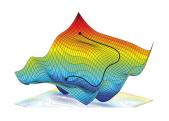
# Introduction to Machine Learning

# **Risk Minimizers**



#### Learning goals

- Know the concepts of the Bayes optimal model (also: risk minimizer, population minimizer)
- Know the Bayes risk
- Know the concept of consistent learners
- Know the concept of the optimal constant model

## **RISK MINIMIZER**

Our goal is to minimize the risk

$$\mathcal{R}_L(f) := \mathbb{E}_{xy}[L(y, f(\mathbf{x}))] = \int L(y, f(\mathbf{x})) d\mathbb{P}_{xy}.$$

for a certain hypothesis  $f(\mathbf{x}) \in \mathcal{H}$  and a loss  $L(y, f(\mathbf{x}))$ .

NB: As  $\mathcal{R}_I$  depends on loss L, we sometimes make this explicit with a subscript if needed, and omit in other cases.

Let us assume we are in an "ideal world":

- ullet The hypothesis space  ${\mathcal H}$  is unrestricted. We can choose any  $f:{\mathcal X} o {\mathbb R}^g.$
- We also assume an ideal optimizer; the risk minimization can always be solved perfectly and efficiently.
- We know  $\mathbb{P}_{xy}$ .

How should f be chosen?

#### **RISK MINIMIZER**

The *f* with minimal risk across all (measurable) functions is called the **risk minimizer**, **population minimizer** or **Bayes optimal model**.

$$f^* = \underset{f:\mathcal{X} \to \mathbb{R}^g}{\arg \min} \mathcal{R}_L(f) = \underset{f:\mathcal{X} \to \mathbb{R}^g}{\arg \min} \mathbb{E}_{xy} \left[ L(y, f(\mathbf{x})) \right]$$
$$= \underset{f:\mathcal{X} \to \mathbb{R}^g}{\arg \min} \int L(y, f(\mathbf{x})) \, d\mathbb{P}_{xy}.$$

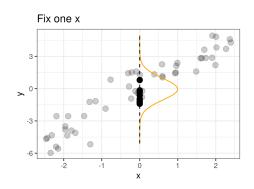
The resulting risk is called Bayes risk

$$\mathcal{R}_{L}^{*} = \inf_{f:\mathcal{X} \to \mathbb{R}^{g}} \mathcal{R}_{L}(f)$$

## **OPTIMAL POINT-WISE PREDICTIONS**

To derive the risk minimizer we usually make use of the following trick:

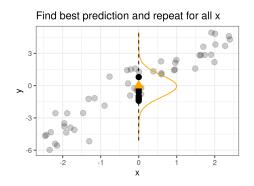
- We can choose  $f(\mathbf{x})$  as we want (unrestricted hypothesis space, no assumed functional form)
- Consequently, for a fixed value  $\mathbf{x} \in \mathcal{X}$  we can select **any** value c we want to predict
- So we construct the **point-wise optimizer** for every  $\mathbf{x} \in \mathcal{X}$ .



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# THEORETICAL AND EMPIRICAL RISK

The risk minimizer is mainly a theoretical tool:

- In practice we need to restrict the hypothesis space  $\mathcal H$  such that we can efficiently search over it.
- In practice we (usually) do not know  $\mathbb{P}_{xy}$ . Instead of  $\mathcal{R}(f)$ , we are optimizing the empirical risk

$$\hat{f} = \operatorname*{arg\,min}_{f \in \mathcal{H}} \mathcal{R}_{emp}(f) = \operatorname*{arg\,min}_{f \in \mathcal{H}} \sum_{i=1}^{n} L\left(y^{(i)}, f\left(\mathbf{x}^{(i)}\right)\right)$$

Note that according to the **law of large numbers** (LLN), the empirical risk converges to the true risk (but beware of overfitting!):

$$\bar{\mathcal{R}}_{\text{emp}}(f) = \frac{1}{n} \sum_{i=1}^{n} L\left(y^{(i)}, f\left(\mathbf{x}^{(i)}\right)\right) \stackrel{n \to \infty}{\longrightarrow} \mathcal{R}(f).$$

# **ESTIMATION AND APPROXIMATION ERROR**

**Goal of learning:** Train a model  $\hat{f}$  for which the true risk  $\mathcal{R}_L\left(\hat{f}\right)$  is close to the Bayes risk  $\mathcal{R}_L^*$ . In other words, we want the **Bayes regret** 

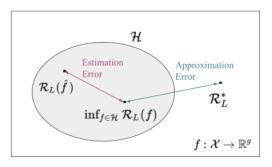
$$\mathcal{R}_L\left(\hat{f}\right)-\mathcal{R}_L^*$$

to be as low as possible.

