

PAC Intro from Understanding ML

Cody Melcher

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Overview

- ▶ Quick conceptual ML

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- ▶ Learning Environment
 1. Framework
 2. ERM (without and with Inductive Bias)

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 1. Necessary for complex tasks, useful for inherent flexibility.
- ▶ Learning Environment
 1. Framework
 2. ERM (without and with Inductive Bias)
- ▶ Formal Definition of the PAC learning model
 1. General bounds

ML Conceptually

- ▶ ML: Automating converting "experience" or data into knowledge.
 1. Necessary for complex tasks, useful for inherent flexibility.

Statistical Learning Framework

► Basic setting:

1. **Domain space** \mathcal{X} ie covariate space.
2. **Label set**: space of possible responses/explanatory variables.
Focus here on $\mathcal{Y} = \{1, 2\}$
3. **Training data** as finite **sequence** of ordered pairs in $\mathcal{X} \times \mathcal{Y}$
4. **Prediction rule** $h : \mathcal{X} \rightarrow \mathcal{Y}$; also called hypothesis or classifier.

Statistical Learning Framework continued

► Notes on basic framework

1. Assume data generated from unknown distribution \mathcal{D} Generally assume iid.
2. Assume perfect/correct but unknown classifier f_i exists and y_i are mapped to by f_i from \mathcal{X} .

3. **Error of h** is

$$L_{\mathcal{D},\{f\}}(h) = P_{x \sim \mathcal{D}}(h(x) \neq f(x)) = \mathcal{D}(\{x : h(x) \neq f(x)\}).$$

Minimize!

4. Note $\mathcal{D} : A \rightarrow [0, 1]$ where $A \subset \mathcal{X}$
5. L called risk or generalization error

Statistical Learning Framework continued

► Empirical Risk Minimization

1. Don't know \mathcal{D} or f . Minimize training error (empirical risk/error) instead.
2. $L_S(h) = \frac{|\{i \in [m]: h(x_i) \neq y_i\}|}{m} \in [0, 1]$
3. Problem: can easily lead to overfitting ie define perfect classifier for training data that is terrible classifier on other data.
4. Solution: restriction on set of possible classifiers \mathcal{H} . Big topic: conditions on \mathcal{H} to guarantee no overfitting.
5. First restriction: $|\mathcal{H}| < \infty$ which leads to...

ERM

- ▶ Realizability Assumption (2.1) etc.
 1. There exists $h^* \in \mathcal{H}$ st $L_{D,f}(h^*) = 0$ ie there exists for every empirical risk minimization problem there exists a perfect classifier.
 2. Possible issue: data drawn from \mathcal{D} could be bad representation of \mathcal{D} . Need to think of error as random.
 3. Let δ be probability sample from \mathcal{D} is non-representative sample so $(1 - \delta)$ is our **confidence parameter** (our sample is representative).
 4. Let ϵ be the **accuracy parameter**
 5. Want to use these to put upper bound on what samples can be realized that lead to classification failure.

ERM cont.

- ▶ Want to figure out bounding $D^m(\{S|_x : L_{(D,f)}(h_s) > \epsilon\})$ ie bound the realized samples of size m mapped by the classifier incorrectly classifies by ϵ .
- ▶ Define the set of bad classifiers as $\mathcal{H}_B = \{h \in \mathcal{H} : L_{D,f}(h) > \epsilon\}$ and let $M = \{S|_x : \exists h \in \mathcal{H}_B, L_S(h) = 0\}$ be the set of misleading samples.
- ▶ Follows that $\{S_x : L_{D,f}(h_s) > \epsilon\} \subset M$ ie set of samples that give a bad classifier are a subset of the misleading samples set.

ERM cont.

- ▶ Can relatedly think of all the possible bad classifiers that manage to give perfect prediction due to a misleading sample; this is equivalent to M .

- ▶ It follows that

$$D^m(\{S|_x : L_{(D,f)}(h_S) > \epsilon\}) \leq D^m(\cup_{h \in H_b} \{S|_x : L_S(h) = 0\})$$

- ▶ By $D(A \cup B) \leq D(A) + D(B)$, it follows that RHS

$$\leq \sum_{h \in H_b} D^m(\{S|_x : L_S(h) = 0\})$$

- ▶ Next due to iid assumption and how we defined ϵ , it follows that $D^m(\{x_i : h(x_i) = y_i\}) = (1 - L_{D,f})^m \leq (1 - \epsilon)^m \leq e^{-\epsilon m}$

ERM cont.

