

# Parallel & Distributed Computing [Y1]

LAB MANUAL/REPORT





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# LAB MANUAL

# **Matrix Operations:**

The parallelized matrix multiplication code uses OpenMP to distribute the computation of matrix elements across multiple threads. Each thread computes individual elements of the result matrix by summing the products of corresponding elements from the input matrices.

When race conditions occur, the results of these operations can become unpredictable due to unsynchronized access to shared data.

#### With the Race condition

```
#include <stdio.h>
#include <omp.h>
#define N 3
void printMatrix(int mat[N][N]){
    for (int i = 0; i < N; i++){
        for (int j = 0; j < N; j++){
            printf("%d ", mat[i][j]);
        printf("\n");
void addMatrices(int A[N][N], int B[N][N], int C[N][N]){
    int sum = 0;
#pragma omp parallel for
    for (int i = 0; i < N; i++){
        for (int j = 0; j < N; j++){
            C[i][j] = A[i][j] + B[i][j];
            sum += C[i][j];
    printf("Sum of all elements in C: %d\n", sum);
}
int main(){
    int A[N][N] = \{\{1, 1, 1\}, \{1, 1, 1\}, \{1, 1, 1\}\};
    int B[N][N] = \{\{2, 2, 2\}, \{2, 2, 2\}, \{2, 2, 2\}\};
    int C[N][N];
    addMatrices(A, B, C);
    printf("Matrix A:\n");
    printMatrix(A);
    printf("\nMatrix B:\n");
    printMatrix(B);
    printf("\nMatrix C (A + B):\n");
    printMatrix(C);
    return 0;
```

```
2 2 2
 2 2 2
 2 2 2
 Matrix C (A + B):
 3 3 3
 3 3 3
 3 3 3
 __(biologist⊛kali)-[~/.../Parallel & Distr
 ibuted Computing/Lab/Report/1.matrix_opera
 tions]
• └─$ cd "/home/biologist/Documents/2UMT/sem
 ester6/Parallel & Distributed Computing/La
 b/Report/1.matrix_operations/" && gcc with
 _race_cond.c -o with_race_cond && "/home/b
 iologist/Documents/2UMT/semester6/Parallel
  & Distributed Computing/Lab/Report/1.matr
 ix_operations/"with_race_cond
 Sum of all elements in C: 27
 Matrix A:
 1 1 1
 1 1 1
 1 1 1
 Matrix B:
 2 2 2
 2 2 2
 2 2 2
 Matrix C (A + B):
 3 3 3
 3 3 3
 3 3 3
```

# Without the Racing condition:

```
#include <stdio.h>
#include <omp.h>
#define N 3
void printMatrix(int mat[N][N]){
#pragma omp parallel for
    for (int i = 0; i < N; i++){
        for (int j = 0; j < N; j++){
            printf("%d ", mat[i][j]);
        printf("\n");
void addMatrices(int A[N][N], int B[N][N], int C[N][N]){
#pragma omp parallel for
    for (int i = 0; i < N; i++){
        for (int j = 0; j < N; j++){
            C[i][j] = A[i][j] + B[i][j];
int main(){
    int A[N][N] = \{\{1, 1, 1\}, \{1, 1, 1\}, \{1, 1, 1\}\};
    int B[N][N] = \{\{2, 2, 2\}, \{2, 2, 2\}, \{2, 2, 2\}\};
    int C[N][N];
    addMatrices(A, B, C);
    printf("Matrix A:\n");
    printMatrix(A);
    printf("\nMatrix B:\n");
    printMatrix(B);
    printf("\nMatrix C (A + B):\n");
    printMatrix(C);
    return 0;
```

```
Matrix A:

1 1 1

1 1 1

1 1 1

Matrix B:

2 2 2

2 2 2

2 2 2

3 3 3

3 3 3

3 3 3

3 3 3
```

**Prefix Sum:** The parallelized prefix sum code leverages OpenMP to divide the array into segments processed by different threads. Each thread calculates partial sums for its segment, and synchronization mechanisms are used to ensure that the cumulative sums from previous segments are correctly included.

