

CPSC 474

Project 2: Compressing a Sparse Matrix

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Pseudocode

Plain English PseudoCode

```
if ("-mFile" is found)
    load the filename that follows
else
    Throw no filename error

Init MPI

if rank is root
    load the file
    while there are lines
        get line and load it into a vector
        save number of lines

for each line
    while there are letters in the line
        parse out the numbers and add them to a master vector

calculate portion that each process will have
calculate # of columns

Broadcast the portion and root values to all processes
resize arrays on all processes to the size of portion
Scatter data to all the processes' resized arrays
Make sure all processes reach the barrier before proceeding

for each local number given to each process
    if number is not 0
        push number to a solution vector
        push number's row and col to solution vector
```

Output individual process solutions

For each non-root process

- send the size of the solution vector to the root rank

- Then send all solution data to the root rank

For the root process

- Gather the sizes of the other processes' solution vectors into a gather sizes array

- Then use the gather sizes array to calculate the offsets of the incoming data in the new array and resize the array to fit all of it.

- Ask the other processes for their solutions and add your own data

- Output the final solutions

Finalize MPI

```

//Detailed PseudoCode
rows = 0
columns = 0
numbers = []
size = 0
rank = 0
rootRank = 0

MPI_Init(argc, argv)
MPI_Comm_rank(MPI_COMM_WORLD, rank)
MPI_Comm_size(MPI_COMM_WORLD, size)

IF rank == rootRank DO:
    // read lines from file to rowVec[]
    rowVec = []
    WHILE line to read in matrixFile DO:
        rowVec[rows] = line
        n = count elements in line separated by comma
        IF n > columns DO:
            columns = n
        ENDIF
        rows += 1
    ENDWHILE
    close matrixFile

    // read numbers from rowVec to numbers[], fill empty cells with 0
    FOR i = 0 to rows - 1 DO:
        n = count elements in rowVec[i] separated by comma
        FOR j = 0 to columns - 1 DO:
            IF j < n DO:
                numbers[i].append(line[j])
            ELSE DO:
                numbers[i].append(0)
            ENDIF
        ENDFOR
    ENDFOR

portion = 0

IF size > 1 DO:
    portion = numbers.size / size
ENDIF

IF rows != 0 DO:
    cols = numbers.size / rows

```

```

localNumbers = []
localNumbers.size = portion

MPI_Bcast(portion, 1, MPI_INT, rootRank, MPI_COMM_WORLD)

MPI_Scatter(numbers, portion, MPI_INT, localNumbers, portion, MPI_INT, rootRank,
MPI_COMM_WORLD)

MPI_Barrier(MPI_COMM_WORLD)

// copy non-zeros from localNumbers to solution[]
solution = []
IF size > 1 DO:
    FOR i = 0 to portion - 1 DO:
        IF localNumbers[i] != 0 DO:
            solution.append(localNumbers[i])
            solution.append((rank * portion + i) % cols))
            solution.append((rank * portion + i) / cols))
        ENDIF
    ENDFOR
ENDIF

IF rank == rootRank DO:
    IF size > 1 DO:
        IF numbers.size % size != 0 DO:
            FOR i = 0 to numbers.size % size - 1 DO:
                IF numbers[portion * size] != 0 DO:
                    solution.append(numbers[portion * size])
                    solution.append((rank * portion + i) % cols))
                    solution.append((rank * portion + i) / cols))
                ENDIF
            ENDFOR
        ENDIF
    ELSE DO:
        IF numbers.size != 0 DO:
            FOR i = 0 to numbers.size - 1 DO:
                IF numbers[i] != 0 DO:
                    solution.append(numbers[i])
                    solution.append((rank * portion + i) % cols))
                    solution.append((rank * portion + i) / cols))
                ENDIF
            ENDFOR
        ENDIF
    ENDIF
ENDIF

gatherSizes = []

