Lab 3: Parallel Smith-Waterman with MPI

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Modified SW mpi.c:

```
67:
```

```
#include <mpi.h> // Use MPI ///ADDED mpi module
Here we include the mpi module to the program.
```

74:

```
#define min(a,b) (((a)<(b) ? (a) : (b)) /* Complete the value
returned when the condition evaluates to false.*/ //if a is smaller
return a, else return b (b is smaller)</pre>
```

We define min to return the smallest value of the 2 inputs.

165:

```
if ( rank==0 ) { /* Use process 0 as the master process
    and have it process the input */ //check that rank is 0
    /**** Error handling for input file ****/
```

285-291:

```
MPI_Bcast( &dim1, 1, MPI_INT, 0, MPI_COMM_WORLD); /* Broadcast dim1 */
///array, length of array, contents of array, root, comm
MPI_Bcast( &dim2, 1, MPI_INT, 0, MPI_COMM_WORLD); /* Broadcast dim2 */
///
MPI_Bcast( &BS, 1, MPI_INT, 0, MPI_COMM_WORLD); /* Broadcast BS */ //
MPI_Bcast( &DELTA, 1, MPI_INT, 0, MPI_COMM_WORLD); /* Broadcast DELTA
*/ ///
MPI_Bcast( sim, AA*AA, MPI_INT, 0, MPI_COMM_WORLD); /* Broadcast
matrix sim */ // AA is size of matrix

nrows = getRowCount(dim1, rank, nprocs); ///getRowCount(int
rowsTotal, int mpiRank, int mpiSize) / rowtotal is dim1, because dim1
is the length of the 1rst sequence
```

We broadcast these variables to all processes from root, and calculate the number of rows per process in case of not being able to distribute the rows among processes evenly.

295:

```
CHECK_NULL ((s1l = (short *) malloc ((short) * (dim1+1))));//create array s1l the size of dim1+1(real seq length), as it will contain the first sequence (dim1 can be exchanged by nrows) /clues from below Create an array of the appropriate size to store the first sequence.
```

303:

```
//MPI_Bcast( s1+1, dim1, MPI_SHORT, 0, MPI_COMM_WORLD );
```

This line divides and sends the contents of s1 into s11 across all the processes.

315:

```
MPI_Bcast( s2+1, dim2, MPI_SHORT, 0, MPI_COMM_WORLD); // broadcasts
contents of array s2 skipping first position and distributes to all
processes (MPI_COMM_WORLD)
```

This line is used to broadcast s2 to all processes.

326:

```
dim1local = (rank ? nrows : dim1); /* condition ? true_case :
false_case */ ///if rank!=0 dim1local=nrows(local row count), else:
dim1local= length of 1rst sequence(dim1)
```

Here we define dimllocal to be nrows if the process is 0, dimllocal will be diml, else, the process will be nrows.

378-383:

```
/* PARALLEL CODE */

for (int jj = 1; jj <= dim2; jj += BS) { /* Strip mining: Define blocks of size BS in the columns */ ///jj+=BS because jumps of size BS 

if (rank != 0) MPI_Recv( &h[0][jj], min(BS, dim2-jj+1), MPI_INT, 
rank-1, 0, MPI_COMM_WORLD, &status); // we receive data from previous 
process and store in ith matrix row -> Size of the buffer = size of 
smallest between BS or dim2-jj+1. Status is a predefined variable 
(MPI_Status, line 153) /tag can be 1 instead of 0

for (i = 1; i <= nrows; i++) /// iterate in nrows range
```

```
for (j = jj; j < min(jj+BS,dim2+1); j++) /* Strip mining: Traverse BS columns within the current block */ // column index beyond current block \rightarrow dim2 as limit of columns (avoid) /min of (next for loop jump (jj loop) vs max size+1), ensures that we are not out of range
```

Here we receive the data sent by the previous process and use it to create the scoring matrix in the following lines

421:

```
if (rank != nprocs-1 ) MPI_Send(&h[nrows][jj], min(BS, dim2-jj+1),
MPI_INT, rank + 1, 0, MPI_COMM_WORLD);
    /* Here we could use a non-blocking send. */ //Tag could be 1 as
well
```

We check that we are not in the last process, if that is true, send the &h matrix on row [nrows] and column [jj] to the next process in order.

