Advanced Machine Learning - Assignment 2

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Exercise 1

Statement:

Consider $\mathcal{H} = \mathcal{H}_1 \cup \mathcal{H}_2$ where

$$\mathcal{H}_1 = \{h_{\theta_1} : \mathbb{R} \to \{0,1\}, \ h_{\theta_1}(x) = \mathbf{1}_{[x \ge \theta_1]}(x) = \mathbf{1}_{[\theta_1,+\infty)}(x), \ \theta_1 \in \mathbb{R}\} \text{ and }$$

$$\mathcal{H}_2 = \{h_{\theta_2} : \mathbb{R} \to \{0, 1\}, \ h_{\theta_2}(x) = \mathbf{1}_{[x < \theta_2]}(x) = \mathbf{1}_{(-\infty, \theta_2)}(x), \ \theta_2 \in \mathbb{R}\}$$

- a) Give an efficient ERM algorithm for learning ${\cal H}$ and compute its complexity for the realizable case.
- b) Give an efficient ERM algorithm for learning ${\cal H}$ and compute its complexity for the agnostic case.
- c) Compute the shattering coefficient $\tau_H(m)$ of the growth function for $m \ge 0$ for hypothesis class \mathcal{H} .
- d) Compare your result with the general upper bound for the growth functions and show that $\tau_H(m)$ obtained at previous point c is not equal to the upper bound.
- e) Does there exist a hypothesis class \mathcal{H} for which $\tau_H(m)$ is equal to the general upper bound (over \mathbb{R} or another domain χ)? If your answer is yes please provide an example, if your answer is no please provide a justification.

Solution:

Point (a): We are in the realizable case \Rightarrow

Given a training set $S = \{(x_1, y_1), (x_2, y_2), ..., (x_m, y_m) \mid x_i \in \mathbb{R}, y_i \in \{0, 1\}, i = \overline{1, m}\}$ exists a function $h_{\theta^*} \in \mathcal{H}$ such that $y_i = h_{\theta^*}(x_i)$.

We consider the following scenarios:

- 1) ----- (only negative examples)

- 4) ++++++++ θ_2^* ----- (case when $h_\theta^* \in \mathcal{H}_2$)

Based on the previous scenarios we can consider the following algorithm:

```
Input: S
Output: h_{\theta_1}^s or h_{\theta_2}^s
Initialization step: \theta_1 = +\infty; \theta_2 = -\infty;
\alpha = arg\min_{i=1,m} x_i; // The index of the first sample in S

// Checking the label of the first sample to choose a hypotesis class (\mathcal{H}_1 or \mathcal{H}_2)

if y_{\alpha} == 0 then

// Set as threshold as the first positive example

\theta_1 = \min_{\substack{i=1,m \\ y_i=1}} x_i if there is such x_i

return h_{\theta_1}

if y_{\alpha} == 1 then

// Set as threshold the last positive example

\theta_2 = \max_{\substack{i=1,m \\ y_i=1}} x_i

return h_{\theta_2}
```

Complexity:

- Initialization step: O(m)

- Finding θ_1 or θ_2 : O(m)

Total Complexity: O(m).

Point (b): We are under the agnostic case \Rightarrow it might not be a function h_{θ} that labels the data (same point might have different labels) or if h_{θ} exists, it might not be in \mathcal{H} .

Let's consider training set $S = \{(x_1, y_1), (x_2, y_2), ..., (x_m, y_m) \mid x_i \in \mathbb{R}, y_i \in \{0, 1\}, i = \overline{1, m}\}$. Based on the fact that we are under agnostic case, we can have $x_i = x_j$ and $y_i \neq y_j$.

