

## **Parallel Computing**

Präsentiert von Jean Sokolov



2

#### **Inhaltsverzeichnis**

- Problemstellung
- Implementation in OpenMP
- Implementation in MPI



#### **PROBLEMSTELLUNG**

**3** 13.07.2023



## **Matrix-Multiplikation**

- Parallelisieren der Kalkulation
- In C++ mit OpenMP und MPI



#### **IMPLEMENTATION IN C++**



# Single-thread

```
1⊕ /*
     Name
                  : m2m serial.cpp
     Version
    Copyright : Your copyright notice
     Description : Matrix matrix multiplication
 9 */
 10 #include <stdio.h>
 11 #include <stdlib.h>
 13 double A[1000][1000];
 14 double B[1000][1000];
 15 double C[1000][1000];
 17⊖ void init m() {
 18 // initialise
 19
        for (int i = 0; i < 1000; ++i) {
 20
             for (int y = 0; y < 1000; ++y) {
                 A[i][y] = rand();
B[i][y] = rand();
 21
 22
 23
24
25 }
26
27⊖ void m2m() {
28
        for (int i = 0; i < 1000; i++) {
 29
            for (int j = 0; j < 1000; j++) {
                 for (int k = 0; k < 1000; k++) {
   C[i][j] += A[i][k] * B[k][j];</pre>
 30
34
                 printf("%.1f \n", C[i][j]);
35
36
37 }
 38
 39⊖ int main(int argc, char *argv[]) {
 40
        init_m();
 41
        m2m();
 42
 43
         return 0;
 44
 45
```



#### IMPLEMENTATION IN OPENMP



#### **Multi-thread**

```
c m2m_serial.cpp
                  1⊝ /*
 3 Name
                : main.cpp
 4 Author
 5 Version
 6 Copyright : Your copyright notice
 7 Description : Matrix matrix multiplication in OpenMP
10 #include <omp.h>
11 #include <stdio.h>
12 #include <stdlib.h>
14 double A[1000][1000];
15 double B[1000][1000];
16 double C[1000][1000];
17
18⊖ void init m() {
19 // initialise matrices
20
        for (int i = 0; i < 1000; ++i) {
21
            for (int y = 0; y < 1000; ++y) {
22
               A[i][y] = rand();
                B[i][y] = rand();
23
24
25
26 }
27
28@ int main(int argc, char *argv[]) {
29
        int i, j, k;
30
        init m();
31
32 #pragma omp parallel for private(i, j, k) shared(A,B,C)
33
        for (i = 0; i < 1000; i++) {
34
            for (j = 0; j < 1000; j++) {
35
                for (k = 0; k < 1000; k++) {
36
                   C[i][j] += A[i][k] * B[k][j];
 37
38
                printf("%.1f \n", C[i][j]);
39
40
41
        return Θ;
42
 43
```



### Laufzeit und Validierung

```
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$ time ./m2m openmp.o > outOpenMP
        0m1,424s
real
        0m7.452s
user
        0m1.700s
sys
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$ time ./m2m serial.o > outSerial
        0m3,655s
real
        0m3,600s
user
        0m0,056s
sys
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$ ls -l
insgesamt 50820
-rwxr-xr-x 1 isd isd 13336 Jun 11 15:50 m2m openmp.o
-rw-r--r-- 1 isd isd 829 Jun 11 15:50 m2m serial.cpp
-rwxr-xr-x 1 isd isd
                        8840 Jun 11 15:50 m2m serial.o
                         892 Jun 11 15:50 main.cpp
-rw-r--r-- 1 isd isd
-rw-r--r-- 1 isd isd 26000000 Jul 12 22:19 outOpenMP
-rw-r--r-- 1 isd isd 26000000 Jul 12 22:19 outSerial
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$ cat outOpenMP | grep 1184512763025494114304.0
1184512763025494114304.0
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$ cat outSerial | grep 1184512763025494114304.0
1184512763025494114304.0
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$ cmp outSerial outOpenMP
outSerial outOpenMP differieren: Byte 3, Zeile 1
isd@debian-eclipse:~/eclipse luna workspace/OpenMP/src$
```



