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FACULTY OF SCIENCE

A Bachelor of Science thesis

Super Learners

and their oracle properties

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Contents

1 Introduction

Our setup closely models what is described in [VDL06] and [LD03]. Let $O_1, \ldots, O_n$ be n-i.i.d. observations distributed according to $P \in \mathcal{P}$ on some measurable space $(\mathcal{O}, \mathcal{A})$ where $O_i \in \mathcal{O}$ for each i and $\mathcal{P}$ is our statistical model. For a parameter set Θ we define the corresponding loss function $L : \mathcal{O} \times \Theta \to [0, \infty)$ as a measurable map such that our goal is to find an estimator $\hat{\theta}$ that minimizes the true risk function $R : \Theta \to \mathbb{R}$ given as

$$R(\theta) = \int L(x, \theta) dP(x) = EL(O_1)$$

The parameter set Θ can be Euclidean, but for the focus of this thesis we will consider it as a collection of functions of the form $\theta : \mathcal{O} \to \mathbb{R}$.

Example 1 (Regression functions Θ). Let $O_1 = (Y_1, X_1), \ldots, O_n = (Y_n, X_n) \in \mathcal{O} = \mathbb{R} \times \mathcal{X}$ be i.i.d. observations distributed according to some $P \in \mathcal{P}$ such that they satisfy the model

$$Y_1 = \theta_0(X_1) + \varepsilon,$$

for an unobservable stochastic error term ε. The goal is to estimate an unknown **regression function** $\theta_0 \in \Theta$ where $\Theta = \{\theta \mid \theta : \mathcal{X} \to \mathbb{R}\}$, is the set of possible regression functions each having $\mathcal{X}$ as their domain. [VDL06]

Example 2 (Parameteric family). Consider the initial setup from example 1. If Y_i is $\mathcal{B}(\mathbb{R}) - \mathcal{B}(\mathbb{R})$ measurable and X_i is $\mathcal{F} - \mathcal{B}(\mathbb{R})$ measurable for some sigma-algebra $\mathcal{F}$ on $\mathcal{X}$, then a **generalized regression model** could be considered as parametrized family of distributions, $\mathcal{P} = \{P_\theta \mid \theta \in \Theta\}$, given that Θ is finite-dimensional.
We can parametrize the conditional probability distributions for Y_1 given $X_1 = x$ as $\mathcal{Q} = \{Q_{\theta(x)} \mid \theta \in \Theta\}$ such that $Q_{\theta(x)}$ is a valid probability distribution on $\mathcal{B}(\mathbb{R})$ for each $x \in X$ and $\theta \in \Theta$. For a given $P_\theta \in \mathcal{P}$ there will exist a $Q_\theta \in \mathcal{Q}$ such that

$$P_\theta(Y \in A \mid X = x) = Q_{\theta(x)}(A) \qquad \text{for all } A \in \mathcal{B}(\mathbb{R}).$$

If we assume that X_1 is distributed according to some H_0 on $\mathcal{X}$, then the distribution P_θ over our observations (the joint over Y and X) will be

$$P_\theta(X \in A, Y \in B) = \int_A Q_{\theta(x)}(B) dH_0(x)$$

for every $A \in \mathcal{F}$ and $B \in \mathcal{B}(\mathbb{R})$.

Example 3 (Logistic regression model). Let $O_1 = (Y_1, X_n), \ldots, O_n = (Y_n, X_n) \in \mathcal{O} = \{0,1\} \times \mathcal{X}$ be i.i.d. observations from some distribution $P_{\theta_0} \in \mathcal{P}$, where Y_i is binary and $\mathcal{X} \subseteq \mathbb{R}^k$. We would like to estimate the parameter function $\theta_0 \in \Theta$

$$\theta_0(x) = E(Y_1 \mid X_1 = x) = P_{\theta_0}(Y_1 = 1 \mid X_1 = x),$$

In logistic regression we assume that $\Theta = \{x \mapsto \text{expit}(\beta x) \mid \beta \in \mathbb{R}^k\}$, so $\theta_0(x) = \text{expit}(\beta_0 x)$, then the goal becomes to estimate the k-dimensional parameter β_0, in this case the $\mathbb{R}^k$ parameter β_0 completely determines θ_0, so Θ is also k-dimensional. The conditional distributions of Y_1 given $X_1 = x$ are Bernoulli distributions and can be parametrized as $\mathcal{Q} = \{\text{Ber}(\text{expit}(\beta x)) \mid \beta \in \mathbb{R}^k\}$. Now from example 2 we know that the statistical model, $\mathcal{P}$, can be parametrized through β, in particular we have

$$\begin{aligned} P_\beta(Y_1 = 1, X_1 \in A) &= \int_A Q_{\theta(x)}(\{1\}) dH_0(x) \\ &= \int_A \text{expit}(\beta x) dH_0(x) \end{aligned}$$

