Task 01:

After generating a random array of size n using the method generate_array which is filled with random integers from 0 to 2n a block of parallel code execution using omp is started in line 73. However, the initial splitting is only supposed to be done exactly once which is why the term #pragma omp single is added to ensure not every thread executes this operation. \

The method quicksort receives an array as well as two indices and sorts the numbers in the given interval by their numeric value. The method divide performs the splitting of the array into two parts using a chosen pivot element and returns the new index of the pivot element which may now be at any index in the given interval. quicksort now recursively calls itself to also sort the new sublists above and below the returned position. To parallelize this omp task is used which creates a new subtask by the next code block with then is executed synchronously by the next available thread.

n	2 PEs	4 PEs	8 PEs
8	0.0000s	0.0001s	0.0000s
80	0.0000s	0.0002s	0.0000s
800	0.0003s	0.0006s	0.0006s
8000	0.0038s	0.0048s	0.0071s
80000	0.0312s	0.0436s	0.0735s
800000	0.4357s	0.7839s	1.4535s
8000000	4.1268s	6.3749s	12.2290s
80000000	172.1905s	130.3264s	92.3354s

Task 02:

To use radix sort to sort an array of integer numbers at each step of the iteration, the algorithm needs to know how many numbers with a 0 and how many numbers with a 1 at the current index are in the array. For that the method one_amount counts the latter and returns it from which the other amount can be calculated using the size of the array. This information can be used to now sort all numbers by putting them in the first or the second part of the list according to that information, while still keeping them in the same order according to all previous iterations. Using the number of elements in the two classes helps in finding the first indices and then just inserting the numbers into a new empty list and updating the indices as needed. The pointer of the new array later is copied to the array to be sorted. \

To parallelize each thread, works on a small subpart of the array to count the class amounts and later to insert them into the new array. A prefix sum has to be used to find the starting indices for each class in each thread.

n	2 PEs	4 PEs	8 PEs
8	0.0015s	0.0019s	0.0032s
80	0.0017s	0.0022s	0.0033s
800	0.0013s	0.0016s	0.0032s
8000	0.0028s	0.0025s	0.0041s
80000	0.0136s	0.0087s	0.0084s

n	2 PEs	4 PEs	8 PEs
800000	0.1426s	0.0964s	0.0949s
8000000	1.5822s	1.1046s	0.8035s
80000000	15.7717s	10.9164s	7.9710s

Task 04:

```
def find_min_pancake_num(initial_config):
   n = size of initial_config
    # general upper bound for pancake problem (proof omitted)
    best = ceil(18 / 11 * n)
    # stack to save a configuration and the amount of steps it took to get there
    s: Stack of pairs (pancake configuration plus an integer number)
    push the pair (initial_configuration, 0) onto s
    while s not empty:
        current_config, num_steps = pop top element from s
        if num_steps > best:
            # stop current iteration if the current iteration can not beat best found solution
            continue
        else:
            parallel for i from 2 to n:
                next_config = flip top i pancakes of current_config
                if next_config is sorted and num_steps + 1 < best:</pre>
                    atomic update best = num_steps + 1
                else:
                    push the pair (next_config, num_steps + 1) onto s
    return best
```

Exercise 1 (Unimplemented parts)

We didn't implement the Parallel Partition successfully because of some technical details. But I will give the pseudocode of the parallel partition below. It may look verbose and complicated, but the key point is to find the correct index to insert the elements. (See line 32. - 34.)

Parallel Partition

```
procedure ParallelPartition(A, n, pivot, offset)
       ⊳ The following 3 arrays are used for storing the count of element in each processors, that are
       smaller/equal/larger to pivot.
3:
       localSmaller \leftarrow Array[1..p]
4:
       localEqual \leftarrow Array[1..p]
5:
       localBigger \leftarrow Array[1..p]
6:
       \triangleright B as a temporary array to store the rearranged array
7:
       B \leftarrow \text{Array}[1..n]
       for tid=1..p in parallel do
8:
9:
          smallerCount \leftarrow 0
10:
          equal \leftarrow 0
          biggerCount \leftarrow 0
11:
12:
          for j=start(tid)...end(tid) do
13:
             if A[j + \text{offset}] < \text{pivot then}
14:
                smallerCount + +
15:
                else if A[j + offset] == pivot then
16:
                  equal + +
17:
                end
18:
                else
19:
                  biggerCount + +
20:
                end
21:
             end
22:
          end
23:
          localSmaller[tid] \leftarrow smallerCount
24:
          localEqual[tid] \leftarrow equal
          localBigger[tid] \leftarrow biggerCount
25:
26:
27:
       prefixSmaller \leftarrow pPrefAdd(localSmaller)
28:
       prefixEqual \leftarrow pPrefAdd(localEqual)
29:
       prefixBigger \leftarrow pPrefAdd(localBigger)
       ▶ Let prefix*[0] be 0
30:
31:
       for tid=1..p in parallel do
          sIdx \leftarrow prefixSmaller[tid - 1]
32:
          eIdx \leftarrow prefixSmaller[p] + prefixEqual[tid-1]
33:
34:
          bIdx \leftarrow prefixSmaller[p] + prefixEqual[p] + prefixBigger[tid - 1]
35:
          for j=start(tid)...end(tid) do
36:
             if A[j + \text{offset}] < \text{pivot then}
                B[\operatorname{sIdx} + +] \leftarrow A[j + \operatorname{offset}]
37:
                else if A[j + offset] == pivot then
38:
                   B[eIdx + +] \leftarrow A[j + offset]
39:
40:
41:
                  B[\text{bIdx} + +] \leftarrow A[j + \text{offset}]
42:
43:
                end
             end
44:
          end
45:
46:
47:
       Copy B back to A[offset..offset + n]
48: end
```

