COMP550 Midterm1 Take Home Part

Siyang Jing

730042022

Definitions:

- $n \in \mathbb{N}^+$, the numer of potions;
- $m \in \mathbb{N}$, the total stench, m is initially 0;
- sequence v[1, 2, ..., n], the (initial) stench factors, where $\forall i \in \{1, 2, ..., n\}, v[i] \in \mathbb{Z}_{50}(i.e.\{0, 1, ..., 49\})$;
- function (,): $\mathbb{Z}_{50} \times \mathbb{Z}_{50} \to \mathbb{Z}_{50}$, the action of pouring potion, defined as (a, b) = a + b, here + is the addition in \mathbb{Z}_{50} ;
- sequence of evaluation should be done sequentially, i.e. we can only pour potions from adjacent one, i.e. (a, b) should only be done if a and b are numbers (in other words expressions that have already been evaluated), e.g. (1, (2, 3)) should be done in the order of (2, 3) and then (1, 5);
- each evaluation (a, b) increments m by $a \times b$, here \times is the regular multiplication, i.e. the effect of pouring potion on the total stench;
- sequence of function evaluations $F_{\sigma}(i, k)$ that sequentially evaluates (,) on the elements of sequence v[i, ..., k] in a certain order, i.e. the order of pouring potion, where $\sigma \in \mathfrak{F}(i, k)$ is some index denoting one possible way of evaluating v[i, ..., k], and $\mathfrak{F}(i, k)$ is the index set for $F_{\sigma}(i, k)$, for example $F_{\mu}(1,3) = (1,(2,3))$, and $F_{\nu}(1,3) = ((1,2),3)$, and etc;
- we notice that the value obtained from the evaluation $F_{\sigma}(i,k)$ does not depend on σ , therefore we define F(i,k) = value of $F_{\sigma}(i,k) = \sum_{l=i}^{l=k} v[l]$;
 - F(i,i) := v[i], which is a simple value, and no evaluation is needed;
- the minimum total stench:
 - $\hat{m} = \min_{\sigma \in \mathfrak{F}(1,n)} \{ m \text{ after sequence of evaluations } F_{\sigma}(1,n) \};$
- the minimum total stench after pouring potions i to k:

$$\hat{m}(i,k) = \min_{\sigma \in \mathfrak{F}(i,k)} \{ m \text{ created by } F_{\sigma}(i,k) \}, \text{ note that } \hat{m}(1,n) = \hat{m};$$

$$\hat{m}(i,i) := 0;$$

• the order of pouring potions i to k, that creates the minimum stench:

$$\hat{F}(i,k) = \underset{\sigma \in \mathfrak{F}(i,k)}{\operatorname{argmin}} \{ m \text{ created by } F_{\sigma}(i,k) \}$$

Statement of the problem:

The aim is to develop an algorithm that,

given
$$v[1,...,n]$$
, finds $\hat{m} = \min_{\sigma \in \mathfrak{F}(1,n)} \{ m \text{ after sequence of evaluations } F_{\sigma}(1,n) \}$,

i.e. finds the order of pouring that creates the minimum total stench.

Statement of the algorithm:

First we notice that, there is a one-one correspondence between

$$\sigma \in \mathfrak{F}(i,k)$$
 and $(j \geqslant i, \mu \in \mathfrak{F}(i,j), \nu \in \mathfrak{F}(j+1,k))$, and, in particular,

$$\forall k > i+1, \sigma \in \mathfrak{F}(i,k), \exists ! j \geqslant i, \mu \in \mathfrak{F}(i,j), \nu \in \mathfrak{F}(j+1,k), s.t. F_{\sigma}(i,k) = (F_{\mu}(i,j), F_{\nu}(j+1,k)),$$

and therefore we have:

$$\hat{m}(i,k) = \min_{\sigma \in \mathfrak{F}(i,k)} \{ m \text{ created by } F_{\sigma}(i,k) \}$$

$$= \min_{i,\mu,\nu} \{ m \text{ created by } (F_{\mu}(i,j), F_{\nu}(j+1,k)) \}$$

$$=\min_{j,\mu,\nu}\{m \text{ created by } F_{\mu}(i,j)+m \text{ created by } F_{\nu}(j+1,k)+F(i,j) imes F(j+1,k)\}$$

$$=\min_{i}\{\min_{\mu,\nu}\{m \text{ created by } F_{\mu}(i,j)+m \text{ created by } F_{\nu}(j+1,k)+F(i,j)\times F(j+1,k)\}\}$$

$$= \min_{j} \{ \min_{\mu} \{ m \text{ created by } F_{\mu}(i,j) \} + \min_{\nu} \{ m \text{ created by } F_{\nu}(j+1,k) \} + F(i,j) \times F(j+1,k) \}$$

$$= \min_{j} \{\hat{m}(i,j) + \hat{m}(j+1,k) + F(i,j) imes F(j+1,k) \}$$

which is the optimal substructure of the problem.