If H_0 has density f w.r.t. Lebesgue measure, we can write

$$P_\beta(Y_1 = 1, X_1 \in A) = \int_A \text{expit}(\beta x) f(x) dm(x)$$

We will now turn our attention to statistical estimators. Statistical literature commonly write that an estimator is stochastic variable taking values in our parameter space $\hat{\theta} \in \Theta$. An estimator is achieved by considering i.i.d. observations $O_1, \ldots, O_2 \in \mathcal{O}$ distributed according to some measure P from some statistical model $\mathcal{P}$. We leave the model unspecified as it can be both parametric or nonparametric. Now let $h : \mathcal{O}^n \to \Theta$ be a measurable map, an estimator created from h is the random variable $T = h(O_1, \ldots, O_n)$. For $\Theta \subseteq \mathbb{R}^k$ the canonical σ-algebra on Θ is the Borel algebra, but when the parameter set is a set of functions, the σ-algebra can only be chosen after careful consideration of constraints on Θ.
In the following section we introduce the terminology "estimator algorithm" which corresponds to the measurable map h from our finite sample observation space to our parameter space.

Definition 1 (Estimator algorithm h)**.** An estimator algorithm is a measurable map $h : \mathcal{O}^n \to \Theta$ for $n \in \mathbb{N}$.

Definition 2 (Statistical Estimator $\hat{\theta}$)**.** Let $O_1, \ldots, O_n \in \mathcal{O}$ be i.i.d. observations distributed according to some $P \in \mathcal{P}$ for a statistical model $\mathcal{P}$ on $\mathcal{O}$. Let $h : \mathcal{O}^n \to \Theta$ be an estimator algorithm. An estimator is the random variable $\hat{\theta} = h(O_1, \ldots, O_n) \in \Theta$.

There is a one-to-one correspondence between the tuples of i.i.d. observations $(O_1, \ldots, O_n) \in \mathcal{O}^n$ and the empirical measures over n observations on $(\mathcal{O}, \mathcal{A})$ defined as

$$P_n(A) = \frac{1}{n} \sum_{i=1}^{n} \delta_{O_i}(A) \qquad \text{for } A \in \mathcal{A}.$$

Note that the empirical measure is a random variable. Thus, we can write $h(P_n)$ as an alternative representation of the estimator $h(O_1, \ldots, O_n)$, by adjusting the notation without introducing ambiguity.

Example 4 (Prediction algorithm). Consider the setup from example 3, where we have i.i.d. observations $O_1 = (Y_1, X_1), \ldots, O_n = (Y_n, X_n)$ such that $Y_i \in \{0, 1\}$ and $\mathcal{X} \in \mathbb{R}^k$ and our goal is to estimate the probability $\theta(x) = P_\theta(Y_1 = 1 \mid X_1 = x)$...

We would now like to consider the scenario where we have a library (set) of learner algorithms, $h_1, \ldots, h_n$. From these algorithms, we can define the set of learners $\{\hat{\theta}_q = h_q(P_n) | 1 \leq q \leq p\}$, where our goal is to find $\hat{\theta}_{\hat{q}}(P_n)$, which denotes the learner that minimizes R and $\hat{q}$ may depend on the observations.
In order to find $\hat{q}$ we have to proceed via cross validation. In cross validation, we randomly split our data into a training set and a test set. Let the random binary vector $S = (S_1, \ldots, S_n) \in \{0, 1\}^n$ be independent of $X_1, \ldots, X_n$ such that $S_i = 0$ indicates that X_i should be in the training set and $S_i = 1$ indicates that X_i belongs to the test set. We can define the empirical distributions over these two subsets, $P^0_{n,S}$ and $P^1_{n,S}$ as

$$P^0_{n,S} = \frac{1}{n_0} \sum_{i:S_i=0} \delta_{X_i}$$
$$P^1_{n,S} = \frac{1}{1 - n_0} \sum_{i:S_i=1} \delta_{X_i}$$

Where n_0 would be the number of S_i's that are marked 0.