We consider the following algorithm:

```
Input: S
Output: h_{\theta_1}^s or h_{\theta_2}^s
Sort S in ascending order of x's.

We obtain S = \{(x_{\sigma(1)}, y_{\sigma(1)}), (x_{\sigma(2)}, y_{\sigma(2)}), \dots, (x_{\sigma(m)}, y_{\sigma(m)})\} with x_{\sigma(1)} \leq x_{\sigma(2)} \leq \dots \leq x_{\sigma(m)}.

Generate a set Z containing values of x without repetition:

Z = \{z_1, z_2, \dots, z_n\} with z_1 = x_{\sigma(1)} < z_2 < \dots z_n = x_{\sigma(m)}

Initialization step:

positive_left = 0; negative_left = 0;
positive_right = # points with label y_i = 1 with i = \overline{1, n};
negative_right = # points with label y_i = 0 with i = \overline{1, n};

if negative_right = 0 then // all samples are positive

\theta_1^s = z_1; return h_{\theta_2^s}

if positive_right = 0 then // all samples are negative

\theta_2^s = z_1; return h_{\theta_2^s}
```

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We consider:
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Loss_1(\theta) = \frac{\# left\_positive_\theta + \# right\_negative_\theta}{n} = error \text{ of } h_{\theta_1} \text{ at threshold } \theta \text{ where } h_{\theta_1} \in \mathcal{H}_1
Loss_2(\theta) = \frac{\# left\_negative_\theta + \# right\_positive_\theta}{n} = error \text{ of } h_{\theta_2} \text{ at threshold } \theta \text{ where } h_{\theta_2} \in \mathcal{H}_2
```

Compute initial Loss₁ and Loss₂ values

```
\theta_1^s = z_1; \theta_2^s = z_1; L_1 = Loss_1(\theta_1^s); L_2 = Loss_2(\theta_2^s)
```