- ▶ Combining the previous inequality with the inequality from two slides back we get, $D^m(\{S|_x : L_{(D,f)}(h_s) > \epsilon\}) \leq |\mathcal{H}|e^{-\epsilon m}$ which provides the upper bound on the realized samples of size m mapped by the classifier incorrectly, which we can summarize as,
- ▶ **Corollary 2.3:** If \mathcal{H} is finite, $\delta \in (0, 1)$ and $\epsilon > 0$ and let m be any integer st. $m \geq \log\left(\frac{|\mathcal{H}|}{\delta}\right)\frac{1}{\epsilon}$. Then for any true classifier f and distribution \mathcal{D} assuming realizability with probability at least $1 - \delta$ over iid sample of size m , we have that for **every** ERM classifier h_s it holds that $L_{D,f}(h_s) \leq \epsilon$.

ERM cont.

- ▶ I.e. for sufficiently large m , classifiers from finite classifier class will be probably (with confidence $1 - \delta$) approximately (up to ϵ wrong) correct.
- ▶ Hence why, PAC learning environment = probably approximately correct.

PAC Learnability Definition

- ▶ (3.1) PAC Learnability: \mathcal{H} is **PAC learnable** if there exists a function $m_H : (0, 1)^2 \rightarrow \mathcal{N}$ and a learning algorithm with the following property: $\forall \epsilon, \delta \in (0, 1), \forall \mathcal{D}$ over \mathcal{X} and for every true classifier $f : \mathcal{X} \rightarrow 0, 1$, and if the realizable assumptions holds, then when running the learning algorithm on $m \geq m_h(\epsilon, \delta)$ iid samples generated by \mathcal{D} and classified by f , the algorithm returns a classifier h st with probability at least $1 - \delta$ that $L_{D,f}(h) \leq \epsilon$

PAC Learnability notes

- ▶ ϵ = how far the classifier can be from the optimal classifier f
- ▶ δ how likely the classifier is to be ϵ close to f
- ▶ notice that m_h determines how many samples needed to be probably accurate; we will focus on finding the minimal m_h ie minimal integer that guarantees probable accuracy.
- ▶ (3.2) $\forall |\mathcal{H}| < \infty, m_h(\epsilon, \delta) \leq \frac{\log(\frac{|\mathcal{H}|}{\delta})}{\epsilon}$

Agnostic PAC

- ▶ Realizability seems unrealistic....solution: agnostic PAC!
- ▶ (3.3) A hypothesis class \mathcal{H} is **agnostic PAC learning** if $\exists m_h : (0, 1)^2 \rightarrow \mathcal{N}$ and a learning algorithm with the following property: $\forall \epsilon, \delta \in (0, 1)$ and $\forall \mathcal{D}$ over $\mathcal{X} \times \mathcal{Y}$ when running the learning algorithm on $m \geq m_h(\epsilon, \delta)$ iid examples generated by \mathcal{D} , a hypothesis h is returned st with probability at least $1 - \delta$,
$$L_D(h) \leq \min_{h' \in \mathcal{H}} (L_D(h') + \epsilon)$$
- ▶ Ie we can guarantee we are only ϵ worse than the best predictor h' from a class of classifiers \mathcal{H} .

Extensions to Agnostic PAC

- ▶ Generally we refer to the agnostic PAC learning environment as PAC learning.
- ▶ Can expand \mathcal{Y} to be larger than $\{0, 1\}$, but still assuming it is finite.
- ▶ Can change risk to be expected squared difference (or something else) to deal with regression problems.
- ▶ Both these extensions change $L_D(h)$, and thus we need to rethink our above definition....

Agnostic PAC with General Loss

- ▶ (3.4) A hypothesis class \mathcal{H} is **agnostic PAC learning** with respect to set Z and loss function $l : H \times Z \rightarrow \mathcal{R}_+$ if $\exists m_h : (0, 1)^2 \rightarrow \mathcal{N}$ and a learning algorithm with the following property: $\forall \epsilon, \delta \in (0, 1)$ and $\forall \mathcal{D}$ over \mathcal{Z} when running the learning algorithm on $m \geq m_h(\epsilon, \delta)$ iid examples generated by \mathcal{D} , a hypothesis h is returned st with probability at least $1 - \delta$, $L_D(h) \leq \min_{h' \in H} (L_D(h') + \epsilon)$
- ▶ where $L_D(h) = E_{(z \sim \mathcal{D})}(l(h, z))$ and $Z = \mathcal{X} \times \mathcal{Y}$ (for our problems though this can be generalized)

TLDR;

- ▶ PAC learning feels natural to stats/math world, learning from "experience" ie data ie we are still trying to approximate a function.
- ▶ New ish to us maybe because PAC learning is about thinking about bounds on what is possible and sorta avoiding asymptotics.
- ▶ Still very general, but allows us to move to VC dimension (what characteristics of \mathcal{H} allow us to deduce it is PAC learnable?), be more rigorous with a lot of the techniques we have learned, etc.