# Отчет

# Задание №2 Решение двумерного уравнения теплопроводности

(неявная схема, MPI+OpenMP) приложение

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# 7. Таблицы и графики, содержащие сведения о размерах сеток, времени решения и эффективности распараллеливания

$$\frac{\partial u}{\partial t} = \frac{\partial}{\partial x} \left( k \left( x, y \right) \frac{\partial u}{\partial x} \right) + \frac{\partial}{\partial y} \left( k \left( x, y \right) \frac{\partial u}{\partial y} \right) + f \left( x, y, t \right), \quad (x, y) \in (0, 1) \times (0, 1), \quad t_{max} = 1$$

$$k(x,y) = 1,$$
  $f(x,y,t) = -exp(-t) \cdot (1 - \pi^2) \cdot (sin(\pi x) + cos(\pi y))$ 

Точное решение:  $u(x,y,t) = exp(-t) \cdot (sin(\pi x) + cos(\pi y)) + 10$ 

### 1) nx = 1000, ny = 1000, nt = 10

| Число | Число | Время выполнения, с | Ускорение | Эффективность, % |
|-------|-------|---------------------|-----------|------------------|
| узлов | ядер  |                     |           |                  |
| 1     | 1     | 21.030              | 1         | 100              |
| 64    | 256   | 0.215               | 24.430    | 38.173           |
| 128   | 512   | 0.460               | 11.411    | 8.915            |

|       | edu-cmc-ski16-052@fen1:~/IMPLICIT_MPI_OPENMP> make accel<br>perl accel.pl |             |                  |                  |            |            |  |  |  |
|-------|---|-------------|------------------|------------------|------------|------------|--|--|--|
| Nodes | Cores   | Time        | Speed Up         | Efficiency, %    | Abs. error | Rel. error |  |  |  |
| 1     | 1   | 21.03000000 | 1                | 100              | 0.01021743 | 0.00105858 |  |  |  |
| 64    | 256   | 0.21519931  | 24.4308404148694 | 38.1731881482334 | 0.01021743 | 0.00105858 |  |  |  |
| 128   | 512   | 0.46070012  | 11.4119787943619 | 8.91560843309527 | 0.01021743 | 0.00105858 |  |  |  |

### 2) nx = 2000, ny = 2000, nt = 10

| Число | Число | Время выполнения, с | Ускорение | Эффективность, % |
|-------|-------|---------------------|-----------|------------------|
| узлов | ядер  |                     |           |                  |
| 1     | 1     | 84.980              | 1         | 100              |
| 64    | 256   | 0.640               | 33.161    | 51.814           |
| 128   | 512   | 0.881               | 24.095    | 18.824           |

| edu-cmc-ski16-052@fen1:~/IMPLICIT_MPI_OPENMP> make accel<br>berl accel.pl |       |             |                  |                  |            |            |  |
|---|-------|-------------|------------------|------------------|------------|------------|--|
| Nodes   | Cores | Time        | Speed Up         | Efficiency, %    | Abs. error | Rel. error |  |
| 1   | 1     | 84.98000000 | 1                | 100              | 0.01026958 | 0.00106405 |  |
| 64  | 256   | 0.64065958  | 33.1611368396302 | 51.8142763119222 | 0.01026958 | 0.00106405 |  |
| 128   | 512   | 0.88168831  | 24.09581680855   | 18.8248568816797 | 0.01026958 | 0.00106405 |  |

# 3) nx = 3000, ny = 3000, nt = 10

| Число | Число | Время выполнения, с | Ускорение | Эффективность, % |
|-------|-------|---------------------|-----------|------------------|
| узлов | ядер  |                     |           |                  |
| 1     | 1     | 192.790             | 1         | 100              |
| 64    | 256   | 1.313               | 36.687    | 57.324           |
| 128   | 512   | 1.565               | 30.784    | 24.050           |

| edu-cmc-ski16-052@fen1:~/IMPLICIT_MPI_OPENMP> make accel<br>perl accel.pl |       |              |                  |                  |            |            |  |  |
|---|-------|--------------|------------------|------------------|------------|------------|--|--|
| Nodes   | Cores | Time         | Speed Up         | Efficiency, %    | Abs. error | Rel. error |  |  |
| 1   | 1     | 192.79000000 | 1                | 100              | 0.01028702 | 0.00106588 |  |  |
| 64  | 256   | 1.31372073   | 36.687782189446  | 57.3246596710094 | 0.01028702 | 0.00106588 |  |  |
| 128   | 512   | 1.56563626   | 30.7846089359223 | 24.0504757311893 | 0.01028702 | 0.00106588 |  |  |

### 4) nx = 4000, ny = 4000, nt = 10

| Число | Число | Время выполнения, с | Ускорение | Эффективность, % |
|-------|-------|---------------------|-----------|------------------|
| узлов | ядер  |                     |           |                  |
| 1     | 1     | 344.460             | 1         | 100              |
| 64    | 256   | 2.486               | 34.630    | 54.110           |
| 128   | 512   | 2.400               | 35.874    | 28.026           |

| edu-cmc-ski16-052@fen1:~/IMPLICIT_MPI_OPENMP> make accel<br>perl accel.pl |       |              |                  |                  |            |            |  |  |
|---|-------|--------------|------------------|------------------|------------|------------|--|--|
| Nodes   | Cores | Time         | Speed Up         | Efficiency, %    | Abs. error | Rel. error |  |  |
| 1   | 1     | 344.46000000 | 1                | 100              | 0.01029575 | 0.00106680 |  |  |
| 64  | 256   | 2.48665432   | 34.6308689983093 | 54.1107328098583 | 0.01029575 | 0.00106680 |  |  |
| 128   | 512   | 2.40047447   | 35.874157828473  | 28.0266858034945 | 0.01029575 | 0.00106680 |  |  |

# 5) nx = 5000, ny = 5000, nt = 10

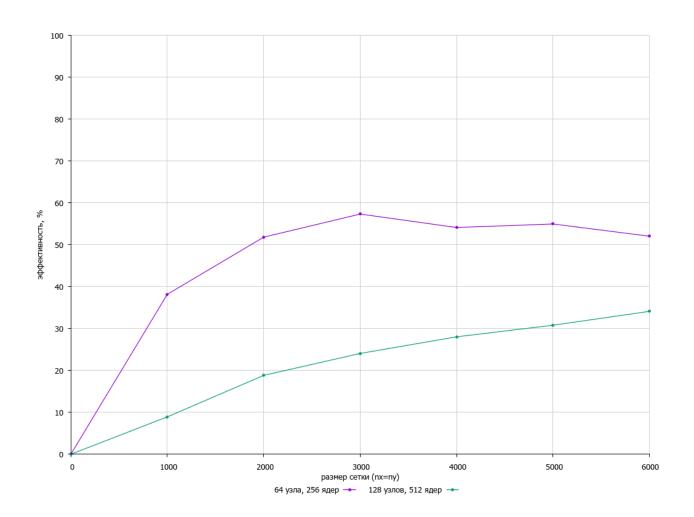
| Число | Число | Время выполнения, с | Ускорение | Эффективность, % |
|-------|-------|---------------------|-----------|------------------|
| узлов | ядер  |                     |           |                  |
| 1     | 1     | 540.010             | 1         | 100              |
| 64    | 256   | 3.838               | 35.174    | 54.959           |
| 128   | 512   | 3.425               | 39.413    | 30.791           |

| edu-cmc-ski16-052@fen1:~/IMPLICIT_MPI_OPENMP> make accel<br>perl accel.pl |       |              |                  |                  |            |            |  |  |
|---|-------|--------------|------------------|------------------|------------|------------|--|--|
| Nodes   | Cores | Time         | Speed Up         | Efficiency, %    | Abs. error | Rel. error |  |  |
| 1   | 1     | 540.01000000 | 1                | 100              | 0.01030099 | 0.00106735 |  |  |
| 64  | 256   | 3.83812213   | 35.1741021852267 | 54.9595346644167 | 0.01030099 | 0.00106735 |  |  |
| 128   | 512   | 3.42531047   | 39.413215585097  | 30.791574675857  | 0.01030099 | 0.00106735 |  |  |

