

# HP-SEE

## Programming with MPI

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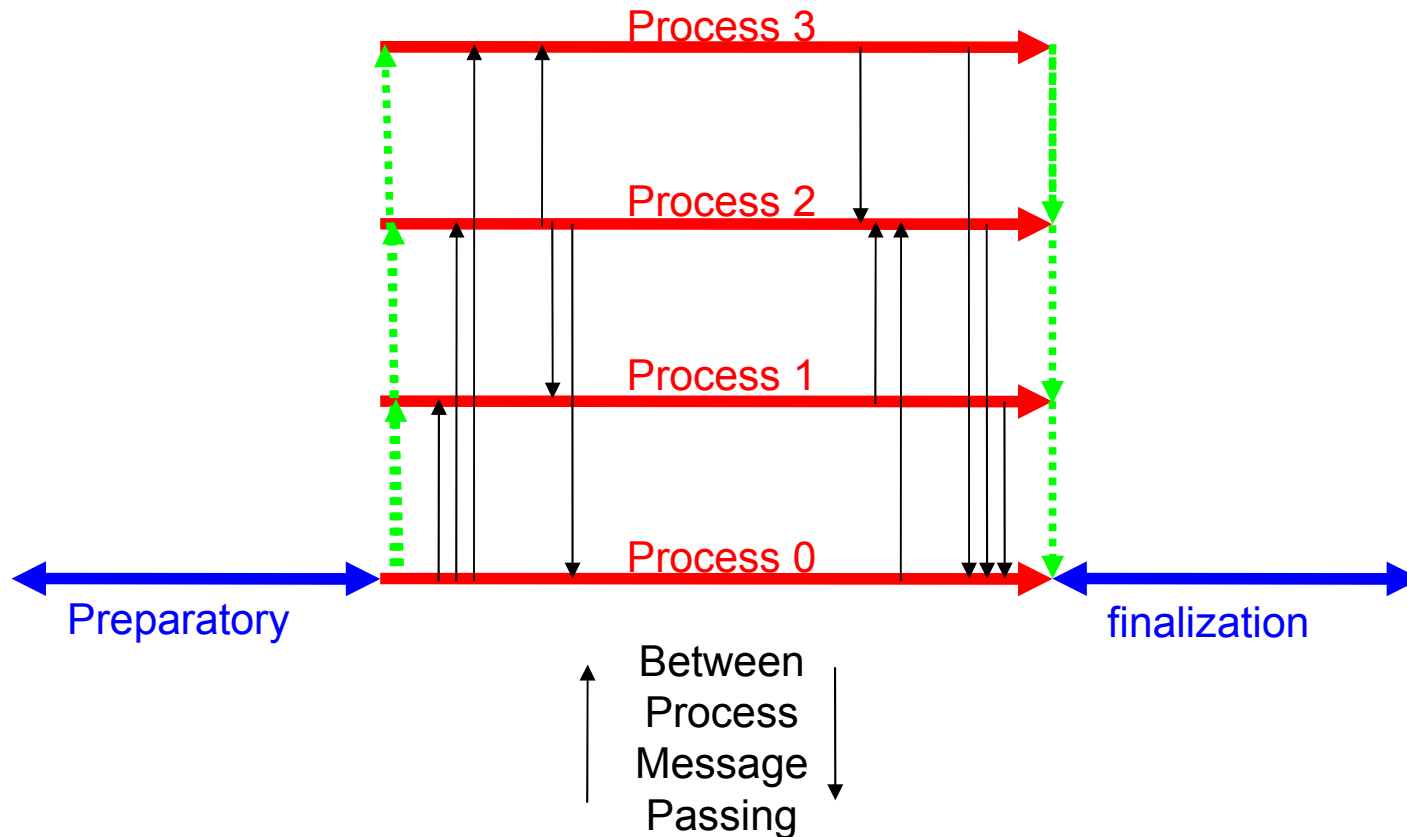
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# Principle of MPI