The Bayes regret can be decomposed as follows:

$$\mathcal{R}_{L}\left(\hat{f}\right) - \mathcal{R}_{L}^{*} = \underbrace{\left[\mathcal{R}_{L}\left(\hat{f}\right) - \inf_{f \in \mathcal{H}} \mathcal{R}_{L}(f)\right]}_{\text{estimation error}} + \underbrace{\left[\inf_{f \in \mathcal{H}} \mathcal{R}_{L}(f) - \mathcal{R}_{L}^{*}\right]}_{\text{approximation error}}$$

## **ESTIMATION AND APPROXIMATION ERROR**



- $\mathcal{R}_L\left(\hat{f}\right)$   $\inf_{f\in\mathcal{H}}\mathcal{R}(f)$  is the **estimation error**. We fit  $\hat{f}$  via empirical risk minimization and (usually) use approximate optimization, so we usually do not find the optimal  $f\in\mathcal{H}$ .
- $\inf_{f \in \mathcal{H}} \mathcal{R}_L(f) \mathcal{R}_L^*$  is the **approximation error**. We need to restrict to a hypothesis space  $\mathcal{H}$  which might not even contain the Bayes optimal model  $f^*$ .

# (UNIVERSALLY) CONSISTENT LEARNERS

**Consistency** is an asymptotic property of a learning algorithm, which ensures the algorithm returns **the correct model** when given **unlimited data**.

Let  $\mathcal{I}: \mathbb{D} \times \Lambda \to \mathcal{H}$  be a learning algorithm<sup>(\*)</sup> that takes a training set  $\mathcal{D}_{\text{train}} \sim \mathbb{P}_{xy}$  of size  $n_{\text{train}}$  and estimates a model  $\hat{f}: \mathcal{X} \to \mathbb{R}^g$ .

The learning method  $\mathcal{I}$  is said to be **consistent** w.r.t. a certain distribution  $\mathbb{P}_{xy}$  if the risk of the estimated model  $\hat{f}$  converges in probability (" $\stackrel{p}{\longrightarrow}$ ") to the Bayes risk  $\mathcal{R}^*$  when  $n_{\text{train}}$  goes to  $\infty$ :

$$\mathcal{R}\left(\mathcal{I}\left(\mathcal{D}_{\mathsf{train}}, \boldsymbol{\lambda}\right)\right) \stackrel{p}{\longrightarrow} \mathcal{R}_{L}^{*} \quad \mathsf{for} \; n_{\mathsf{train}} \to \infty.$$

 $^{(*)}$   $\lambda \in \Lambda$  denotes hyperparameters of the learning algorithm.

# (UNIVERSALLY) CONSISTENT LEARNERS

Consistency is defined w.r.t. a particular distribution  $\mathbb{P}_{xy}$ . But since we usually do not no  $\mathbb{P}_{xy}$ , consistency does not offer much help to choose an algorithm for a particular task.

More interesting is the stronger concept of **universal consistency**: An algorithm is universally consistent if it is consistent for **any** distribution.

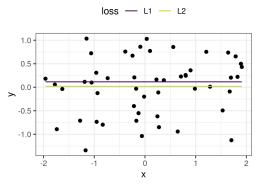
In Stone's famous consistency theorem from 1977, the universal consistency a weighted average estimator as KNN was proven. Many other ML model have since then been proven to be universally consistent (SVMs, ANNs, etc.).

**Note** that universal consistency is obviously a desirable property - however, (universal) consistency does not tell us anything about convergence rates ...

## **OPTIMAL CONSTANT MODEL**

While the risk minimizer gives us the (theoretical) optimal solution, the **optimal constant model** (also: featureless predictor) gives us an computable empirical lower baseline solution.

The constant model is the model  $f(\mathbf{x}) = \theta$  that optimizes the empirical risk  $\mathcal{R}_{\text{emp}}(\theta)$ .



# RISK MINIMIZER AND OPTIMAL CONSTANT

In the following chapters, we will derive the risk minimizers and optimal constant models for different loss functions.

Name	Risk Minimizer	Optimal Constant
L2	$f^*(\mathbf{x}) = \mathbb{E}_{y x}[y \mid \mathbf{x}]$	$\hat{f}(\mathbf{x}) = \frac{1}{n} \sum_{i=1}^{n} y^{(i)}$
L1	$f^*(\mathbf{x}) = med_{y x}[y \mid \mathbf{x}]$	$\hat{f}(\mathbf{x}) = med(y^{(i)})$
0-1	$h^*(\mathbf{x}) = \operatorname{argmax}_{l \in \mathcal{Y}} \mathbb{P}(y = l \mid \mathbf{x})$	$\hat{h}(\mathbf{x}) = mode\left\{y^{(i)}\right\}$
Brier	$\pi^*(\mathbf{x}) = \mathbb{P}(y = 1 \mid \mathbf{x} = \mathbf{x})$	$\hat{\pi}(\mathbf{x}) = \frac{1}{n} \sum_{i=1}^{n} y^{(i)}$
Bernoulli (on probs)	$\pi^*(\mathbf{x}) = \mathbb{P}(y = 1 \mid \mathbf{x} = \mathbf{x})$	$\hat{\pi}(\mathbf{x}) = \frac{1}{n} \sum_{i=1}^{n} y^{(i)}$
Bernoulli (on scores)	$f^*(\mathbf{x}) = \log\left(\frac{\mathbb{P}(y=1 \mid \mathbf{x})}{1 - \mathbb{P}(y=1 \mid \mathbf{x})}\right)$	$\hat{f}(\mathbf{x}) = \log \frac{n_{+1}}{n_{-1}}$