```
for i \leftarrow 2 to n in
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// updating positive and negative counts on both sides
    // for each change of threshold
    if y_i == 0 then
        negative_left = negative_left +1
        negative_right = negative_right -1
    if y_i == 1 then
        positive_left = positive_left +1
        positive_right = positive_right -1
    // recomputing Loss values for each {\cal H} class for the new threshold
    L_1^i = Loss_1(z_i)
    L_2^i = Loss_2(z_i)
    // selecting threshold with the best error for the both {\cal H} classes
    if L_1^i < L_1 then
        \theta_1^s = z_i
       L_1 = L_1^i
    if L_2^i < L_2 then
        \theta_2^s = z_i
        L_2 = L_2^i
if L_1 < L_2 then
    return h_{\theta_i^s}
```

Complexity:

return $h_{\theta_s^s}$

- 1. sorting: $O(m \cdot log(m))$
- 2. computing Z: O(m)
- 3. initialization step: O(m)
- 4. computing loss = O(1)
- 5. selecting best threshold: O(m)

Total Complexity: $O(m \cdot log(m))$

Point (c): We know that $\tau_{\mathcal{H}}(m) = \max_{C \subseteq X: |C| = m} |\mathcal{H}_C|$ and $\mathcal{H} = \mathcal{H}_1 \cup \mathcal{H}_2$ $\Rightarrow \tau_{\mathcal{H}}(m) = \tau_{\mathcal{H}_1 \cup \mathcal{H}_2}(m) = \max_{C \subseteq X: |C| = m} |\mathcal{H}_{1C}| + |\mathcal{H}_{2C}| - |\mathcal{H}_{1C} \cap \mathcal{H}_{2C}|.$

$$\Rightarrow \tau_{\mathcal{H}}(m) = \tau_{\mathcal{H}_1 \cup \mathcal{H}_2}(m) = \max_{C \subseteq X \cdot |C| = m} |\mathcal{H}_{1C}| + |\mathcal{H}_{2C}| - |\mathcal{H}_{1C} \cap \mathcal{H}_{2C}|.$$

We consider $C = \{c_1, c_2, ..., c_m\}$ a set of m points with $c_i \le c_j$.

For \mathcal{H}_{1C} we can have at most m + 1 different functions:

$$a_1 < c_1 < a_2 < c_2 < \dots < a_m < c_m < a_{m+1} \Longrightarrow |\mathcal{H}_{1C}| = \left| \{h_{a_1}, h_{a_2}, \dots \ h_{a_m}, h_{a_{m+1}} \} \right| = m+1. \ [1]$$

 h_{a_1} labels points c_1, c_2, \ldots, c_m with labels $(1, 1, \ldots, 1, 1)$.

 h_{a_2} labels points c_1, c_2, \ldots, c_m with labels $(0, 1, \ldots, 1, 1)$.

 h_{a_m} labels points c_1, c_2, \dots, c_m with labels $(0, 0, \dots, 0, 1)$.

 $h_{a_{m+1}}$ labels points c_1, c_2, \dots, c_m with labels $(0, 0, \dots, 0, 0)$.

For \mathcal{H}_{2C} we have at most m + 1 different functions:

$$b_1 < c_1 < b_2 < c_2 < \dots < b_m < c_m < b_{m+1} \Rightarrow |\mathcal{H}_{2C}| = |\{h_{b_1}, h_{b_2}, \dots h_{b_m}, h_{b_{m+1}}\}| = m+1.$$
 [2]

 h_{b_1} labels points c_1, c_2, \ldots, c_m with labels $(0, 0, \ldots, 0, 0)$.

 h_{b_2} labels points c_1, c_2, \ldots, c_m with labels $(1, 0, \ldots, 0, 0)$.

 h_{b_m} labels points c_1, c_2, \dots, c_m with labels $(1, 1, \dots, 1, 0)$.

 $h_{b_{m+1}}$ labels points c_1, c_2, \ldots, c_m with labels $(1, 1, \ldots, 1, 1)$.

We can observe that h_{a_1} is equivalent to $h_{b_{m+1}}$ and $h_{a_{m+1}}$ is equivalent to h_{b_1} . [3]

From [1], [2] and [3] we have $\tau_{\mathcal{H}_1 \cup \mathcal{H}_2}(m) = m + 1 + m + 1 - 2 = 2m$.

Point (d): Lemma Sauer-Shelah-Perles

Let \mathcal{H} be a hypothesis class with $VC\dim(\mathcal{H}) \leq d < \infty$. Then, for all m, we know that:

$$\tau_{\mathcal{H}}(m) \leq \sum_{i=0}^{d} C_{m}^{i}$$

In particular, if m > d + 1 then $\tau_{\mathcal{H}}(m) \le (\frac{em}{d})^d = O(m^d)$

