

- **#pragma omp task**

One thread creates a task to write the tasks onto sticky notes.

All threads to execute the sticky notes.

```
#include <iostream>
#include <string>
#include <omp.h>

using namespace std;
int main() {
    omp_set_num_threads( 2 );
#pragma omp parallel
{
    #pragma omp task
    cout<<"A"<<endl;
    #pragma omp task
    cout<<"B"<<endl;
}

    return 0;
}
```

Que- Why do we get 4 things printed when we only have print statements in 2 tasks?

Ans: Each of the two threads was assigned two tasks, resulting in a total of four tasks on the sticky notes

- **#pragma omp taskwait**

Causes all tasks to wait until they are completed.

If you care about order, do this:

```
#include <iostream>
#include <string>
#include <omp.h>

using namespace std;
int main() {
    omp_set_num_threads(2);
#pragma omp parallel
{
    #pragma omp task
    cout<<"A"<<endl;
    #pragma omp taskwait
    #pragma omp task
    cout<<"B"<<endl;
    #pragma omp taskwait
}

    return 0;
}
```

Factorial using tasks

```
#include <iostream>
#include <omp.h>
using namespace std;

long factorial(int n) {
    if (n == 0 || n == 1)
        return 1;

    long result;

    #pragma omp task shared(result)

    result = n * factorial(n - 1);

    #pragma omp taskwait

    return result;
}

int main() {
    int n;
    cout << "Enter a number: ";
    cin >> n;
    long result;
    #pragma omp parallel
    {
        #pragma omp single
        result = factorial(n);
    }
    cout << "Factorial of " << n << " is " << result << endl;
    return 0;
}
```

- `pragma omp parallel for collapse(2)`

Specifying the **COLLAPSE** clause allows you to parallelize multiple loops in a nest without introducing nested parallelism.

collapse(2): The collapse clause tells OpenMP to combine the two nested loops into one large loop. The 2 in `collapse(2)` means that OpenMP will collapse the two innermost loops into a single loop with a single iteration space.

Example Without collapse(2)

```
#include <iostream>
#include <omp.h>
using namespace std;
int main() {
    const int N = 4;
    int matrix[N][N] = {0};

    #pragma omp parallel for
    for (int i = 0; i < N; i++) {
        for (int j = 0; j < N; j++) {
            matrix[i][j] = i + j;
        }
    }
    for (int i = 0; i < N; i++) {
        for (int j = 0; j < N; j++) {
            cout << matrix[i][j] << " ";
        }
        cout << endl;
    }

    return 0;
}
```

- Only the outer loop (i loop) is parallelized.
- The inner loop (j loop) runs sequentially within each thread.

Example With collapse(2)

```
#include <iostream>
#include <omp.h>
using namespace std;
int main() {
    const int N = 4;
    int matrix[N][N] = {0};
    #pragma omp parallel for collapse(2)
    for (int i = 0; i < N; i++) {
        for (int j = 0; j < N; j++) {
            matrix[i][j] = i + j;
        }
    }
    for (int i = 0; i < N; i++) {
        for (int j = 0; j < N; j++) {
            cout << matrix[i][j] << " ";
        }
        cout << endl;
    }

    return 0;
}
```

What is Scheduling in OpenMP

Scheduling is a method in OpenMP to distribute iterations to different threads in for loop.

you can use `#pragma omp parallel for` directly without scheduling, it is equal to `#pragma omp parallel for schedule(static,1)`

Static

`#pragma omp parallel for schedule (static, chunk-size)`

If you do not specify chunk-size variable, OpenMP will divide iterations into chunks that are approximately equal in size and it distributes chunks to threads **in order**

If you specify chunk-size variable, the iterations will be divided into $\text{iter_size} / \text{chunk_size}$ chunks.

Example

```
#include <iostream>
#include <omp.h>
using namespace std;
int main()
{
    #pragma omp parallel for schedule(static, 3)
    for (int i = 0; i < 20; i++)
    {
        cout<<"Thread "<< omp_get_thread_num()<<" is
        running number \n"<< i;
    }
    return 0;
}
```

20 iterations will be divided into 7 chunks (6 with 3 iters, 1 with 2 iters).



Que: But what if $\text{iter_size} / \text{chunk_size}$ is larger than the number of threads in your computer, or number of threads you specified in `omp_set_num_threads(thread_num)`?

Ans: OpenMP will still split task into 7 chunks, but distributes the chunks to threads **in a circular order**, like the following figure shows



Dynamic

`#pragma omp parallel for schedule(dynamic, chunk-size)`

OpenMP will still split task into `iter_size/chunk_size` chunks, but distribute trunks to threads dynamically without any specific order.

Example

```
#include <iostream>
#include <omp.h>
using namespace std;
int main()
{
    #pragma omp parallel for schedule(dynamic, 3)
    for (int i = 0; i < 20; i++)
    {
        cout<<"Thread "<< omp_get_thread_num()<<" is running number \n"<< i;
    }
    return 0;
}
```

20 iterations will be divided into threads but not in order any chunk not done any thread can take it to execute.

