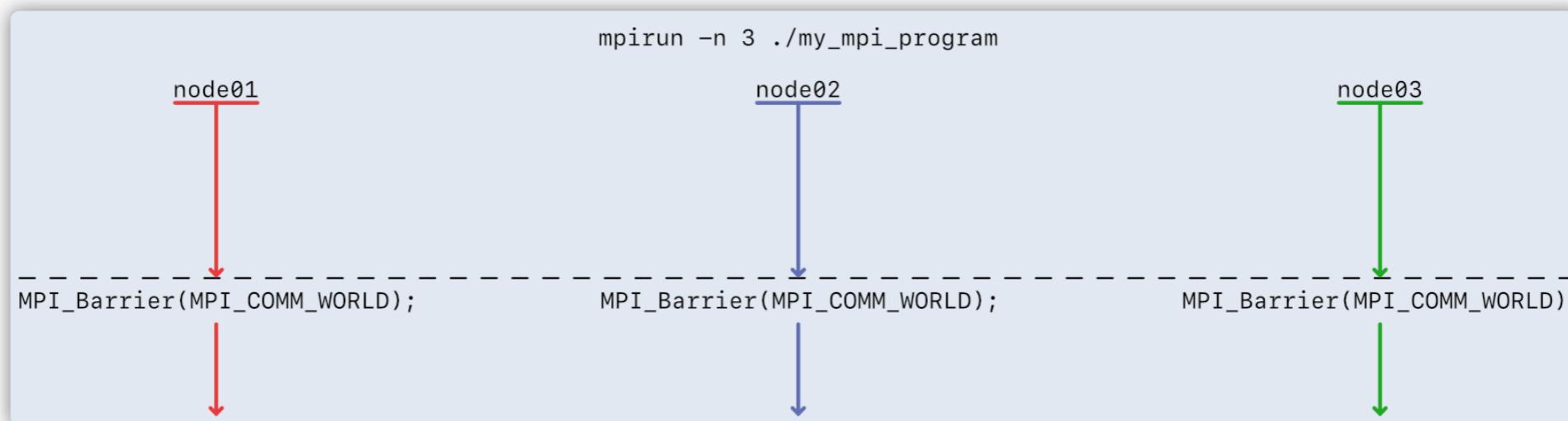


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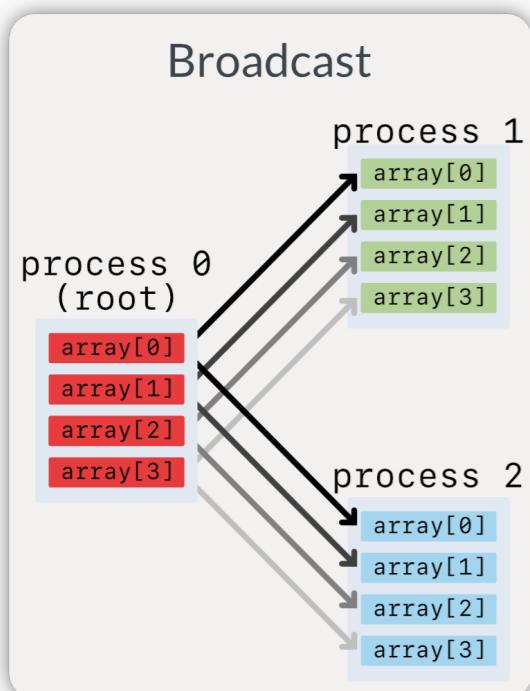


Collective operations

- The first set of communication functions we will look at are *collective operations*
- Collective: all processes must be involved in the operation (as opposed to *point-to-point* communications)
- Examples (this list is not exhaustive!):
 - Broadcast a variable from one process to all processes (Broadcast)
 - Distribute elements of an array on one process to multiple processes (Scatter)
 - Collect elements of arrays scattered over processes into a single process (Gather)
 - Sum a variable over all processes (Reduction)

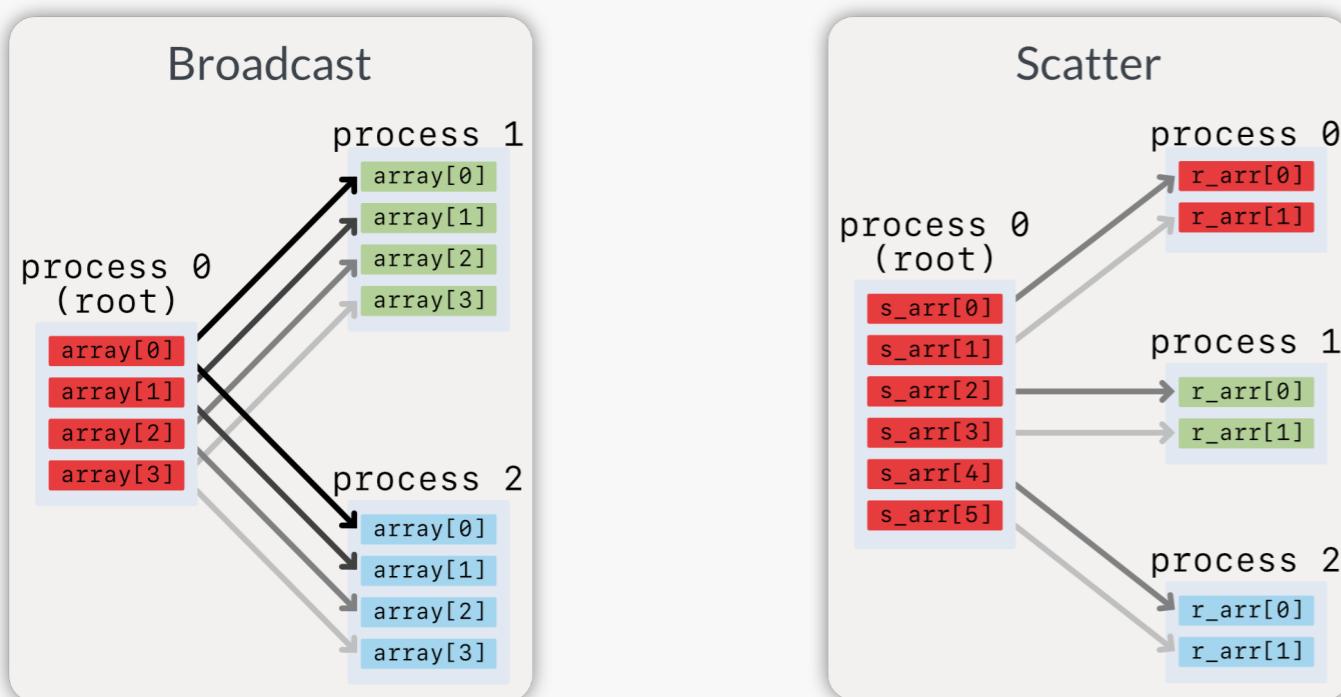
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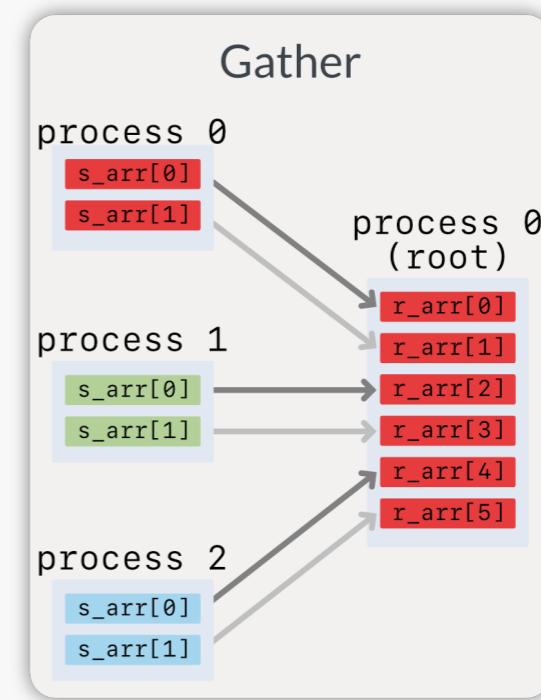
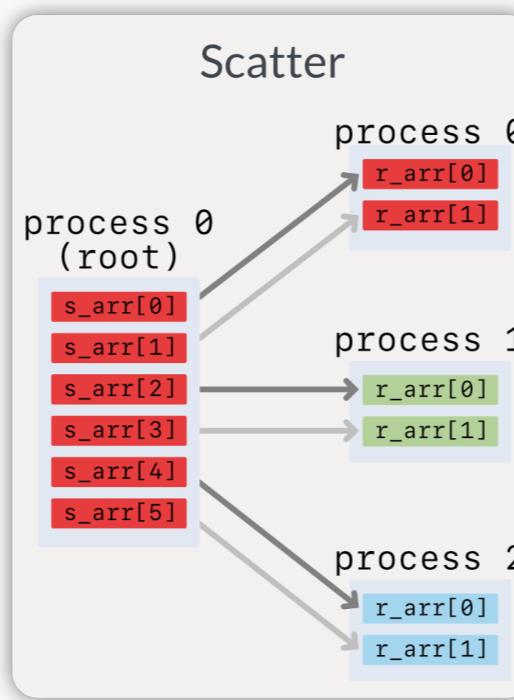
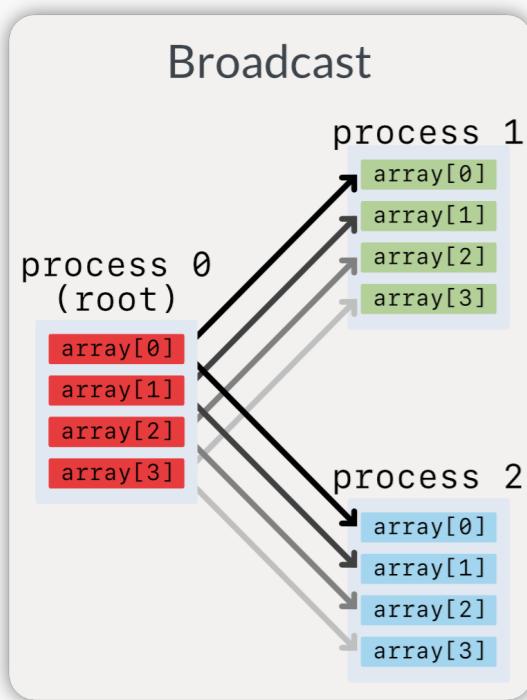
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Collective operations: Broadcast

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- The `MPI_Datatype` is important since MPI uses it to estimate the size in bytes that need to be transferred
- Full list of types available in MPI documentation. E.g. see: <https://www.mpich.org/static/docs/latest/www3/Constants.html>

Collective operations: Scatter

- Scatter:

```
MPI_Scatter(  
    const void *sendbuf, int sendcount, MPI_Datatype sendtype,  
    void *recvbuf, int recvcount, MPI_Datatype recvtype,  
    int root, MPI_Comm comm  
>);
```

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```

- `sendcount` is the number of elements to be sent to *each* process
- `sendbuf` is only relevant in the root process
- *Example:* distribute a 12-element array from process 0, assuming 3 processes in total (including root)

```
double arr_all[12]; /* ← this only needs to be defined on process with rank = 0 */  
double arr_proc[4];  
MPI_Scatter(arr_all, 4, MPI_DOUBLE, arr_proc, 4, MPI_DOUBLE, 0, MPI_COMM_WORLD);
```

Collective operations: Scatter

- Scatter:

```
MPI_Scatter(  
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```

- *Example:* distribute each element of a 4-element array to 4 processes in total (including root)

```
double arr[4]; /* ← this only needs to be defined on process with rank = 0 */  
double element;  
MPI_Scatter(arr, 1, MPI_DOUBLE, &element, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
```

Collective operations: Gather

- Gather:

```
MPI_Gather(  
    const void *sendbuf, int sendcount, MPI_Datatype sendtype,  
    void *recvbuf, int recvcount, MPI_Datatype recvtype, int root,  
    MPI_Comm comm  
)
```

Collective operations: Gather

- Gather:

```
MPI_Gather(  
    const void *sendbuf, int sendcount, MPI_Datatype sendtype,  
    void *recvbuf, int recvcount, MPI_Datatype recvtype, int root,  
    MPI_Comm comm  
)
```

- `recvcount` is the number of elements to be received by *each* process

Collective operations: Gather

- Gather:

```
MPI_Gather(  
    const void *sendbuf, int sendcount, MPI_Datatype sendtype,  
    void *recvbuf, int recvcount, MPI_Datatype recvtype, int root,  
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)
```

- `recvcount` is the number of elements to be received by *each* process
- `recvbuf` is only relevant in the root process

Collective operations: Gather

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```
MPI_Gather(  
    const void *sendbuf, int sendcount, MPI_Datatype sendtype,  
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)
```