We know that VC dim($\mathcal{H}_{thresholds}$) = 1, based on that fact it's trivial to prove that $VC \dim(\mathcal{H})$ = $VC\dim(\mathcal{H}_1\cup\mathcal{H}_2)=2.$

At the previous point, we proved that $\tau_H(m) = 2m$ [1]

Considering $VCdim(\mathcal{H}) = 2 = d$ we have the following cases:

Case 1: m > 3, we have that $\tau_{\mathcal{H}}(m) \leq \left(\frac{em\ d}{d}\right) \implies \tau_{\mathcal{H}}(m) \leq \left(\frac{em\ 2}{d}\right) \stackrel{[1]}{\Longrightarrow} 2m \leq \frac{e^2\ m^2}{4} \iff$ $8m \le e^2 m^2 \mid \cdot \frac{1}{m}$ (because we know that m > 0) \iff $8 \le e^2 m \iff \frac{8}{e^2} \le m \iff 1.082 \dots \le m$ (this holds for every m > 3) [2].

Case 2:
$$m = 1 \Rightarrow \tau_{\mathcal{H}}(m) \leq \sum_{i=0}^{2} C_{m}^{i} \Rightarrow \tau_{\mathcal{H}}(m) \leq C_{m}^{0} + C_{m}^{1} + C_{m}^{2} \Rightarrow \tau_{\mathcal{H}}(m) \leq 1 + m + \frac{m(m-1)}{2} = 1 + 1 + 0 = 2 \stackrel{[1]}{\Rightarrow} 2 \cdot m = 2 \cdot 1 = 2 \leq 2 \quad [3].$$

Case 3:
$$m = 2 \Rightarrow \tau_{\mathcal{H}}(m) \leq \sum_{i=0}^{2} C_{m}^{i} \Rightarrow \tau_{\mathcal{H}}(m) \leq C_{m}^{0} + C_{m}^{1} + C_{m}^{2} \Rightarrow \tau_{\mathcal{H}}(m) \leq 1 + m + \frac{m(m-1)}{2} = 1 + 2 + 1 = 4 \stackrel{[1]}{\Rightarrow} 2 \cdot m = 2 \cdot 2 = 4 \leq 4 \quad [4].$$

Case 4:
$$m = 3 \Rightarrow \tau_{\mathcal{H}}(m) \le \sum_{i=0}^{2} C_{m}^{i} \Rightarrow \tau_{\mathcal{H}}(m) \le C_{m}^{0} + C_{m}^{1} + C_{m}^{2} \Rightarrow \tau_{\mathcal{H}}(m) \le 1 + m + \frac{m(m-1)}{2} = 1 + 3 + 3 = 7 \stackrel{[1]}{\Rightarrow} 2 \cdot m = 2 \cdot 3 = 6 \le 7$$
 [5].

From [2], [3], [4], [5] we obtain that $\tau_{\mathcal{H}}(m) \leq \sum_{i=0}^{2} C_m^i$ as expected from Lemma Sauer-Shelah-Perles, to a more general form we have that 2m (our result obtained earlier) $\leq C_m^0 + C_m^1 + C_m^2 = 1 + m + \frac{m(m-1)}{2} = \frac{m^2 + m + 2}{2} \iff 2m \leq \frac{m^2 + m + 2}{2}$ for m > 0, which is correct, but is a high limit from the one computed.

Point (e): We know that $\tau_{\mathcal{H}}(m) = \max_{C \subseteq X} |\mathcal{H}_C|$ and from the *Lemma Sauer-Shelah-Perles* we know that $\tau_{\mathcal{H}}(m) \le \sum_{i=0}^d C_m^i$. Let's consider the definition from the lecture of $\mathcal{H}_{\text{thresholds}}$ such that

$$\mathcal{H}_{\text{thresholds}} = \left\{ h_a : \mathbb{R} \to \{0,1\} \mid h_a(x) = \begin{cases} 1, & x < a \\ 0, & \text{otherwise} \end{cases} \right\}$$

We know that $VC \dim(\mathcal{H}_{thresholds}) = 1$, based on that fact we get:

$$\tau_{\mathcal{H}_{\text{thresholds}}}(m) \le \sum_{i=0}^{1} C_m^i = C_m^0 + C_m^1 = 1 + m [1]$$

Let's consider $C = \{c_1, c_2, ..., c_m\}$ a set of m points, with $c_i < c_j$.

If we take $a_1 < c_1 < a_2 < c_2 < ... < c_m < a_{m+1}$ then \mathcal{H}_C will have $|\mathcal{H}_C| = |\{h_{a_1}, h_{a_2}, ..., h_{a_{m+1}}\}| = m+1$ (there will be at most m+1 functions, from *Lecture 8*). [2]

From [1] and [2] we conclude that $\mathcal{H}_{\text{thresholds}}$ has the growth function $(\tau_{\mathcal{H}}(m))$ equal to the general upper bound.

Exercise 2

Statement:

Consider a modified version of the AdaBoost algorithm that runs for exactly three rounds as follows:

- the first two rounds run exactly as in AdaBoost (at round 1 we obtain distribution $\mathbf{D}^{(1)}$, weak classifier h_1 with error ε_1 , at round 2 we obtain distribution $\mathbf{D}^{(2)}$, weak classifier h_2 with error ε_2).
