Parallel Programming Final Project

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1 PROJECT1

1.1 MPI_Allgather

Use MPI_Send and MPI_Recv to implement the MPI_ALLGATHER function. For each process, it first send it data to all of the rest process using MPI_Send. For simplicity, each process send to itself by **memcpy**. Then each process recive data from other processes by MPI_Recv. We test the algorithm by generating send_buffer and recv_buffer of length 10000, and number of process is set to 4.

Using my_MPI_Allgather:

Total elapsed time: 0.000029

Using MPI_Allgather:

Total elapsed time: 0.000010

1.2 Gemm

The Matrix is required to be parallizaed in the tile unit. Which is cannon algorithm. Cannon's algorithm is a distributed algorithm for matrix multiplication for two-dimensional meshes. It is especially suitable for computers laid out in an N × N mesh. Consider two n × n matrices A and B partitioned into p blocks. A(i, j) and B(i, j) (0i, jp) of size $(n/\sqrt{p})(n/\sqrt{p})$ each. The initial step of the algorithm regards the alignment of the matrixes. Align the blocks of A and B in such a way that each process can independently start multiplying its local submatrices. This is done by shifting all submatrices A(i, j) to the left (with wraparound) by i steps and all submatrices B(i, j) up (with wraparound) by j steps. Perform local block multiplication. Each block of A moves one step left and each block of B moves one step up (again with wraparound). Perform next block multiplication, add to partial result, repeat until all blocks have been multiplied.

In our experiment, we randomly generated 1024x1024 matrix and test the time elapsed during multiplication.

Large matrix multiplication random generate 1024x1024 matrix...

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Multiplication size: (1024, 1024) x (1024, 1024).

Total elapsed time: 17.719502

Convolution operation Using a 4 \star 4 kernel to do the

pooling operation for the matrix.

Matrix size: (10214, 1024), Kernel size (4, 4).

Total elapsed time: 0.121011

Pooling operation Using a 4 \star 4 kernel to do the

convolution operation for the matrix

Matrix size: (1024, 1024), kernel size (4, 4).

Total elapsed time: 0.107230

1.3 Wordcount

In this problem we use map<string, int> as a dict to map a word to its frequency. We define the word as a sequence which is made up of only english letters(lowercase or uppercase). The common function of this two algorithms is performing wordcounting, given the file reader and the block size to read. For this function, we extract all the suitable word and add the lowercase of the word to the dict. The big file problem deals with a single file of size 820kb, while the small file problem deals with 100 file of size ranging from 0kb to 400 kb.

1.3.1 big file. For big file, we divide the whole data block into **p** pieces(p is the total number of process), and assign each process to one of its block. After getting the word frequency dictionary of all processes, we need to merge all the dictionaries into a final dictionary. Each process first send the size of the dictionary, then for each word, it first send the length of the word, then send the word and the frequency. The reason to send the size of dictionary and the length of the word is MPI_Recv need the block size before it receive the actual word. Finally, we sort the final dictionary to get the few most frequent words.

The output is shown as follows.

wordcount big file the most frequent words top 20

<word></word>	<freq></freq>	
the	8233	
of	3895	
and	3159	
to	2752	
that	2456	
a	2312	
in	2298	
he	2005	
was	1857	
had	1694	
she	1638	

her	1619	
with	1467	
his	1327	
it	876	
on	868	
him	836	
not	825	
for	770	
they	723	
=======		

Total elapsed time: 2.327309

1.3.2 *small file.* The idea of the small file wordcount is same as the big file wordcount, except that it distribute the small files to all the process. For each process, it perform wordcount of each file and merge all frequencies locally. Then it merge all the dictionaries into a final dictionary.

