Boosting

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Outline

Introduction

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Boosting - General Idea

Boosting is a technique that trains models which each have high Bias – low Variance – and it creates a particular linear combination of these models which is a better fit for the data (this "boosted" model having increased Variance).

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- ► Ideally, the simple models, that have high Bias, allow for very efficient training, keeping computation times small.
- Two popular algorithms for this general idea are AdaBoost and "Gradient Boosted Decision Trees" – which are used in a tool called XGBoost.

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For binary classification, write the empirical 0-1 loss function as $\mathcal{L}_{\mathcal{S}}^{01}$; it is defined as follows. Given a prediction function $f:\mathbb{R}^d \to \{1,-1\}$, the loss function on data $\mathcal{S}=\{(\mathbf{x}_i,y_i)\}_{i=1}^n$ is the probability that $f(\mathbf{x}_i)\neq y_i$. That is,

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Population loss: probability that $f(\mathbf{x}) \neq y$ when (\mathbf{x},y) is drawn from the population.

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A **learner** or **learning algorithm** is a procedure or rule that finds (or trains) f_S from given training data S.

Recall how each empirical loss function $\mathcal{L}_{\mathcal{S}}$ is the average of a "per-example" loss over the set \mathcal{S} . In the 0-1 case, a per-example loss of

$$\boldsymbol{\ell}^{01}(f,(\mathbf{x},y)) = \begin{cases} 0, & \text{if } f(\mathbf{x}) = y \\ 1, & \text{else} \end{cases}$$

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In this average, each point in S is weighted equally. However, given a probability vector $\mathbf{p} = (p_1, p_2, \dots, p_n)$, we could weight points in S according to \mathbf{p} :

$$p_1 \ell^{01}(f, (\mathbf{x}_1, y_1)) + p_2 \ell^{01}(f, (\mathbf{x}_2, y_2)) + \ldots + p_n \ell^{01}(f, (\mathbf{x}_n, y_n)).$$

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We'll write $\mathcal{L}_{\mathcal{S}}^{01}(f, \mathbf{p})$ for this weighted loss function. Note that a similar weighted loss exists for any "standard" empirical loss function.¹

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We'll write $\mathcal{L}^{01}_{S}(f,\mathbf{p})$ for this weighted loss function. Note that a similar weighted loss exists for any "standard" empirical loss function.1 Below, we begin with a learner that does okay, but $\mathcal{L}^{01}_{\mathcal{S}}(f_{\mathcal{S}})$ may not be that

"small"; however, through weighted loss and adaptively changing p, we build

one with much better 0-1 loss.

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For γ , with $0<\gamma<\frac{1}{2}$, we call a learning algorithm a γ -weak learner if, for any $0<\delta<1$, there is a positive integer N so that whenever $|\mathcal{S}|\geq N$ then the learning algorithm determines $f_{\mathcal{S}}$ so that $\mathcal{L}^{01}_{\mathcal{S}}(f_{\mathcal{S}})\leq \frac{1}{2}-\gamma$.

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Note that, for a sufficiently large set of training data S, a γ -weak learner will produce a prediction function f_S that is incorrect less than half of the time. That is, for small γ , this is a bit better than using a coin flip.

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Decision stumps can be viewed as weak learners. More precisely, for data where the label should have a single sign (+1 or -1) for all $\mathbf x$ that have some coordinate x_j that is in a fixed interval, decision stumps are γ -weak learners for some $0<\gamma<\frac{1}{6}$.

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Hence, it is common to use decision trees with small max depth as the base model in a boosting algorithm.

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One chooses a number T of rounds for the AdaBoost algorithm. In round t, with $1 \le t \le T$, say that f_t is the basic model that is trained in that round. That round will also have a corresponding coefficient w_t . The final (boosted) model is given by

$$f(\mathbf{x}) = \operatorname{sign}(\mathbf{w}_1 f_1(\mathbf{x}) + \mathbf{w}_2 f_2(\mathbf{x}) + \ldots + \mathbf{w}_T f_T(\mathbf{x})).$$

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The procedure for finding f_t and w_t is the following. In each round t, there is a probability vector $\mathbf{p}^{(t)} \in \mathbb{R}^n$. In round 1, we give $\mathbf{p}^{(1)}$ uniform weights; that is, $\mathbf{p}^{(1)} = (1/n, 1/n, \dots, 1/n)$.

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Once done with T rounds, the boosted model is the one given by linear combination above.