# 6) nx = 6000, ny = 6000, nt = 10

| Число | Число | Время выполнения, с | Ускорение | Эффективность, % |
|-------|-------|---------------------|-----------|------------------|
| узлов | ядер  |                     |           |                  |
| 1     | 1     | 776.620             | 1         | 100              |
| 64    | 256   | 5.830               | 33.301    | 52.034           |
| 128   | 512   | 4.448               | 43.649    | 34.100           |

| edu-cmc-ski16-052@fen1:~/IMPLICIT_MPI_OPENMP> make accel<br>perl accel.pl |       |              |                  |                  |            |            |  |  |
|---|-------|--------------|------------------|------------------|------------|------------|--|--|
| Nodes   | Cores | Time         | Speed Up         | Efficiency, %    | Abs. error | Rel. error |  |  |
| 1   | 1     | 776.62000000 | 1                | 100              | 0.01030449 | 0.00106771 |  |  |
| 64  | 256   | 5.83015086   | 33.30188268919   | 52.0341917018593 | 0.01030449 | 0.00106771 |  |  |
| 128   | 512   | 4.44807063   | 43.6492619272999 | 34.1009858807031 | 0.01030449 | 0.00106771 |  |  |



#### 9. Дополнительные материалы

#### Компиляция и запуск

#### Makefile

```
compile_XL_03:
mpixlcxx_r implicit_single.cpp -03 -qhot -qstrict
                                                            -o single.x:
mpixlcxx_r implicit_mpi.cpp
                             -03 -qhot -qstrict -qsmp=omp -o mpi.x;
compile_XL_04:
mpixlcxx_r implicit_single.cpp -04 -qhot -qstrict
                                                            -o single.x;
                             -04 -qhot -qstrict -qsmp=omp -o mpi.x;
mpixlcxx_r implicit_mpi.cpp
compile_XL_05:
mpixlcxx_r implicit_single.cpp -05 -qhot -qstrict
mpixlcxx_r implicit_mpi.cpp     -05 -qhot -qstrict -qsmp=omp -o mpi.x;
mpisubmit.bg -n 1 ./single.x data --stdout=1.out --stderr=1.err
run_omp_2:
mpisubmit.bg -n 2 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=2.out --stderr=2.err
run omp 4:
mpisubmit.bg -n 4 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=4.out --stderr=4.err
mpisubmit.bg -n 8 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=8.out --stderr=8.err
run_omp_16:
mpisubmit.bg -n 16 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=16.out --stderr=16.err
run_omp_32:
mpisubmit.bg -n 32 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=32.out --stderr=32.err
mpisubmit.bg -n 64 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=64.out --stderr=64.err
run_omp_128:
mpisubmit.bg -n 128 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=128.out --stderr=128.err
mpisubmit.bg -n 256 -w 00:10:00 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=256.out --stderr=256.err
run_omp_512:
mpisubmit.bg -n 512 -w 00:05:00 ./mpi.x data -mode SMP -env OMP_NUM_THREADS=4 --stdout=512.out --stderr=512.err
cancel:
llcancel -u edu-cmc-ski16-052
rm *.out *.err core* 2>/dev/null || echo > /dev/null
llq -u edu-cmc-ski16-052
accel:
perl accel.pl
run1_16_03:
make cancel
make clean
make compile_XL_03
make run_omp_1
for i in 2 4 8 16; do make run_omp_$$i; done;
run1 16 04:
make cancel
```

```
make clean
make compile_XL_04
make run_omp_1
for i in 2 4 8 16; do make run_omp_$$i; done;
run1_16_05:
make cancel
make clean
make compile_XL_05
make run_omp_1
for i in 2 4 8 16; do make run_omp_$$i; done;
run1_32_03:
make cancel
make clean
make compile_XL_03
make run_omp_1
for i in 2 4 8 16 32; do make run_omp_$$i; done;
run1_32_04:
make cancel
make clean
make compile_XL_04
make run_omp_1
for i in 2 4 8 16 32; do make run_omp_$$i; done;
run1_32_05:
make cancel
make clean
make compile_XL_05
make run_omp_1
for i in 2 4 8 16 32; do make run_omp_$$i; done;
run1_64_03:
make cancel
make clean
make compile_XL_03
make run_omp_1
for i in 2 4 8 16 32 64; do make run_omp_$$i; done;
run1_64_04:
make cancel
make clean
make compile_XL_04
make run_omp_1
for i in 2 4 8 16 32 64; do make run_omp_$$i; done;
run1_64_05:
make cancel
make clean
make compile_XL_05
make run_omp_1
for i in 2 4 8 16 32 64; do make run_omp_$$i; done;
run1_128_03:
make cancel
make clean
make compile_XL_03
make run_omp_1
for i in 2 4 8 16 32 64 128; do make run_omp_$$i; done;
run1_128_04:
make cancel
make clean
make compile_XL_04
make run_omp_1
for i in 2 4 8 16 32 64 128; do make run_omp_$$i; done;
run1_128_05:
make cancel
make clean
```

```
make compile_XL_05
make run_omp_1
for i in 2 4 8 16 32 64 128; do make run_omp_$$i; done;
run1_256_03:
make cancel
make clean
make compile_XL_03
make run_omp_1
for i in 2 4 8 16 32 64 128 256; do make run_omp_$$i; done;
run1_256_04:
make cancel
make clean
make compile_XL_04
make run_omp_1
for i in 2 4 8 16 32 64 128 256; do make run_omp_$$i; done;
run1_256_05:
make cancel
make clean
make compile_XL_05
make run_omp_1
for i in 2 4 8 16 32 64 128 256; do make run_omp_$$i; done;
run1_512_03:
make cancel
make clean
make compile_XL_03
make run_omp_1
for i in 2 4 8 16 32 64 128 256 512; do make run_omp_$$i; done;
run1_512_04:
make cancel
make clean
make compile_XL_04
make run_omp_1
for i in 2 4 8 16 32 64 128 256 512; do make run_omp_$$i; done;
run1_512_05:
make cancel
make clean
make compile_XL_05
make run_omp_1
for i in 2 4 8 16 32 64 128 256 512; do make run_omp_$$i; done;
```

Скрипт вывода времени работы, ускорения и эффективности вычислений для серии запусков (парсер выходных файлов):

```
accel.pl
```

```
use strict;
use warnings;
my coeff = 80;
$coeff = $ARGV[0] if defined $ARGV[0];
my @files = ();
files[_] = (2 ** _] for 0..10;
my %time = ();
my %accel = ();
my %param = ();
my %abs_err = ();
my %rel_err = ();
for my $name(@files){
           my $fh;
            if(name == 1){
                       open($fh, "<", $name.".out") or last;</pre>
            }else{
                       open($fh, "<", $name.".out") or next;
            while(<$fh>){
                       if(/^Time:\s+(.+)$/){
                                   time{mame} = 1;
                       }elsif(/^Abs. error:\s+(.+)$/){
                                   abs_err{name} = 1;
                       }elsif(/^Rel. error:\s+(.+)$/){
                                   $rel_err{$name} = $1;
                       if(/^0:\s+(.+):\s+(.+)$/){
                                   $param{$1} = {} unless exists($param{$1});
                                   $param{$1}->{$name} = $2;
           }
            close($fh);
}
if(keys %time){
           return unless defined $time{1};
            for(keys %time){
                        $accel{$_} = $time{1} / $time{$_};
                        \c = 1;
            printf("%-10s %-15s %-25s %-25s %-25s %-20s %-20s \n", "Nodes", "Cores", "Time", "Speed Up", "Efficiency, %", "Abs. error", '
            for(sort {$a <=> $b} keys %time){
                       if(\$\_ == 1){
                                 printf("%-10s %-15s %-25s %-25s %-20s %-20s \n", $_, $_, $time{$_}, $accel{$_}, ($accel{$_} / $_ * 100), $abs_e
                                  printf("%-10s \%-15s \%-25s \%-25s \%-25s \%-20s \%-20s \%-30s \%-
           }
```

```
}else{
   print "No tasks completed\n";
print "n";
if(keys %param){
    for my $i(sort {$a cmp $b} keys %param){
       return unless defined $param{$i}{1};
       my $h = $param{$i};
       my $show = 0;
        for(keys %{$h}){
           next if h->\{_\} == 0;
            if(h->\{1\} / h->\{\{\}\} < \ * \ *coeff / 100)\{
                show = 1;
       }
        if($show){
           printf("%-10s %-25s %-25s", "np", $i, "%");
           print "\n";
            for(sort {$a <=> $b} keys %{$h}){
                printf("%-10s %-25s %-25s\n", $_, $h->{$_}, $h->{1} / $h->{$_});
           print "\n";
       }
   }
}
```