## Working with PROCESSES

### MESSAGE PASSING INTERFACE



# Test MPI Source Code



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```
#include <mpi.h>
#include <stdio.h>
#include <stdlib.h>
#include <math.h>
typedef struct
{ double thred, start, stop;
} metad; metad * array, procs;
int main(int argc, char **argv)
{ int rank, size, i, j, k;
  double sinsin, starttime, stoptime;
  long int li, Nruntime; Nruntime=9999999;
MPI_Init(&argc, &argv); // init MPI
MPI_Comm_rank(MPI_COMM_WORLD, &rank);
MPI_Comm_size(MPI_COMM_WORLD, &size);
// print inputs & take time in mother process
if (rank==0)
{ starttime = MPI_Wtime();
  array = malloc(size*sizeof(metad));
  for (i=0; i<size; i++)
  { array[i].thred=0;
    array[i].start=0; array[i].stop =0;
  }
}
```

```
MPI_Scatter(&array[rank],3,MPI_DOUBLE,
           &procs,3,MPI_DOUBLE,0,MPI_COMM_WORLD);
// run processes inner loops
procs.start = MPI_Wtime();
for (j=0; j<16/size; j++)
  for (k=0; k<Nruntime; k++)
    { sinsin=sin(k); }
procs.thred= rank;
procs.stop = MPI_Wtime();
MPI_Gather(&procs,3,MPI_DOUBLE,array,3,MPI_DOUBLE,0,MPI_COMM_WORLD);
// print process runtime vector in mother process
if(rank==0)
{ for (i=0; i<size; i++)
    printf("Iter %d process %f Runtime %f \n",
          i, array[i].thred, array[i].stop - array[i].start);
  stoptime = MPI_Wtime();
  printf("time = %f \n",stoptime-startime);
} MPI_Finalize(); // close MPI
return 0; }
```



# Comments on the Source

## Working with processes

Each process manages its own central memory area

Loop on processes is missing ~ automatically via **mpirun**

Part of preparation & finalization in mother process

## Procedures to distribute and collect data

- MPI\_Scatter and MPI\_Gather  
    distribute & collect slots of data
- MPI\_Bcast for single values  
    distribute the same data with synchronization
- MPI\_Send and MPI\_Recv  
    process-to-process etc.



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# Preparation of Test

## Compilation

```
build.sh = mpicc test.c -o test
```

## Script

```
run.sh = (/usr/bin/time mpirun -np $2 ./$1)
```

## Execution

```
./run.sh test <no_process>
```

## Tested in hardware

Dell Inspiron ~ 1 processor 2 cores



# MPI Test Example