The result is shown as follows

wordcount small files

the most frequent words top 20

<Word> <Frea> ______ the 485 is 218 208 to of 156 150 а for 134 122 as file 117 in 116 and 114 you 110 99 jpeg if 97 are 93 this 82 image 79 or 71 that 71 70 LISE be 68

Total elapsed time: 0.117455

2 PROJECT2

2.1 Monte Carlo

The idea of Monte Carlo is randomly throw a niddle into a 1x1 square area several times(denoted as n). and count times it fall into the circle area(r=1) inside the square area(denoted as m). and compute π by

$$\pi_{estimate} = \frac{m}{n}$$

We just parallelize all the individual trials. The result of our experiment is shown as follows

```
Exercise 1: Monte Carlo
------
total trial time=10000000
pi = 3.14145
```

2.2 Quick Sort

The idea of quick sort is parallel the quick sort in recursive form In our experiment we generate a big array of size 100000, and we assign random value each of it element. after quicksort we check its correctness and extimate the time elapsed.

```
allocting a array of length 1000000...

Total elapsed time: 0.046374

checking quick sort result..
```

2.3 Page Rank

For Pagerank algorithm we need to implement two classes, Graph and SubGraph. Graph class contain all the nodes of a whole graph. SubGraph class contain partial nodes of the whole graph. Beside, it contain the map from global index of father graph to local index in the subgraph. The general idea of pagerank is to first divide it into several subgraphs and then parallelize each iteration. The main body of the code is shown as follows.

```
for (int t = 0; t < iterations; t++){
    cout << "iter : " << t << " running" << endl;
    // #pragma omp parallel
    {
        #pragma omp for
        for (int tid = 0; tid < threads; tid++){
            // do pagerank on the local sub-Graph
            sgs[tid].updatePageRankLocal();
        }
        syncGraph(sgs, threads);
    }
    graphMerge(g, threads, sgs);
}</pre>
```

- 2.3.1 graphPartition. To make sure each sub graph have as most inside link as possible. We first count degree of each node, choose few nodes with most degree as root, and perform DFS to get the rest of subgraph.
- 2.3.2 updatePageRankLocal. Since for each subgraph, there exist edge linked to other subgraph. we only perform pagerank locally, which means we don't accept pagerank from outside the subgraph.
- 2.3.3 syncGraph. At the end of each iteration all subgraph must be synchronized to get the real pagerank result.
- 2.3.4 graphMerge. At the end of program, we need to merge all subgraphs into the original graph.

The number of node is set as 1024000, each node have degree ranging from 1 to 10. The threads is set as 100. we first check the correctness of our algorithm by set the initial pagerank of each node as 1. If it is correct the pagerank of each node would stay unchanged(Assume that the edges in the graph are bi-directional). the result is shown as follows

Initiallzing graph of node=1024000...

generate edge count of different nodes ranges from 1 to 10 executing page rank $\,$

iter : 0 running
iter : 1 running
iter : 2 running
iter : 3 running
iter : 4 running

Total elapsed time: 77.663224s

The elapsed time is too long for few iterations..

3 PROJECT3

3.1 Build Hadoop distributed system

First, you need to prepare four virtual machines named master, slave1, slave2, and slave3. The network of four virtual machines is configured as bridge mode, and static IP is allocated.

192.168.3.77 master 192.168.3.74 slave1 192.168.3.76 slave2 192.168.3.78 slave3

The four machines then exchange the SSH public key with each other, allowing them to ping each other. Then install Java8-Open-JKD and Hadoop and add them to the environment variables. Finally, configure the Hadoop configuration files, including core-site.xml, .hdfs-site.xml, yarn-site.xml, mapred-site.xml. Enter command "jps" test the system

hadoop@master:~/hadoop-3.3.1/etc/hadoop\$ jps

4288 Jps 3009 NameNode 3923 NodeManager 3736 ResourceManager

3422 SecondaryNameNode

3198 DataNode

The configuration was successful.

3.2 Mapper

The Mapper file combine the input file name with every piece of temperature data in that file

<filename, temperature>

3.3 Reducer

The Reducer reduce to the max temperature among data with same filename.