# CSE 708 Seminar

# Subset Sum Count (Using OpenMP in C)

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- 2) Sequential Solution
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- 4) Slurm Script
- 5) Parallel Solution
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#### 1) Problem Statement

Determine the count of the subsets within an array whose sum equals a given target sum.

```
array[] = { 3, 1, 2, 4, 5} Target Sum = 6
subsets = {3, 1, 2}, {4, 2}, {1,5}
Subset Sum Count = 3
```

Constraint: All array elements are whole numbers.

#### 2) Sequential Solution

```
if ( array[i] > j ): DP[i][j] = DP[i-1][j]
else: DP[i][j] = DP[i-1][j] + DP[i-1][j-array[i]]
```

#### 2) Sequential Solution

```
if ( array[i] > j ): DP[i][j] = DP[i-1][j]
else: DP[i][j] = DP[i-1][j] + DP[i-1][j-array[i]]
```

```
static int subsetSum(int array[], int array_size, int sum){
    int DP[][] = new int[array_size + 1][sum + 1];
   DP[0][0] = 1;
    for(int i = 1; i <= sum; i++)</pre>
     DP[0][i] = 0;
    for(int i = 1; i <= array_size; i++){</pre>
        for(int j = 0; j <= sum; j++){</pre>
            // DP Formula
            if (array[i-1] > j)
              DP[i][j] = DP[i-1][j];
              DP[i][j] = DP[i-1][j] + DP[i-1][j-array[i-1]];
    return DP[array_size][sum];
```

## 3) MPI vs OpenMP

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MPI

**OpenMP** 

</>>

Cores

Nodes.

Tasks/

Processes/

**Instances** 

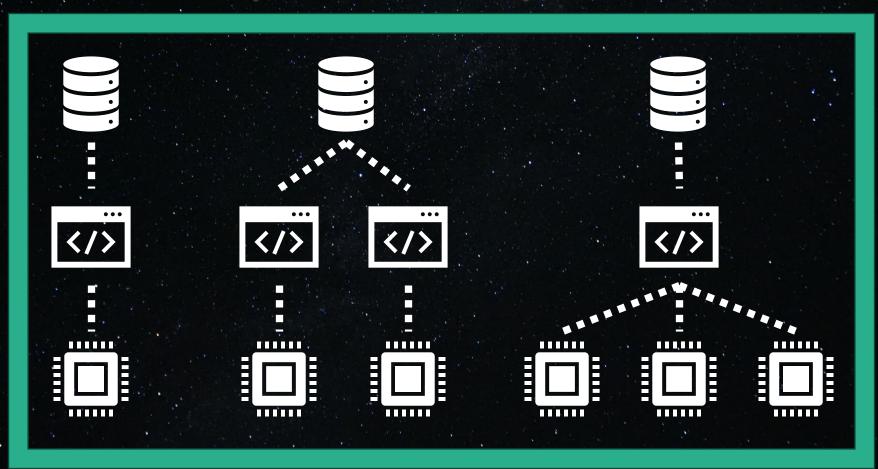
# 3] MPI vs OpenMP

Hybrid = MPI + OpenMP

Nodes.