- in the third round we compute for each i = 1, 2, ..., m:

$$D^{3}(i) = \begin{cases} \frac{D^{(1)}(i)}{Z}, & \text{if } h_{1}(x_{i}) \neq h_{2}(x_{i}), \\ 0, & \text{otherwise} \end{cases}$$

where Z is a normalization factor such that $\mathbf{D}^{(3)}$ is a probability distribution.

- obtain weak classifier h_3 with error ε_3 .
- output the final classifier $h_{final}(x) = sign(h_1(x) + h_2(x) + h_3(x))$.

Assume that at each round t = 1, 2, 3 the weak learner returns a weak classifier h_t for which the error ε_t satisfies $\varepsilon_t \le 1/2 - \gamma_t$, $\gamma_t > 0$.

- a) What is the probability that the classifier h_1 (selected at round 1) will be selected again at round 2? Justify your answer.
- b) Consider $\gamma = \min \{\gamma_1, \gamma_2, \gamma_3\}$. Show that the training error of the final classifier h_{final} is at most $\frac{1}{2} \frac{3}{2}\gamma + 2\gamma^3$ and show that this is strictly smaller than $\frac{1}{2} \gamma$.

Solution:

Point (a): We know that $\varepsilon_t \leq \frac{1}{2} - \gamma$, $\gamma > 0$ for t = 1, 2, 3For t = 1 let's assume that $\varepsilon_t = 0$ (our estimator does not make any mistakes, x_i is correctly classified, with $i = \overline{1, m}$) $\Rightarrow \Pr_{i > D^{(t)}}[h_t(x_i) \neq y_i] = \sum_{i=1}^m D^{(t)} \cdot \mathbf{1}_{[h_t(x_i) \neq y_i]} = 0$

Based on that, we have the distribution at iteration t + 1:

$$D^{(t+1)}(i) = \frac{D^{(t)}(i) \cdot e^{-w_t}}{Z_{t+1}} = \frac{D^{(t)}(i) \cdot \sqrt{\frac{\varepsilon_t}{1-\varepsilon_t}}}{Z_{t+1}} \quad [1]$$

where
$$Z_{t+1} = \sum_{i=1}^{m} D^{(t)}(i) \cdot e^{-w_t h_t(x_i) y_i} = \sum_{i=1}^{m} D^{(t)}(i) \cdot e^{-w} = \sum_{i=1}^{m} D^{(t)}(i) \cdot \sqrt{\frac{\varepsilon_t}{1-\varepsilon_t}} = \sqrt{\frac{\varepsilon_t}{1-\varepsilon_t}} \cdot \sum_{i=1}^{m} D^{(t)}(i) \xrightarrow{\frac{D^{(t)} probability}{distribution}} \sqrt{\frac{\varepsilon_t}{1-\varepsilon_t}} \cdot 1 = \sqrt{\frac{\varepsilon_t}{1-\varepsilon_t}} [2].$$

From [1] and [2] we have
$$D^{(t+1)}(i) = \frac{D^{(t)}(i) \cdot \sqrt{\frac{\ell_t}{1-\ell_t}}}{\sqrt{\frac{\ell_t}{1-\ell_t}}} = D^{(t)}(i)$$
, for $i = \overline{1, m}$.

This implies that if the first estimator has $\varepsilon_t = 0 \Rightarrow D^{(t+1)} = D^{(t)} \Rightarrow$ the estimator chosen at t+1 will be the same as the one at iteration t (the sample importance has not change) $\Rightarrow Pr[h_t = h_{t+1} \mid \varepsilon_t = 0] = 1$ [3]

For the general case when we have $\varepsilon_t \in (0, \frac{1}{2})$ we consider $D^{(t)}$ the probability distribution of importance to contain correctly classified and misclassified examples.

We can write this in a more formal manner as:

$$\sum_{i=1}^{m} D^{(t)}(i) \cdot e^{-w_{t}h_{t}(x_{i})y_{i}} = \sum_{i=1}^{m} D^{(t)}(i) \cdot e^{-w_{t}} \cdot \mathbf{1}_{[h_{t}(x_{i})=y_{i}]} + \sum_{i=1}^{m} D^{(t)}(i) \cdot e^{w_{t}} \cdot \mathbf{1}_{[h_{t}(x_{i})\neq y_{i}]}$$

We want to prove that $D^{(t)} \neq D^{(t+1)}$ because from that we have $h_t \neq h_{t+1}$.

For the misclassified samples we have that: $\sum_{i=1}^{m} D^{(t)}(i) \cdot e^{-w_t h_t(x_i) y_i} \cdot \mathbf{1}_{[h_t(x_i) \neq y_i]} = \sum_{i=1}^{m} D^{(t)}(i) \cdot e^{w_t} \cdot \mathbf{1}_{[h_t(x_i) \neq y_i]} = \sum_{i=1}^{m} D^{(t)}(i) \cdot \mathbf{1}_{[h_t(x_i) \neq y_i]} \cdot e^{w_t} = \varepsilon_t \cdot e^{w_t} = \varepsilon_t \cdot \sqrt{\frac{1-\varepsilon_t}{\varepsilon_t}} = \sqrt{\varepsilon_t (1-\varepsilon_t)} \quad [4]$