### Пример использования:

make run1\_64\_03 (очистка очереди, компиляция, запуск на 1, 2, 4, 8, 16, 32, 64 ядра) make accel (вывод результатов)

#### 10. Приложение

#### Листинг программы

#### 1) Последовательная версия

```
implicit single.cpp
#include <iostream>
#include <iomanip>
#include <cassert>
#include <cstdlib>
#include <cmath>
#include <ctime>
#include <mpi.h>
using namespace std;
// -----
char* filename;
// -----
double a1, b1;
double a2, b2;
double t_max;
int nx, ny, nt;
double dx, dy, dt;
double dx2, dy2;
double dt_1_2;
double* x;
double* y;
double* t;
double** u;
double** u_half;
// -----
double *A_i, *B_i, *C_i;
double *A_j, *B_j, *C_j;
// -----
double* F_i;
double* F_j;
double* X_i;
double* X_j;
void print_grid();
void print_x();
void print_y();
void print_t();
void print_u();
void print_u_half();
void print_u_exact(int);
void print_error();
void free_memory();
// BEGIN----- EXACT SOLUTION ------
double u_exact(double x, double y, double t){
   // return exp(-t) * sin(M_PI * x) * cos(M_PI * y);
   return exp(-t) * (sin(M_PI * x) + cos(M_PI * y)) + 10;
// END----- EXACT SOLUTION ------
```

```
// BEGIN----- F(x, y, t) ------
double F(double x, double y, double t){
  // return exp(-t) * sin(M_PI * x) * cos(M_PI * y) * (2 * M_PI * M_PI - 1);
  return - exp(-t) * (sin(M_PI * x) + cos(M_PI * y)) * (1 - M_PI * M_PI);
// END----- F(x, y, t) -------
// BEGIN----- READ DATA -----
void read_data(){
  FILE* f;
  f = fopen(filename, "r");
  if(f == NULL){
     cout << "Cannot open file \"" << filename << "\"" << endl;</pre>
      exit(1);
  int count = fscanf(f, "%lf %lf %lf %lf %lf %d %d %d", &a1, &b1, &a2, &b2, &t_max, &nx, &ny, &nt);
  if(count < 8){
     cout << "Wrong data format in file \"" << filename << "\"" << endl << endl;
     cout << "Usage: a1 b1 a2 b2" << endl;</pre>
                 t_max" << endl;
     cout << "
     cout << "
                 nx ny nt" << endl;</pre>
     cout << endl;</pre>
     cout << "Example: 0 1 0 1" << endl;</pre>
               1" << endl;
     cout << "
     cout << "
                  10 10 5" << endl;
     cout << endl;</pre>
     exit(1);
  }
  fclose(f);
// END----- READ DATA ------
\label{eq:const_double *_A, const_double *_B, const_double *_B, const_double *_C, double *f, double *x) {} \\
  double A[n];
  double B[n];
  double C[n];
  for(int i = 0; i < n; i++){
     A[i] = _A[i];
     B[i] = _B[i];
C[i] = _C[i];
  // -----
  double coeff;
  for(int i = 1; i < n; i++){
     coeff = A[i] / B[i-1];
     B[i] -= coeff * C[i-1];
     f[i] -= coeff * f[i-1];
  x[n-1] = f[n-1] / B[n-1];
  for(int i = n-2; i \ge 0; i--){
     x[i] = (f[i] - C[i] * x[i+1]) / B[i];
// END----- THOMAS ALGORITHM ------
```

```
// BEGIN----- SET TRIAG COEFF -----
void set_triag_coeff(){
   A_i = new double[ny];
   B_i = new double[ny];
   C_i = new double[ny];
   for(int i = 0; i < ny; i++){</pre>
      A_i[i] = -1.0 / dy2;
      B_i[i] = 2.0 / dt + 2.0 / dy2;
      C_i[i] = -1.0 / dy2;
   }
   A_i[0] = 0;
   C_i[ny-1] = 0;
   A_j = new double[nx];
   B_j = new double[nx];
   C_j = new double[nx];
   for(int i = 0; i < nx; i++){
      A_j[i] = -1.0 / dx2;
      B_j[i] = 2.0 / dt + 2.0 / dx2;
C_j[i] = -1.0 / dx2;
   A_{j}[0] = 0;
   C_j[nx-1] = 0;
// END----- SET TRIAG COEFF -----
// BEGIN----- SET GRID -----
void set_grid(){
   read_data();
   assert(a1 < b1);
   assert(a2 < b2);
   assert(t_max > 0);
   assert(nx > 0);
   assert(ny > 0);
   assert(nt > 0);
   dx = (b1 - a1) / nx;
   dy = (b2 - a2) / ny;
   dt = t_max / nt;
   dx2 = dx * dx;
   dy2 = dy * dy;
   dt_1_2 = dt / 2.0;
   try{
      x = new double[nx+1];
      y = new double[ny+1];
      t = new double[nt+1];
      u = new double*[ny+1];
      u_half = new double*[ny+1];
      for(int i = 0; i <= ny; i++){
          u[i] = new double[nx+1];
          u_half[i] = new double[nx+1];
          for(int j = 0; j \le nx; j++){
             u[i][j] = 0;
             u_half[i][j] = 0;
      }
   }catch(std::bad_alloc){
      cout << "Cannot allocate memory" << endl;</pre>
```

```
exit(1):
  }
  for(int i = 0; i \le ny; i++){
     y[i] = a2 + i * dy;
  for(int j = 0; j \le nx; j++){
     x[j] = a1 + j * dx;
  }
  for(int k = 0; k \le nt; k++){
     t[k] = k * dt;
}
// END----- SET GRID ------
// BEGIN----- INITIAL CONDITIONS -----
double g_0(int j, int i){
  return u_exact(x[j], y[i], t[0]);
void set_initials(){
  for(int i = 0; i <= ny; i++){
     for(int j = 0; j <= nx; j++){
        u[i][j] = g_0(j, i);
  }
}
// END----- INITIAL CONDITIONS -----
// BEGIN----- BOUNDARY CONDITIONS -----
void set_boundaries(double** _u, double _t){
  for(int i = 0; i <= ny; i++){
      _u[i][0] = u_exact(x[0], y[i], _t);
      _u[i][nx] = u_exact(x[nx], y[i], _t);
  }
  for(int j = 0; j \le nx; j++){
     u[0][j] = u_exact(x[j], y[0], _t);
      _u[ny][j] = u_exact(x[j], y[ny], _t);
// END----- BOUNDARY CONDITIONS -----
// BEGIN-----
              ------ SOLVE ------
void solve(){
  set_triag_coeff();
  F_i = new double[ny-1];
  F_j = new double[nx-1];
  X_i = new double[ny-1];
  X_j = \text{new double[nx-1]};
  set_initials();
  clock_t t1 = clock();
  for(int k = 0; k < nt; k++){
     double t_half = t[k] + dt_1_2;
     double t_k = t_half + dt_1_2;
      // BEGIN----- x-direction ------
     set_boundaries(u_half, t_half);
     for(int i = 1; i < ny; i++){
        for(int j = 1; j < nx; j++){
           F_{j}[j-1] = (u[i+1][j] - 2 * u[i][j] + u[i-1][j]) / dy2 + F(x[j], y[i], t_half) + 2 * u[i][j] / dt;
```

```
if(j == 1){
               F_{j[j-1]} += u_{exact}(x[j-1], y[i], t_{half}) / dx2;
            else if(j == nx-1){
               F_j[j-1] += u_exact(x[j+1], y[i], t_half) / dx2;
         }
         solveMatrix(nx-1, A_j, B_j, C_j, F_j, X_j);
         for(int j = 0; j < nx-1; j++){
            u_half[i][j+1] = X_j[j];
      }
      // END----- x-direction ------
      // BEGIN----- y-direction ------
      set_boundaries(u, t_k);
      for(int j = 1; j < nx; j++){
         for(int i = 1; i < ny; i++){
           F_i[i-1] = (u_half[i][j-1] - 2 * u_half[i][j] + u_half[i][j+1]) / dx2 + F(x[j], y[i], t_half) + 2 * u_half[i][j]
         F_i[0] += u_exact(x[j], y[0], t_k) / dy2;
         F_i[ny-2] += u_exact(x[j], y[ny], t_k) / dy2;
         solveMatrix(ny-1, A_i, B_i, C_i, F_i, X_i);
         for(int i = 0; i < ny-1; i++){
            u[i+1][j] = X_i[i];
      }
      // END------ y-direction ------
   clock_t t2 = clock();
   cout << "Time: " << (double)(t2 - t1) / (double)CLOCKS_PER_SEC << endl << endl;</pre>
// END----- SOLVE ------
int main(int argc, char** argv){
   cout << fixed << setprecision(8);</pre>
   if(argc < 2){
      cout << "Usage: ./a.out data" << endl;</pre>
      exit(1);
   }
   filename = argv[1];
   set_grid();
   solve();
   print_error();
   free_memory();
   return 0;
}
// BEGIN-----
             ------ PRINT GRID ------
void print_x(){
   cout << "x: ";
```

```
for(int j = 0; j \le nx; j++){
       cout << x[j] << " ";
   cout << endl << endl;</pre>
}
void print_y(){
   cout << "y: ";
   for(int i = 0; i \le ny; i++){
       cout << y[i] << " ";
   cout << endl << endl;</pre>
}
void print_t(){
   cout << "t: ";
   for(int k = 0; k \le nt; k++){
       cout << t[k] << " ";
   cout << endl << endl;</pre>
}
void print_grid(){
   print_x();
   print_y();
   print_t();
// END----- PRINT GRID -----
// BEGIN------ PRINT SOLUTION ------
void print_u(){
   cout << "u:" << endl;</pre>
   for(int i = 0; i <= ny; i++){
       for(int j = 0; j \le nx; j++){
           cout << setw(12) << setprecision(8) << u[i][j] << " ";</pre>
       cout << endl;</pre>
   }
   cout << endl;</pre>
}
void print_u_half(){
   cout << "u_half:" << endl;</pre>
   for(int i = 0; i <= ny; i++){
       for(int j = 0; j \le nx; j++){
           \verb|cout| << \verb|setw(12)| << \verb|setprecision(9)| << \verb|u_half[i][j]| << " ";
       cout << endl;</pre>
   }
   cout << endl;</pre>
}
void print_u_exact(int k){
   for(int i = 0; i <= ny; i++){
       for(int j = 0; j \le nx; j++){
           cout << u_exact(x[j], y[i], t[k]) << " ";
       cout << endl;</pre>
   }
```

```
cout << endl;</pre>
}
// END----- PRINT SOLUTION -----
// BEGIN-----PRINT ERROR ------
void print_error(){
  double max_abs_error = 0;
  double max_rel_error = 0;
  for(int i = 1; i < ny; i++){
     for(int j = 1; j < nx; j++){
        double sln = u[i][j];
        double sln_e = u_exact(x[j], y[i], t[nt]);
        if(abs(sln_e) < 1e-10 \mid | abs(sln) < 1e-10){
           continue;
        double abs_error = abs(sln - sln_e);
        double rel_error = abs_error / abs(sln_e);
        max_abs_error = max(max_abs_error, abs_error);
        max_rel_error = max(max_rel_error, rel_error);
     }
  }
  cout << "Abs. error: " << setprecision(8) << max_abs_error << endl;</pre>
  cout << "Rel. error: " << setprecision(8) << max_rel_error << endl;</pre>
}
// END-----PRINT ERROR ------
// BEGIN------ FREE MEMORY ------
void free_memory(){
  // -----
  delete[] x;
  delete[] y;
  delete[] t;
  // -----
  delete[] A_i;
  delete[] B_i;
  delete[] C_i;
  delete[] A_j;
  delete[] B_j;
  delete[] C_j;
  // -----
  for(int i = 0; i <= ny; i++){
     delete[] u[i];
     delete[] u_half[i];
  }
  delete[] u;
  delete[] u_half;
  delete[] F_i;
  delete[] F_j;
  delete[] X_i;
  delete[] X_j;
  // -----
// END----- FREE MEMORY ------
```