- `recvcount` is the number of elements to be received by *each* process
- `recvbuf` is only relevant in the root process
- *Example:* collect a 9-element array at process 0, by concatenating 3 elements from each of 3 processes in total (including root)

```
double arr_all[9]; /* ← this only needs to be defined on process with rank = 0 */  
double arr_proc[3];  
MPI_Gather(arr_proc, 3, MPI_DOUBLE, arr_all, 3, MPI_DOUBLE, 0, MPI_COMM_WORLD);
```

Collective operations: Gather

- Gather:

```
MPI_Gather(  
    const void *sendbuf, int sendcount, MPI_Datatype sendtype,  
    void *recvbuf, int recvcount, MPI_Datatype recvtype, int root,  
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double arr_all[9]; /* ← this only needs to be defined on process with rank = 0 */  
double arr_proc[3];  
MPI_Gather(arr_proc, 3, MPI_DOUBLE, arr_all, 3, MPI_DOUBLE, 0, MPI_COMM_WORLD);
```

- *Example*: collect a 4-element array at process 0, by concatenating an element from each of 4 processes in total (including root)

```
double arr[4]; /* ← this only needs to be defined on process with rank = 0 */  
double element;  
MPI_Gather(&element, 1, MPI_DOUBLE, arr, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
```

Collective operations: Reduction

- Reduction:

```
MPI_Reduce(  
    const void *sendbuf, void *recvbuf, int count,  
    MPI_Datatype datatype, MPI_Op op, int root,  
    MPI_Comm comm  
)
```

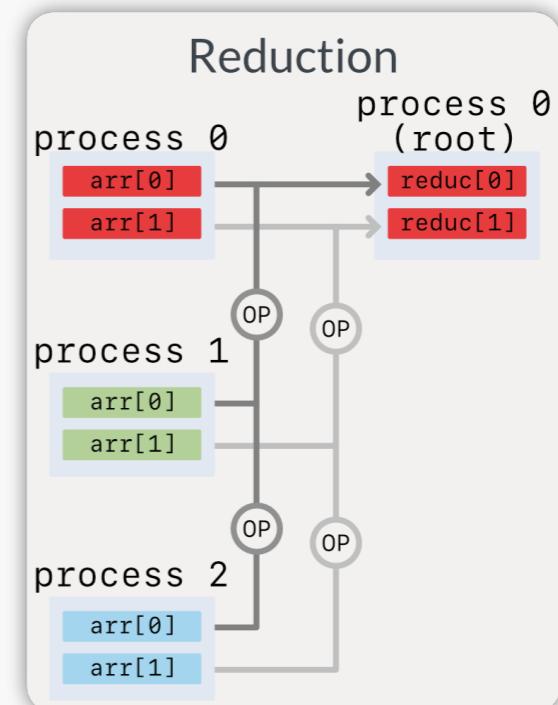
Collective operations: Reduction

- Reduction:

```
MPI_Reduce(  
    const void *sendbuf, void *recvbuf, int count,  
    MPI_Datatype datatype, MPI_Op op, int root,  
    MPI_Comm comm  
)
```

- Notes:

- `MPI_Op` is an operation, e.g. `MPI_SUM`, `MPI_PROD`, etc.
- The correct result of the operation depends on specifying the datatype correctly
- `count` is the number of elements of the arrays and is the same for send and receive, e.g. in the example on the right, `count = 2`
- The operation is over all processes in `comm`



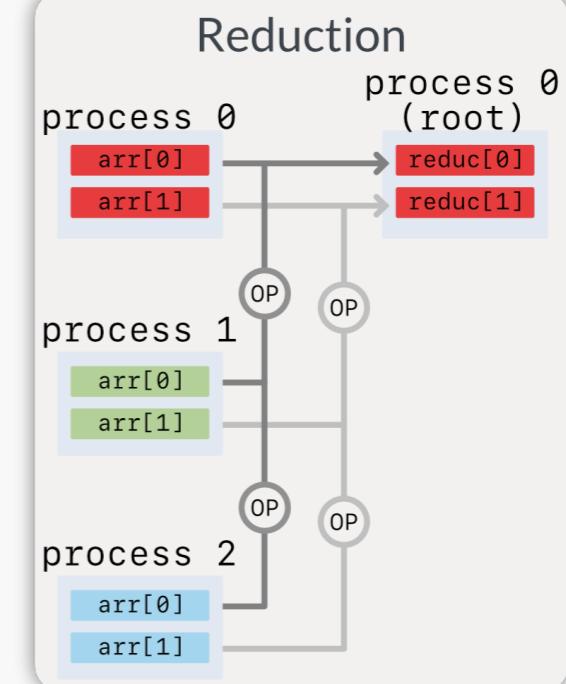
Collective operations: Reduction

- Reduction:

```
MPI_Reduce(  
    const void *sendbuf, void *recvbuf, int count,  
    MPI_Datatype datatype, MPI_Op op, int root,  
    MPI_Comm comm  
)
```

- Example:* Sum each element of a 3-element array over all processes

```
double s_arr[3];  
double r_arr[3]; /* ← only needs to      *  
                  * be defined on root */  
MPI_Reduce(s_arr, r_arr, 3, MPI_DOUBLE,  
           MPI_SUM, 0, MPI_COMM_WORLD);
```



Collective operations: Reduction

- Reduction:

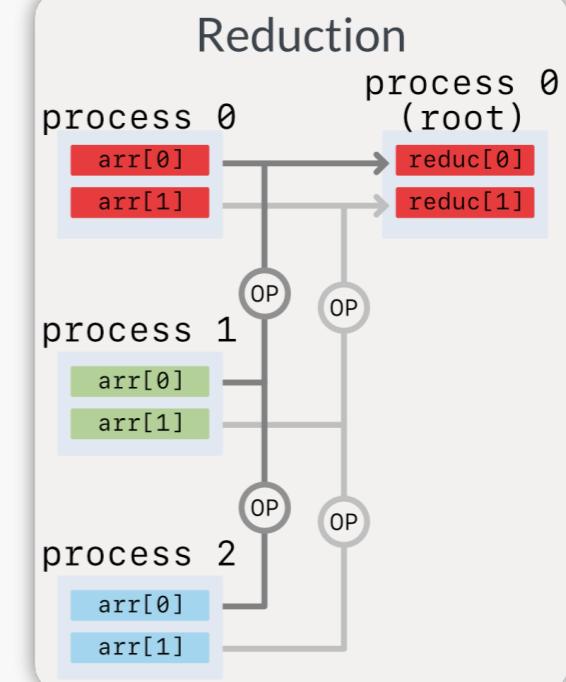
```
MPI_Reduce(  
    const void *sendbuf, void *recvbuf, int count,  
    MPI_Datatype datatype, MPI_Op op, int root,  
    MPI_Comm comm  
)
```

- Example:* Sum each element of a 3-element array over all processes

```
double s_arr[3];  
double r_arr[3]; /* ← only needs to      *  
                   *      be defined on root */  
MPI_Reduce(s_arr, r_arr, 3, MPI_DOUBLE,  
           MPI_SUM, 0, MPI_COMM_WORLD);
```

- Example:* Sum variable `var` over all processes

```
double var;  
double sum; /* ← only needs to      *  
               *      be defined on root */  
MPI_Reduce(&var, &sum, 1, MPI_DOUBLE,  
           MPI_SUM, 0, MPI_COMM_WORLD);
```



Collective operations

Some additional notes on variants of the collectives we have covered

- `MPI_Scatterv()` and `MPI_Gatherv()`
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→ which buffer depends on the specific MPI function
 - Instructs MPI to use the **same** buffer for receive and send
 - E.g. below, the sum will be placed in `var` of the root process (process with `rank = 0`):

```
if(rank != 0) {
    MPI_Reduce(&var, null, 1, MPI_DOUBLE, MPI_SUM, 0, MPI_COMM_WORLD);
} else {
    MPI_Reduce(MPI_IN_PLACE, &var, 1, MPI_DOUBLE, MPI_SUM, 0, MPI_COMM_WORLD);
}
```

Point-to-point communication

- Communications that involve transfer of data between two processes

Point-to-point communication

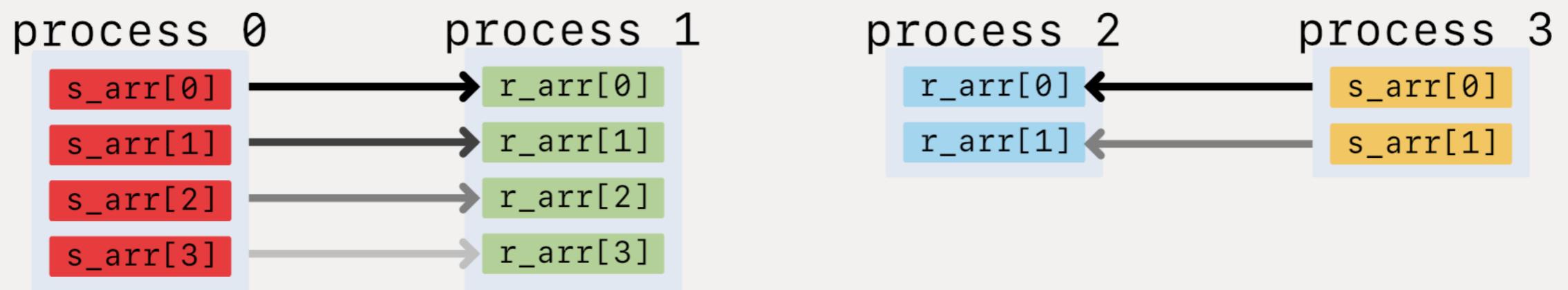
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 - Posting a receive which does not have a matching send

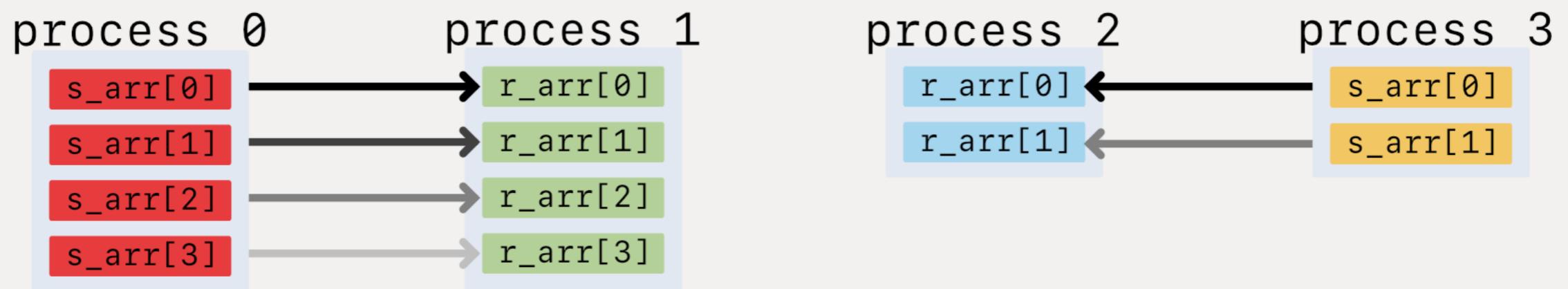
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Two point-to-point communications are depicted above
↳ between i) process 0 and 1 and between ii) process 2 and 3

Point-to-point communication

- Send/Receive

```
MPI_Send(void *buf, int n, MPI_Datatype type, int dest, int tag, MPI_Comm comm)
MPI_Recv(void *buf, int n, MPI_Datatype type, int srce, int tag, MPI_Comm comm, MPI_Status *status)
```

Point-to-point communication

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- Use of `MPI_ANY_SOURCE` in `MPI_Recv()` means "accept data from any source"
- `status` can be used to query the result of the receive (e.g. how many elements were received). We will use `MPI_STATUS_IGNORE` in place of `status`, which ignores the status

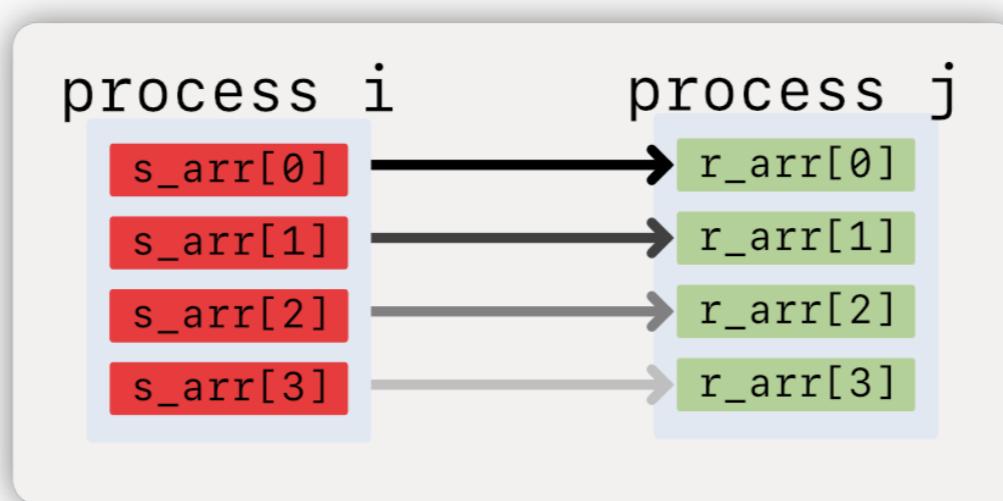
Point-to-point communication

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Point-to-point communication