In similar manner we have the normalization factor:

$$\sum_{i=1}^{m} D^{(t)}(i) \cdot e^{-w_{t}h_{t}(x_{i})y_{i}} = \varepsilon_{t} \cdot \sqrt{\frac{1-\varepsilon_{t}}{\varepsilon_{t}}} + (1-\varepsilon_{t}) \cdot \sqrt{\frac{\varepsilon_{t}}{1-\varepsilon_{t}}} = 2\sqrt{\varepsilon_{t}(1-\varepsilon_{t})} [5]$$

By dividing [4] to [5] we get:

$$\sum_{i=1}^{m} D^{(t+1)}(i) \cdot \mathbf{1}[y_i \neq h_t(x_i)] = \frac{\sqrt{\varepsilon_t(1-\varepsilon_t)}}{2\sqrt{\varepsilon_t(1-\varepsilon_t)}} = \frac{1}{2}$$

We can observe how in $D^{(t+1)}$ the misclassified samples by h_t represent half of importance from the probability distribution \Rightarrow if $\varepsilon_t \in (0, \frac{1}{2}) \Rightarrow D^{(t)} \neq D^{(t+1)}$ (because the misclassified samples will weight more in the next iteration) $\Rightarrow h^{(t)} \neq h^{(t+1)}$

$$\Rightarrow Pr[h^{(t)} = h^{(t+1)} | \varepsilon_t \in (0, \frac{1}{2})] = 0$$
 [6]

From [3] and [6]
$$\Rightarrow Pr[h^{(t)} = h^{(t+1)}] = \begin{cases} 1, & \varepsilon_t = 0 \\ 0, & \text{otherwise} \end{cases}$$

Point (b): Assuming that the training error of h_{final} is at most $\frac{1}{2} - \frac{3}{2}\gamma + 2\gamma^2$ we will show that is strictly smaller than $\frac{1}{2} - \gamma$.

We have that
$$\frac{1}{2} - \frac{3}{2}\gamma + 2y^3 \le \frac{1}{2} - \gamma \iff \frac{3}{2}y - 2\gamma^3 \ge \gamma \iff \frac{1}{2}\gamma - 2\gamma^3 \ge 0 \iff \frac{1}{2}\gamma(1 - 4\gamma^2) \ge 0 \iff \frac{1}{2}\gamma(1 - 2\gamma)(1 + 2\gamma) \ge 0$$
 [1].

Based on the fact we know $\varepsilon_t \in (0, \frac{1}{2}) \Longrightarrow \gamma \in (0, \frac{1}{2})$ [2].

From [1] and [2] we conclude that $\frac{1}{2}y > 0$, $1 - 2\gamma > 0$, $1 + 2\gamma > 0 \Rightarrow \frac{1}{2}\gamma \cdot (1 - 2\gamma)(1 + 2\gamma) \ge 0 \Rightarrow \frac{1}{2}\gamma - \frac{3}{2}\gamma + 2\gamma^2 \le \frac{1}{2} - \gamma$

Exercise 3

Statement:

Let Σ be a finite alphabet and let $\chi = \Sigma^m$ be a sample space of all strings of length m over Σ . Let \mathcal{H} be a hypothesis space over χ , where $\mathcal{H} = \{h_w : \Sigma^m \to \{0,1\}, \ w \in \Sigma^*, \ 0 < |w| \le m$, s.t $h_w(x) = 1$ if w is a substring of x.

- a. Give an upper bound (any upper bound that you can come up) of the VCdimension of \mathcal{H} in terms of $|\Sigma|$ and m.
- b. Give an efficient algorithm for finding a hypothesis h_w consistent with a training set in the realizable case. What is the complexity of your algorithm?

Example: let $\Sigma = \{a, b, c\}$, m = 4 and the training set $S = \{(aabc, 1), (baca, 0), (bcac, 0), (abba, 1)\}$. The output of the algorithm should be h_{ab} .

Solution:

Point (a): From *Lecture 6* we know that $VC \dim(\mathcal{H}) \leq \log_2 |\mathcal{H}|$.

Let's consider $\Sigma = \{a, b\}$ this implies that the sample space

$$\chi = \{a, b\}^m = \{\underbrace{aaa...a}_{m}, \underbrace{baa...a}_{m}, \underbrace{bba...a}_{m}, ..., \underbrace{bbb...b}_{m}\} \text{ with } |\chi| = 2^m$$

$$\mathcal{H} = \{h_w : \Sigma^m \to \{0,1\}, \ w \in \Sigma^*, \ 0 < |w| \le m, \text{ s.t } h_h(x) = 1 \text{ if w is a substring of x} \}$$