Exercise 3

3.1

The **insert** operation can result in concurrent write if it is not carefully designed, given the case where $\mathbf{index}(A[i]) == \mathbf{index}(A[j])$.

The other parts are guaranteed to be thread-safe.

3.2

For a parallel version of $\mathbf{copy}(A, B)$, we will need the help of another array C.

```
procedure Parallel Copy(A, B)
         \triangleright C stores the size of each bucket in B
3:
         C \leftarrow \text{Array}[1..\text{nb}]
4:
         \mathbf{pfill}(C,0)
         for i=1..nb in parallel do
5:
6:
            C[i] \leftarrow \operatorname{len}(B[i])
7:
         end
        P \leftarrow \mathbf{pPrefAdd}(C)
8:
        for i=1..nb-1 in parallel do
9:
            beg \leftarrow P[0] as 0 beg \leftarrow P[i-1] + 1
10:
11:
            end \leftarrow P[i]
12:
            for j=beg..end in parallel do
13:
14:
               A[j] \leftarrow B[i][j - \text{beg}]
15:
            end
16:
         end
17: end
```

3.2.1 Comments

- For the beg and end in the above algorithm, we note that (P[i] (P[i-1] + 1) + 1) is equal to C[i], namely len(B[i]).
- I don't see the necessity of using **pfill**(). It's likely used for avoiding data corruption.

3.2.2 Complexity Analysis

Operation	Time Complexity	Processors Required
Get the lengths	O(1)	nb
Parallel Prefix Sum	$O(\log \mathrm{nb})$	nb/2
Insert	O(1)	n
Overall	$O(\log \mathrm{nb})$	n

For fast Parallel Bucket Sort, we need enough number of buckets. However, when it comes to the parallel copy operation, a lower nb is preferred. Herein lies a tradeoff.

Exercise 5

(a)

Given an index i, to find the final position j of a_i in the merged list, we can use Binary Search to find the biggest element b_k in B that is smaller than a_i (or the smallest element in B that is bigger

than a_i), which requires $O(\log n)$ comparisons. (See section **Parallel Merge sort** in the lecture slide for example) Then, j = i + k, as there are i + k elements in total that are smaller than a_i .

(b)

Denote by **computeInsertPos**(A, B, i) the algorithm described in (a), which returns the final position j to insert our $a_i = A[i]$.

```
1: \mathbf{procedure} \ \mathsf{Parallel\_Merge}(A,B)
2: C \leftarrow \mathsf{Array}[1..2n]
3: \mathbf{for} \ i = 1..n in parallel \mathbf{do}
4: \flat \ j_1 \neq j_2 as the elements are unique
5: j_1 \leftarrow \mathbf{computeInsertPos}(A,B,i)
6: j_2 \leftarrow \mathbf{computeInsertPos}(B,A,i)
7: C[j_1] \leftarrow A[i]
8: C[j_2] \leftarrow B[i]
9: \mathbf{end}
10: \mathbf{end}
```

(b.1) Comments

- If the elements are not unique and $j_1 == j_2$ by accident, we can set j_2 as $j_1 + 1$ to avoid data race.
- The total work T_1 is apparently $O(n\log n)$ and the span T_∞ is $O(\log n)$.

(c)

This task seems impossible from the software perspective. With O(n) processors, we can compare a_i with $b_{1..n}$ in O(1). (but don't we need to broadcast a_i to all the processors?)

That said, it's still not possible to reduce the results in O(1). Consider the following 2 cases:

```
1. \max_{\{k: a_i \ge b_k\}} k;
2. \sum_{\kappa \in \{k: a_i \ge b_i\}} 1.
```

In both cases, the reduction takes $O(\log \#\{k: a_i \ge b_k\})$. O(1) is only made possible if we have some hardware "hacks" or some magic.

If such a hack exists, we can use $n \cdot O(n) = O(n^2)$ processors for merging two lists in O(1).