#### With the Race condition:

```
Prefix sum: 1 3 6 10 15 21 28 36 45 55
  __(biologist⊛kali)-[~/.../Parallel & Distribut
 ed Computing/Lab/Report/2.prefix_sum]
• └$ gcc -fopenmp race.c -o race
  ___(biologist⊛kali)-[~/.../Parallel & Distribut
 ed Computing/Lab/Report/2.prefix_sum]
• └$ ./race
 Prefix sum: 1 3 6 10 15 21 28 36 45 55
 __(biologist⊛kali)-[~/.../Parallel & Distribut
 ed Computing/Lab/Report/2.prefix_sum]
• └$ gcc -fopenmp race.c -o race
  ___(biologist⊛kali)-[~/.../Parallel & Distribut
 ed Computing/Lab/Report/2.prefix_sum]
• └$ ./race
 Prefix sum: 1 3 6 10 15 21 28 36 45 56
    -(biologist®kali)-[~/…/Parallel & Distribut|
 ed Computing/Lab/Report/2.prefix_sum]
• └─$ gcc -fopenmp race.c -o race
  ___(biologist⊛kali)-[~/.../Parallel & Distribut∎
 ed Computing/Lab/Report/2.prefix_sum]
• └$ ./race
 Prefix sum: 1 3 6 10 15 21 28 36 45 55
 __(biologist⊛kali)-[~/.../Parallel & Distribut
 ed Computing/Lab/Report/2.prefix_sum]
• └$ gcc -fopenmp race.c -o race
 __(biologist⊛kali)-[~/.../Parallel & Distribut
 ed Computing/Lab/Report/2.prefix_sum]
• └-$ ./race
 Prefix sum: 1 3 6 10 15 21 28 36 45 54
```

# Without the Racing condition

#### Code:

```
#include <stdio.h>
#include <omp.h>
int main()
    int arr[] = \{1, 2, 3, 4, 5, 6, 7, 8, 9, 10\};
    int prefix_sum[10];
    prefix_sum[0] = arr[0]; // First element remains the same
#pragma omp parallel for
    for (int i = 1; i < 10; i++)
        prefix_sum[i] = prefix_sum[i - 1] + arr[i];
    printf("Prefix sum: ");
    for (int i = 0; i < 10; i++)
        printf("%d ", prefix_sum[i]);
    printf("\n");
    return 0;
```

```
Distributed Computing/Lab/Report]

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```

#### **Recurrence Sum:**

In the parallelized recurrence sum code, each thread computes its partial sum of elements based on the recurrence relation. To prevent race conditions, critical sections or other synchronization mechanisms are employed to ensure that concurrent updates to shared variables are properly coordinated. This coordination guarantees the correctness of the computed sum.

#### With the Race condition

#### Code:

```
#include <stdio.h>
#include <omp.h>
int main(){
    int n;
    printf("Enter the number of terms in the sequence: ");
    scanf("%d", &n);
    int sequence[n];
    int start_value;
    printf("Enter the starting value of the sequence: ");
    scanf("%d", &start_value);
    sequence[0] = start_value;
#pragma omp parallel for
    for (int i = 1; i < n; i ++){
        sequence[i] = sequence[i - 1] + 3;
    printf("Sequence: ");
    for (int i = 0; i < n; i ++){
        printf("%d ", sequence[i]);
    printf("\n");
    return 0;
```

```
Enter the number of terms in the sequence: 9
Enter the starting value of the sequence: 4
Sequence: 4 7 10 13 16 19 22 25 28

—(biologist@kali)-[~/.../Parallel & Distribut ed Computing/Lab/Report/3.recurrance_sum]

—$ cd "/home/biologist/Documents/2UMT/semeste r6/Parallel & Distributed Computing/Lab/Report/3.recurrance_sum/" && gcc Norace.c -o Norace && "/home/biologist/Documents/2UMT/semester6/P arallel & Distributed Computing/Lab/Report/3.r ecurrance_sum/"Norace
Enter the number of terms in the sequence: 9
Enter the starting value of the sequence: 4
Sequence: 4 8 12 16 20 24 28 32 36
```

# Without the Racing condition Code:

```
#include <stdio.h>
#include <omp.h>
int main(){
    int n;
    printf("Enter the number of terms in the sequence: ");
    scanf("%d", &n);
    int sequence[n];
    int start_value;
    printf("Enter the starting value of the sequence: ");
    scanf("%d", &start_value);
    sequence[0] = start_value;
#pragma omp parallel
        int thread_id = omp_get_thread_num();
        int local_start = (n * thread_id) / omp_get_num_threads();
        int local_end = (n * (thread_id + 1)) / omp_get_num_threads();
        for (int i = local_start + 1; i < local_end; i++){</pre>
#pragma omp critical
                sequence[i] = sequence[i - 1] + 3;
    printf("Sequence: ");
   for (int i = 0; i < n; i++){
        printf("%d ", sequence[i]);
    printf("\n");
    return 0;
```

```
Enter the number of terms in the sequence: 9
Enter the starting value of the sequence: 4
Sequence: 4 7 10 13 16 19 22 25 28

—(biologist@kali)-[~/.../Parallel & Distributed Computing/Lab/Report/3.recurrance_sum]

