

FEDERAL STATE AUTONOMOUS EDUCATIONAL  
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ITMO UNIVERSITY

Report

MPI. Assignments 12 – 13

Parallel algorithms for the analysis and synthesis of data

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# 1 Assignments

## 1.1 Assignment 12. MPI. Delayed interactions. Scheme of an iterative method with exchange over a ring topology using pending requests.

### 1.1.1 Formulation of the problem

MPI allows for non-blocking operations to form whole packets of requests for communication operations `MPI_SEND_INIT` and `MPI_RECV_INIT`, which are started by the `MPI_START` or `MPI_STARTALL` functions.

Checking for completion of execution is performed by conventional means using the functions of the **WAIT** and **TEST** families.

Find and fix errors in `ASSIGNMENT12.C`, add the for loop. When should you use a loop?

### 1.1.2 Example of launch parameters and output. Detailed description of solution

Code for **assignment 12** is [here](#).

Compilation example: `MPIC++ -O ./CPF/12.O ASSIGNMENT12.C`

Launch example: `MPIRUN -OVERSUBSCRIBE -NP 10 ./CPF/12.O`

```
aptmess@improfeo: ~/ITMO/parallel_algorithms/HT/hw_mpi
(base) aptmess@improfeo:~/ITMO/parallel_algorithms/HT/hw_mpi$ mpic++ -o ./cpf/12.o Assignment12.c
(base) aptmess@improfeo:~/ITMO/parallel_algorithms/HT/hw_mpi$ mpirun --oversubscribe -np 10 ./cpf/12.o
1 (previous) -> 2 (current) -> 3 (next)
7 (previous) -> 8 (current) -> 9 (next)
2 (previous) -> 3 (current) -> 4 (next)
3 (previous) -> 4 (current) -> 5 (next)
0 (previous) -> 1 (current) -> 2 (next)
8 (previous) -> 9 (current) -> 0 (next)
9 (previous) -> 0 (current) -> 1 (next)
4 (previous) -> 5 (current) -> 6 (next)
5 (previous) -> 6 (current) -> 7 (next)
6 (previous) -> 7 (current) -> 8 (next)
(base) aptmess@improfeo:~/ITMO/parallel_algorithms/HT/hw_mpi$ _
```

Let's move to the the code and explain how it works.

```

1  #include <iostream>
2  #include "mpi.h"
3
4  using namespace std;
5  int main(int argc, char **argv)
6  {
7      int rank, size, prev, next;
8      float sbuf[2], rbuf[2];
9
10     MPI_Init(&argc, &argv);
11
12     MPI_Request reqs[4];
13     MPI_Status stats[4];
14
15     MPI_Comm_size(MPI_COMM_WORLD, &size);
16     MPI_Comm_rank(MPI_COMM_WORLD, &rank);
17
18     prev = rank - 1;
19     next = rank + 1;
20
21     if (rank == 0) prev = size - 1;
22     if (rank == size - 1) next = 0;
23
24     MPI_Recv_init(&rbuf[0], 1, MPI_FLOAT, prev, 5, MPI_COMM_WORLD, &reqs[0]);
25     MPI_Recv_init(&rbuf[1], 1, MPI_FLOAT, next, 6, MPI_COMM_WORLD, &reqs[1]);
26     MPI_Send_init(&sbuf[0], 1, MPI_FLOAT, prev, 6, MPI_COMM_WORLD, &reqs[2]);
27     MPI_Send_init(&sbuf[1], 1, MPI_FLOAT, next, 5, MPI_COMM_WORLD, &reqs[3]);
28
29     for (int i = 0; i < 2; i++) sbuf[i] = float(rank);
30
31     MPI_Startall(4, reqs);
32     MPI_Waitall(4, reqs, stats);
33
34     cout << rbuf[0] << " (previous)" << " -> " << rank << " (current)" << " -> " << rbuf[1] << " (next)" << '\n' << endl;
35
36     for (int i = 0; i < 4; i++) MPI_Request_free(&reqs[i]);
37
38     MPI_Finalize();
39 }