Tasks/
Processes/
Instances

Cores



## 3) MPI vs OpenMP

#### MPI.

```
#SBATCH --nodes=4

#SBATCH --ntasks-per-node=2

#SBATCH --cpus-per-task=1
```

#### **OpenMP**

```
#SBATCH --nodes=1
#SBATCH --ntasks-per-node=1
#SBATCH --cpus-per-task=3
```

#### Hybrid

```
#SBATCH --nodes=4

#SBATCH --ntasks-per-node=2

#SBATCH --cpus-per-task=3
```

#### 3] MPI vs OpenMP

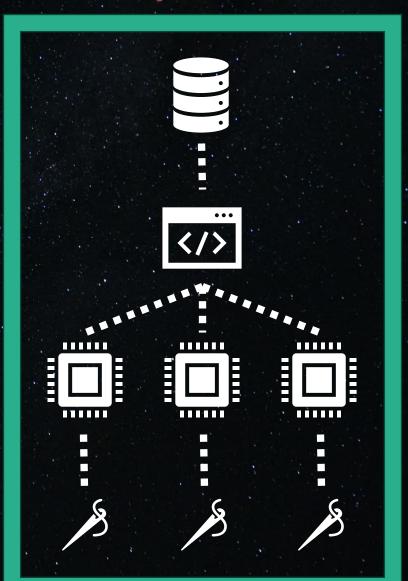
**OpenMP** 

**Nodes** 

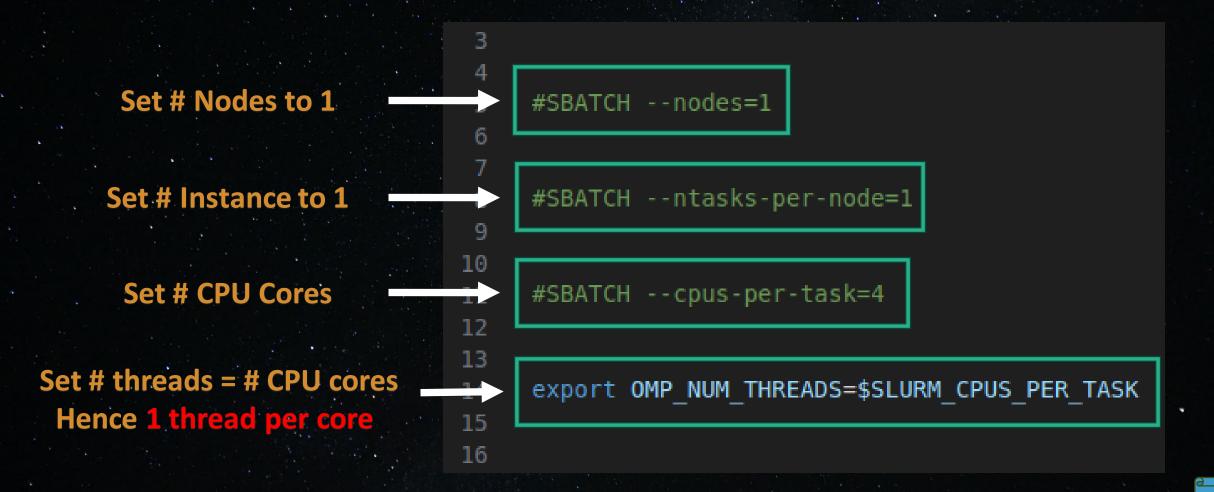
Tasks / Processes / Instances

**Cores (Hardware)** 

Threads (Software)



# 4) Slurm Script



# 4) Slurm Script

**OpenMP Configuration** 

Reserve all nodes & cores (Pizza Box) for exclusive use

Set # threads = # CPU cores Hence 1 thread per core

**Enables support for OpenMP pragmas and directives** 

```
    slurm.sh 
    x

708 > open_mp_c_execution > parallel > \( \sigma \) slurm.sh
       #!/bin/bash
       #SBATCH --nodes=1
       #SBATCH --ntasks-per-node=1
       #SBATCH --cpus-per-task=64
       #SBATCH --time=00:10:00
       #SBATCH --partition=general-compute
       #SBATCH --gos=general-compute
  11
       #SBATCH --cluster=ub-hpc
  12
       #SBATCH --reservation=ubhpc-future
  13
       #SBATCH --job-name="ssc 64 threads"
  15
       #SBATCH --output=output openmp-64.txt
       #SBATCH --exclusive
       module load intel
       export OMP NUM THREADS=$SLURM CPUS PER TASK
  23
  24
                      -o compiled file ssc openmp.c
       qcc - fopenmp
       ./compiled file
```