—$ cd "/home/biologist/Documents/2UMT/semester6/Parallel & Distributed Computing/Lab/Report/3.recurrance_sum/" && gcc Norace.c -o Norace && "/home/biologist/Documents/2UMT/semester6/Parallel & Distributed Computing/Lab/Report/3.recurrance_sum/"Norace
Enter the number of terms in the sequence: 9
Enter the starting value of the sequence: 4
Sequence: 4 7 10 13 16 19 22 25 28
```

# **Depth First Search (DFS):**

DFS is a graph traversal algorithm that explores as far as possible along each branch before backtracking. It's used to traverse or search for nodes in a graph or tree structure.

When race conditions occur, the results of these operations can become unpredictable due to unsynchronized access to shared data.

# With the Race condition

```
#include <stdio.h>
#include <omp.h>
#define V 11
#define NUM_THREADS 4 // Number of threads to use
char vertexLabels[V] = {'A', 'B', 'C', 'D', 'E', 'F', 'G', 'H', 'I', 'J', 'K'};
int visited[V] = {0}; // Array to track visited vertices
int parent[V];
int graph[V][V] = {
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {1, 1, 0, 0, 1, 0, 0, 0, 0, 0, 0},
    {0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0},
    {0, 0, 1, 1, 0, 1, 0, 0, 0, 0, 0},
    {0, 0, 0, 1, 1, 0, 1, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 1, 0, 1, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 1, 0},
void DFS(int v){
    visited[v] = 1;
    for (int i = 0; i < V; i++){
        if (graph[v][i] && !visited[i]){
            parent[i] = v;
            DFS(i);
int main(){
    omp_set_num_threads(NUM_THREADS);
#pragma omp parallel for
    for (int v = 0; v < V; v++){
        if (!visited[v]){
            DFS(v);
    printf("\nEdges of the spanning tree:\n");
    for (int i = 1; i < V; i++){
        printf("(%c, %c)\n", vertexLabels[parent[i]], vertexLabels[i]);
    return 0;
```

```
__(biologist⊛kali)-[~/.../Parallel & Distri
buted Computing/Lab/Report/5.DFS]
└$ ./race
Edges of the spanning tree:
(C, B)
(A, C)
(F, D)
(C, E)
(G, F)
(A, G)
(K, H)
(H, I)
(A, J)
(J, K)
__(biologist⊛kali)-[~/.../Parallel & Distri
buted Computing/Lab/Report/5.DFS]
└$ ./race
Edges of the spanning tree:
(C, B)
(E, C)
(A, D)
(F, E)
(G, F)
(A, G)
(K, H)
(H, I)
(A, J)
(J, K)
—(biologist⊛kali)-[~/…/Parallel & Distri
buted Computing/Lab/Report/5.DFS]
└$ ./race
Edges of the spanning tree:
(C, B)
(A, C)
(F, D)
(D, E)
(G, F)
(A, G)
(K, H)
(H, I)
(A, J)
(J, K)
```

# Without the Racing condition

```
#include <stdio.h>
#include <omp.h>
#define V 11
#define NUM_THREADS 4 // Number of threads to use
char vertexLabels[V] = {'A', 'B', 'C', 'D', 'E', 'F', 'G', 'H', 'I', 'J', 'K'};
int visited[V] = {0}; // Array to track visited vertices
int parent[V];
int graph[V][V] = {
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {1, 1, 0, 0, 1, 0, 0, 0, 0, 0, 0},
    {0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0},
    {0, 0, 1, 1, 0, 1, 0, 0, 0, 0, 0},
    {0, 0, 0, 1, 1, 0, 1, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 1, 0, 1, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 1, 0},
void DFS(int v){
    visited[v] = 1:
    for (int i = 0; i < V; i ++){
        if (graph[v][i] && !visited[i]){
            parent[i] = v;
            DFS(i);
int main(){
    omp_set_num_threads(NUM_THREADS);
#pragma omp parallel for
    for (int v = 0; v < V; v++){
#pragma omp critical
       if (!visited[v]){
            DFS(v);
    printf("\nEdges of the spanning tree:\n");
    for (int i = 1; i < V; i + 1)
        printf("(%c, %c)\n", vertexLabels[parent[i]], vertexLabels[i]);
    return 0;
```

```
—(biologist⊛kali)-[~/…/Parallel & Distri
buted Computing/Lab/Report/5.DFS]
_$ ./Nrace
Edges of the spanning tree:
(A, C)
(E, D)
(C, E)
(D, F)
(F, G)
(G, H)
(H, I)
(H, K)
__(biologist⊛kali)-[~/.../Parallel & Distri
buted Computing/Lab/Report/5.DFS]
_$ ./Nrace
Edges of the spanning tree:
(C, B)
(A, C)
(E, D)
(D, F)
(F, G)
(G, H)
(H, I)
```

# **Breadth First Search (BFS):**

#### With the Race condition

DFS is a graph traversal algorithm that explores as far as possible along each branch before backtracking. It's used to traverse or search for nodes in a graph or tree structure.

When race conditions occur, the results of these operations can become unpredictable due to unsynchronized access to shared data.

