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Analysis and Design of Algorithms

Homework 2

Question -1:

Let the three candies types “circus peanuts”, “Heath bars”, and “Cioccolateria Gardini chocolate truffles” be represented by integers 0, 1 and 2 respectively.

Sub-problems:

i - denotes the i^{th} of the animal

j - denotes the candy type

Let the animals be $a_0, a_1, \dots, a_i, \dots, a_{n-1}$ (n is the number of animals)

Let $\forall i \in [0, n)$ and $\forall j \in [0, 2]$, $\text{scores}[i][j]$ denote the optimal answer for the game with animals $a_i, a_{i+1}, \dots, a_{n-1}$ and the candy in the hand of the player is ' j '

Recurrence:

1. Base Case: $i == n$ (no animal remaining to swap candy)
 $\text{scores}[i][j] = 0, \forall j \in [0, 2]$
2. The player and animal have the same candy
 $\text{scores}[i][j] = 1 + \text{scores}[i + 1][j]$
3. The player and animal do not have the same candy
 $\text{scores}[i][j] = \max(-1 + \text{scores}[i + 1][\text{candies}[i]], \text{scores}[i + 1][j])$ {here $\text{candies}[i]$ is the candy present with the i^{th} animal}

Final Solution:

$\text{scores}[0][0]$ is our final answer

The player tries to swap his candy with all the animals from a_0 to a_{n-1} ; with the initial candy present with him to be of type 0 i.e. “circus peanuts”.

Correctness of Recurrence:

For the i^{th} animal, one of the following cases, are possible,

1. Player and the Animal have the same candy
2. Player and the Animal do not have the same candy

Case 1: Player and the Animal have the same candy

For this case the player is free to either swap or not swap the candy with the animal.

If the player wishes to swap then $\text{score}[i][j]$ would become $1 + \text{score}[i+1][j]$.

If the player doesn't want to swap then the $\text{score}[i][j]$ would remain the same.

Clearly the best choice for the player is to swap the candy and increase the score.

That is if $\text{candies}[i] == j$ then $\text{score}[i][j] = 1 + \text{score}[i+1][j]$

Case 2: Player and the Animal do not have the same candy

For this case the player is again free to either swap or not swap the candy with the animal.

If the player wishes to swap then $\text{score}[i][j]$ would become $-1 + \text{score}[i+1][\text{candies}[i]]$.

If the player doesn't want to swap then the $\text{score}[i][j]$ would remain the same.

Here the score is not just dependent on the effect of the swap process on the score, it is also dependent on the score the player gains after he receives the i^{th} animal candy in swapping.

So, the optimal answer in this case is to take the maximum score of both the cases

That is if $\text{candies}[i] != j$ then $\text{score}[i][j] = \max(\text{score}[i+1][j], -1 + \text{score}[i+1][\text{candies}[i]])$

Pseudocode:

```

FUNCTION CandyScoreRoutine(candies, n, i, candy, scores):
    IF i = n:
        RETURN 0
    ENDIF

    IF scores[i][candy] != -inf:
        RETURN scores[i][candy]
    ENDIF

    IF candies[i] = candy:
        scores[i][candy] <- 1 + CandyScoreRoutine(candies, n, i + 1, candy, scores)
    ELSE:
        scores[i][candy] <- max(-1 + CandyScoreRoutine(candies, n, i+1, candies[i], scores),
CandyScoreRoutine(candies, n, i+1, candy, scores))
    ENDIF

    RETURN scores[i][candy]

ENDFUNCTION

FUNCTION CandyScore(candies):
    types <- 3
    startCandy <- 0
    n <- len(candies)

    scores <- [[-inf for _ in range(types)] for _ in range(n)]

```

```
RETURN CandyScoreRoutine(candies, n, 0, startCandy, scores)
```

```
ENDFUNCTION
```

Runtime :

$O(n)$:- We have $3 * n$ states in our DP, each of which is visited only once, hence the total run time of our algorithm is $O(n)$.