```

### Assignment12 code

Firstly i wrote a whole program that can be compiled and works correctly as a ring topology as we can see on picture. As fixing errors in lines 24 – 27 i have add by symbol & address of variable RBUF and SBUF (also i have created float arrays with the same names), also do the same for REQS. In lines 26 – 27 we can see that there is no values in array SBUF, so for loop is for mapping float value of rank to each piece of array (line 29). Inside for loop there is no need to start sending and recieving messages on each iteration - we should do it only ones as i did it in lines 31 – 32. After that i displayed information in ring topology and used for loop for function MPI\_REQUEST\_FREE to free inactive persistent requests created with either MPI\_RECV\_INIT or MPI\_SEND\_INIT and friends. All that i said is the answer on question when should we use loop: free inactive requests and may be for initialzing data to send. That's all, program works correctly and errors are fixed.

## 1.2 Assignment 13. MPI. Collective process interactions. Barrier.

### 1.2.1 Formulation of the problem

Find out which process will perform the multiplication of two  $500 \times 500$  square matrices faster.

Complete the code ASSIGNMENT13.C. You can use the necessary code from the previous assignments.

### 1.2.2 Example of launch parameters and output. Detailed description of solution

Code for assignment 13 is [here](#).

Compilation example: `MPIC++ -O ./CPF/13.O ASSIGNMENT13.C`

Launch example: `MPIRUN -OVERSUBSCRIBE -NP 5 ./CPF/13.O 500`

```
(base) aptmess@improfeo:~/ITM0/parallel_algorithms/HT/hw_mpi$ mpirun --oversubscribe -np 5 ./cpf/13.o 500
process 2, execution time=2.03871
process 4, execution time=2.37145
process 3, execution time=2.41065
process 0, execution time=2.42486
process 1, execution time=2.48914
(base) aptmess@improfeo:~/ITM0/parallel_algorithms/HT/hw_mpi$ _
```

Let's move to the the code and explain how it works.

```

1  #include <mpi.h>
2  #include <stdio.h>
3  #include <ctime>
4  #include <cstdlib>
5  #include <iostream>
6
7  using namespace std;
8
9  int main(int argc, char **argv)
10 {
11     int matrix_size = atoi(argv[1]);
12     int rank, size;
13     MPI_Init(&argc, &argv);
14
15     double start_time, end_time;
16     MPI_Comm_size(MPI_COMM_WORLD, &size);
17     MPI_Comm_rank(MPI_COMM_WORLD, &rank);
18
19     //matrix initialization by each process
20
21     int A[matrix_size][matrix_size], B[matrix_size][matrix_size], C[matrix_size][matrix_size];
22
23     for (int i = 0; i < matrix_size; i++)
24     {
25         for (int j = 0; j < matrix_size; j++)
26         {
27             A[i][j] = 1 + rand() % 10;
28             B[i][j] = 1 + rand() % 10;
29             C[i][j] = 0;
30         }
31     }
32
33     MPI_Barrier(MPI_COMM_WORLD); //barrier process synchronization
34
35     // start timing for each process
36
37     start_time = MPI_Wtime();
38
39     // matrix multiplication
40
41     for (int i = 0; i < matrix_size; i++) {
42         for (int j = 0; j < matrix_size; j++) {
43             for (int k = 0; k < matrix_size; k++) {
44                 C[i][j] += A[i][k] * B[k][j];
45             }
46         }
47     }
48
49     // end of timing
50
51     end_time = MPI_Wtime();
52
53     // output the execution time of matrix multiplication at each process
54
55     cout << "process " << rank ;
56     cout << ", execution time=" << (end_time - start_time);
57     cout << '\n' << endl;
58     MPI_Finalize();
59 }

```

Assignment13 code

What should we need was written in file by comments, by job was to fill the gaps as code. Firstly i initialize matrices  $A$  and  $B$  with the same seed for each process with random integers from 1 to 10 and matrix  $C$  by null values. Next there are a barrier function which is waiting when all processes initialize their matrices, after that we start timing for eahc process, do known matrix multiplication algorithm, end timing and show execution time of matrix multiplication at each process and understand which process performed faster. In our example it is process with rank 2. Program works correctly and algorithm is explained.

## 1.3 Appendix

The link to the sourse code which is placed on my [github](#).