#### 5) Parallel Solution

Sequential Code Snippet

Parallel OpenMP
Code Snippet

```
115
       for (int i = 1; i <= n; i++){
           for (int j = 0; j \leftarrow sum; j++){
116
                if (a[i-1] \rightarrow j){
117
                    dp[i][j] = dp[i - 1][j];
118
119
                else{
120
121
                    dp[i][j] = dp[i - 1][j] + dp[i - 1][j - a[i - 1]];
122
123
124
```

```
for (int i = 1; i <= n; i++) {
126
          #pragma omp parallel for
127
          for (int j = 0; j <= sum; j++) {
128
              if (a[i - 1] > j){
129
                   #pragma omp atomic write
130
131
                  dp[i][j] = dp[i - 1][j];
132
              else {
134
                   #pragma omp atomic write
                   dp[i][j] = dp[i - 1][j] + dp[i - 1][j - a[i - 1]];
135
136
137
138
```

#### 5) Parallel Solution

```
// Row Wise Iteration

// For (int i = 1; i <= n; i++) {

// #pragma omp parallel for

for (int j = 0; j <= sum; j++) {

// Parallel column execution for a given row i

// Resumes sequential execution

// Resumes sequential execution</pre>
```

```
for (int i = 1; i <= n; i++) {
    #pragma omp parallel for
    for (int j = 0; j <= sum; j++) {
        int thread_id = omp_get_thread_num();
        thread_table[i][j] = thread_id;

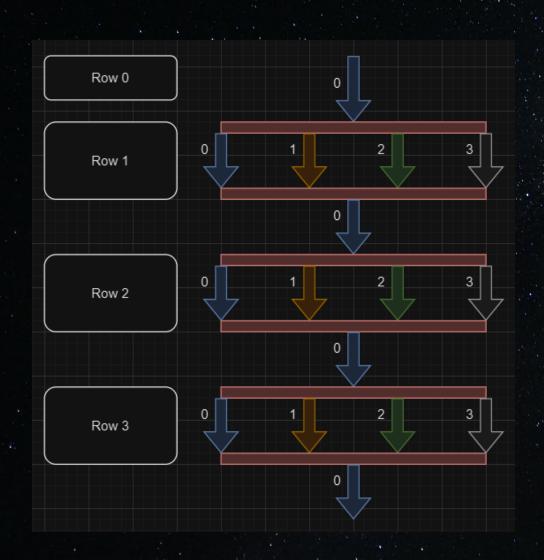
        if (a[i - 1] > j){
            #pragma omp atomic write
            dp[i][j] = dp[i - 1][j];
        }
        else {
            #pragma omp atomic write
            dp[i][j] = dp[i - 1][j] + dp[i - 1][j - a[i - 1]];
        }

140
        }

141       }
```

```
Max Threads = 4
     Input Array = [3, 1, 2, 4, 5]
     Target Sum = 6
     Thread Table:
     Row 0: 0 0 0 0 0 0 0
     Row 1: 0 0 1 1 2 2 3
     Row 2: 0 0 1 1 2 2 3
     Row 3: 0 0 1 1 2 2 3
     Row 4: 0 0 1 1 2 2 3
11
     Row 5: 0 0 1 1 2 2 3
12
13
     DP Table:
14
     Row 0: 1 0 0 0 0 0 0
15
     Row 1: 1 0 0 1 0 0 0
     Row 2: 1 1 0 1 1 0 0
17
     Row 3: 1 1 1 2 1 1 1
     Row 4: 1 1 1 2 2 2 2 2
     Row 5: 1 1 1 2 2 3 3
21
     Subset Sum Count = 3
22
```