To develop an algorithm from this optimal substructure, we either

- first recursively call the function to compute $\hat{m}(i,j)$ and $\hat{m}(j+1,k)$, and then iterate over j to find $\hat{m}(i,k)$, which describes a top-down recursive algorithm,
- or we can first compute all the $\hat{m}(i, i)$, then $\hat{m}(i, i+1)$, and then $\hat{m}(i, i+2)$, till $\hat{m}(1, n)$ to get the desired solution, which describes a bottom-up algorithm.

Either way, we need to memoize/store the values of $\hat{m}(i, j)$ that have been calculated to avoid repetitive calculation of the same value, and we also need calculate and store the values of F(i, j) beforehand to avoid repetition.

The time cost for bottom-up solution is $\Theta(n^3)$, and the space cost is $\Theta(n^2)$, there is no worst case, best case, or average case for the algorithm, since in any situation, we calculate everything.

Below is the pseudo code for the bottom-up algorithm:

The input v is a sequence (1D array), which stores the sequence v described above.

The m 2D array stores $\hat{m}(i, j)$ described above in m[i, j], and m[i, i] = 0.

The f 2D array stores the value of the function evaluation F(i,j) at f[i,j], and f[i,i] = v[i].

The output is $m[1, n] = \hat{m}(1, n) = \hat{m}$, the minimum total stench.

Algorithm Potion-Adding-Order(v)

```
n=v.\mathrm{length} let m[1...n,1...n] and f[1...n,1...n] be new tables for i=1 to n m[i,i]=0
```

```
f[i,i] = v[i]
   for j = i + 1 to n
       f[i, j] = f[i, j-1] + v[j] \mod 50
   end
end
for l=2 to n
   for i = 1 to n - l + 1
       i = i + l - 1
       m[i,j] = \infty
       for k = i to i - 1
           q = m[i, k] + m[k+1, j] + f[i, k] \times f[k+1, j]
           if q < m[i, j]
               m[i, j] = q
           end
       end
   end
end
return m[1, n]
```

Argument for Optimality (for the bottom up solution):

First, I argue that no matter what method is employed, we need to find $\hat{m}(i,k)$ for every i,k.

Suppose we did not find $\hat{m}(i, k)$, i.e. for $\sigma \in \mathfrak{F}(i, k)$, we don't know which σ corresponds to the $F_{\sigma}(i, k)$ that produces $\hat{m}(i, k)$.

```
Then we don't know \min_{\sigma} \{\hat{m}(1,i-1) + m \text{ created by } F_{\sigma}(i,k) + F(1,i-1) \times F(i,k)\},
which equals \min_{\mu,\nu} \{m \text{ created by } F_{\mu}(1,i-1) + m \text{ created by } F_{\nu}(i,k) + F(1,i-1) \times F(i,k)\}.
Then we don't know \min_{j} \{\min_{\mu,\nu} \{m \text{ created by } F_{\mu}(1,j-1) + m \text{ created by } F_{\nu}(j,k) + F(1,j-1) \times F(j,k)\}\},
which equals \min_{\sigma \in \mathfrak{F}(i,k)} \{m \text{ created by } F_{\sigma}(1,k)\}, which is simply \hat{m}(1,k).
```

By the same logic, we reason that we don't know $\hat{m}(1,n)$, i.e. even if we get some value, correct or not, we don't know whether it is the minimum, whether it is the answer.

Second, I argue that in finding each $\hat{m}(i,k)$, we need to iterate through all the values j between i and k to calculate $\hat{m}(i,j) + \hat{m}(j+1,k) + F(i,j) \times F(j+1,k)$.

Suppose we did not go through j_0 , then we don't know $\min_j \{\hat{m}(i,j) + \hat{m}(j+1,k) + F(i,j) \times F(j+1,k)\}$, since this j_0 we did not go through could be exactly the one that produces the above minimum. Therefore we don't know $\hat{m}(i,k)$.

From the argument above, we can see that we need $\Theta(n^3)$ calculations of $\hat{m}(i, j) + \hat{m}(j + 1, k) + F(i, j) \times F(j+1, k)$, for each $i \leq j < k$, no matter the form/implementation of the algorithm.

In our bottom-up solution which also stores all calculated $\hat{m}(i,j)$ and F(i,j), we have constant time cost for calculating $\hat{m}(i,j) + \hat{m}(j+1,k) + F(i,j) \times F(j+1,k)$. When we need to calculate this expression, since the lag between i and j, or j+1 and k, is less than the lag between i and k, we will have already calculated and stored the values of $\hat{m}(i,j)$, $\hat{m}(j+1,k)$, F(i,j), and F(j+1,k), and therefore the calculation is constant. Since the calculations are necessarily needed, and the cost for each calculation is optimal (constant), the whole algorithm is optimal.