If we were to use a non-blocking send (nb SW mpi.c):

```
if (rank != nprocs-1 ) {
    MPI_Request request;
    MPI_Isend(&h[nrows][jj], min(BS, dim2-jj+1), MPI_INT, rank + 1, 0,
MPI_COMM_WORLD, &request);//every row is BS length, when last row, BS
is too long (you are already in the last BS range)
    MPI_Wait(&request, MPI_STATUS_IGNORE);
}
}
/* Here we could use a non-blocking send. */ //MPI_Isend+MPI_wait+
request \to MPI_Send data and count??
```

In this case we would use MPI_Isend and MPI_Wait to avoid blocking the process

432:

```
MPI_Allreduce(localres, globalres, 1, MPI_2INT, MPI_MAXLOC, MPI_COMM_WORLD); /// we store the output in globalres array and use MAXLOC operation to find max value
```

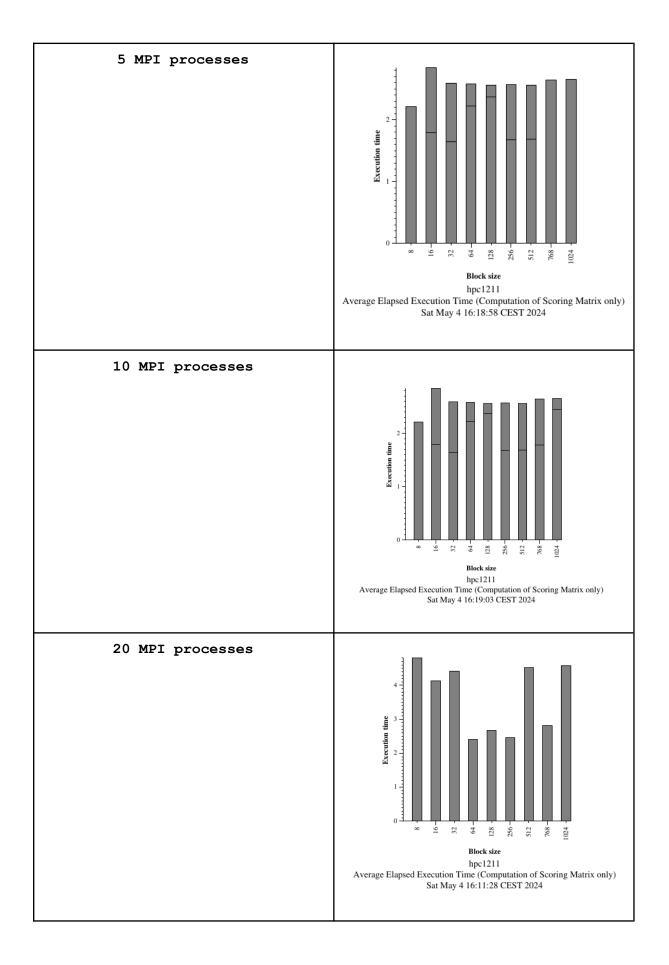
Allreduce is used to gather elements across all processes, then compute the maximum value across all processes and identify the process that contains it.

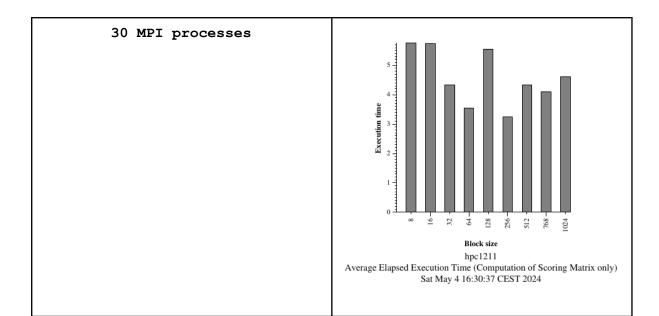
508:

```
if (rank==0) free(s1); /* Only the master process allocated it */ //master process\rightarrow rank 0, if rank is 0, end
```

This line is used to end the program and free the memory when the last process (0) is reached.

Blocksize SW mpi.c:





You can observe that the execution time increases with the more processors you have.

Also it can be observed that the blocking size with smaller execution time overall is block size 256, which makes it the optimal value as a blocking factor.

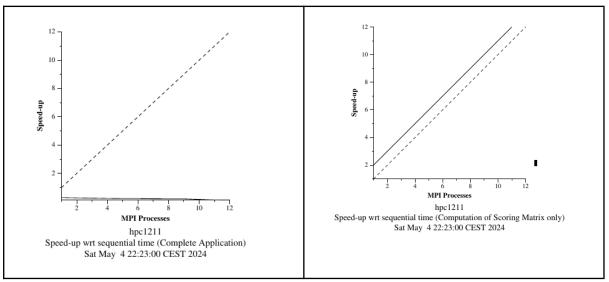
Strong SW mpi.c:

There appears to be an error when executing this program.

According to the error exit of the sbatch submission it appears to be a syntax error.

```
(standard_in) 2: syntax error
(standard_in) 1: syntax error
(standard_in) 1: syntax error
(standard_in) 2: syntax error
(standard_in) 1: syntax error
(standard_in) 1: syntax error
(standard_in) 1: syntax error
(standard_in) 2: syntax error
(standard_in) 1: syntax error
(standard_in) 1: syntax error
```

Even with this errors the code created the following 2 plots



The graph belonging to the computation of the scoring matrix presents no problem and seems to be plotted correctly, while the one belonging to the complete application seems to be wrongly plotted, this may be due to the errors mentioned above.