### 5) Parallel Solution



```
Max Threads = 4
     Input Array = [3, 1, 2, 4, 5]
     Target Sum = 6
     Thread Table:
     Row 0: 0 0 0 0 0 0 0
     Row 1: 0 0 1 1 2 2 3
     Row 2: 0 0 1 1 2 2 3
     Row 3: 0 0 1 1 2 2 3
     Row 4: 0 0 1 1 2 2 3
11
     Row 5: 0 0 1 1 2 2 3
12
13
     DP Table:
     Row 0: 1 0 0 0 0 0 0
15
     Row 1: 1 0 0 1 0 0 0
17
     Row 2: 1 1 0 1 1 0 0
     Row 3: 1 1 1 2 1 1 1
     Row 4: 1 1 1 2 2 2 2 2
     Row 5: 1 1 1 2 2 3 3
21
     Subset Sum Count = 3
22
```

#### A) Standard Execution (Amdahl's Law):

- Total size of input data remains same
- Increase the number of cores
- With more cores, each core operates on lesser data

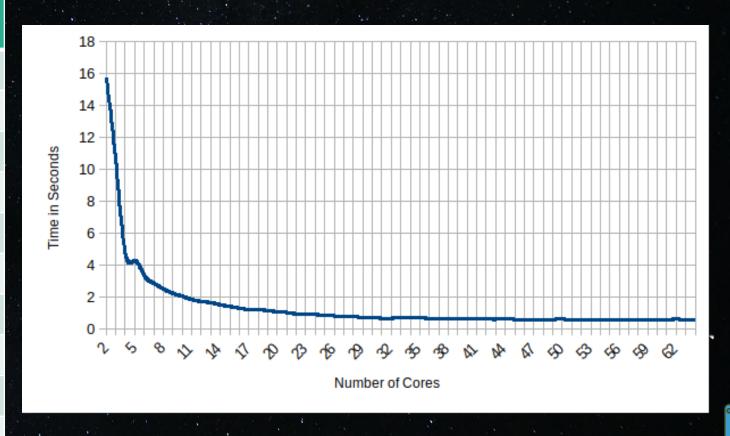
#### B) Scaled Execution (Gustafson's Law):

- Fix amount of data that each core operates on
- Increase the number of cores
- With more cores, the total size of the input data should get increased, because data per core is constant

#### C) Sequential Execution

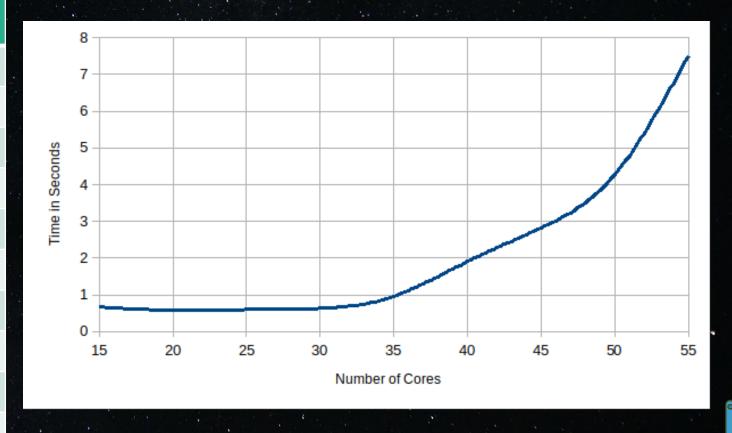
#### A) Standard Execution (Amdahl's Law):

Cores	Data/Core (Col/Core)	Input Size (Total Col)	Time (in seconds)	
1	100 M	100 M	31.46356	
2	50 M	100 M	15.75797	
3	33.3 M	100 M	10.53631	
4	25 M	100 M	4.781726	
7	14.3 M	100 M	2.885004	
15	6.67 M	100 M	1.430131	
25	4 M	100 M	0.890899	
35	2.85 M	100 M	0.716509	
45	2.2M	100 M	0.617453	
64	1.56 M	100 M	0.589691	