#### 2) Параллельная версия

```
implicit mpi.cpp
#include <iostream>
#include <iomanip>
#include <cassert>
#include <cstdlib>
#include <cmath>
#include <mpi.h>
#include "thomas_x.cpp"
#include "thomas_y.cpp"
#include "thomas_y_helper.h"
using namespace std;
// -----
char* filename;
// -----
double a1, b1;
double a2, b2;
double t_max;
int nx, ny, nt;
double dx, dy, dt;
double dx2, dy2;
double dt_1_2;
double* x;
double* y;
double* t;
double** u;
double** u_half;
// -----
double *A_i, *B_i, *C_i;
double *A_j, *B_j, *C_j;
// -----
int mpirank;
int mpisize;
bool rank_0;
bool rank_np;
int src_up, src_dn;
int dst_up, dst_dn;
int snd_up_tag, snd_dn_tag;
int rcv_up_tag, rcv_dn_tag;
MPI_Datatype row, row_1, column;
// -----
int ny_per_node;
int ny_tail;
int my_ny;
int row_0_ny;
int row_n_ny;
double** u_up;
double** u_down;
double** u_left;
double** u_right;
// -----
void print_grid();
```

```
void print_x();
void print_y();
void print_t();
void print_u();
void print_u_half();
void print_u_exact(int);
void print_error();
void free_memory();
// -----
// BEGIN------ EXACT SOLUTION ------
double u_exact(double x, double y, double t){
  // return exp(-t) * sin(M_PI * x) * cos(M_PI * y);
  return exp(-t) * (sin(M_PI * x) + cos(M_PI * y)) + 10;
// END----- EXACT SOLUTION ------
// BEGIN------ F(x, y, t) ------
double F(double x, double y, double t){
  // return exp(-t) * sin(M_PI * x) * cos(M_PI * y) * (2 * M_PI * M_PI - 1);
  return - exp(-t) * (sin(M_PI * x) + cos(M_PI * y)) * (1 - M_PI * M_PI);
// END----- F(x, y, t) ------
// BEGIN----- READ DATA ------
void read_data(){
  FILE* f;
  f = fopen(filename, "r");
  if(f == NULL){
      cout << "Cannot open file \"" << filename << "\"" << endl;</pre>
      exit(1);
  7
  int count = fscanf(f, "%lf %lf %lf %lf %lf %d %d %d", &a1, &b1, &a2, &b2, &t_max, &nx, &ny, &nt);
  if(count < 8){
      cout << "Wrong data format in file \"" << filename << "\"" << endl << endl;</pre>
     cout << "Usage: a1 b1 a2 b2" << endl;</pre>
     cout << "
                 t_max" << endl;
     cout << "
                 nx ny nt" << endl;</pre>
     cout << endl;</pre>
     cout << "Example: 0 1 0 1" << endl;</pre>
     cout << "
                  1" << endl;
     cout << "
                   10 10 5" << endl;
     cout << endl;</pre>
     exit(1);
  }
  fclose(f);
// END----- READ DATA -----
// BEGIN------ SET TRIAG COEFF ------
void set_triag_coeff(){
  A_i = new double[ny];
  B_i = new double[ny];
  C_i = new double[ny];
  for(int i = 0; i < ny; i++){
      A_i[i] = -1.0 / dy2;
      B_i[i] = 2.0 / dt + 2.0 / dy2;
      C_i[i] = -1.0 / dy2;
```

```
C_i[ny-1] = 0;
   A_j = new double[nx];
   B_j = new double[nx];
   C_j = new double[nx];
   for(int i = 0; i < nx; i++){
       A_j[i] = -1.0 / dx2;
      B_{j}[i] = 2.0 / dt + 2.0 / dx2;
      C_{j}[i] = -1.0 / dx2;
   C_j[nx-1] = 0;
}
// END----- SET TRIAG COEFF -----
// BEGIN----- SET GRID -----
void set_grid(){
   read_data();
   assert(a1 < b1);</pre>
   assert(a2 < b2);</pre>
   assert(t_max > 0);
   assert(nx > 0);
   assert(ny > 0);
   assert(nt > 0);
   dx = (b1 - a1) / nx;
   dy = (b2 - a2) / ny;
   dt = t_max / nt;
   dx2 = dx * dx;
   dy2 = dy * dy;
   dt_1_2 = dt / 2.0;
   // -----
   ny_per_node = ((ny + 1) + 2 * (mpisize-1)) / mpisize;
   ny_tail = ((ny + 1) + 2 * (mpisize-1)) % mpisize;
   my_ny = ny_per_node - 1;
   if(mpirank < ny_tail)</pre>
      my_ny++;
   // -----
   try{
      x = new double[nx+1];
      y = new double[my_ny+1];
      t = new double[nt+1];
      u = new double*[my_ny+1];
      u_half = new double*[my_ny+1];
      for(int i = 0; i <= my_ny; i++){
          u[i] = new double[nx+1];
          u_half[i] = new double[nx+1];
          for(int j = 0; j <= nx; j++){
             u[i][j] = 0;
             u_half[i][j] = 0;
      }
   }catch(std::bad_alloc){
      cout << "Cannot allocate memory" << endl;</pre>
   }
   double y0 = 0;
```

```
if(mpisize == 1){
     y0 = a2;
   }else if(!rank_0){
     int _y0 = 0;
      if(ny_tail){
        _y0 = (mpirank <= ny_tail) ? mpirank : ny_tail;</pre>
     y0 = a2 + ((ny_per_node-2) * (mpirank) + _y0) * dy;
  }
   for(int i = 0; i <= my_ny; i++){</pre>
     y[i] = y0 + i * dy;
   // -----
   for(int j = 0; j \le nx; j++){
     x[j] = a1 + j * dx;
   for(int k = 0; k \le nt; k++){
     t[k] = k * dt;
}
// END----- SET GRID -----
                    ------ INITIAL CONDITIONS ------
double g_0(int j, int i){
   return u_exact(x[j], y[i], t[0]);
void set_initials(){
   int i_1 = (rank_0) ? 0 : 1;
   int i_2 = (rank_np) ? (my_ny) : (my_ny-1);
   for(int i = i_1; i <= i_2; i++){
     for(int j = 0; j \le nx; j++){
        u[i][j] = g_0(j, i);
  }
}
// END----- INITIAL CONDITIONS -----
// BEGIN----- BOUNDARY CONDITIONS -----
void set_boundaries(double** _u, int k){
   if(!rank_0){
      #pragma omp parallel for
      for(int j = 0; j \le nx; j++){
        _u[my_ny][j] = u_down[k][j];
  }
   if(!rank_np){
     #pragma omp parallel for
     for(int j = 0; j \le nx; j++){
         u[0][j] = u_up[k][j];
     }
  }
   #pragma omp parallel for
   for(int i = 0; i <= my_ny; i++){