 $\iff \mathcal{H} = \{h_w : \{a,b\}^m \to \{0,1\}, \ w \in \{a,b\}^*, \ 0 < |w| \le m, \text{ s.t } h_h(x) = 1 \text{ if w is a substring of x} \}$

$$\iff \mathcal{H} = \begin{cases} h_w : \{a, b\}^m \to \{0, 1\} \mid w \in \{a, b, aa, ab, ba, bb, \dots, \underbrace{bb \dots b}_{\text{m times}} \end{cases} \\ h_w(x) = 1, \text{ if w is substring of x} \end{cases}$$

We have that $|\mathcal{H}| = \left| \{ h_a, h_b, h_{aa}, h_{ab}, h_{ba}, h_{bb}, \dots, h_{bb \dots b} \} \right| = 2 + 4 + 8 + \dots + 2^m = \sum_{i=1}^m 2^i = 2 \cdot \frac{2^m - 1}{2 - 1} = 2^{m+1} - 2.$

To generalize, if
$$|\Sigma| = k \in \mathbb{N}^*$$
, $\Sigma = \{c_i \mid i \in \overline{1, k}\} \Longrightarrow \chi = \{c_i \mid i \in \overline{1, k}\}^m \Longrightarrow |\chi| = k^m = |\Sigma|^m$.

$$|\mathcal{H}| = \left| \{ h_{c_1}, h_{c_2}, \dots h_{c_k}, h_{c_1 c_1}, h_{c_1 c_2}, \dots, h_{c_k c_k}, \dots \underbrace{h_{c_k c_k}, \dots h_{c_k c_k}}_{\text{m times}} \} \right| \\ \Longrightarrow k^1 + k^2 + k^3 + \dots k^m = \sum_{i=1}^m k^i = k \cdot \frac{k^{m-1}}{k-1} = k \cdot \frac{k^{m-1}}{$$

$$\frac{k^{m+1}-k}{k-1}$$
 , where $k=|\Sigma|$

$$\Rightarrow |\mathcal{H}| = \frac{|\Sigma|^{m+1} - |\Sigma|}{|\Sigma| - 1}$$

$$\Rightarrow VC \dim(\mathcal{H}) \leq \log_2 |\mathcal{H}| = \log_2 \frac{|\Sigma|^{m+1} - |\Sigma|}{|\Sigma| - 1}$$

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Point (b): We are under the realizable case \Rightarrow Given a training set S = \{(x_1, y_1), (x_2, y_2), ..., (x_n, y_n) \mid x_i \in \Sigma^m, y_i \in \{0, 1\}, i = \overline{1, n}\} exists a function h_w \in \mathcal{H} with w \in \Sigma^*, 0 < |w| \le m such that y_i = h_{w^*}(x_i). Let's consider the following algorithm for our hypothesis class \mathcal{H}:
```

```
Input: S
Output: h_{w^s}
positive words = [word for (word,y) in S if y == 1]
negative_words = [word for (word,y) in S if y == 0]
P = length(positive words)
if P == 0 then
   w_s = \text{word} \notin \text{negative\_words}, with | word | = m
   return h_{w_a}
positive_substrings = dict() // initialize empty dict
negative_substrings = set() // initialize empty set
// We need to find the substrings that appear in all
// p positive words
for word in positive_words do
   substrings = generate all substrings(word)
   for substring in substrings do
    positive_substrings[substring] += 1
   end
end
// Next we need to generate all substrings from negative words
// Such that the selected W it's not contained in them
for word in negative words do
   substrings = generate_all_substrings(word)
   for substring in substrings do
       negative_substrings.update(substring)
   end
end
// Finally, we select W as the first substring that appears in all
// positive examples and it's not in the set of negative substrings
for substring, counts in positive substrings.items() do
   if substring not in negative substrings and count == P then
       return h_{substring}
end
```

Complexity:

- 1. Initialization step: O(n)
- 2. Iteration through positive and negative words: O(n)
- 3. Generationg all substrings for a word of size m: $O(m^2)$
- 4. Finding the W_S : $O(n^2 \cdot m^4)$
 - iterating through $n \cdot m^2$ substrings: $O(n \cdot m^2)$
 - searching substring in maximum $n \cdot m^2$ possible negative substrings: $O(n \cdot m^2)$

Total Complexity: $O(n^2 \cdot m^4)$