```
#include <stdio.h>
#include <omp.h>
#include <stdlib.h>
#define V 11
#define NUM_THREADS 4 // Number of threads to use
char vertexLabels[V] = {'A', 'B', 'C', 'D', 'E', 'F', 'G', 'H', 'I', 'J', 'K'};
int visited[V] = {0}; // Array to track visited vertices
int parent[V];
int graph[V][V] = {
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {1, 1, 0, 0, 1, 0, 0, 0, 0, 0, 0},
    {0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0},
    {0, 0, 1, 1, 0, 1, 0, 0, 0, 0, 0},
    {0, 0, 0, 1, 1, 0, 1, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 1, 0, 1, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 1, 0},
};
void BFS(int start_vertex)
    int queue[V];
    int front = 0, rear = 0;
    queue[rear++] = start_vertex;
    visited[start_vertex] = 1;
    while (front ≠ rear){
        int current_vertex = queue[front++];
        for (int i = 0; i < V; i \leftrightarrow){
            if (graph[current_vertex][i] && !visited[i]){
                queue[rear++] = i;
                visited[i] = 1;
                parent[i] = current_vertex;
int main(){
    omp_set_num_threads(NUM_THREADS);
#pragma omp parallel for
    for (int v = 0; v < V; v++)
        if (!visited[v]){
            BFS(v);
    printf("\nEdges of the spanning tree:\n");
    for (int i = 1; i < V; i \leftrightarrow){
        printf("(%c, %c)\n", vertexLabels[parent[i]], vertexLabels[i]);
    return 0;
```

# output:

```
Edges of the spanning tree:
(C, B)
(E, C)
(A, D)
(D, E)
(G, F)
(A, G)
(G, H)
(H, I)
(A, J)
(J, K)
—(biologist⊛kali)-[~/…/Parallel & Distribut
ed Computing/Lab/Report/4.BFS]
└$ ./race
Edges of the spanning tree:
(C, B)
(A, C)
(A, D)
(D, E)
(G, F)
(A, G)
(G, H)
(H, I)
(A, J)
(H, K)
```

Without the Racing condition

```
#include <stdio.h>
#include <omp.h>
#include <stdlib.h>
#define V 11
#define NUM_THREADS 4 // Number of threads to use
char vertexLabels[V] = {'A', 'B', 'C', 'D', 'E', 'F', 'G', 'H', 'I', 'J', 'K'}; int visited[V] = \{0\}; // Array to track visited vertices
int parent[V];
int graph[V][V] = {
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0},
    {1, 1, 0, 0, 1, 0, 0, 0, 0, 0, 0},
    {0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0},
    {0, 0, 1, 1, 0, 1, 0, 0, 0, 0, 0},
    {0, 0, 0, 1, 1, 0, 1, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 0, 1, 0, 0, 0},
    {0, 0, 0, 0, 0, 1, 1, 0, 1, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1},
    {0, 0, 0, 0, 0, 0, 0, 1, 0, 1, 0},
void BFS(int start_vertex){
    int queue[V];
    int front = 0, rear = 0;
    queue[rear++] = start_vertex;
    visited[start_vertex] = 1;
    while (front \neq rear){
        int current_vertex = queue[front++];
        for (int i = 0; i < V; i++){
             if (graph[current_vertex][i] && !visited[i]){
                 queue[rear++] = i;
                 visited[i] = 1;
                 parent[i] = current_vertex;
int main(){
    omp_set_num_threads(NUM_THREADS);
#pragma omp parallel for
    for (int v = 0; v < V; v + +){
#pragma omp critical
        if (!visited[v]){
             BFS(v);
    printf("\nEdges of the spanning tree:\n");
    for (int i = 1; i < V; i \leftrightarrow){
        printf("(%c, %c)\n", vertexLabels[parent[i]], vertexLabels[i]);
    return 0;
```

```
Edges of the spanning tree:
(C, B)
(A, C)
(E, D)
(C, E)
(E, F)
(F, G)
(F, H)
(H, I)
(K, J)
(H, K)
┌──(biologist®kali)-[~/.../Parallel & Distribut
ed Computing/Lab/Report/4.BFS]
└$ ./Nrace
Edges of the spanning tree:
(C, B)
(A, C)
(E, D)
(C, E)
(E, F)
(F, G)
(F, H)
(H, I)
(K, J)
(H, K)
```