      _u[i][0] = u_left[k][i];
      _u[i][nx] = u_right[k][i];
   }
}
              ----- BOUNDARY CONDITIONS -----
// BEGIN----- MPI_PREPARE ------
void mpi_prepare(){
   MPI_Type_contiguous(nx, MPI_DOUBLE, &row);
```

```
MPI_Type_commit(&row);
   MPI_Type_contiguous(nx-1, MPI_DOUBLE, &column);
   MPI_Type_commit(&column);
   MPI_Type_vector(nx-1, 1, mpisize, MPI_DOUBLE, &row_1);
   MPI_Type_commit(&row_1);
   src_dn = dst_dn = mpirank+1;
   src_up = dst_up = mpirank-1;
   snd_up_tag = 1000 * dst_up + 2;
   snd_dn_tag = 1000 * dst_dn + 0;
   rcv_up_tag = 1000 * mpirank + 0;
   rcv_dn_tag = 1000 * mpirank + 2;
// END------ MPI_PREPARE ------
// BEGIN------ PREPARE_BOUNDARIES ------
void prepare_boundaries(){
   u_up = new double*[(nt+1)*2];
   u_down = new double*[(nt+1)*2];
   u_left = new double*[(nt+1)*2];
   u_right = new double*[(nt+1)*2];
   for(int k = 0; k \le nt*2; k++){
      u_up[k] = new double[nx+1];
      u_down[k] = new double[nx+1];
      u_left[k] = new double[my_ny+1];
      u_right[k] = new double[my_ny+1];
   for(int k = 0; k \le nt; k++){
      for(int j = 0; j < nx; j++){
          u_up[k*2][j] = u_exact(x[j], y[0], t[k]);
         u_down[k*2][j] = u_exact(x[j], y[my_ny], t[k]);
         if(k < nt){
             u_up[k*2+1][j] = u_exact(x[j], y[0], t[k]+dt_1_2);
             u_down[k*2+1][j] = u_exact(x[j], y[my_ny], t[k]+dt_1_2);
      }
      for(int i = 0; i <= my_ny; i++){
          u_left[k*2][i] = u_exact(x[0], y[i], t[k]);
         u_right[k*2][i] = u_exact(x[nx], y[i], t[k]);
         if(k < nt){
             u_{\text{left}[k*2+1][i]} = u_{\text{exact}(x[0], y[i], t[k]+dt_1_2)};
             u_right[k*2+1][i] = u_exact(x[nx], y[i], t[k]+dt_1_2);
      }
   }
}
// END----- PREPARE_BOUNDARIES -----
void solve(){
   mpi_prepare();
   // -----
   set_triag_coeff();
   double* F_j = new double[nx-1];
   double* X_j = new double[nx-1];
   prepare_boundaries();
   set_initials();
```

```
// -----
double** A_nx_i = new double*[nx];
double** B_nx_i = new double*[nx];
double** C_nx_i = new double*[nx];
double** F_nx_i = new double*[nx];
double** X_nx_i = new double*[nx];
for(int i = 0; i < nx; i++){
   A_nx_i[i] = new double[my_ny];
   B_nx_i[i] = new double[my_ny];
   C_nx_i[i] = new double[my_ny];
   F_nx_i[i] = new double[my_ny];
   X_nx_i[i] = new double[my_ny+1];
}
// -----
Helper* helper = new Helper(mpirank, mpisize, nx, my_ny, int(y[0] / dy));
MPI_Barrier(MPI_COMM_WORLD);
double t1 = MPI_Wtime();
for(int k = 0; k < nt; k++){
   double t_half = t[k] + dt_1_2;
   double t_k = t_half + dt_1_2;
   // BEGIN----- x-direction ------
   MPI_Status st1, st2;
   MPI_Request req_snd_up, req_snd_dn;
   MPI_Request req_rcv_up, req_rcv_dn;
   if(mpisize != 1){
       if(!rank_0){
          MPI_Irecv(&u[0][1], 1, row, src_up, rcv_up_tag, MPI_COMM_WORLD, &req_rcv_up);
       }
       if(!rank_np){
          MPI_Irecv(&u[my_ny][1], 1, row, src_dn, rcv_dn_tag, MPI_COMM_WORLD, &req_rcv_dn);
   }
   set_boundaries(u_half, 2*k+1);
   if(mpisize != 1){
       if(!rank 0){
          MPI_Isend(&u[1][1], 1, row, dst_up, snd_up_tag, MPI_COMM_WORLD, &req_snd_dn);
       if(!rank_np){
          MPI_Isend(&u[my_ny-1][1], 1, row, dst_dn, snd_dn_tag, MPI_COMM_WORLD, &req_snd_up);
   }
   for(int i = 2; i < my_ny-1; i++){
       u_half = thomas_x(u, A_j, B_j, C_j, F_j, X_j, k, i, t_half);
   }
   if(!rank_0){
       MPI_Wait(&req_rcv_up, &st1);
   u_{half} = thomas_x(u, A_j, B_j, C_j, F_j, X_j, k, 1, t_{half});
   if(!rank_np){
       MPI_Wait(&req_rcv_dn, &st2);
   7
   u_half = thomas_x(u, A_j, B_j, C_j, F_j, X_j, k, my_ny-1, t_half);
```

```
// BEGIN------ y-direction ------
      set_boundaries(u, k*2+2);
      #pragma omp parallel for
      for(int _i = 0; _i < my_ny; _i++){</pre>
         A_nx_i[1][_i] = A_i[_i];
         B_nx_i[1][_i] = B_i[_i];
         C_nx_i[1][_i] = C_i[_i];
      }
      #pragma omp parallel for
      for(int j = 1; j < nx; j++){
         for(int i = 1; i < my_ny; i++){</pre>
            if(rank_0){
            F_nx_i[j][0] += u_up[2*k+2][j] / dy2;
         }else if(rank_np){
            F_nx_i[j][my_ny_2] += u_down[2*k+2][j] / dy2;
      }
      X_nx_i = thomas_y(A_nx_i, B_nx_i, C_nx_i, F_nx_i, X_nx_i, helper);
      #pragma omp parallel for
      for(int j = 1; j < nx; j++){
         for(int i = 0; i < my_ny-1; i++){</pre>
            u[i+1][j] = X_nx_i[j][i];
      }
      // END----- y-direction ------
   }
   MPI_Barrier(MPI_COMM_WORLD);
   double t2 = MPI_Wtime();
   if(rank_0){
      cout << "nprocs: " << mpisize << endl;</pre>
      cout << "Time: " << (t2 - t1) << endl << endl;</pre>
   }
   MPI_Barrier(MPI_COMM_WORLD);
// END-----
                   int main(int argc, char** argv){
   cout << fixed << setprecision(8);</pre>
   if(argc < 2){
      cout << "Usage: ./a.out data" << endl;</pre>
      exit(1);
   }
   filename = argv[1];
   MPI_Init(&argc, &argv);
   MPI_Comm_rank(MPI_COMM_WORLD, &mpirank);
   MPI_Comm_size(MPI_COMM_WORLD, &mpisize);
   rank_0 = (mpirank == 0);
   rank_np = (mpirank == mpisize-1);
   set_grid();
```