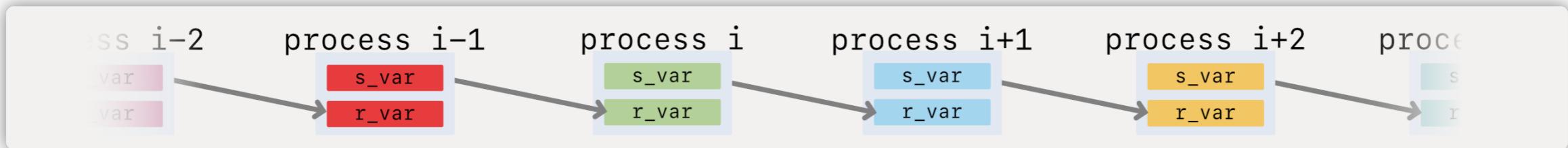
- Send/Receive; a trivial example



```
if(rank == i) {  
    MPI_Send(s_arr, 4, MPI_DOUBLE, j, 0, MPI_COMM_WORLD);  
}  
if(rank == j) {  
    MPI_Recv(r_arr, 4, MPI_DOUBLE, i, MPI_ANY_TAG, MPI_COMM_WORLD, MPI_STATUS_IGNORE);  
}
```

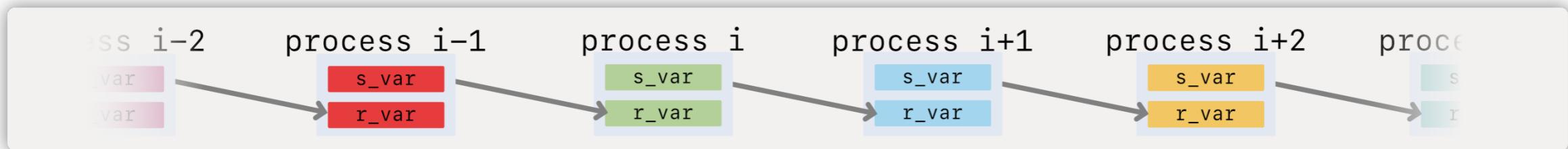
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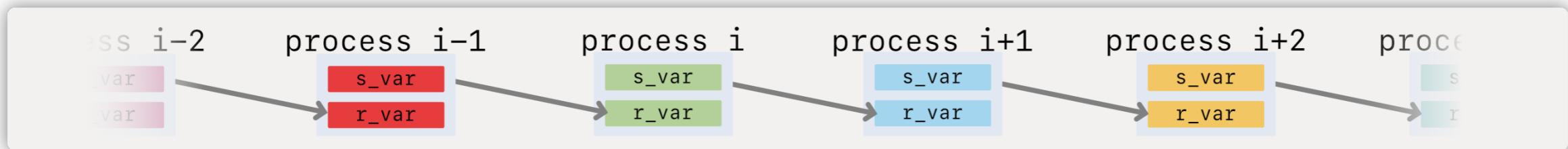


- This will **not** work:

```
MPI_Send(&s_var, 1, MPI_DOUBLE, rank+1, 0, MPI_COMM_WORLD);
MPI_Recv(&r_var, 1, MPI_DOUBLE, rank-1, MPI_ANY_TAG, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
```

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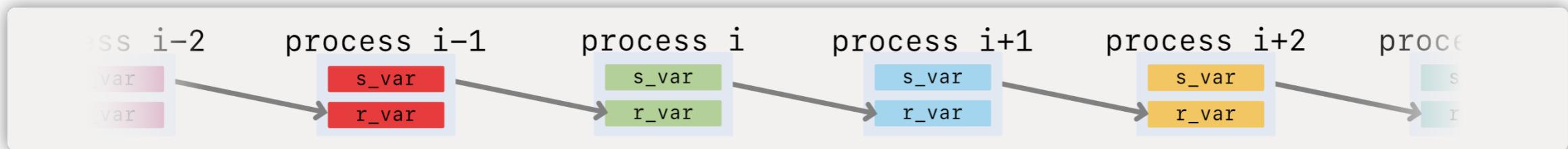
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- It results in a **deadlock**:
 - an `MPI_Recv()` can only be posted once an `MPI_Send()` completes
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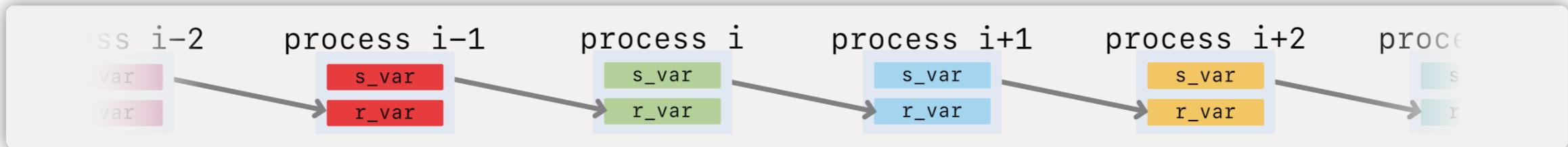
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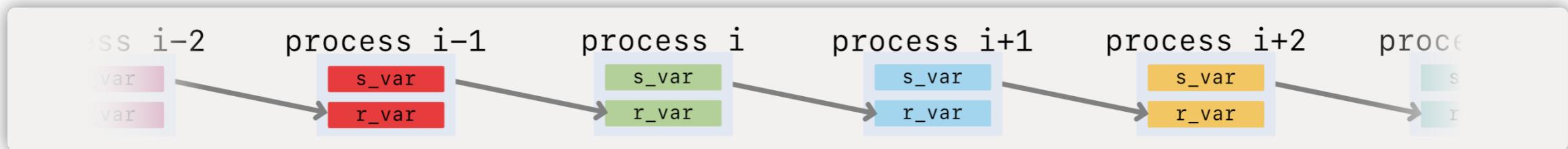
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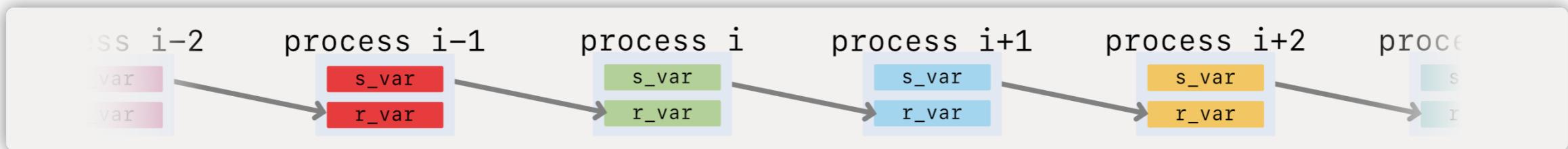
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 - Serializes communications that may be done faster in parallel
 - Inelegant, obscure, and error-prone

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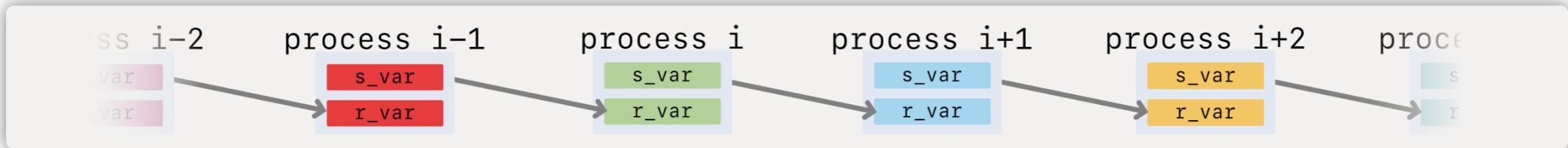


- A more efficient and elegant solution is to use `MPI_Sendrecv()`:

```
MPI_Sendrecv(void *sendbuf, int sendcount, MPI_Datatype sendtype, int dest, int sendtag,
            void *recvbuf, int recvcount, MPI_Datatype recvtype, int srce, int recvtag,
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```

- For the depicted example:

```
MPI_Sendrecv(&s_var, 1, MPI_DOUBLE, rank+1, 0,
             &r_var, 1, MPI_DOUBLE, rank-1, MPI_ANY_TAG,
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 - The function `MPI_Wait()` is used to block until the operation has complete

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MPI_Isend(sendbuf, ... , request);
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- `MPI_Sendrecv_replace()`
 - Like `MPI_Sendrecv()` but with a single `buf` rather than separate `sendbuf` and `recvbuf`
 - ↳ The receive message overwrites the send message

Exercises

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- Note that if you have modified `ex${n}.c` correctly, the job should complete in **less than one minute**

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 - ex05: Use of `MPI_Scatter()` and `MPI_Gather()`
 - ex06: Use of `MPI_Scatter()` and `MPI_Reduce()`
 - ex07: Use of `MPI_Send()` and `MPI_Recv()`
 - ex08: Use of `MPI_Sendrecv()`

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 - ex06: Use of `MPI_Scatter()` and `MPI_Reduce()`
 - ex07: Use of `MPI_Send()` and `MPI_Recv()`
 - ex08: Use of `MPI_Sendrecv()`
- All exercises have been tested with OpenMPI and the GNU Compiler. Please use:

```
module load gompi
```

for all exercises.

Exercises

Ex01

- Modify `ex01.c` to call `MPI_Comm_size()` and `MPI_Comm_rank()` with the appropriate arguments

```
int nproc, rank;  
/*  
 * TODO: call `MPI_Comm_size()` and `MPI_Comm_rank()` with the  
 * appropriate arguments  
 */  
MPI_Comm_size(/* TODO */);  
MPI_Comm_rank(/* TODO */);
```

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Ex01

- Modify `ex01.c` to call `MPI_Comm_size()` and `MPI_Comm_rank()` with the appropriate arguments

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MPI_Comm_rank(/* TODO */);
```

- To compile, type `make`. Remember to first load the appropriate module (`gompi`):

```
[user@front01 ex01]$ module load gompi  
[user@front01 ex01]$ make  
mpicc -c ex01.c  
mpicc -o ex01 ex01.o  
[user@front01 ex01]$
```

Exercises

Ex01

- A job script has been prepared to run ex01:

```
[user@front01 ex01]$ cat sub-01.sh
#!/bin/bash
#SBATCH --job-name=01
#SBATCH --nodes=2
#SBATCH --ntasks=8
#SBATCH --output=ex01-output.txt
#SBATCH --time=00:01:00

module load gompi
mpirun ./ex01
```

Exercises

Ex01

- A job script has been prepared to run ex01:

```
[user@front01 ex01]$ cat sub-01.sh
#!/bin/bash
#SBATCH --job-name=01
#SBATCH --nodes=2
#SBATCH --ntasks=8
#SBATCH --output=ex01-output.txt
#SBATCH --time=00:01:00

module load gompi
mpirun ./ex01
```

- 2 nodes, 8 processes, meaning 4 processes per node

Exercises

Ex01

- A job script has been prepared to run `ex01`:

```
[user@front01 ex01]$ cat sub-01.sh
#!/bin/bash
#SBATCH --job-name=01
#SBATCH --nodes=2
#SBATCH --ntasks=8
#SBATCH --output=ex01-output.txt
#SBATCH --time=00:01:00

module load gompi
mpirun ./ex01
```

- 2 nodes, 8 processes, meaning 4 processes per node
- program output will be redirected to file `ex01-output.txt`

Exercises

Ex01

- A job script has been prepared to run `ex01`:

```
[user@front01 ex01]$ cat sub-01.sh
#!/bin/bash
#SBATCH --job-name=01
#SBATCH --nodes=2
#SBATCH --ntasks=8
#SBATCH --output=ex01-output.txt
#SBATCH --time=00:01:00

module load gompi
mpirun ./ex01
```

- 2 nodes, 8 processes, meaning 4 processes per node
- program output will be redirected to file `ex01-output.txt`
- requests 1 minute. If not done by then, the scheduler will kill the job

Exercises

Ex01

- Submit the job script:

```
[user@front01 ex01]$ sbatch sub-01.sh
Submitted batch job 69711
[user@front01 ex01]$
```

Exercises

Ex01

- Submit the job script:

```
[user@front01 ex01]$ sbatch sub-01.sh
Submitted batch job 69711
[user@front01 ex01]$
```

- Check status of job:

```
[user@front01 ex01]$ squeue -u $USER
JOBID PARTITION      NAME      USER ST      TIME   NODES NODELIST(REASON)
69712      cpu        01      user    R      0:00      2 cn[01-02]
[user@front01 ex01]$
```

Exercises

Ex01

- Submit the job script:

