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```
#include <stdio.h>
#include <stdlib.h>
#include <omp.h>

int main(int argc, char *argv[]) {
    int nenv = 3;
    omp_set_num_threads(nenv); // set number of threads
    printf("nenv: %d\n", nenv);

    int chunk = 5;
    omp_set_schedule(omp_sched_static, chunk);
    // omp_set_schedule(omp_sched_dynamic, chunk);
    // omp_set_schedule(omp_sched_guided, chunk);
    printf("chunk size: %d\n", chunk);

    int i = 0;
    int n = 17;
    int a[n];
    int t[nenv];

    #pragma omp parallel for schedule(runtime)
    for (i=0; i<n; i++) {
        a[i] = omp_get_thread_num(); // chosen thread per iteration
        t[omp_get_thread_num()]++; // parallel increment
    }

    printf("a (schedule): ");
    for (i=0; i<n; i++) {
        printf("%d ", a[i]);
    }
    printf("\n");

    printf("t (counter): ");
    for (i=0; i<nenv; i++) {
        printf("%d ", t[i]);
    }
    printf("\n");
}
```

The variable `a` stores the selected thread number for each parallel iteration, while `t` stores a non-atomic counter that all threads with the same ID increment. Unless no two threads are assigned the same iteration, the final value of `t` will be non-deterministic as each `var++` operation is in fact a read-modify-write operation:

```
movl -4(%rbp), %eax # load var into eax
addl $1, %eax       # increment eax by 1
movl %eax, -4(%rbp) # store eax back into var
```

See Tables 1 and 2 for the values of α and t for different scheduling strategies.

[illegible]

Table 2: Values of array t for different scheduling strategies - keep in mind that these values are not reproducible / deterministic.

case / t	0	1	2
static, 0	74307862	7	1806905557
static, 1	8591638	7	1872621781
dynamic, 1	6150416	18	1875062992
dynamic, 2	40737057	1	1840476368
guided, 5	51370273	1	1829843168

2 Exercise 2

Table 3: Duration of independent tasks we want to schedule optimally.

Task ID	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16
Task duration	1	2	1	2	1	2	1	2	1	2	1	2	1	2	4	3	3

2.1 Optimal Schedule

Assuming the given tasks can be executed independently and the goal is to minimize the total execution time, with 4 workers available, the optimal schedule can be determined through the "Longest Processing Time" (LPT) rule. The LPT rule schedules tasks in decreasing order of processing time, which is optimal for minimizing the total completion time. This is often a good heuristic for this type of problem and tends to provide good results.

First, we sort the tasks by duration in descending order, then we assign the first n tasks to the n workers in a round-robin fashion.

```
def schedule_tasks(task_durations):
    sorted_tasks = sorted(enumerate(task_durations), key=lambda x: x[1])

    worker_availability = [0] * 4

    scheduled_tasks = []
    for task_id, duration in sorted_tasks:
        # find the worker who becomes available the soonest
        min_time = min(worker_availability)
        worker_index = worker_availability.index(min_time)

        # update worker availability
        worker_availability[worker_index] += duration

        # Record the assignment
        scheduled_tasks.append((worker_index, task_id, min_time, min_time + duration))

    return scheduled_tasks

task_durations = [1, 2, 1, 2, 1, 2, 1, 2, 1, 2, 1, 2, 1, 2, 4, 3, 3]
scheduled_tasks = schedule_tasks(task_durations)

for worker, task_id, start_time, end_time in scheduled_tasks:
    print(f"Worker {worker}: Task {task_id}, Start: {start_time}, End: {end_time}")
```

2.2 Schedule `static,3`

2.3 Schedule `dynamic,2`

3 Exercise 3

3.1 Fix the problems with this OpenMP code

4 Exercise 4

4.1 What is the output of the three different versions?

4.2 How often is the function `omp_tasks` called?

5 Exercise 5

5.1 Parallelize the pixel computation

5.2 Running time analysis

5.3 Influence of schedule parameter

6 Exercise 6

6.1 Parallelize the filter computation

6.2 Strong scaling analysis

6.3 Weak scaling analysis

7 Exercise 7

7.1 Convert OpenMP code to CUDA

7.2 Running time analysis

7.3 Impact of block size

7.4 Running time: CPU vs GPU code

8 Addendum: Raw Data

1168	1	1	0.0603872
1168	1	1	0.0607409
1168	1	1	0.0600319
1168	2	1	0.196807
1168	2	1	0.2452
1168	2	1	0.19003
1168	4	1	3.45923
1168	4	1	3.90704
1168	4	1	3.45583
1168	8	1	5.395
1168	8	1	5.45436
1168	8	1	4.53896
1168	16	1	10.7055
1168	16	1	10.5507
1168	16	1	10.2593
1168	24	1	17.3402
1168	24	1	18.5362
1168	24	1	17.2604
1168	32	1	26.1056
1168	32	1	25.1663
1168	32	1	27.9486

Figure 1: Raw output from "filter strong" job.

1168	1	1	0.060196
1168	1	1	0.0609
1168	1	1	0.060195
1168	2	2	0.401089
1168	2	2	0.635222
1168	2	2	1.18221
1168	4	4	14.4383
1168	4	4	13.3359
1168	4	4	9.2267
1168	8	8	44.0875
1168	8	8	44.8141
1168	8	8	42.5354

Figure 2: Raw output from "weak scaling" job. Timed out on *slurmstepd* due to time out / time limit.

90	1	0.110155
90	1	0.109749
90	1	0.109885
90	2	0.056617
90	2	0.056599
90	2	0.056612
90	4	0.045880
90	4	0.045966
90	4	0.045863
90	8	0.031120
90	8	0.031132
90	8	0.031170
90	16	0.018182
90	16	0.018227
90	16	0.018220
90	24	0.013238
90	24	0.013257
90	24	0.013180
90	32	0.014816
90	32	0.017296
90	32	0.014814
1100	1	16.306608
1100	1	16.316588
1100	1	16.284397
1100	2	8.175213
1100	2	8.178992
1100	2	8.170321
1100	4	6.621239
1100	4	6.678632
1100	4	6.639713
1100	8	4.557337
1100	8	4.554004
1100	8	4.586490
1100	16	2.447131
1100	16	2.448894
1100	16	2.447200
1100	24	1.731222
1100	24	1.718731
1100	24	1.718424
1100	32	1.312658
1100	32	1.313263
1100	32	1.320209

Figure 3: Raw output from "juliap" job.

"static"	1100	16	2.450491
"static"	1100	16	2.448260
"static"	1100	16	2.449136

Figure 4: Raw output from "juliap2" job.