// END----- x-direction ------

```
solve();
   print_error();
   free_memory();
   MPI_Finalize();
   return 0;
}
// BEGIN----- PRINT GRID ------
void print_x(){
   MPI_Barrier(MPI_COMM_WORLD);
   if(rank_0){
       cout << "x: ";
       for(int j = 0; j \le nx; j++){
           cout << x[j] << " ";
       cout << endl << endl;</pre>
   }
   MPI_Barrier(MPI_COMM_WORLD);
}
void print_y(){
   MPI_Barrier(MPI_COMM_WORLD);
   for(int r = 0; r < mpisize; r++){
       if(r == mpirank){
           cout << "rank: " << r << endl;</pre>
           cout << "my_ny: " << my_ny << endl;</pre>
           cout << "y: ";
           for(int i = 0; i <= my_ny; i++){</pre>
               cout << y[i] << " ";
           cout << endl << endl;</pre>
       }
       MPI_Barrier(MPI_COMM_WORLD);
   MPI_Barrier(MPI_COMM_WORLD);
void print_t(){
   MPI_Barrier(MPI_COMM_WORLD);
   if(rank_0){
       cout << "t: ";
       for(int k = 0; k \le nt; k++){
           cout << t[k] << " ";
       cout << endl << endl;</pre>
   }
   MPI_Barrier(MPI_COMM_WORLD);
}
void print_grid(){
   print_x();
   print_y();
```

```
print_t();
}
// END----- PRINT GRID -----
// BEGIN------ PRINT SOLUTION ------
void print_u(){
   MPI_Barrier(MPI_COMM_WORLD);
   if(mpirank == 0){
       cout << "u:" << endl;</pre>
   }
   MPI_Barrier(MPI_COMM_WORLD);
   for(int r = 0; r < mpisize; r++){
       if(r == mpirank){
           for(int i = 0; i <= my_ny; i++){</pre>
               for(int j = 0; j \le nx; j++){
                  cout << setw(12) << setprecision(8) << u[i][j] << " ";</pre>
              cout << endl;</pre>
           }
           cout << endl;</pre>
       }
       MPI_Barrier(MPI_COMM_WORLD);
   MPI_Barrier(MPI_COMM_WORLD);
}
void print_u_half(){
   MPI_Barrier(MPI_COMM_WORLD);
   cout << setprecision(8);</pre>
   if(mpirank == 0){
       cout << "u_half:" << endl;
   MPI_Barrier(MPI_COMM_WORLD);
   for(int r = 0; r < mpisize; r++){
       if(r == mpirank){
           for(int i = 0; i <= my_ny; i++){
              for(int j = 0; j \le nx; j++){
                  cout << setw(12) << setprecision(9) << u_half[i][j] << " ";</pre>
              cout << endl;</pre>
           cout << endl;</pre>
       }
       MPI_Barrier(MPI_COMM_WORLD);
   }
   MPI_Barrier(MPI_COMM_WORLD);
}
void print_u_exact(int k){
   MPI_Barrier(MPI_COMM_WORLD);
   cout << setprecision(8);</pre>
   if(mpirank == 0){
       cout << "u_half:" << endl;</pre>
   }
```

```
MPI_Barrier(MPI_COMM_WORLD);
   for(int r = 0; r < mpisize; r++){
       if(r == mpirank){
           for(int i = 0; i <= my_ny; i++){
              for(int j = 0; j \le nx; j++){
                  cout << setw(12) << setprecision(9) << u_exact(x[j], y[i], t[k]) << " ";</pre>
              cout << endl;</pre>
           }
           cout << endl;</pre>
       MPI_Barrier(MPI_COMM_WORLD);
   }
   MPI_Barrier(MPI_COMM_WORLD);
}
// END----- PRINT SOLUTION -----
// BEGIN------ PRINT ERROR ------
void print_error(){
   double max_abs_error = 0;
   double max_rel_error = 0;
   int i1, i2;
   int j1 = 0, j2 = nx;
   if(mpirank == 0){
       i1 = 0;
       i2 = my_ny_1;
   }else if(mpirank == mpisize-1){
       i1 = 1;
       i2 = my_ny;
   }else{
       i1 = 1;
       i2 = my_ny_1;
   for(int i = i1; i <= i2; i++){
       for(int j = j1; j \le j2; j++){
           double sln = u[i][j];
           double sln_e = u_exact(x[j], y[i], t[nt]);
           if(abs(sln_e) < 1e-10 \mid | abs(sln) < 1e-9){
              continue;
           double abs_error = abs(sln - sln_e);
           double rel_error = abs_error / abs(sln_e);
           max_abs_error = max(max_abs_error, abs_error);
           max_rel_error = max(max_rel_error, rel_error);
       }
   }
   double* abs_err = new double[mpisize];
   double* rel_err = new double[mpisize];
   MPI_Gather(&max_abs_error, 1, MPI_DOUBLE, abs_err, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
   MPI_Gather(&max_rel_error, 1, MPI_DOUBLE, rel_err, 1, MPI_DOUBLE, 0, MPI_COMM_WORLD);
   if(mpirank == 0){
       double max_abs_error = 0;
       double max_rel_error = 0;
       for(int i = 0; i < mpisize; i++){</pre>
```

```
max_abs_error = max(abs_err[i], max_abs_error);
        max_rel_error = max(rel_err[i], max_rel_error);
     }
     cout << "Abs. error: " << setprecision(8) << max_abs_error << endl;</pre>
     cout << "Rel. error: " << setprecision(8) << max_rel_error << endl;</pre>
  }
}
// END------ PRINT ERROR ------
// BEGIN------ FREE MEMORY ------
void free_memory(){
  delete[] x;
  delete[] y;
  delete[] t;
  for(int i = 0; i <= my_ny; i++){</pre>
     delete[] u[i];
     delete[] u_half[i];
  delete[] u;
  delete[] u_half;
// END----- FREE MEMORY ------
```

```
thomas x.cpp
#ifndef _THOMAS_X_
#define _THOMAS_X_
extern int nx;
extern double** u_left;
extern double** u_right;
extern double dx2:
extern double dy2;
extern double* x;
extern double* y;
extern double dt;
extern double F(double x, double y, double t);
// BEGIN----- THOMAS ALGORITHM -----
void solveMMatrix(const int n, const double *_A, const double *_B, const double *_C, double *f, double *x){
   // -----
   double A[n];
   double B[n];
   double C[n];
   for(int i = 0; i < n; i++){
       A[i] = _A[i];
       B[i] = _B[i];
       C[i] = _C[i];
   }
   // -----
   double coeff;
   for(int i = 1; i < n; i++){
       coeff = A[i] / B[i-1];
       B[i] -= coeff * C[i-1];
       f[i] -= coeff * f[i-1];
   x[n-1] = f[n-1] / B[n-1];
   for(int i = n-2; i \ge 0; i--){
       x[i] = (f[i] - C[i] * x[i+1]) / B[i];
}
// END----- THOMAS ALGORITHM -----
extern double** u_half;
double** thomas_x(double** u, double* A_j, double* B_j, double* C_j, double* F_j, double* X_j, int k, int i, double t_half){
   #pragma omp parallel for
   for(int j = 1; j < nx; j++){
       F_{j}[j-1] = (u[i+1][j] - 2 * u[i][j] + u[i-1][j]) / dy2 + F(x[j], y[i], t_half) + 2 * u[i][j] / dt;
   F_{j}[0] += u_{ft}[2*k+1][i] / dx2;
   F_{j[nx-2]} += u_{right[2*k+1][i]} / dx2;
   solveMMatrix(nx-1, A_j, B_j, C_j, F_j, X_j);
   \hbox{\#pragma omp parallel for }
   for(int j = 0; j < nx-1; j++){
       u_half[i][j+1] = X_j[j];
   return u_half;
}
#endif
```

```
thomas y.cpp
#ifndef THOMAS_CPP_
#define THOMAS_CPP_
#include <iostream>
#include <cstring>
#include <mpi.h>
#include "thomas_y_helper.h"
using namespace std;
extern int mpirank;
extern int mpisize;
extern bool rank_0;
extern bool rank_np;
extern int my_ny;
extern int nx;
extern int src_up, src_dn;
extern int dst_up, dst_dn;
extern int snd_up_tag, snd_dn_tag;
extern int rcv_up_tag, rcv_dn_tag;
extern MPI_Datatype row_1, column;
void solveMatrix(int n, double *a, double *b, double *c, double *f, double *x){
   double m;
    int i_prev;