```
[user@front01 ex01]$ sbatch sub-01.sh
Submitted batch job 69711
[user@front01 ex01]$
```

- Check status of job:

```
[user@front01 ex01]$ squeue -u $USER
JOBID PARTITION      NAME      USER ST      TIME   NODES NODELIST(REASON)
69712      cpu        01      user    R      0:00      2 cn[01-02]
[user@front01 ex01]$
```

- The job runs very quickly. You may see no output above if the job has finished:

```
[user@front01 ex01]$ squeue -u $USER
JOBID PARTITION      NAME      USER ST      TIME   NODES NODELIST(REASON)
[user@front01 ex01]$
```

Exercises

Ex01

- If done, the file `ex01-output.txt` should have been created
- Inspect the file:

```
[user@front01 ex01]$ cat ex01-output.txt
This is rank = 3 of nproc = 8 on node: cn01
This is rank = 0 of nproc = 8 on node: cn01
This is rank = 1 of nproc = 8 on node: cn01
This is rank = 2 of nproc = 8 on node: cn01
This is rank = 7 of nproc = 8 on node: cn02
This is rank = 5 of nproc = 8 on node: cn02
This is rank = 6 of nproc = 8 on node: cn02
This is rank = 4 of nproc = 8 on node: cn02
[user@front01 ex01]$
```

Exercises

Ex01

- If done, the file `ex01-output.txt` should have been created
- Inspect the file:

```
[user@front01 ex01]$ cat ex01-output.txt
This is rank = 3 of nproc = 8 on node: cn01
This is rank = 0 of nproc = 8 on node: cn01
This is rank = 1 of nproc = 8 on node: cn01
This is rank = 2 of nproc = 8 on node: cn01
This is rank = 7 of nproc = 8 on node: cn02
This is rank = 5 of nproc = 8 on node: cn02
This is rank = 6 of nproc = 8 on node: cn02
This is rank = 4 of nproc = 8 on node: cn02
[user@front01 ex01]$
```

- Note the order is nondeterministic; whichever process reaches the print statement first prints

Exercises

Ex01

- If done, the file `ex01-output.txt` should have been created
- Inspect the file:

```
[user@front01 ex01]$ cat ex01-output.txt
This is rank = 3 of nproc = 8 on node: cn01
This is rank = 0 of nproc = 8 on node: cn01
This is rank = 1 of nproc = 8 on node: cn01
This is rank = 2 of nproc = 8 on node: cn01
This is rank = 7 of nproc = 8 on node: cn02
This is rank = 5 of nproc = 8 on node: cn02
This is rank = 6 of nproc = 8 on node: cn02
This is rank = 4 of nproc = 8 on node: cn02
[user@front01 ex01]$
```

- Note the order is nondeterministic; whichever process reaches the print statement first prints
- We can use synchronization to serialize the print statements and ensure the correct order

Exercises

Ex02

- `ex02.c` is similar to `ex01.c`
- A for-loop is included over the print statement:

```
/*
 * TODO: add an MPI_Barrier() to ensure the ranks print in-order
 */
for(int i=0; i<nproc; i++) {
    if(rank == i)
        printf(" This is rank = %d of nproc = %d on node: %s\n", rank, nproc, hname);
}
```

Exercises

Ex02

- `ex02.c` is similar to `ex01.c`
- A for-loop is included over the print statement:

```
/*
 * TODO: add an MPI_Barrier() to ensure the ranks print in-order
 */
for(int i=0; i<nproc; i++) {
    if(rank == i)
        printf(" This is rank = %d of nproc = %d on node: %s\n", rank, nproc, hname);
}
```

- Study the code and carefully place an `MPI_Barrier()` so that the print statements will be executed in rank ordered

Exercises

Ex02

- `ex02.c` is similar to `ex01.c`
- A for-loop is included over the print statement:

```
/*
 * TODO: add an MPI_Barrier() to ensure the ranks print in-order
 */
for(int i=0; i<nproc; i++) {
    if(rank == i)
        printf(" This is rank = %d of nproc = %d on node: %s\n", rank, nproc, hname);
}
```

- Study the code and carefully place an `MPI_Barrier()` so that the print statements will be executed in rank ordered
- As before, when done:
 - use `make` to compile and
 - `sbatch sub-02.sh` to submit the prepared job script

Exercises

Ex02

- Inspect the file `ex02-output.txt`:

```
[user@front01 ex02]$ cat ex02-output.txt
This is rank = 0 of nproc = 8 on node: cn01
This is rank = 1 of nproc = 8 on node: cn01
This is rank = 2 of nproc = 8 on node: cn01
This is rank = 3 of nproc = 8 on node: cn01
This is rank = 4 of nproc = 8 on node: cn02
This is rank = 5 of nproc = 8 on node: cn02
This is rank = 6 of nproc = 8 on node: cn02
This is rank = 7 of nproc = 8 on node: cn02
[user@front01 ex02]$
```

Exercises

Ex02

- Inspect the file `ex02-output.txt`:

```
[user@front01 ex02]$ cat ex02-output.txt
This is rank = 0 of nproc = 8 on node: cn01
This is rank = 1 of nproc = 8 on node: cn01
This is rank = 2 of nproc = 8 on node: cn01
This is rank = 3 of nproc = 8 on node: cn01
This is rank = 4 of nproc = 8 on node: cn02
This is rank = 5 of nproc = 8 on node: cn02
This is rank = 6 of nproc = 8 on node: cn02
This is rank = 7 of nproc = 8 on node: cn02
[user@front01 ex02]$
```

- The print statements should appear in rank order

Exercises

Ex03

- This exercise demonstrates `MPI_Bcast()`
- A file `data.txt` is included:

```
[user@front01 ex03]$ cat data.txt  
3.14159265359  
[user@front01 ex03]$
```

Exercises

Ex03

- This exercise demonstrates `MPI_Bcast()`
- A file `data.txt` is included:

```
[user@front01 ex03]$ cat data.txt  
3.14159265359  
[user@front01 ex03]$
```

- In `ex03.c`, the root process (process with `rank = 0`) calls the `readline()` function to read the single line from the file

Exercises

Ex03

- This exercise demonstrates `MPI_Bcast()`
- A file `data.txt` is included:

```
[user@front01 ex03]$ cat data.txt  
3.14159265359  
[user@front01 ex03]$
```

- In `ex03.c`, the root process (process with `rank = 0`) calls the `readline()` function to read the single line from the file
- Your task is to use a `MPI_Bcast()` to broadcast the variable `readline()` to all processes:

```
/*  
 * TODO: add an MPI_Bcast() with the appropriate arguments to  
 * broadcast variable `val' from the root process to all processes  
 */  
MPI_Bcast(/* TODO */);
```

Exercises

Ex03

- If done correctly, `ex03-output.txt` should include the following output:

```
[user@front01 ex03]$ cat ex03-output.txt
This is rank = 7 of nproc = 8 on node: cn02 | Got from root var = 3.141593
This is rank = 6 of nproc = 8 on node: cn02 | Got from root var = 3.141593
This is rank = 5 of nproc = 8 on node: cn02 | Got from root var = 3.141593
This is rank = 4 of nproc = 8 on node: cn02 | Got from root var = 3.141593
This is rank = 3 of nproc = 8 on node: cn01 | Got from root var = 3.141593
This is rank = 0 of nproc = 8 on node: cn01 | Got from root var = 3.141593
This is rank = 2 of nproc = 8 on node: cn01 | Got from root var = 3.141593
This is rank = 1 of nproc = 8 on node: cn01 | Got from root var = 3.141593
[user@front01 ex03]$
```

With the order being undetermined also in this case

Exercises

Ex04

- This exercise demonstrates the scatter operation
- The file `data.txt` now includes eight lines:

```
[user@front01 ex04]$ cat data.txt  
0.42816572487  
0.57721566490  
0.66274341935  
0.69314718056  
1.41421356237  
1.61803398875  
2.71828182846  
3.14159265359  
[user@front01 ex04]$
```

Exercises

Ex04

- This exercise demonstrates the scatter operation
- The file `data.txt` now includes eight lines:

```
[user@front01 ex04]$ cat data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
[user@front01 ex04]$
```

- The root process reads the eight lines into the eight-element double precision array `vals[]` using the function `readlines()`

```
/*
 * root process: read `n` lines of "data.txt" into array `vars[]`
 */
int nelems = 8;
double vars[nelems];
if(rank == 0) {
    char fname[] = "data.txt";
    readlines(nelems, vars, fname);
}
```

Exercises

Ex04

- Your task is to scatter the array `vals[]` so that each of the eight processes receives one element of the array into variable `var`:

```
double var;
/*
 * TODO: use an MPI_Scatter() to distribute the elements of `vars[]'
 * to the processes, one element for each process. Assume the number
 * of processes is the same as the number of elements.
 */
MPI_Scatter(/* TODO */);
```

Exercises

Ex04

- Your task is to scatter the array `vals[]` so that each of the eight processes receives one element of the array into variable `var`:

```
double var;
/*
 * TODO: use an MPI_Scatter() to distribute the elements of `vars[]'
 * to the processes, one element for each process. Assume the number
 * of processes is the same as the number of elements.
 */
MPI_Scatter(/* TODO */);
```

- Once done, compile (`make`) and submit the job script (`sbatch sub-04.sh`)

Exercises

Ex04

- Your task is to scatter the array `vals[]` so that each of the eight processes receives one element of the array into variable `var`:

```
double var;
/*
 * TODO: use an MPI_Scatter() to distribute the elements of `vars[]'
 * to the processes, one element for each process. Assume the number
 * of processes is the same as the number of elements.
 */
MPI_Scatter(/* TODO */);
```

- Once done, compile (`make`) and submit the job script (`sbatch sub-04.sh`)
- If done correctly, you should see the following in `ex04-output.txt`:

```
[user@front01 ex04]$ cat ex04-output.txt
This is rank = 2 of nproc = 8 on node: cn01 | Got from root var = 0.662743
This is rank = 3 of nproc = 8 on node: cn01 | Got from root var = 0.693147
This is rank = 1 of nproc = 8 on node: cn01 | Got from root var = 0.577216
This is rank = 4 of nproc = 8 on node: cn02 | Got from root var = 1.414214
This is rank = 5 of nproc = 8 on node: cn02 | Got from root var = 1.618034
This is rank = 6 of nproc = 8 on node: cn02 | Got from root var = 2.718282
This is rank = 0 of nproc = 8 on node: cn01 | Got from root var = 0.428166
This is rank = 7 of nproc = 8 on node: cn02 | Got from root var = 3.141593
[user@front01 ex04]$
```

Exercises

Ex04

- Your task is to scatter the array `vals[]` so that each of the eight processes receives one element of the array into variable `var`:

```
double var;
/*
 * TODO: use an MPI_Scatter() to distribute the elements of `vars[]'
 * to the processes, one element for each process. Assume the number
 * of processes is the same as the number of elements.
 */
MPI_Scatter(/* TODO */);
```

- Once done, compile (`make`) and submit the job script (`sbatch sub-04.sh`)
- If done correctly, you should see the following in `ex04-output.txt`:

```
[user@front01 ex04]$ cat ex04-output.txt
This is rank = 2 of nproc = 8 on node: cn01 | Got from root var = 0.662743
This is rank = 3 of nproc = 8 on node: cn01 | Got from root var = 0.693147
This is rank = 1 of nproc = 8 on node: cn01 | Got from root var = 0.577216
This is rank = 4 of nproc = 8 on node: cn02 | Got from root var = 1.414214
This is rank = 5 of nproc = 8 on node: cn02 | Got from root var = 1.618034
This is rank = 6 of nproc = 8 on node: cn02 | Got from root var = 2.718282
This is rank = 0 of nproc = 8 on node: cn01 | Got from root var = 0.428166
This is rank = 7 of nproc = 8 on node: cn02 | Got from root var = 3.141593
[user@front01 ex04]$
```

- Pro tip:** you can sort the output by piping through `sort`, i.e. `cat ex04-output.txt | sort`

Exercises

Ex05

- This exercise demonstrates the gather operation
- The exercise starts like ex04:
 - The root process reads the eight values from `data.txt` and stores them in `vars[]`

Exercises

Ex05

- This exercise demonstrates the gather operation
- The exercise starts like ex04:
 - The root process reads the eight values from `data.txt` and stores them in `vars[]`
- We would like:

Exercises

Ex05

- This exercise demonstrates the gather operation
- The exercise starts like ex04:
 - The root process reads the eight values from `data.txt` and stores them in `vars[]`
- We would like:
 - The elements of `vars[]` to be scattered, one element to each of eight processes (same as ex04)

Exercises

Ex05

- This exercise demonstrates the gather operation
- The exercise starts like ex04:
 - The root process reads the eight values from `data.txt` and stores them in `vars[]`
- We would like:
 - The elements of `vars[]` to be scattered, one element to each of eight processes (same as ex04)
 - Each process to divide its element, stored in `var`, by two

Exercises

Ex05

- This exercise demonstrates the gather operation
- The exercise starts like ex04:
 - The root process reads the eight values from `data.txt` and stores them in `vars[]`
- We would like:
 - The elements of `vars[]` to be scattered, one element to each of eight processes (same as ex04)
 - Each process to divide its element, stored in `var`, by two
 - The process' `var` variables to be gathered back into `vars[]` of the root process

Exercises

Ex05

- This exercise demonstrates the gather operation
- The exercise starts like ex04:
 - The root process reads the eight values from `data.txt` and stores them in `vars[]`
- We would like:
 - The elements of `vars[]` to be scattered, one element to each of eight processes (same as ex04)
 - Each process to divide its element, stored in `var`, by two
 - The process' `var` variables to be gathered back into `vars[]` of the root process