#### **IMPLEMENTATION IN MPI**

**10** 13.07.2023



#### **Statische Prozesszahl**

```
10 #include "mpi.h"
11 #include <iostream>
12 using namespace std;
14 #define num 1000
15 #define USE_MPI_IN_PLACE 0
17 double A[num][num];
18 double B[num][num];
19 double C[num][num];
21 void init_m() {
        for (int i = 0; i < num; ++i) {
           for (int y = 0; y < num; ++y) {
   A[i][y] = rand();</pre>
                B[i][y] = rand();
       }
29 }
32⊖ int main(int argc, char *argv[]) {
      int n, rank, size, i, j, k;
        MPI::Init(argc, argv);
        size = MPI::COMM WORLD.Get size();
        rank = MPI::COMM_WORLD.Get_rank();
41
            if (rank==0) printf("Matrix size not divisible by number of processors\n");
            MPI Finalize();
43
            exit(-1);
44
45
46
        n = num/size:
        MPI::COMM_WORLD.Bcast(B, num*num, MPI::DOUBLE, 0);
        if (rank == 0)
            MPI::COMM_WORLD.Scatter(A, num*n, MPI::DOUBLE, MPI_IN_PLACE, num*n, MPI::DOUBLE, 0);
            MPI::COMM_WORLD.Scatter(A, num*n, MPI::DOUBLE, A[rank*n], num*n, MPI::DOUBLE, 0);
        for (i = n*rank; i < n*(rank+1); i++) {</pre>
            for (j = 0; j < num; j++) {
                for (k = 0; k < num; k++) {
                   C[i][j] += A[i][k] * B[k][j];
                printf("%.1f \n", C[i][j]);
61
            MPI::COMM WORLD.Gather(MPI IN PLACE, num*n, MPI::DOUBLE, C[n*rank], num*n, MPI::DOUBLE, 0):
63
            MPI::COMM_WORLD.Gather(C[n*rank], num*n, MPI::DOUBLE, C, num*n, MPI::DOUBLE, 0);
64
65
66
        MPI::Finalize():
67
        return 0;
68 }
```



### **Dynamische Prozesszahl**

```
#include "mpi.h"
     #include <iostre
     using namespace std;
     #define num 1000
     #define USE_MPI_IN_PLACE 0
     double A[num][num];
    double B[num][num];
double C[num][num];
  18⊖ void init_m() {
    // initialise

for (int i = 0; i < num; ++i) {

    for (int y = 0; y < num; ++y) {

        A[i][y] = rand();

        B[i][y] = rand();
 29@ int main(int argc, char *argv[]) {
30    int n, rank, size, i, j, k, proc;
          MPI::Init(argc, argv);
size = MPI::COMM_WORLD.Get_size();
rank = MPI::COMM_WORLD.Get_rank();
          while (num%proc!=0) {
    //if (rank==0) printf("Matrix size not divisible by number of processors\n");
              proc--;
              rrr_comm rew_comm;
MPI_comm_split(MPI_cOMM_WORLD, old_rankproc, old_rank, &new_comm);
int new_rank, new_size;
MPI_comm_rank(new_comm, &new_rank);
               MPT Comm new comm
              MPI_Comm_size(new_comm, &new_size);
MPI_Barrier(new_comm);
              //if (rank==055proc!=size)
//printf("Matrix cannot be divided by specified process count. Continuing with %i processes instead.\n", new_size);
               //printf("#%i here\n", old rank);
               n = num/proc;
MPI Bcast(B, num*num, MPI::DOUBLE, 0, new comm);
                        MPI_Scatter(A, num*n, MPI::DOUBLE, MPI_IN_PLACE, num*n, MPI::DOUBLE, 0, new_comm);
                        MPI_Scatter(A, num*n, MPI::DOUBLE, A[new_rank*n], num*n, MPI::DOUBLE, 0, new_comm);
                   //printf("%.1f \n", C[i][j]);
               if (new_rank<proc){
                        MPI_Gather(MPI_IN_PLACE, num*n, MPI::DOUBLE, C[n*new_rank], num*n, MPI::DOUBLE, 0, new_comm);
                   else
MPI_Gather(C[n*new_rank], num*n, MPI::DOUBLE, C, num*n, MPI::DOUBLE, 0, new_comm);
          //printf("#%i done \n", rank);
MPI::Finalize();
          return 0:
```

