12ML:: BASICS

Data

 $\mathcal{X} \subset \mathbb{R}^p$: p-dimensional feature / input space Usually we assume $\mathcal{X} \equiv \mathbb{R}^p$, but sometimes, dimensions may be bounded (e.g., for categorical or non-negative features.)

 $\mathcal{Y} \subset \mathbb{R}^g$: target space

e.g.: $\mathcal{Y}=\mathbb{R}$, $\mathcal{Y}=\{0,1\}$, $\mathcal{Y}=\{-1,1\}$, $\mathcal{Y}=\{1,\ldots,g\}$ with g classes

 $x = (x_1, \dots, x_p)^T \in \mathcal{X}$: feature vector

 $y \in \mathcal{Y}$: target / label / output

 $\mathbb{D}_n = (\mathcal{X} \times \mathcal{Y})^n \subset \mathbb{D}$: set of all finite data sets of size n

 $\mathbb{D} = \bigcup_{n \in \mathbb{N}} (\mathcal{X} \times \mathcal{Y})^n$: set of all finite data sets

 $\mathcal{D} = ((\mathsf{x}^{(1)}, \mathsf{y}^{(1)}), \dots, (\mathsf{x}^{(n)}, \mathsf{y}^{(n)})) \in \mathbb{D}_n$: data set with n observations

 $\mathcal{D}_{\mathsf{train}}, \ \mathcal{D}_{\mathsf{test}} \subset \mathcal{D}$: data for training and testing (often: $\mathcal{D} = \mathcal{D}_{\mathsf{train}} \dot{\cup} \; \mathcal{D}_{\mathsf{test}}$)

 $(x^{(i)}, y^{(i)}) \in \mathcal{X} \times \mathcal{Y} : i$ -th observation or instance

 \mathbb{P}_{xy} : joint probability distribution on $\mathcal{X} \times \mathcal{Y}$

Model and Learner

Model / hypothesis: $f: \mathcal{X} \to \mathbb{R}^g$, $\mathsf{x} \mapsto f(\mathsf{x})$ is a function that maps feature vectors to predictions, often parametrized by $oldsymbol{ heta} \in \Theta$ (then we write f_{θ} , or, equivalently, $f(x \mid \theta)$).

 $\Theta \subset \mathbb{R}^d$: parameter space

 $\boldsymbol{\theta} = (\theta_1, \theta_2, ..., \theta_d) \in \Theta$: model parameters Some models may traditionally use different symbols.

 $\mathcal{H} = \{f: \mathcal{X} \to \mathbb{R}^g \mid f \text{ belongs to a certain functional family}\}$:

hypothesis space

Set of functions defining a specific model class to which we restrict our learning task

Learner $\mathcal{I}: \mathbb{D} \times \mathbf{\Lambda} \to \mathcal{H}$ takes a training set $\mathcal{D}_{\mathsf{train}} \in \mathbb{D}$ and produces a model $f:\mathcal{X} \to \mathbb{R}^g$, its hyperparameters set to $\boldsymbol{\lambda} \in \boldsymbol{\Lambda}$.

For a parametrized model this can be adapted to $\mathcal{I}: \mathbb{D} imes \mathbf{\Lambda} o \Theta$

 $\Lambda \subset \mathbb{R}^{foo}$: hyperparameter space

 $\boldsymbol{\lambda} = (\lambda_1, \lambda_2, ..., \lambda_{foo})^T \in \boldsymbol{\Lambda}$: model hyperparameters

 $\pi_k = \mathbb{P}(y = k)$: **prior probability** for class kIn case of binary labels we might abbreviate $\pi = \mathbb{P}(y = 1)$.

 $\pi_k(x) = \mathbb{P}(y = k \mid x)$: **posterior probability** for class k, given xIn case of binary labels we might abbreviate $\pi(x) = \mathbb{P}(y = 1 \mid x)$.

 $\mathcal{L}(m{ heta})$ and $\ell(m{ heta}) = \log \mathcal{L}(m{ heta})$: likelihood and log-likelihood for parameter $oldsymbol{ heta}$

These are based on a statistical model.

 $\epsilon = y - f(x)$ or $\epsilon^{(i)} = y^{(i)} - f(x^{(i)})$: (i-th) **residual** in regression

yf(x) or $y^{(i)}f(x^{(i)})$: margin for (i-th) observation in binary classification (with $\mathcal{Y} = \{-1, 1\}$).

 \hat{y} , \hat{f} , $\hat{\pi}_k(x)$, $\hat{\pi}(x)$ and $\hat{\theta}$

The hat symbol denotes **learned** functions and parameters.

oss and Risk

 $L: \mathcal{Y} \times \mathbb{R}^g o \mathbb{R}:$ loss function

L(y, f(x)) quantifies the "quality" of the prediction f(x) for a single observation x.

 $\mathcal{R}_{\mathsf{emp}}:\mathcal{H} o \mathbb{R}: \mathsf{empirical} \mathsf{\; risk}$

The ability of a model f to reproduce the association between x and ythat is present in the data \mathcal{D} can be measured by the summed loss:

$$\mathcal{R}_{emp}(f) = \sum_{i=1}^{n} L\left(y^{(i)}, f\left(\mathbf{x}^{(i)}\right)\right)$$

Since f is usually defined by **parameters** θ , this becomes:

$$\mathcal{R}_{emp}: \mathbb{R}^d
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Learning then amounts to empirical risk minimization – figuring out which model \hat{f} has the smallest summed loss:

$$\hat{f} = rg \min_{oldsymbol{ heta} \in \Theta} \mathcal{R}_{\mathsf{emp}}(oldsymbol{ heta}).$$

Components of Learning

Learning = Hypothesis space + Risk + Optimization.

Hypothesis space: Defines (and restricts!) what kind of model f can be learned from the data.

Examples: linear functions, decision trees

Risk: Quantifies how well a model performs on a given data set. This allows us to rank candidate models in order to choose the best one.

Examples: squared error, negative (log-)likelihood

Optimization: Defines how to search for the best model, i.e., the model with the smallest risk, in the hypothesis space.

Examples: gradient descent, quadratic programming

Regression Losses

Basic idea (L2 loss / squared error):

 $L(y, f(x)) = (y - f(x))^2$ or $L(y, f(x)) = 0.5(y - f(x))^2$

Convex and differentiable

Tries to reduce large residuals (loss scaling quadratically)

Optimal constant model: $\hat{f}(x) = \text{mean of } y|x$

Basic idea (L1 loss / absolute error):

L(y, f(x)) = |y - f(x)|

Convex and more robust

Non-differentiable for y = f(x), optimization becomes harder

Optimal constant model: $\hat{f}(x) = \text{median of } y | x$

Classification

Assume we are given a classification problem:

 $\mathsf{x} \in \mathcal{X}$ feature vector $y \in \mathcal{Y} = \{1, \dots, g\}$ categorical output variable (label) $\mathcal{D} = ((\mathsf{x}^{(1)}, \mathsf{y}^{(1)}), \dots, (\mathsf{x}^{(n)}, \mathsf{y}^{(n)}))$ observations of x and y

Classification usually means to construct g discriminant functions:

 $f_1(x), \ldots, f_g(x)$, so that we choose our class as $h(x) = arg \max_{k \in \{1,...,g\}} f_k(x)$

Linear Classifier: If the functions $f_k(x)$ can be specified as linear functions, we will call the classifier a *linear classifier*.

Binary classification: If only 2 classes exist, we can use a single discriminant function $f(x) = f_1(x) - f_2(x)$.