```
/* TODO: use an MPI_Scatter() to distribute the elements of `vars[]'  
 * to the processes, one element for each process. Assume the number  
 * of processes is the same as the number of elements. Same as in  
 * exercise ex04.  
 */  
MPI_Scatter(/* TODO */);  
  
/* Divide by two on each rank */  
var = var*0.5;  
  
/* TODO: use an MPI_Gather() to collect `var' from each rank to the  
 * array `vars[]' on the root process  
 */  
MPI_Gather(/* TODO */);
```

Exercises

Ex05

- At the end, the root process prints all elements of `vars[]`

```
/*
 * root process: print the elements of `vars[]` obtained via the
 * `MPI_Gather()`
 */
if(rank == 0)
    for(int i=0; i<nlems; i++)
        printf(" vars[%d] = %lf\n", i, vars[i]);
```

Exercises

Ex05

- At the end, the root process prints all elements of `vars[]`

```
/*
 * root process: print the elements of `vars[]` obtained via the
 * `MPI_Gather()`
 */
if(rank == 0)
    for(int i=0; i<nlems; i++)
        printf(" vars[%d] = %lf\n", i, vars[i]);
```

- Inspect the output `ex05-output.txt` and ensure the result is correct:

```
[user@front01 ex05]$ cat ex05-output.txt
vars[0] = 0.214083
vars[1] = 0.288608
vars[2] = 0.331372
vars[3] = 0.346574
vars[4] = 0.707107
vars[5] = 0.809017
vars[6] = 1.359141
vars[7] = 1.570796
[user@front01 ex05]$
```

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)
- We would like:

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)
- We would like:
 - The root process to read all elements into an array `vars[]`

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)
- We would like:
 - The root process to read all elements into an array `vars[]`
 - The elements to be scattered to all processes
 - ↳ Each process should receive `nelems_loc = nelems / nproc` elements

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)
- We would like:
 - The root process to read all elements into an array `vars[]`
 - The elements to be scattered to all processes
 - ↳ Each process should receive `nelems_loc = nelems / nproc` elements
 - Each process to sum its local elements, storing the result into `sum_loc`

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)
- We would like:
 - The root process to read all elements into an array `vars[]`
 - The elements to be scattered to all processes
 - ↳ Each process should receive `nelems_loc = nelems / nproc` elements
 - Each process to sum its local elements, storing the result into `sum_loc`
 - To use a reduction operation to obtain the grand total over all 2520 elements

Exercises

Ex06

- This exercise demonstrates the reduction operation
- `data.txt` is now a file with 2520 random numbers, one per line (`nelems = 2520`)
- We would like:
 - The root process to read all elements into an array `vars[]`
 - The elements to be scattered to all processes
 - ↳ Each process should receive `nelems_loc = nelems / nproc` elements
 - Each process to sum its local elements, storing the result into `sum_loc`
 - To use a reduction operation to obtain the grand total over all 2520 elements
- Note that in `ex06.c` we explicitly check that the number of processes divides the number of elements in `data.txt`:

```
/*
 * Abort if the number of processes does not divide `nelems' exactly
 */
int nelems = 2520;
if(nelems % nproc != 0) {
    fprintf(stderr, " nelems = %d not divisible by nproc = %d\n", nelems, nproc);
    MPI_Abort(MPI_COMM_WORLD, 1);
}
```

Exercises

Ex06

- Your `TODOs` are to complete the scatter and reduction operations:

```
/*
 * TODO: use an MPI_Scatter() to distribute the elements of `vars[]'
 * to each process's `vars_loc[]` array
 */
MPI_Scatter(/* TODO */);

/*
 * `sum_loc` holds the sum over each process's local elements
 */
double sum_loc = 0;
for(int i=0; i<nelems_loc; i++)
    sum_loc += vars_loc[i];

/*
 * TODO: use an MPI_Reduce() to sum `sum_loc` over all processes and
 * store in `sum` of the root process
 */
double sum;
MPI_Reduce(/* TODO */);
```

Exercises

Ex06

- The root process prints the result at the end. If correct, you should see:

```
[user@front01 ex06]$ cat ex06-output.txt
Used 8 processes, sum = 1266.960662
[user@front01 ex06]$
```

Exercises

Ex06

- The root process prints the result at the end. If correct, you should see:

```
[user@front01 ex06]$ cat ex06-output.txt
Used 8 processes, sum = 1266.960662
[user@front01 ex06]$
```

- Note that `ex06.c` allows running with any number of processes which divide 2520 exactly

Exercises

Ex06

- The root process prints the result at the end. If correct, you should see:

```
[user@front01 ex06]$ cat ex06-output.txt
Used 8 processes, sum = 1266.960662
[user@front01 ex06]$
```

- Note that `ex06.c` allows running with any number of processes which divide 2520 exactly
- Try, for example, modifying `sub-06.sh` to use 40 processes (20 per node)

Exercises

Ex06

- The root process prints the result at the end. If correct, you should see:

```
[user@front01 ex06]$ cat ex06-output.txt  
Used 8 processes, sum = 1266.960662  
[user@front01 ex06]$
```

- Note that `ex06.c` allows running with any number of processes which divide 2520 exactly
- Try, for example, modifying `sub-06.sh` to use 40 processes (20 per node)
- You should see an identical sum:

```
[user@front01 ex06]$ cat ex06-output.txt  
Used 40 processes, sum = 1266.960662  
[user@front01 ex06]$
```

Exercises

Ex07

- This exercise demonstrates `MPI_Send()` and `MPI_Recv()`
- `data.txt` is the same as in `ex04`, with eight elements:

```
[user@front01 ex07]$ cat data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
[user@front01 ex07]$
```

Exercises

Ex07

- This exercise demonstrates `MPI_Send()` and `MPI_Recv()`
- `data.txt` is the same as in `ex04`, with eight elements:

```
[user@front01 ex07]$ cat data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
[user@front01 ex07]$
```

- All elements are read by the root process into array `vars[]`

Exercises

Ex07

- This exercise demonstrates `MPI_Send()` and `MPI_Recv()`
- `data.txt` is the same as in `ex04`, with eight elements:

```
[user@front01 ex07]$ cat data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
[user@front01 ex07]$
```

- All elements are read by the root process into array `vars[]`
- The elements are then scattered, one to each process, and stored in variable `var0`

Exercises

Ex07

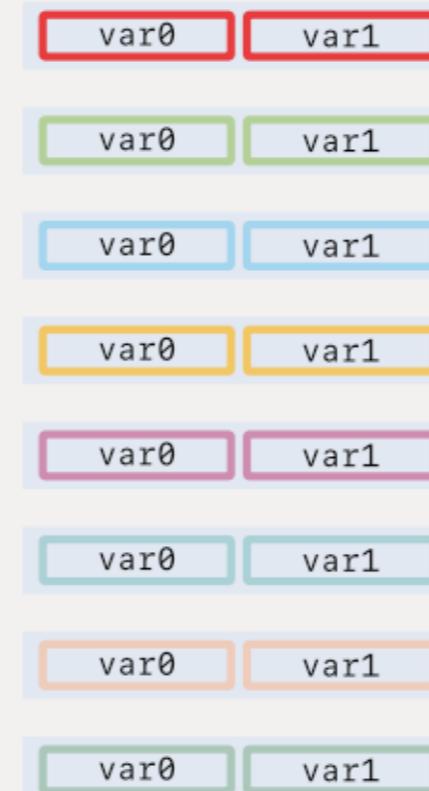
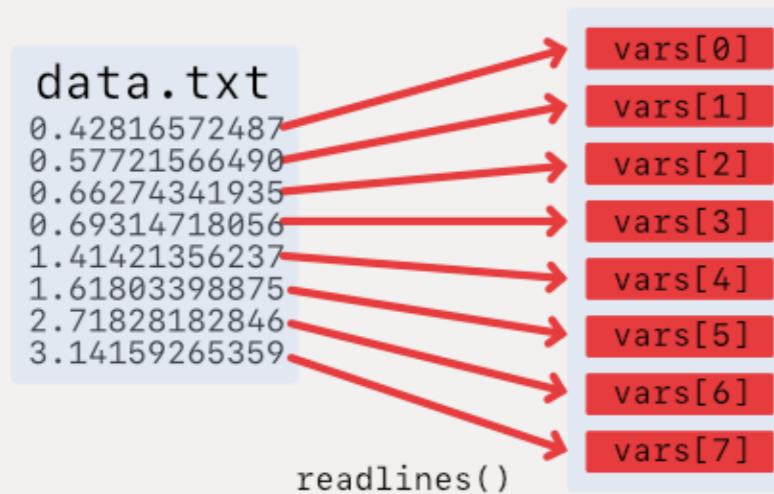
- This exercise demonstrates `MPI_Send()` and `MPI_Recv()`
- `data.txt` is the same as in `ex04`, with eight elements:

```
[user@front01 ex07]$ cat data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
[user@front01 ex07]$
```

- All elements are read by the root process into array `vars[]`
- The elements are then scattered, one to each process, and stored in variable `var0`
- Using `MPI_Send()` and `MPI_Recv()`, We would like that:
 - All processes with even ranks store in variable `var1` the value of `var0` corresponding to their next odd rank
 - All processes with odd ranks store in variable `var1` the value of `var0` corresponding to their previous even rank

Exercises

Ex07



Exercises

Ex07

```
data.txt  
0.42816572487  
0.57721566490  
0.66274341935  
0.69314718056  
1.41421356237  
1.61803398875  
2.71828182846  
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