Question-2:

Sub-problems:

i - denotes the i^{th} of the house

j - denotes the number of bakeries

Let the houses be $h_0, h_1, \dots, h_i, \dots, h_{n-1}$ (n is the number of houses)

Let $\forall i \in [0, n)$ and $\forall j \in (0, k]$, $\text{sumDistances}[i][j]$ denote the optimal answer for the houses h_0, h_1, \dots, h_i having j spots for the bakeries distributed among themselves.

RecurrenceMarks:

1. Base Cases:

a. $i=-1$ and $j=0$, i.e. all k spots have been picked.

$$\text{sumDistances}[i][j] = 0$$

b. $i=-1$ and $j \neq 0$, i.e j spots for bakeries have not been picked.

$$\text{sumDistances}[i][j] = \text{inf}$$

c. $i \neq -1$ and $j=0$, i.e i houses are not considered while picking spots for the bakeries

$$\text{sumDistances}[i][j] = \text{inf}$$

2. $\forall i \in [0, n)$ and $j > 0$

$$\text{sumDistances}[i][j] = \min_{k=0 \dots i} (\text{sumDistances}[k-1][j-1] + \text{sumMedianDistance}[k][i]),$$

where $\text{sumMedianDistance}[k][i]$ is the sum of distances walked by people of houses k to i , each day to visit bakery present in a house h_b with $k \leq b \leq i$.

Final Solution :

$$\text{sumDistances}[n-1][k],$$

$(n-1, k)$ means we are picking k bakeries from houses 0 to $n-1$, for positioning our bakeries.

Correctness of Recurrence :

For a range of houses, h_0, h_1, \dots, h_i with k spots to be picked up, we can break the problem into two parts, picking $k-1$ spots in the range h_0, h_1, \dots, h_{j-1} and picking 1 spot in the range h_j, h_{j+1}, \dots, h_i (for some j such that $0 \leq j \leq i$).

Then the sum of distances to the nearest bakery in the range $(0, i)$ with k bakeries, i.e $\text{sumDistance}[i][k]$ would become $\text{distance}[j][i] + \text{sumDistance}[j-1][k-1]$, where $\text{distance}[j][i]$ is the minimum sum of all the distances to the nearest bakery for the range of houses h_j, h_{j+1}, \dots, h_i .

Therefore the $\text{sumDistance}[i][k] = \min_{j=0 \dots i} (\text{distance}[j][i] + \text{sumDistance}[j-1][k-1])$

Similarly, $\text{sumDistance}[j-1][k-1]$ can be computed optimally.

So, we can deduce that for an appropriate unique value of j , the quantity $\{ \text{distance}[j][i] + \text{sumDistance}[j-1][k-1] \}$ would be minimum and hence, $\text{sumDistance}[i][k]$ would be computed optimally as the least minimum sum of distances.

Pseudocode:

```
FUNCTION getMinimumDistanceRoutine(distance, sumDistances, n, k):
    IF n = -1 AND k = 0:
        RETURN 0
    ELSEIF n = -1 OR k = 0:
        RETURN inf
    ENDIF
    IF sumDistances[n][k] != inf:
        RETURN sumDistances[n][k]
    ENDIF

    for i in range(n, -1, -1):
        sumDistances[n][k] <- min(sumDistances[n][k], distance[i][n] +
getMinimumDistanceRoutine(distance, sumDistances, i-1, k-1))
    ENDFOR

    RETURN sumDistances[n][k]

ENDFUNCTION

FUNCTION getMinimumDistance(houses, n, k):
    sumDistances <- [[inf for _ in range(k+1)] for _ in range(n)]

    distance <- [[0 for _ in range(n)] for _ in range(n)]

    for i in range(n):
        for j in range(i+1, n):
            medianHouse <- houses[(i+j)//2]
            for l in range(i, j+1):
                distance[i][j] += abs(medianHouse - houses[l])
            ENDFOR
        ENDFOR
    ENDFOR

    RETURN getMinimumDistanceRoutine(distance, sumDistances, n-1, k)

ENDFUNCTION
```

Runtime :

$O(n^3)$ -

The matrix sumMedianDistance is filled in $O(n^3)$ computations.