```

```

displacements = []
solSize = solution.size

IF rank == rootRank DO:
  FOR i = 0 to size - 1 DO:
    IF i != rootRank DO:
      gatherSizes.append(0)
      MPI_Recv(gatherSizes[i], 1, MPI_INT, i, 0, MPI_COMM_WORLD, MPI_STATUS_IGNORE)
    ELSE DO:
      gatherSizes.append(solSize)
    ENDIF
  ENDFOR

  sum = 0
  FOR i = 0 to gatherSizes.size - 1 DO:
    displacements.append(sum)
    sum += gatherSizes[i]
  ENDFOR

  finalAnswer = []
  finalAnswer.size = sum

  MPI_Gatherv(solution, solution.size, MPI_INT, finalAnswer, gatherSizes,
displacements, MPI_INT, rootRank, MPI_COMM_WORLD)

  print(finalAnswer)
ELSE DO:
  // send the sizes of the solution buffer from each process
  MPI_Send(solSize, 1, MPI_INT, rootRank, 0, MPI_COMM_WORLD)
  // then gather the results into one buffer
  MPI_Gatherv(solution, solution.size, MPI_INT, NULL, NULL, NULL, MPI_INT,
rootRank, MPI_COMM_WORLD)

ENDIF

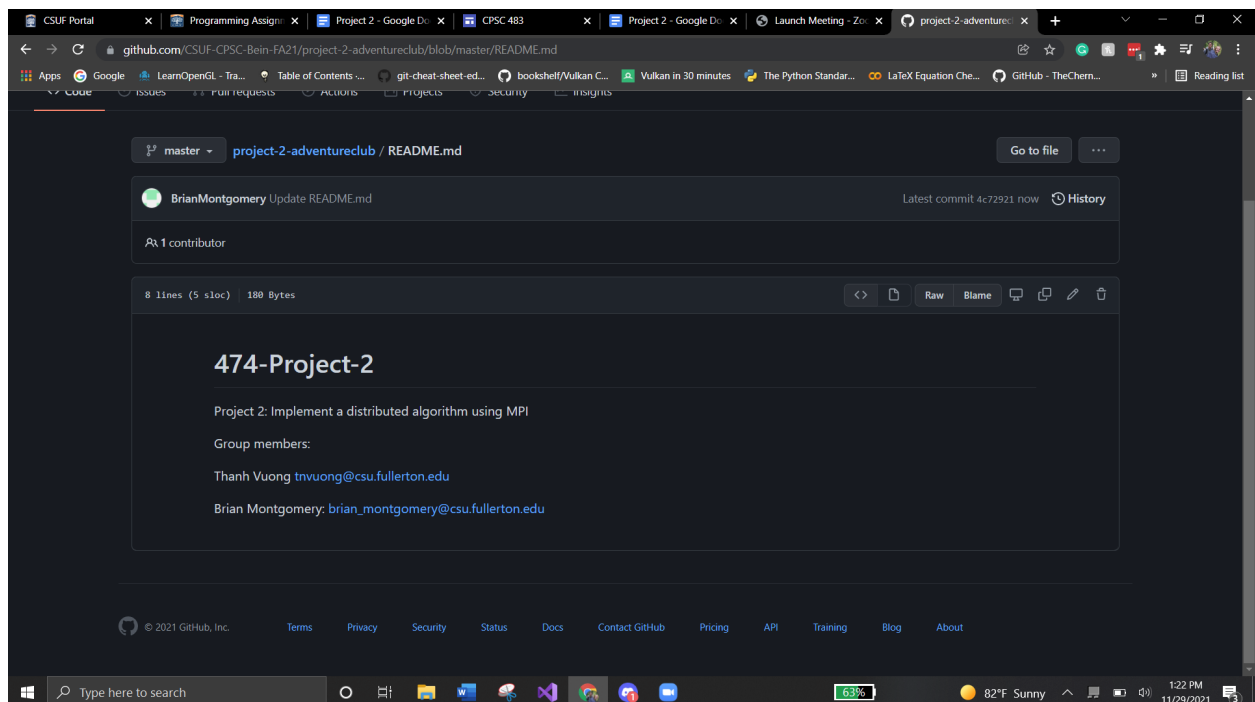
MPI_Finalize()

```

How to Run the Code

1. Have MSMPI MS-MPI SDK and Redistributions installed on a Windows machine.
2. Download the GitHub repository from <https://github.com/CSUF-CPSC-Bein-FA21/project-2-adventureclub/tree/master>
3. Open command prompt and navigate to the repository's /x64/Debug folder.
4. Run the command `mpiexec -n 4 ./mpi.exe -mFile matrix.txt`
 - This command runs the program executable with 4 processes and uses the input of matrix.txt which is also found in the /x64/Debug folder.

Screenshots



```
C:\Users\brian\Documents\Fall 2021\474 Parallel and Distrbuted Computing\Assignments\MPI\x64\Debug>mpiexec -n 4 ./mpi.exe -mFile matrix.txt
Rank: 0 Solution:
Rank: 1 Solution: (value = 2, column = 8, row = 7),
Rank: 2 Solution: (value = 3, column = 17, row = 9),
Rank: 3 Solution: (value = 6, column = 6, row = 12), (value = 8, column = 10, row = 13), (value = 1, column = 15, row = 13),
Final Solution: (value = 2, column = 8, row = 7), (value = 3, column = 17, row = 9), (value = 6, column = 6, row = 12), (value = 8, column = 10, row = 13), (value = 1, column = 15, row = 13),
C:\Users\brian\Documents\Fall 2021\474 Parallel and Distrbuted Computing\Assignments\MPI\x64\Debug>
```

Input used in this screenshot:

0, 0
0, 0
0, 0
0, 0
0,0
0,0
0,0,0,0,0,0,0,0,2,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0
0,0
0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,3,0,0,0,0,0,0,0,0,0
0,0
0,0
0,0,0,0,0,0,6,0
0,0,0,0,0,0,0,0,0,0,8,0,0,0,0,0,1,0,0,0,0,0,0,0,0,0,0,0
0,0