var0	var1
------	------

var0	var1
------	------

var0	var1
------	------

var0	var1
------	------

var0	var1
------	------

var0	var1
------	------

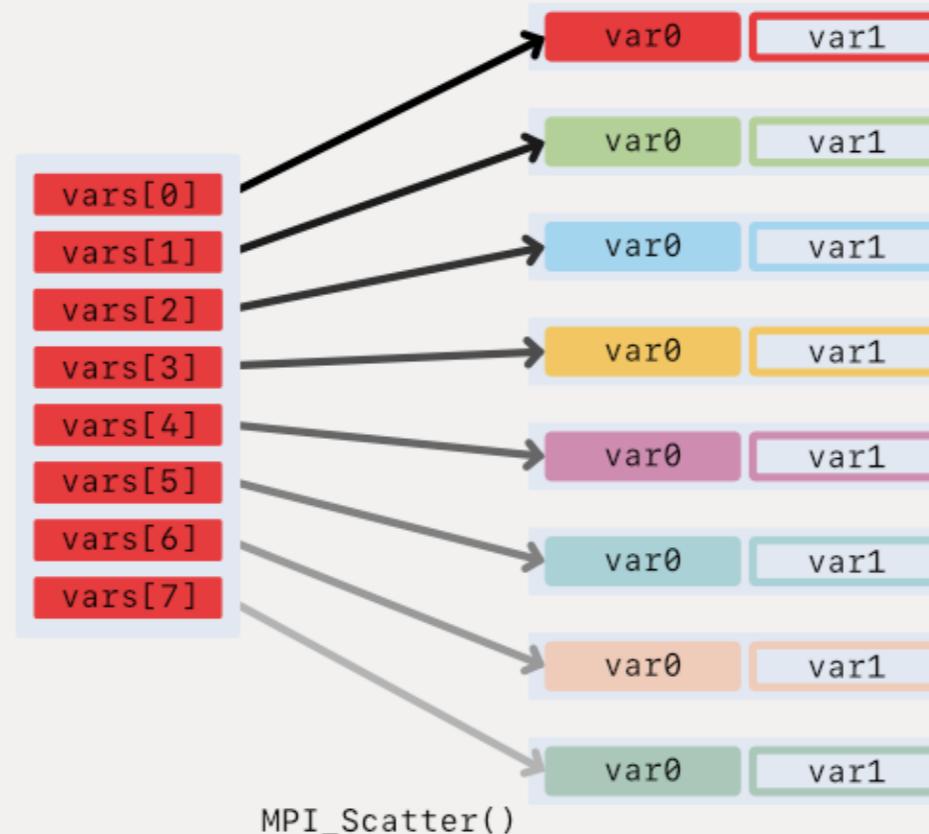
var0	var1
------	------

var0	var1
------	------

Exercises

Ex07

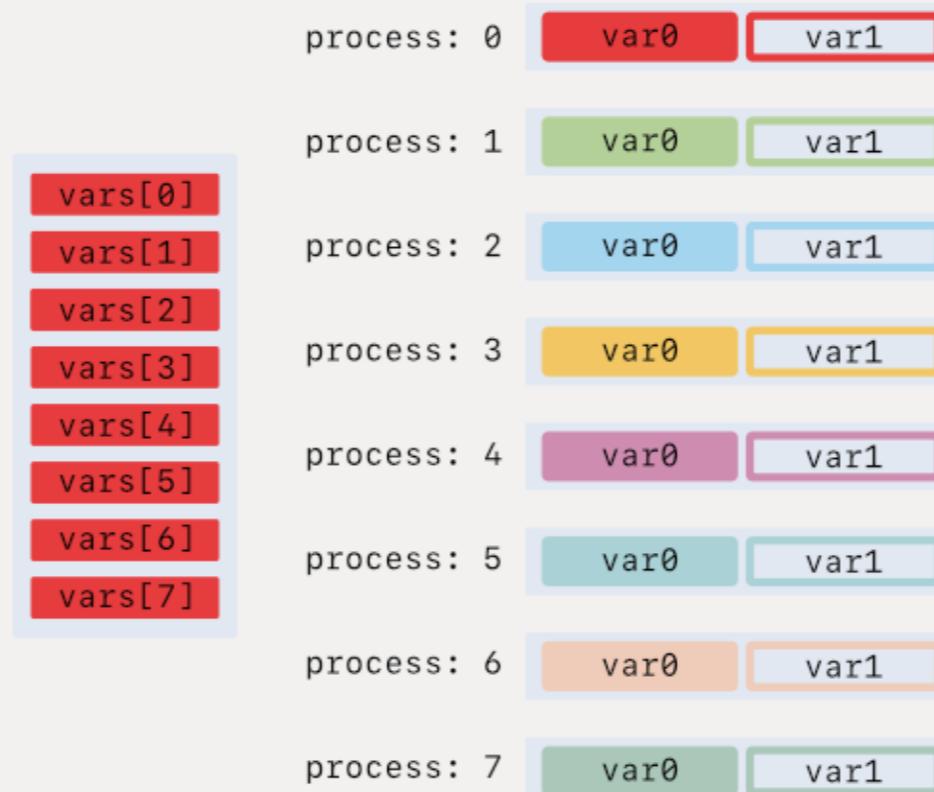
```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```



Exercises

Ex07

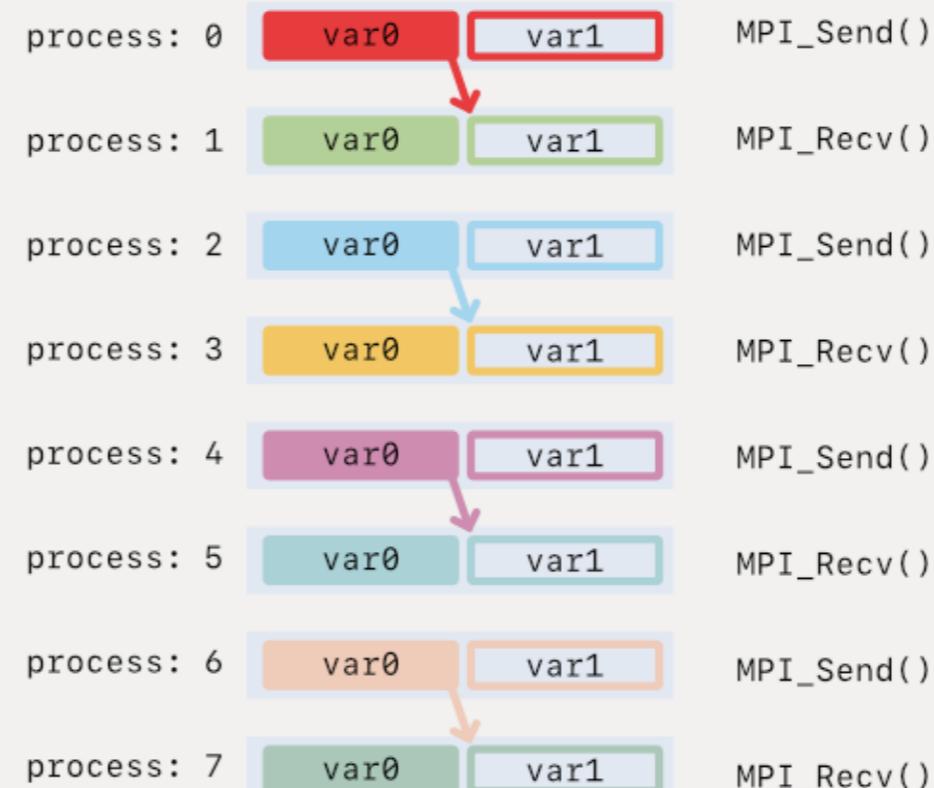
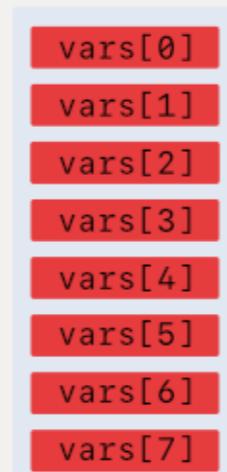
```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```



Exercises

Ex07

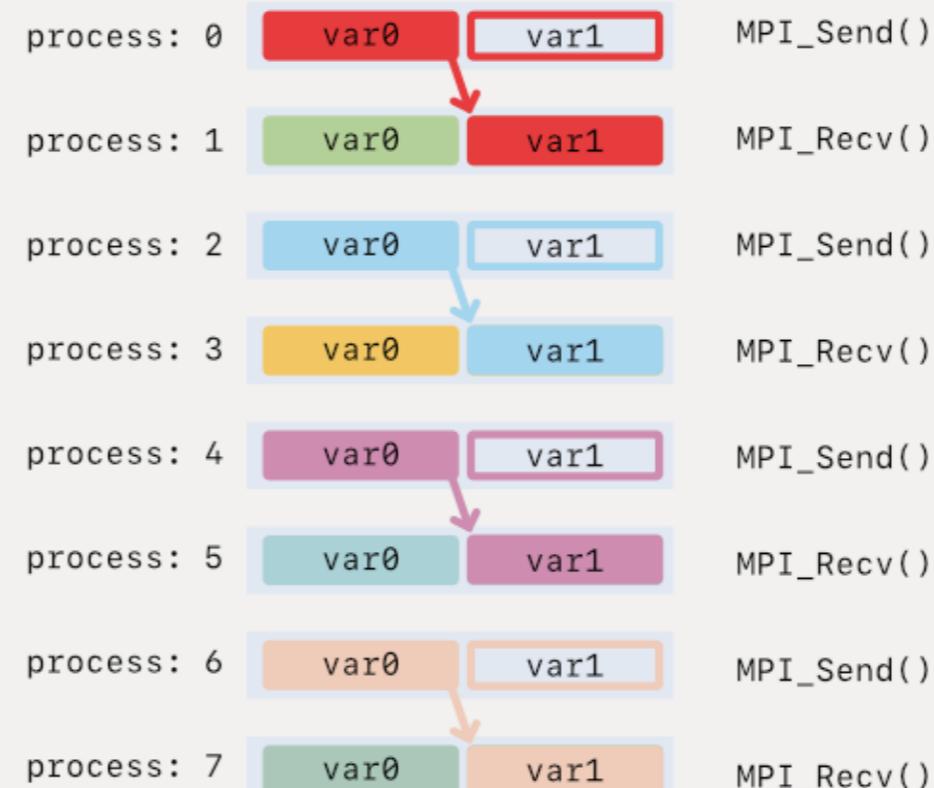
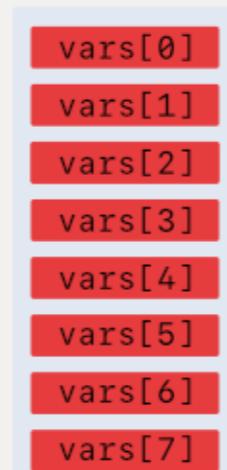
```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```



Exercises

Ex07

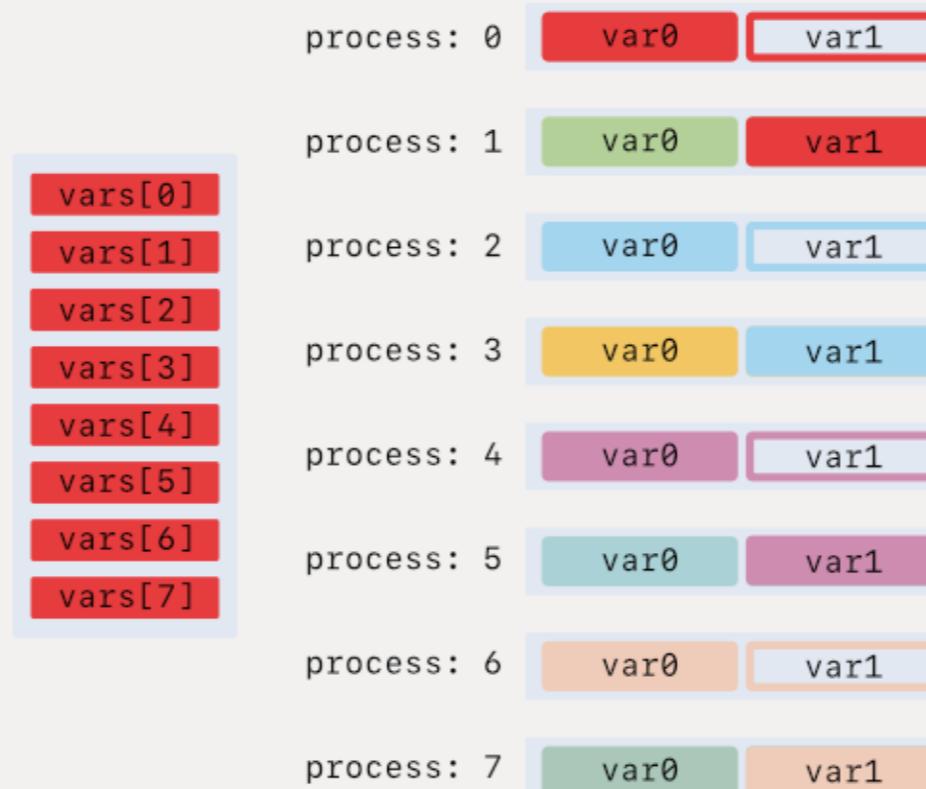
```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```



Exercises

Ex07

```
data.txt  
0.42816572487  
0.57721566490  
0.66274341935  
0.69314718056  
1.41421356237  
1.61803398875  
2.71828182846  
3.14159265359
```