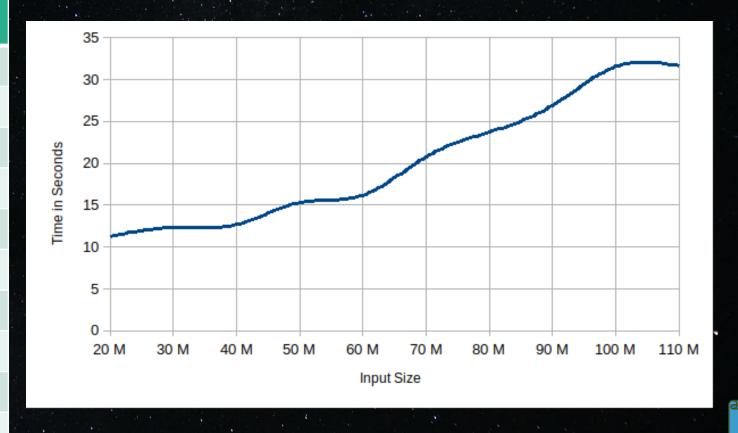
#### B) Scaled Execution (Gustafson's Law):

Cores	Data/Core (Col/Core)	Input Size (Total Col)	Time (in seconds)
10	2 M	20 M	0.624523
15	2 M	30 M	0.675528
20	2 M	40 M	0.586507
25	2 M	50 M	0.607871
30	2 M	60 M	0.6315
35	2 M	70 M	0.96475
40	2 M	80 M	1.916973
45	2 M	90 M	2.823243
50	2 M	100 M	4.27448
55	2 M	110 M	7.51923



#### C) Sequential Execution

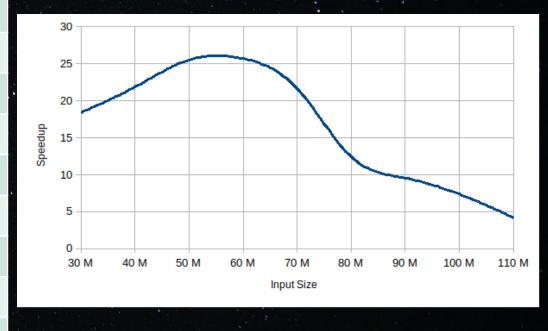
Cores	Data/Core (Col/Core)	Input Size (Total Col)	Time (in seconds)
1	20 M	20 M	11.30215
1	30 M	30 M	12.34482
1	40 M	40 M	12.6843
1	50 M	50 M	15.32529
1	60 M	60 M	16.19791
1	70 M	70 M	20.78737
1	80 M	80 M	23.74192
1	90 M	90 M	26.91424
1	100 M	100 M	31.57325
1	110 M	110 M	31.6582



#### Speedup

Formula: Speedup = Tseq/Tp
Tseq is the execution time of sequential algorithm
Tp is the execution time of parallel algorithm with
p processors

Input Size (Total Col)	seq	Tseq	Cores	Data/Core (Col/Core)	Тр	Speedup
20 M	1	11.3	10	2 M	0.62	18.2258
30 M	1	12.34	15	2 M	0.67	18.4179
40 M	1	12.68	20	2 M	0.58	21.8620
50 M	1	15.32	25	2 M	0.6	25.5334
60 M	1	16.19	30	2 M	0.63	25.6984
70 M	1	20.78	35	2 M	0.96	21.6458
80 M	1	23.74	40	2 M	1.91	12.4293
90 M	1	26.91	45	2 M	2.82	9.54255
100 M	1	31.57	50	2 M	4.27	7.3944
110 M	1	31.65	55	2 M	7.51	4.2143



## 7) Reference

- 1) GFG: <a href="https://www.geeksforgeeks.org/count-of-subsets-with-sum-equal-to-x/">https://www.geeksforgeeks.org/count-of-subsets-with-sum-equal-to-x/</a>
- 2) Princeton University BootCamp:

  <a href="https://princetonuniversity.github.io/PUbootcamp/sessions/parallel-programming/Intro PP bootcamp 2018.pdf">https://princetonuniversity.github.io/PUbootcamp/sessions/parallel-programming/Intro PP bootcamp 2018.pdf</a>
  - 3) Dr. Jones Lectures on OpenMP