    for(int i = 1; i < n; i++){
       i_prev = i-1;
       m = a[i] / b[i_prev];
       b[i] -= m * c[i_prev];
       f[i] -= m * f[i_prev];
   x[n-1] = f[n-1] / b[n-1];
    for(int i = n-2; i \ge 0; i--){
       x[i] = (f[i] - c[i] * x[i+1]) / b[i];
}
void down(double** A, double** B, double** C, double** F_i, double** d_l, int i_down){
    double coeff;
    int i_next;
    for(int i = 0; i <= i_down; i++){</pre>
       i_next = i+1;
       // -----
       coeff = B[1][i];
       B[1][i] /= coeff;
       C[1][i] /= coeff;
        if(!rank_0)
           d_1[1][i] /= coeff;
       int J;
        #pragma omp parallel for private(J)
       for (J = 1; J < nx; J++){
           F_i[J][i] /= coeff;
```

```
coeff = A[1][i];
       B[1][i_next] -= C[1][i] * coeff;
       A[1][i] -= B[1][i] * coeff;
       if(!rank_0)
           d_l[1][i_next] -= d_l[1][i] * coeff;
       #pragma omp parallel for private(J)
       for(J = 1; J < nx; J++){
           F_i[J][i_next] -= F_i[J][i] * coeff;
       // -----
   }
   coeff = B[1][i_down];
   B[1][i_down] /= coeff;
   A[1][i_down] /= coeff;
   C[1][i_down] /= coeff;
   if(!rank_0)
       d_l[1][i_down] /= coeff;
   #pragma omp parallel for
   for(int J = 1; J < nx; J++){
       F_i[J][i_down] /= coeff;
}
void up(double** A, double** B, double** C, double** F_i, double** d_l, double** d_r, int i_down){
   double coeff;
   d_r[1][i_down-1] = C[1][i_down-1];
   C[1][i_down-1] = 0;
   int i_next;
   for(int i = i_down-2; i >= 0; i--){
       i_next = i+1;
       coeff = C[1][i];
       C[1][i] -= B[1][i_next] * coeff;
       if(!rank_0)
           d_1[1][i] -= d_1[1][i_next] * coeff;
       d_r[1][i] -= d_r[1][i_next] * coeff;
       #pragma omp parallel for
       for(int J = 1; J < nx; J++){
           F_i[J][i] -= F_i[J][i_next] * coeff;
   }
}
double** thomas_y(double** A_i, double** B_i, double** C_i, double** F_i, double** X_i, Helper* helper){
   // -----
   int M = helper->M;
   int H = helper->H;
   int W = helper->W;
   int i_0 = helper->i_0;
   int i_down = helper->i_down;
   double** d_l = helper->d_l;
```

```
double** d_r = helper->d_r;
helper->flush();
int* block_size = helper->block_size;
int* block_start = helper->block_start;
int* block_fin = helper->block_fin;
int* block_shift = helper->block_shift;
int block_size_max = helper->block_size_max;
double A[nx][H];
double C[nx][H];
double F[nx][H];
double triple_dn[nx][3];
double triple_up[nx][3];
double sln[mpisize][block_size_max][H];
MPI_Request* req_sln_rcv = helper->req_sln_rcv;
MPI_Request* req_sln_snd = helper->req_sln_snd;
MPI_Status* stat_sln = helper->stat_sln;
// -----
if(!rank_np){
      MPI_Irecv(helper->triple_dn, nx+2, MPI_DOUBLE, src_dn, rcv_dn_tag * 1000, MPI_COMM_WORLD, &(helper->req_triple_down_rcv_dn_tag * 1000, MPI_COMM_WORLD, &(he
for(int r = 0; r < mpisize; r++){
      MPI_Irecv(sln[r], block_size[r] * H, MPI_DOUBLE, r, (r+1), MPI_COMM_WORLD, &req_sln_rcv[r]);
       \label{eq:mpi_line} \mbox{MPI\_Irecv(\&C[1][r], 1, row\_1, r, r * 4, MPI\_COMM\_WORLD, \&helper->req\_diags\_rcv[r + mpisize]); } 
      MPI_Irecv(&F[1][r], 1, row_1, r, r * 5, MPI_COMM_WORLD, &helper->req_diags_rcv[r + mpisize * 2]);
}
//------
//-----
double* a_down_all = helper->a_down_all;
double* c_down_all = helper->c_down_all;
double* f_down_all = helper->f_down_all;
//-----
d_1[1][0] = A_i[1][0];
d_r[1][i_down] = B_i[1][i_down];
down(A_i, B_i, C_i, F_i, d_l, i_down);
up(A_i, B_i, C_i, F_i, d_l, d_r, i_down);
#pragma omp parallel for
for(int J = 1; J < nx; J++){
      helper->triag_snd_buf[J] = F_i[J][0];
helper->triag_snd_buf[nx] = d_1[1][0];
helper->triag_snd_buf[nx+1] = d_r[1][0];
if(!rank 0){
      MPI_Isend(helper->triag_snd_buf, nx+2, MPI_DOUBLE, dst_up, snd_up_tag * 1000, MPI_COMM_WORLD, &(helper->req_triags_snd)
int J;
#pragma omp parallel for private(J)
for(J = 1; J < nx; J++){
      for(int i = 0; i < M; i++){
             C_{i}[J][i] = C_{i}[1][i];
              d_1[J][i] = d_1[1][i];
```

```
d_r[J][i] = d_r[1][i];
   }
}
//-----
if(!rank_np){
   MPI_Wait(&(helper->req_triple_down_rcv), &(helper->stat_triple_down));
double A_i_1_down = A_i[1][i_down];
double B_i_1_down = B_i[1][i_down];
#pragma omp parallel for private(J)
for(J = 1; J < nx; J++){
   double coef;
   double* h = helper->triple_dn;
   // -----
   triple_dn[J][0] = h[nx];
   triple_dn[J][1] = h[nx+1];
   triple_dn[J][2] = h[J];
   A_{i[J][i_down]} = A_{i_1down}
   B_i[J][i_down] = B_i_1_down;
   coef = C_i[J][i_down];
   B_i[J][i_down] -= triple_dn[J][0] * coef;
   F_i[J][i_down] -= triple_dn[J][2] * coef;
   triple_dn[J][1] *= -coef;
   // -----
   coef = B_i[J][i_down];
   A_i[J][i_down] /= coef;
   B_i[J][i_down] /= coef;
   C_i[J][i_down] /= coef;
   F_i[J][i_down] /= coef;
   f_down_all[J-1] = F_i[J][i_down];
   if(!rank_0){
      d_l[J][i_down] /= coef;
      a_down_all[J-1] = d_l[J][i_down];
   }
   triple_dn[J][1] /= coef;
   c_down_all[J-1] = triple_dn[J][1];
   d_r[J][i_down] /= coef;
   // -----
for(int i = 0; i < mpisize; i++){</pre>
   MPI_Isend(a_down_all, 1, column, i, mpirank * 3, MPI_COMM_WORLD, &helper->req_diags_snd[i]);
   MPI_Isend(c_down_all, 1, column, i, mpirank * 4, MPI_COMM_WORLD, &helper->req_diags_snd[i + mpisize]);
   MPI_Isend(f_down_all, 1, column, i, mpirank * 5, MPI_COMM_WORLD, &helper->req_diags_snd[i + mpisize * 2]);
}
MPI_Waitall(mpisize * 3, helper->req_diags_rcv, helper->stat_recv);
//-----
#pragma omp parallel for private(J)
for(J = block_start[mpirank]; J < block_fin[mpirank]; J++){</pre>
   double B[H];
   double X[H];
   for(int i = 0; i < H; i++){
      B[i] = 1;
```

```
solveMatrix(H, A[J], B, C[J], F[J], X);
        for(int i = 0; i < H; i++){
            sln[mpirank][J - block_shift[mpirank]][i] = X[i];
   }
    for(int r = 0; r < mpisize; r++){
        MPI_Isend(sln[mpirank], block_size[mpirank] * H, MPI_DOUBLE, r, (mpirank+1), MPI_COMM_WORLD, &req_sln_snd[r]);
   int numdone;
    int _mpisize = mpisize;
    while(_mpisize--){
        MPI_Waitany(mpisize, req_sln_rcv, &numdone, stat_sln);
        int r = numdone;
        #pragma omp parallel for
        for(int J = block_start[r]; J < block_fin[r]; J++){</pre>
            if(rank_0){
                X_i[J][i_down] = sln[r][J-block_shift[r]][0];
                for(int i = M-2; i >= 0; i--){
                    X_i[J][i] = F_i[J][i] - X_i[J][i_down] * d_r[J][i];
            }else{
                double _l = sln[r][J-block_shift[r]][mpirank-1];
                double _r = sln[r][J-block_shift[r]][mpirank];
                X_i[J][i_down] = _r;
                for(int i = M-2; i >= 0; i--){
                    X_i[J][i] = F_i[J][i] - (d_l[J][i] * _l + d_r[J][i] * _r);
            }
       }
   }
    return X_i;
}
```