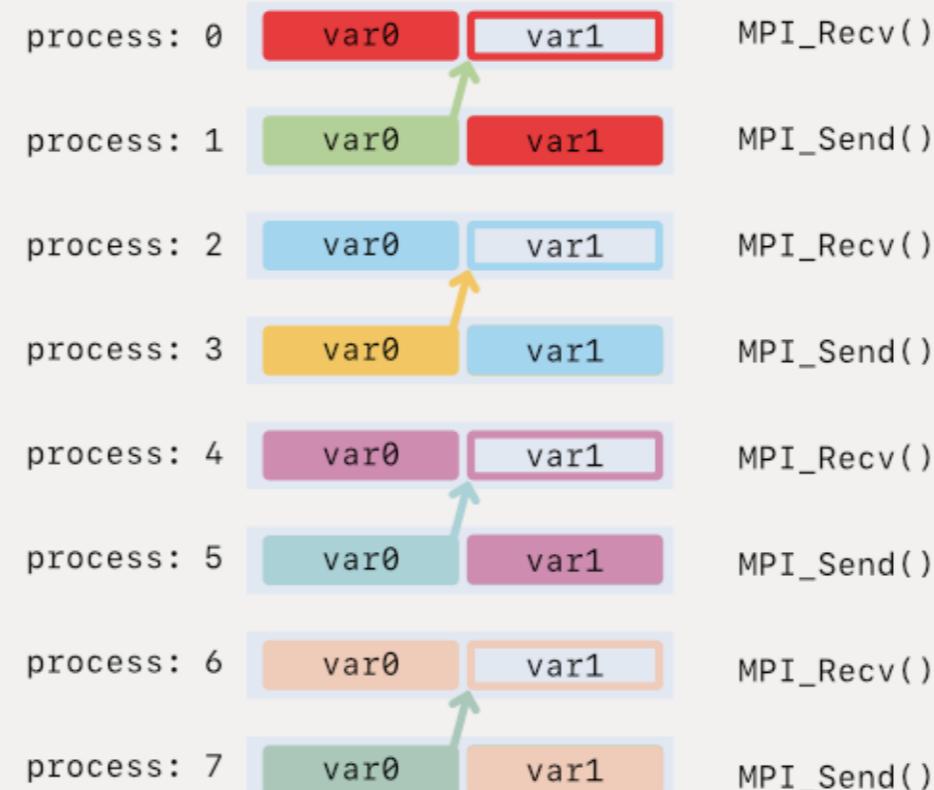


Exercises

Ex07

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

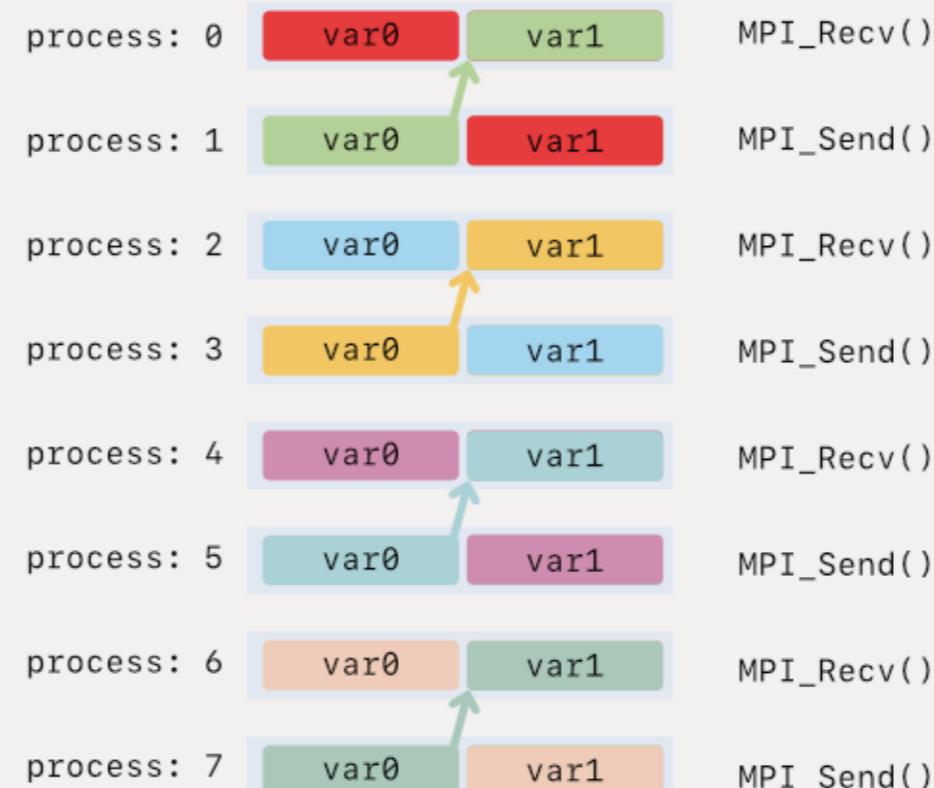


Exercises

Ex07

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

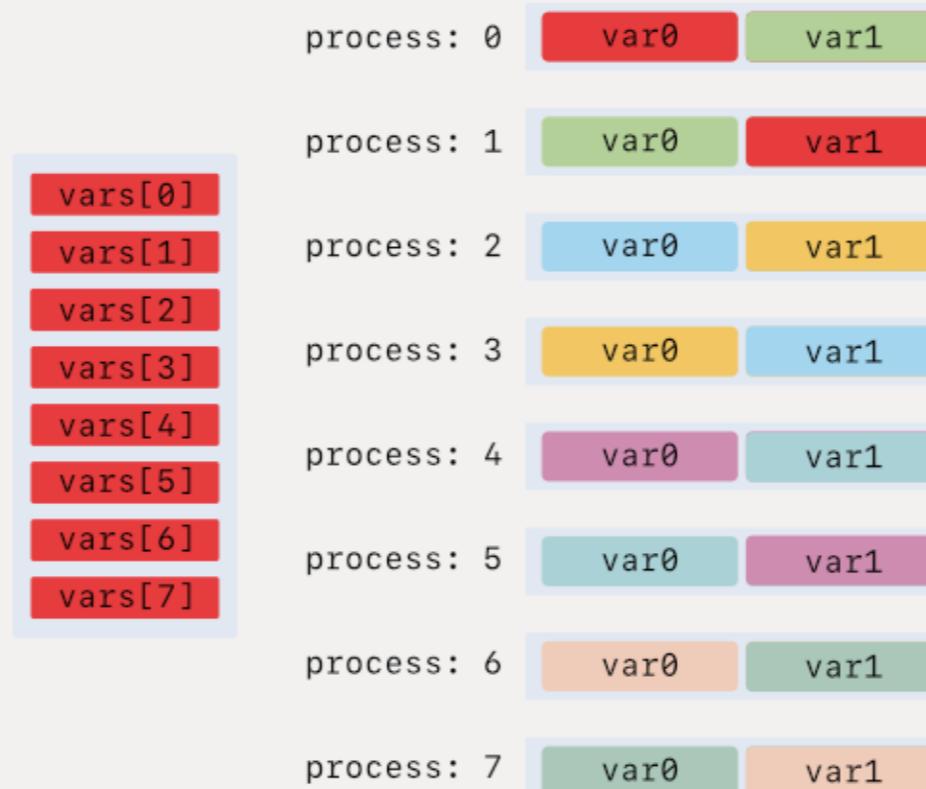
vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]



Exercises

Ex07

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```



Exercises

Ex07

- The `TODO`s are for completing the arguments of the `MPI_Recv()`s and `MPI_Send()`s

```
double var1;
/*
 * TODO: Use `MPI_Send()` and `MPI_Recv()` appropriately, so that
 * `var0` of each even rank is copied into `var1` of the next odd
 * rank
 */
if(rank % 2 == 0) {
    MPI_Send(/* TODO */);
} else {
    MPI_Recv(/* TODO */);
}

/*
 * TODO: Use `MPI_Send()` and `MPI_Recv()` appropriately, so that
 * `var0` of each odd rank is copied into `var1` of the previous
 * even rank
 */
if(rank % 2 == 1) {
    MPI_Send(/* TODO */);
} else {
    MPI_Recv(/* TODO */);
}
```

Exercises

Ex07

- The correct output should look like this:

```
[user@front01 ex07]$ cat ex07-output.txt |sort
This is rank = 0 of nproc = 8 on node: cn01 | var0 = 0.428166    var1 = 0.577216
This is rank = 1 of nproc = 8 on node: cn01 | var0 = 0.577216    var1 = 0.428166
This is rank = 2 of nproc = 8 on node: cn01 | var0 = 0.662743    var1 = 0.693147
This is rank = 3 of nproc = 8 on node: cn01 | var0 = 0.693147    var1 = 0.662743
This is rank = 4 of nproc = 8 on node: cn02 | var0 = 1.414214    var1 = 1.618034
This is rank = 5 of nproc = 8 on node: cn02 | var0 = 1.618034    var1 = 1.414214
This is rank = 6 of nproc = 8 on node: cn02 | var0 = 2.718282    var1 = 3.141593
This is rank = 7 of nproc = 8 on node: cn02 | var0 = 3.141593    var1 = 2.718282
[user@front01 ex07]$
```

Exercises

Ex08

- This exercise demonstrates the use of `MPI_Sendrecv()`
- The same `data.txt` with eight elements is used as before
- The elements are read by the root process and scattered to all processes as before

Exercises

Ex08

- This exercise demonstrates the use of `MPI_Sendrecv()`
- The same `data.txt` with eight elements is used as before
- The elements are read by the root process and scattered to all processes as before
- Now each process has three variables:
 - `var_proc` contains the element received from the scatter
 - `var_next` is to contain `var_proc` of the next process
 - `var_prev` is to contain `var_proc` of the previous process

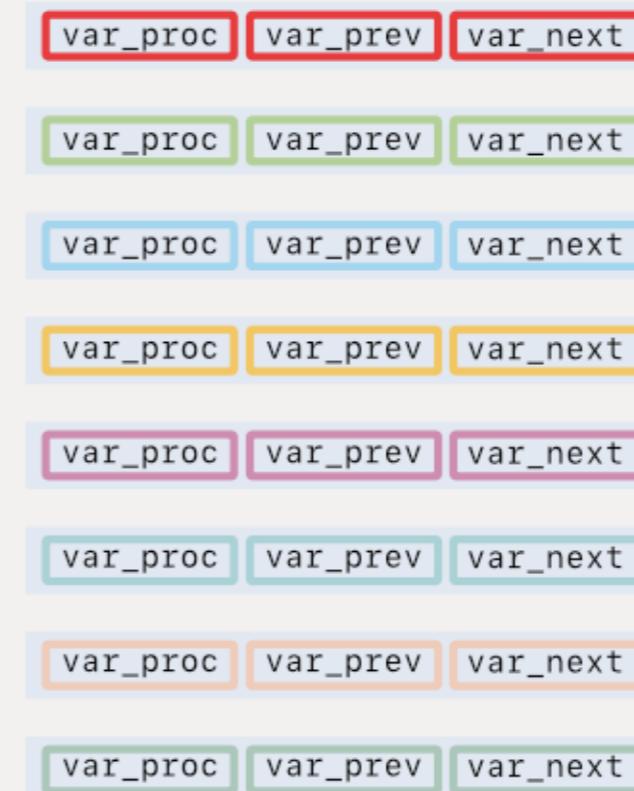
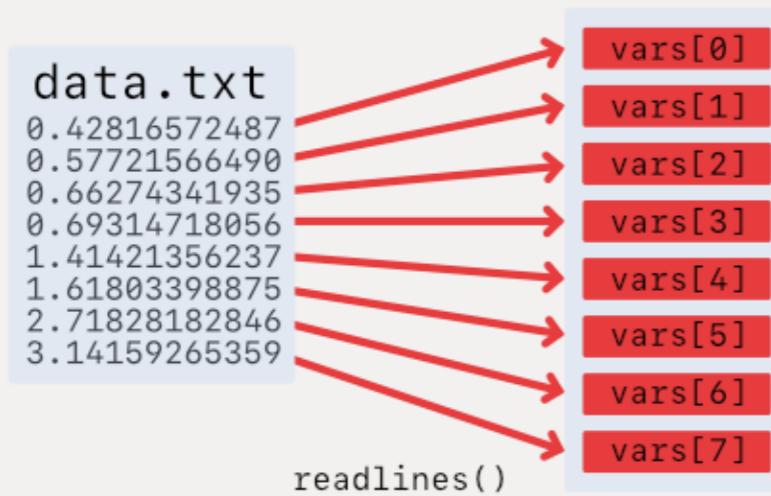
Exercises

Ex08

- This exercise demonstrates the use of `MPI_Sendrecv()`
- The same `data.txt` with eight elements is used as before
- The elements are read by the root process and scattered to all processes as before
- Now each process has three variables:
 - `var_proc` contains the element received from the scatter
 - `var_next` is to contain `var_proc` of the next process
 - `var_prev` is to contain `var_proc` of the previous process
- Use two `MPI_Sendrecv()` to achieve this