13.07.2023 **12** 



### Laufzeit und Validierung

```
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debug$ time sudo mpirun -np 8 ./m2m MPI > out3
real
        0m1,053s
        0m6,612s
user
        0m1,140s
sys
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debuq$ time sudo mpirun -np 7 ./m2m MPI > out3
        0m1.125s
real
        0m6,404s
user
        0m2.240s
SVS
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debug$ time sudo mpirun -np 6 ./m2m MPI > out3
        0m1,178s
real
        0m6,728s
user
        0m2.284s
SVS
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debug$ time sudo mpirun -np 3 ./m2m MPI > out3
        0m2,486s
real
        0m7,284s
user
        0m4,672s
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debug$ ls -l out3
-rw-r--r-- 1 isd isd 26000000 Jul 12 22:23 out3
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debug$ cat out3 | grep 1184512763025494114304.0
1184512763025494114304.0
isd@debian-eclipse:~/eclipse ptp workspace/m2m MPI/Debug$
```



### **Speichereffiziente Variation**

```
7 #include "mpi.h"
 8 #include <iostream>
 9 using namespace std;
11 #define num 1000
12 #define USE MPI IN PLACE 0
13
14 double A[num][num];
15 double B[num][num];
16 double C[num][num];
17 double tmpB[num][num];
19 MPI Status status;
21 void init m() {
22 // initialise
      for (int i = 0; i < num; ++i) {
           for (int y = 0; y < num; ++y) {
24
               A[i][v] = rand();
25
26
               B[i][y] = rand();
27
28
       }
29 }
30
31 void rotate b matrix(){
       for (int i = 0; i < num; ++i) {
33
           for (int y = 0; y < num; ++y) {
34
               tmpB[num-1-y][i]=B[i][y];
35
36
37
       for (int i = 0; i < num; ++i) {
38
           for (int y = 0; y < num; ++y) {
39
               B[i][y]=tmpB[i][y];
40
41
42 }
```



### **Speichereffiziente Variation**

```
440 int main(int argc, char *argv[]) {
45
       int n, rank, size, i, j, k, proc;
46
       init m();
47
       rotate b matrix();
48
49
       MPI::Init(argc, argv);
       MPI_Barrier(MPI_COMM_WORLD);
51
       size = MPI::COMM WORLD.Get size();
       rank = MPI::COMM WORLD.Get rank();
53
55
       while (num%proc!=0) {
            //if (rank==0) printf("Matrix size not divisible by number of processors\n");
57
58
59
       int old rank = rank;
60
           MPI Comm new comm;
61
            MPI_Comm_split(MPI_COMM_WORLD, old_rank<=proc, old_rank, &new_comm);</pre>
62
            int new rank, new size;
           MPI_Comm_rank(new_comm, &new_rank);
63
            MPI Comm size(new comm, &new size);
65
           MPI Barrier(new comm);
66
67
            n = num/proc;
68
            //printf("packet size=%i here\n using %i slaves\n", n, proc);
69
            if (new rank==0){
70
               for (int i = 1; i<=proc; i++){
71
                   for (int j = 0; j < num; j++){
                       for (int k = n*(i-1); k < n*i; k++) {
                           MPI_Send(&B[num-1-j], num, MPI_DOUBLE, i, 0, new_comm);
74
75
                       MPI_Send(&A[j], num, MPI_DOUBLE, i, 0, new_comm);
76
77
78
79
            if (new rank<=proc && new rank != 0){
               for (int i = n*(new_rank-1); i < n*new_rank; i++) {</pre>
80
81
                   MPI_Recv(&A[i], num, MPI_DOUBLE, 0, 0, new_comm, &status);
82
                    for (int j = 0; j < num; j++) {
                       MPI_Recv(&B[num-1-j], num, MPI_DOUBLE, 0, 0, new_comm, &status);
83
84
                        for (int k = 0; k < num; k++) {
85
                           C[i][j] += A[i][k] * B[num-1-j][k];
86
87
                       printf("%.1f \n", C[i][j]);
88
89
90
91
            MPI Barrier(new comm);
92
       MPI::Finalize();
93
        return Θ;
94 }
95
```



### Quellcode

 https://github.com/JeanSokolov/Parallel Computing



17

#### Vielen Dank für Ihre Aufmerksamkeit