```
$ ./run.sh test 4
```

```
input nopr=4
```

```
Iter 0 process 0 Runtime 3.508056
```

```
Iter 1 process 1 Runtime 3.685100
```

```
Iter 2 process 2 Runtime 3.748083
```

```
Iter 3 process 3 Runtime 3.768022
```

```
time = 3.768220
```

```
7.26 user 0.13 system 0:04.83 elapsed 152% CPU
```

```
(0 avgtext + 0 avgdata 14448 maxresident)k
```

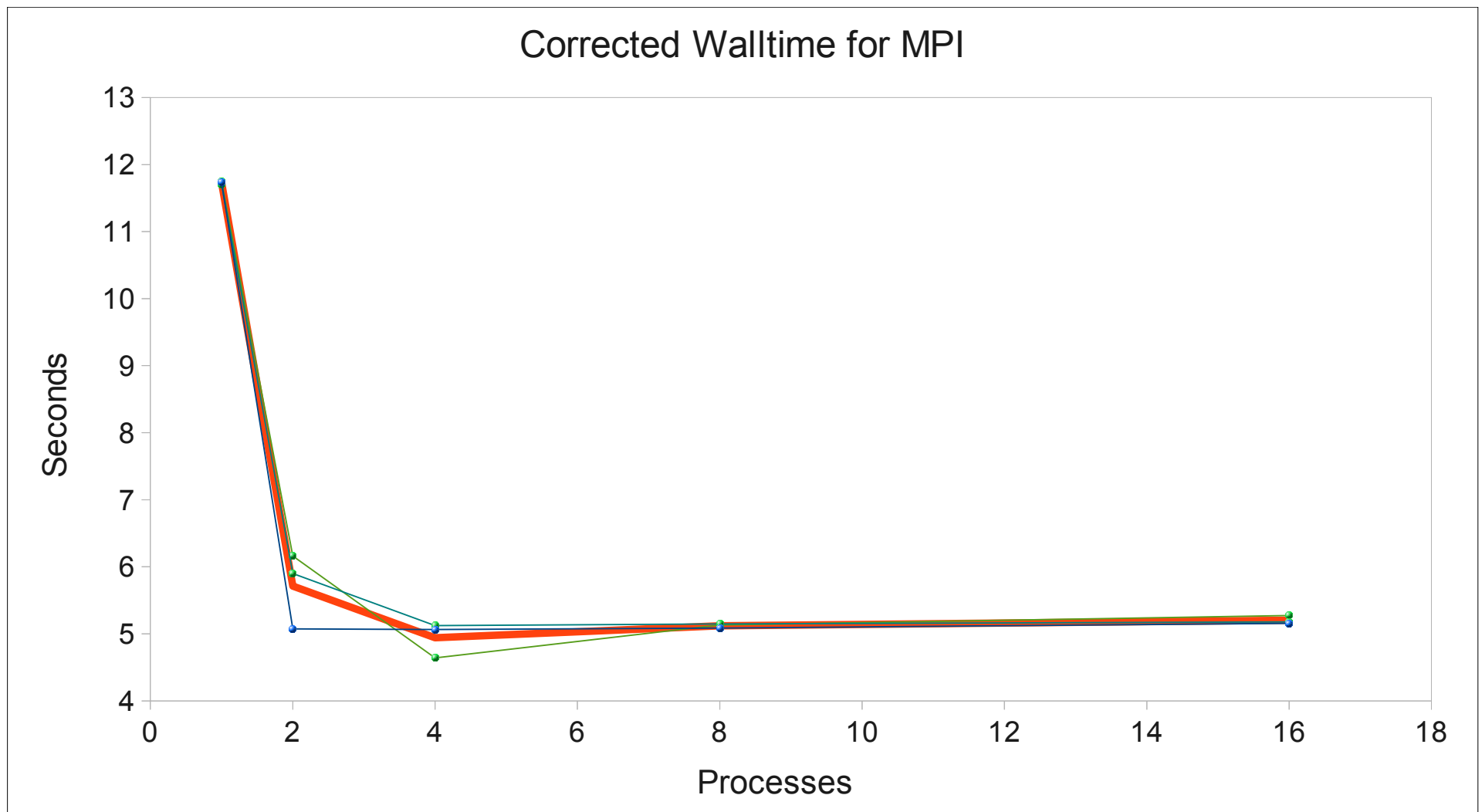
```
103 inputs + 680 outputs (19 major + 5707 minor) pagefaults 0 swaps
```

# Test MPI Walltime



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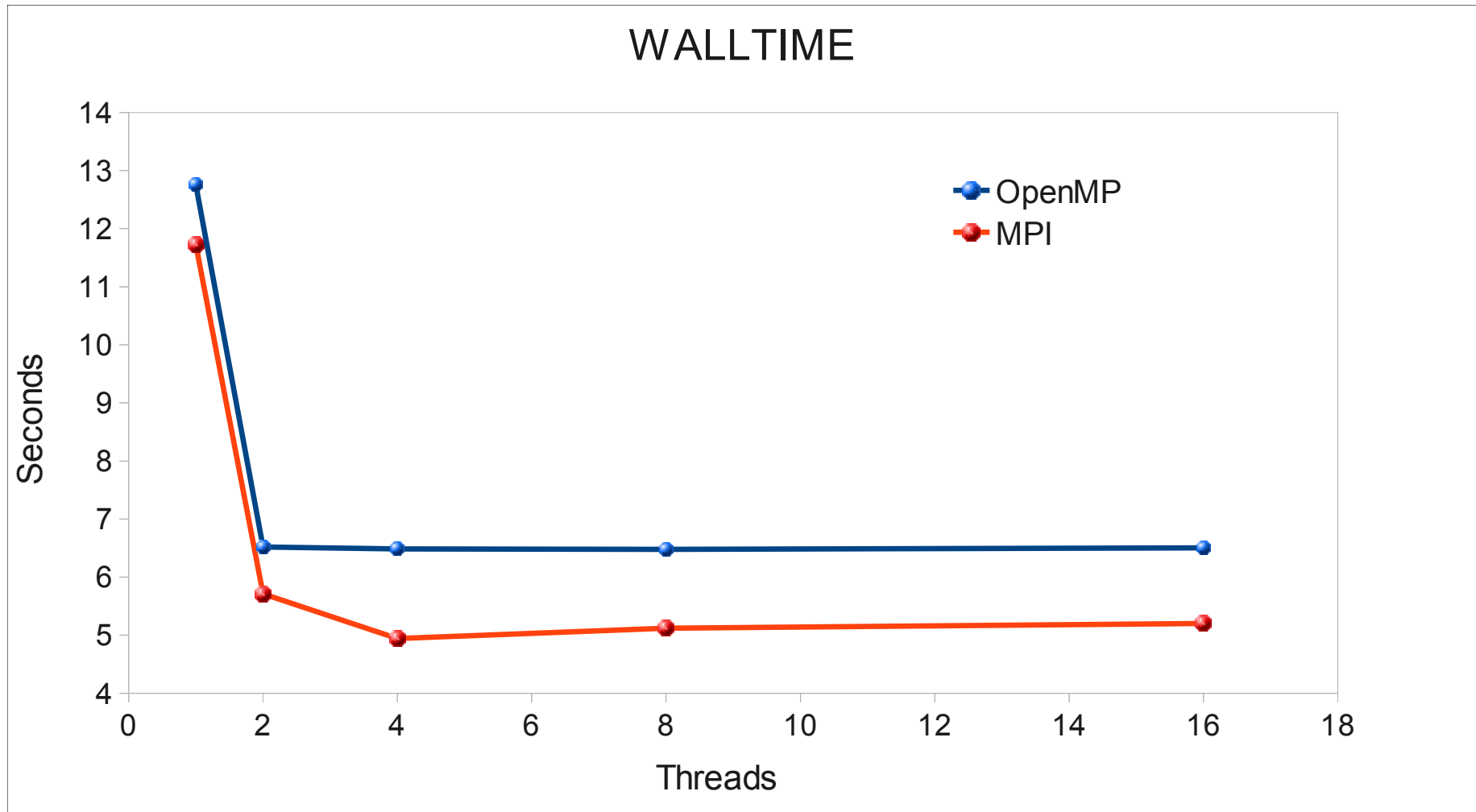


# Compared with OpenMP



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# Comments on Performance

The case of four processes in two cores

Started simultaneously

Shared the time of CPU (total of 152%)

Similar runtime for all processes

Performance ~ number of cores

Compared with OpenMP

- OpenMP requires shared memory for all cores
- MPI may run in distributed hardware ...



# Conclusions on MPI

## More difficult to be implemented

- Need for radical modifications in the code
  - replacement of loops
  - use of message passing routines
  - dealing with distributed data files

## But ... More easy to run

- Does not require shared memory architecture
- May run in distributed hardware (cluster / grid)



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# References

<http://www.open-mpi.org/>

<http://software.intel.com/>

<http://www.openchannelsoftware.com/>