Example 5 (Random splits). For $n = 9$ observations one could for example define the distribution of the random vector S as

$$P(S = (0, 0, 0, 0, 0, 0, 1, 1, 1)) = \frac{1}{3}$$
$$P(S = (0, 0, 0, 1, 1, 1, 0, 0, 0)) = \frac{1}{3}$$
$$P(S = (1, 1, 1, 0, 0, 0, 0, 0, 0)) = \frac{1}{3},$$

i.e. 3-fold cross-validation.
In general for n observations we have 2^n ways of choosing which observations should be in the training set and in the validation set. It might not be desirable to define the discrete probabilities for S over $\{0, 1\}^n$ simply as $\frac{1}{2^n}$ for each possible combination of training/validation data, since that would also include the combination where $n_1 = 0$. To ensure that we always have $n_1 > 0$, then let n_1 be given, then we see that there are $\binom{n}{n_1}$ ways of choosing both the validation and training set. We can therefore define the distribution of S as

$$P(S = s) = \binom{n}{n_1}^{-1} \qquad \text{for each } s \in \{0, 1\}^n \text{ where } \sum_i s_i = n_1$$

Definition 3 (True risk of q'th learner averaged over splits)**.** Given the data $O_1, \ldots, O_n \in \mathcal{O}$ and a set of learners $\{\theta_q(P^0_{n,S}) \mid 1 \leq q \leq p\}, p \in \mathbb{N}$ applied to our training data $P^0_{n,S}$. The risks of these learners averaged over some split-variable S is given as a function of q

$$q \mapsto E_S \int L(x, \theta_q(P^0_{n,S}))dP(x) = E_S R(\theta_q(P^0_{n,S}))$$

Where P is the true distribution for our data X.

Definition 4 (Oracle selector)**.** The oracle selector is a function $\tilde{q} : \mathcal{O}^n \to \{1, \ldots, p\}$ which finds the learner that minimizes the true risk given our data $O_1, \ldots, O_n \in O$.

$$\tilde{q}(O_1, \ldots, O_n) = \arg\min_{1 \leq q \leq p} E_S R(\theta_q(P^0_{n,S}))$$

Where $P^0_{n,s}$ is the empirical distribution over the training set of $O_1, \ldots, O_n$ as specified by some split-variable S.

In similar manner to the above the defintions, we can define the cross-validation risk and the cross-validation selector for our learners

Definition 5 (Cross-validation risk of i'th learner averaged over splits)**.** Given the data $O_1, \ldots, O_n \in \mathcal{O}$ and a set of learners $\{\theta_q(P^n_{n,S}) \mid 1 \leq q \leq p\}, p \in \mathbb{N}$. The cross-validation risks of these learners averaged over some split-variable S is given as a function of q

$$q \mapsto E_S \int L(x, \theta_q(P^0_{n,S}))dP^1_{n,s}(x) = E_S \hat{R}(\theta_q(P^0_{n,S}))$$

Where $P^1_{n,S}$ is the empircal distribution over the validation of $O_1, \ldots, O_n$. We write $\hat{R}$ for the empirical risk over the validation set.

Definition 6 (Cross-validation selector)**.** The cross-validation selector is a function $\hat{q} : \mathcal{O}^n \to \{1, \ldots, p\}$ which finds the learner that minimizes the cross-validation risk given our data $O_1, \ldots, O_n \in \mathcal{O}$.

$$\hat{q}(O_1, \ldots, O_n) = \arg\min_{1 \leq q \leq p} E_S \hat{R}(\theta_q(P^0_{n,S}))$$

Where $\hat{R}$ is the empirical risk over the validation set and $P^0_{n,s}$ is the empirical distribution over the training set of $O_1, \ldots, O_n$ as specified by some split-variable S.

We are interested in the risk difference between the cross-validation selector and and the oracle selector, we remark that the optimal risk is attained at the true value θ_0

$$R(\theta_0) = \int L(x, \theta_0)dP(x),$$

and clearly it is the case that $R(\theta_0) \leq R(\theta)$ for any learner θ of θ_0. Given a set of learners we define the centered conditional risk as the difference

$$\begin{aligned}\Delta_S(\theta_{\hat{q}}, \theta_0) &= R(\theta_{\hat{q}}(P^0_{n,S})) - R(\theta_0) \\ &= E_S \int L(x, \theta_{\hat{q}}(P^0_{n,S})) - L(x, \theta_0)dP(x)\end{aligned}$$

The following result is due to [LD03]:

Theorem 7 (Asymptotic equality)**.** *The cross validation selector $\hat{q}$ performs asymptotically as well as the oracle selector $\tilde{q}$ in the sense that*

$$\frac{\Delta_S(\theta_{\hat{q}}, \theta_0)}{\Delta_S(\theta_{\tilde{q}}, \theta_0)} \to 1 \qquad \textit{in probability for } n \to \infty$$

2 The discrete super learner, dSL

2.1 Finite sample properties

3 The ensemble super learner, eSL

4 Simulation results

5 Discussion

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