The matrix sumDistances has $n*k$ states, each of which requires n computations to be filled.

Hence the whole sumDistances is filled in $O(n^2 * k)$ computations.

Therefore total Complexity is $O(n^3 + n^2 * k) = O(n^3)$ [$k \leq n$]

Question-3:

Sub-problems:

Let n denote the number of drops.

Let $\forall i \in [0, n)$ and $\forall j \in [i, n)$, $\text{energy}[i][j]$ denote the maximum energy gained when drops d_i, d_{i+1}, \dots, d_j are merged to form a single drop

Recurrence:

1. Base Case: $i=j$
 $\text{energy}[i][i] = 0$, energy gained by just a single drop is 0 i.e the initial energy a drop possesses is 0.
2. $\forall j > i$
 $\text{energy}[i][j] = \max(\text{energy}[i][j-1] + \text{combinedDrops}[i][j-1]^2 + \text{drops}[j]^2, \text{energy}[i+1][j] + \text{drops}[i]^2 + \text{combinedDrops}[i+1][j]^2)$
3. $\forall j < i$
 $\text{energy}[i][j] = 0$

Final Solution:

$\text{energy}[0][n-1]$ is the required maximum energy, when drops d_0, d_1, \dots, d_{n-1} are merged.

Correctness of Recurrence :

Let $\text{energy}[i][j]$ denotes the energy gained when all the drops from d_i to d_j are combined.

We can combine the drops i to j in primarily two ways:

1. Combine the drops i to $j-1$ and then combine this combined drop with the j^{th} drop.
In this the total energy gained till this point would be $\text{energy}[i][j-1]$ (energy gained while combining drops i to $j-1$) and $\text{combinedDrops}[i][j-1]^2$ (volume of the combined drop from i to $j-1$) + $\text{drops}[j]^2$, which is-
 $\text{energy}[i][j-1] + (\text{combinedDrops}[i][j-1]^2 + \text{drops}[j]^2)$
2. Combine the drops $i+1$ to j and then combine this combined drop with the i^{th} drop
Drawing on the lines of the method followed in the 1st point, energy gained in this case would be $(\text{drops}[i]^2 + \text{combinedDrops}[i+1][j]^2) + \text{energy}[i+1][j]$

Now, the maximum energy gained when combining drops from d_i to d_j would be

$$\text{energy}[i][j] = \max(\text{energy}[i][j-1] + (\text{combinedDrops}[i][j-1]^2 + \text{drops}[j]^2), (\text{drops}[i]^2 + \text{combinedDrops}[i+1][j]^2) + \text{energy}[i+1][j])$$

Recursively applying the same logic on smaller parts, we can have all the possible ways to merge drops d_i to d_j , the maximum desired energy would be the max of all possible energies at each level of merging.

Pseudocode :

```

FUNCTION getMaxEnergy(drops):
    n <- len(drops)

    combinedDrops <- [[0 for _ in range(n)] for _ in range(n)]
    energy <- [[0 for _ in range(n)] for _ in range(n)]

    for i in range(n):
        combinedDrops[i][i] <- drops[i]
    ENDFOR

    for i in range(1, n):
        for j in range(n-i):
            r <- j
            c <- j+i

            combinedDrops[r][c] <- drops[r] + combinedDrops[r+1][c]

            energy1 <- ( drops[r]^2 + combinedDrops[r+1][c]^2 ) + energy[r+1][c]
            energy2 <- energy[r][c-1] + ( drops[c]^2 + combinedDrops[r][c-1]^2 )

            energy[r][c] <- max(energy1, energy2)

        ENDFOR
    ENDFOR

    maxEnergy <- energy[0][n-1]

    RETURN maxEnergy

ENDFUNCTION

```

Runtime :

$O(n^2)$:

There are n^2 entries in our energy matrix, each of which gets filled only once, which is $O(n^2)$ computations to fill the energy matrix.