#endif

```
thomas y helper.h
#ifndef MPI_Helper_H
#define MPI_Helper_H
#include <mpi.h>
class Helper{
public:
   Helper(int _mpirank, int _mpisize, int _nx, int _my_ny, int _i_0){
      // -----
      mpirank = _mpirank;
      nx = _nx;
      my_ny = _my_ny;
      // -----
      M = my_ny_1;
      H = mpisize;
      W = H + 1;
      i_0 = _i_0;
      i_down = M-1;
      req_diags_rcv = new MPI_Request[mpisize*3];
      req_diags_snd = new MPI_Request[mpisize*3];
      stat_recv = new MPI_Status[mpisize*3];
      // -----
      req_sln_rcv = new MPI_Request[mpisize];
      req_sln_snd = new MPI_Request[mpisize];
      stat_sln = new MPI_Status[mpisize];
      // -----
      B = new double[H];
      X = new double[H];
      triag_snd_buf = new double[nx+2];
      triple_dn = new double[nx+2];
      // -----
      d_1 = new double*[nx];
      d_r = new double*[nx];
      for(int i = 0; i < nx; i++){
          d_l[i] = new double[my_ny-1];
          d_r[i] = new double[my_ny-1];
      // -----
      block_size = new int[mpisize];
      block_start = new int[mpisize];
      block_fin = new int[mpisize];
      block_shift = new int[mpisize];
      for(int r = 0; r < mpisize; r++){
          block_size[r] = nx / mpisize;
      block_size[mpisize-1] += nx % mpisize;
      block_start[0] = 1;
      for(int r = 1; r < mpisize; r++){
         block_start[r] = 0;
          for(int i = 0; i < r; i++){
             block_start[r] += block_size[i];
```

```
}
   block_fin[0] = block_size[0];
   for(int r = 1; r < mpisize; r++){
       block_fin[r] = block_start[r] + block_size[r];
   block_shift[0] = 0;
   for(int r = 1; r < mpisize; r++){
       block_shift[r] = block_start[r];
   block_size_max = 0;
   for(int r = 0; r < mpisize; r++){
       block_size_max = std::max(block_size_max, block_size[r]);
   // -----
   a_down_all = new double[nx-1];
   c_down_all = new double[nx-1];
   f_down_all = new double[nx-1];
   // -----
}
~Helper(){
   // -----
   delete[] req_diags_rcv;
   delete[] req_diags_snd;
   delete[] stat_recv;
   // -----
   delete[] req_sln_rcv;
   delete[] req_sln_snd;
   delete[] stat_sln;
   // -----
   delete[] triag_snd_buf;
   delete[] triple_dn;
   for(int i = 0; i < nx; i++){
       delete[] d_l[i];
       delete[] d_r[i];
   }
   delete[] d_1;
   delete[] d_r;
   // -----
   delete[] block_size;
   delete[] block_start;
   delete[] block_fin;
   delete[] block_shift;
   // -----
   delete[] a_down_all;
   delete[] c_down_all;
   delete[] f_down_all;
}
void flush(){
   for(int J = 1; J < nx; J++){
       for(int i = 0; i < M; i++){
          d_1[J][i] = d_r[J][i] = 0;
       }
   }
}
// -----
```

```
int nx;
   int my_ny;
   // -----
   int M;
   int H;
   int W;
   // -----
   int i_0;
   int i_down;
   // -----
   double** A;
   double* B;
   double** C;
   double** F;
   double* X;
   // -----
   MPI_Request* req_diags_snd;
   MPI_Request* req_diags_rcv;
   MPI_Status* stat_recv;
   // -----
   MPI_Request* req_sln_rcv;
   MPI_Request* req_sln_snd;
   MPI_Status* stat_sln;
   // -----
   double* triag_snd_buf;
   double* triple_dn;
   // -----
   double** d_l;
   double** d_r;
   // -----
   MPI_Request req_triple_down_rcv;
   MPI_Request req_triags_snd;
   MPI_Status stat_triple_down;
   // -----
   int* block_size;
   int* block_start;
   int* block_fin;
   int* block_shift;
   int block_size_max;
   // -----
   double* a_down_all;
   double* c_down_all;
   double* f_down_all;
   // -----
private:
   int mpirank;
   int mpisize;
```

#endif