Exercises

Ex08



Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

var_proc
var_prev
var_next

var_proc
var_prev
var_next

var_proc
var_prev
var_next

var_proc
var_prev
var_next

var_proc
var_prev
var_next

var_proc
var_prev
var_next

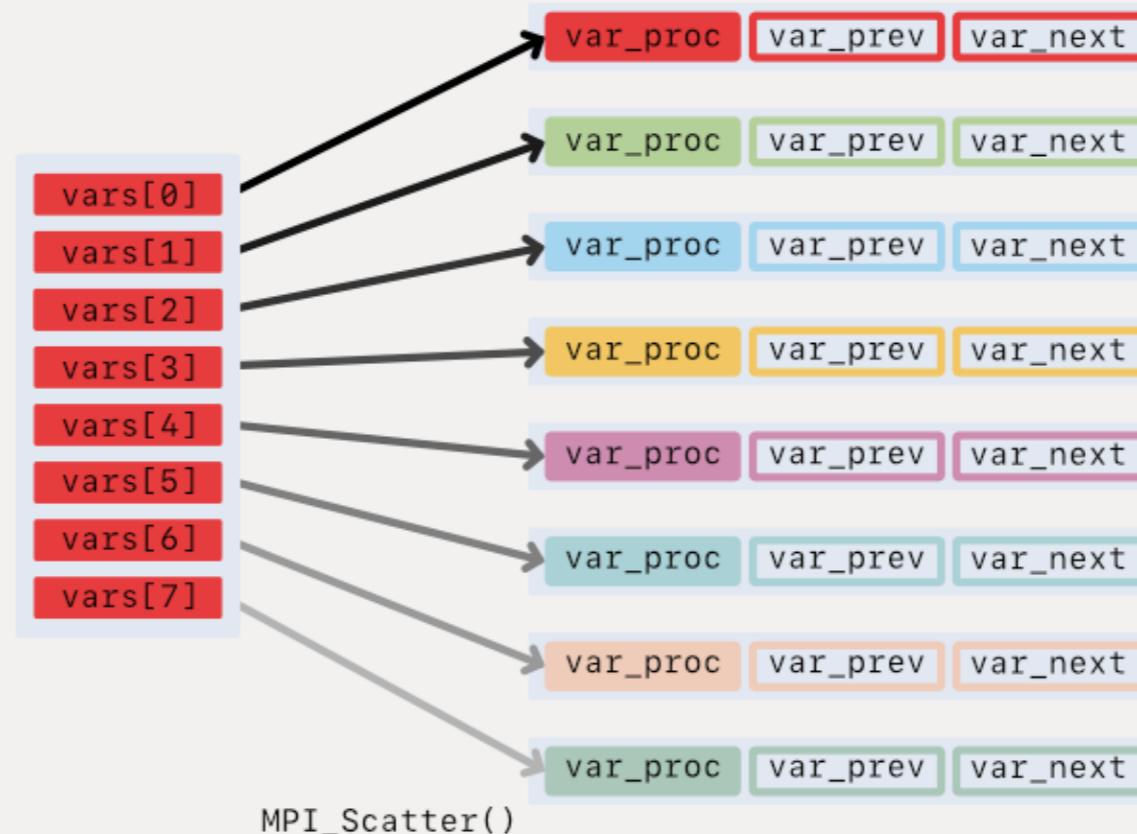
var_proc
var_prev
var_next

var_proc
var_prev
var_next

Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```



Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

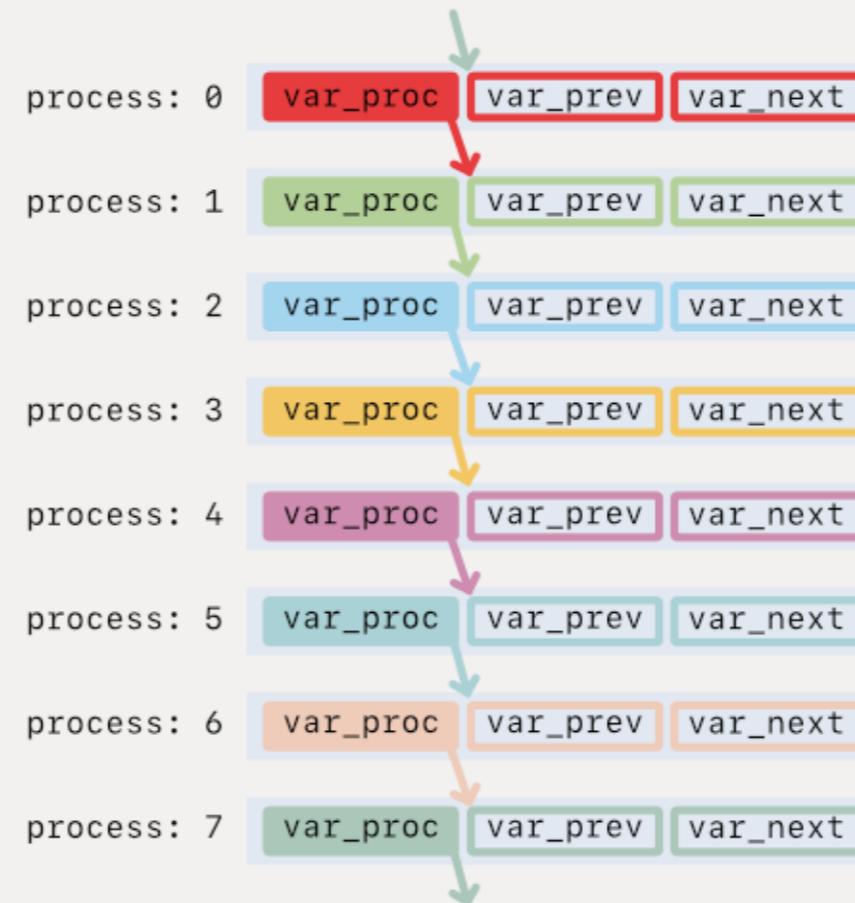
process: 0 var_proc var_prev var_next
process: 1 var_proc var_prev var_next
process: 2 var_proc var_prev var_next
process: 3 var_proc var_prev var_next
process: 4 var_proc var_prev var_next
process: 5 var_proc var_prev var_next
process: 6 var_proc var_prev var_next
process: 7 var_proc var_prev var_next

Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

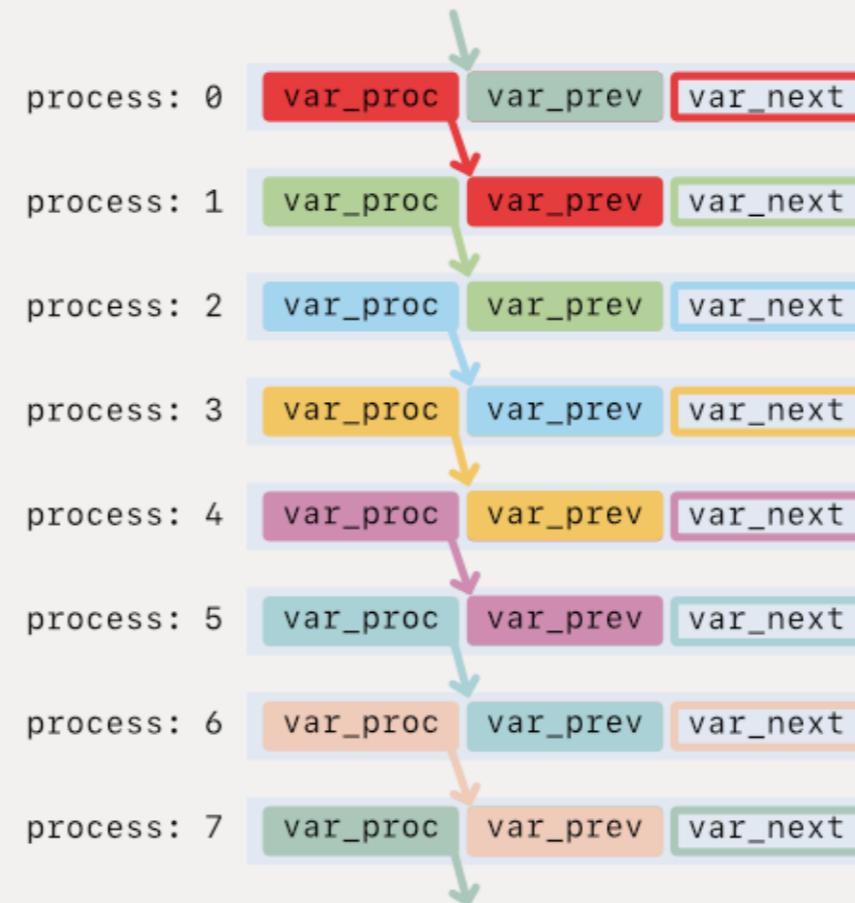


Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]



Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

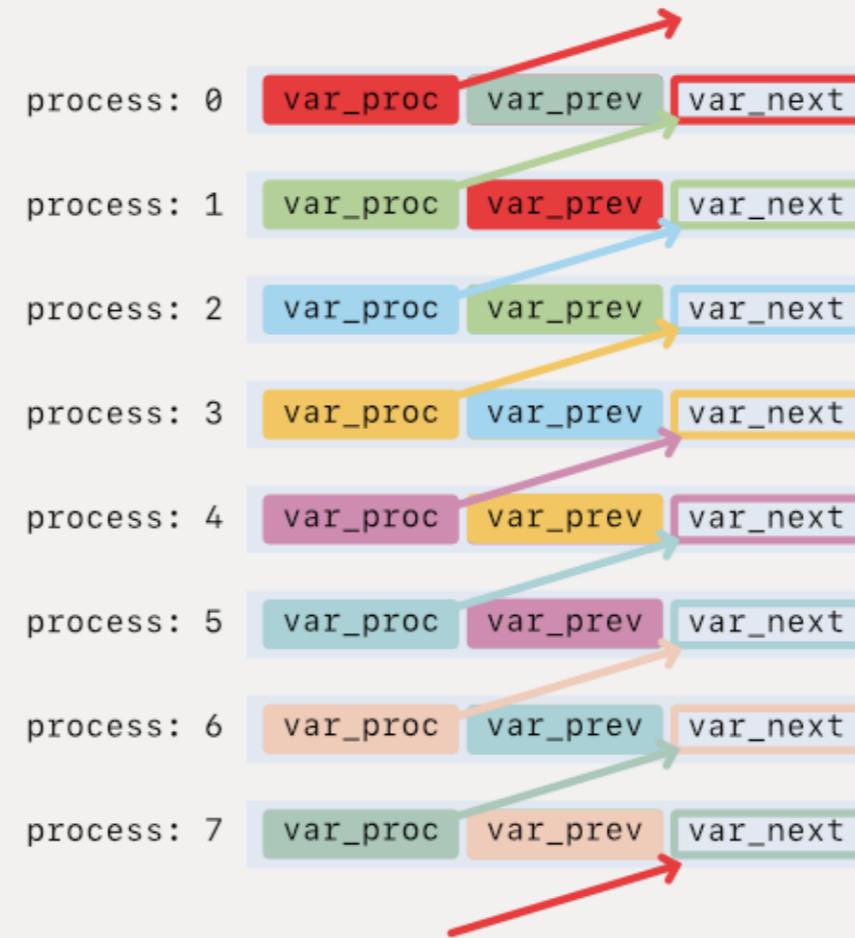
process: 0	var_proc	var_prev	var_next
process: 1	var_proc	var_prev	var_next
process: 2	var_proc	var_prev	var_next
process: 3	var_proc	var_prev	var_next
process: 4	var_proc	var_prev	var_next
process: 5	var_proc	var_prev	var_next
process: 6	var_proc	var_prev	var_next
process: 7	var_proc	var_prev	var_next

Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]



Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]



Exercises

Ex08

```
data.txt
0.42816572487
0.57721566490
0.66274341935
0.69314718056
1.41421356237
1.61803398875
2.71828182846
3.14159265359
```

vars[0]
vars[1]
vars[2]
vars[3]
vars[4]
vars[5]
vars[6]
vars[7]

process: 0 var_proc var_prev var_next
process: 1 var_proc var_prev var_next
process: 2 var_proc var_prev var_next
process: 3 var_proc var_prev var_next
process: 4 var_proc var_prev var_next
process: 5 var_proc var_prev var_next
process: 6 var_proc var_prev var_next
process: 7 var_proc var_prev var_next

Exercises

Ex08

- The `TODO`s are for completing the arguments of the `MPI_Sendrecv()`s

```
double var_next, var_prev;
/*
 * TODO: Use `MPI_Sendrecv()` appropriately, so that `var_proc` of
 * each rank is copied into `var_prev` of the next rank. Assume
 * periodicity of ranks, i.e. if the sender is the last process
 * (`rank = nproc - 1') then send to the first process (`rank =
 * 0')
 */
MPI_Sendrecv(/* TODO */);

/*
 * TODO: Use `MPI_Sendrecv()` appropriately, so that `var_proc` of
 * each rank is copied into `var_next` of the previous rank. Assume
 * periodicity of ranks, i.e. if the sender is the first process
 * (`rank = 0') then send to the last process (`rank = nproc - 1')
 */
MPI_Sendrecv(/* TODO */);
```

Exercises

Ex08

- The correct output should look like this:

```
[user@front01 ex08]$ cat ex08-output.txt | sort
This is rank = 0 of nproc = 8 on node: cn01  | var_proc = 0.428166  var_prev = 3.141593  var_next = 0.577216
This is rank = 1 of nproc = 8 on node: cn01  | var_proc = 0.577216  var_prev = 0.428166  var_next = 0.662743
This is rank = 2 of nproc = 8 on node: cn01  | var_proc = 0.662743  var_prev = 0.577216  var_next = 0.693147
This is rank = 3 of nproc = 8 on node: cn01  | var_proc = 0.693147  var_prev = 0.662743  var_next = 1.414214
This is rank = 4 of nproc = 8 on node: cn02  | var_proc = 1.414214  var_prev = 0.693147  var_next = 1.618034
This is rank = 5 of nproc = 8 on node: cn02  | var_proc = 1.618034  var_prev = 1.414214  var_next = 2.718282
This is rank = 6 of nproc = 8 on node: cn02  | var_proc = 2.718282  var_prev = 1.618034  var_next = 3.141593
This is rank = 7 of nproc = 8 on node: cn02  | var_proc = 3.141593  var_prev = 2.718282  var_next = 0.428166
[user@front01 ex08]$
```