Notes for Elements of Statistical Learning

Sebastian Claici

December 2, 2024

Chapter 2: Overview of Supervised Learning

Statistical Decision Theory

The goal is to minimize the expected prediction error:

$$EPE(f) = E (Y - f(X))^{2}$$

$$= \int [y - f(x)]^{2} p(x, y) dx dy$$
(1)

If we break down the expectation as $E_{X,Y} = E_X E_{Y|X=x}$ we can rewrite this as

$$\begin{aligned} \text{EPE}(f) &= \mathbf{E}_X \mathbf{E}_{Y|X=x} (Y - f(X))^2 \\ &= \int_X \int_Y [y - f(x)]^2 p(y|X = x) p(x) \, \mathrm{d}y \, \mathrm{d}x \\ &= \int_X p(x) \left(\int_Y [y - f(x)]^2 p(y|X = x) \, \mathrm{d}y \right) \mathrm{d}x \end{aligned}$$

We have moved the dependence on p(x) outside the inner expectation. Since f is unconstrained, we can solve for the optimal f pointwise. That is:

$$\underset{f}{\operatorname{arg\,min}} \operatorname{EPE}(f) = \underset{c}{\operatorname{arg\,min}} \int_{Y} [y - c]^{2} p(y|X = x) \, \mathrm{d}y$$

Differentiating wrt c and using the fact that

$$\int_{Y} y \ p(y|X=x) \, \mathrm{d}y = \mathrm{E}(Y|X=x)$$

gives us (2.13) in the book.

Nearest-neighbor methods try to model the regression function directly by averaging predictions around the query point x. To drive this point home, we can show that $NN(x) \to x$ as the number of training points $N \to \infty$.

To sketch this proof out, assume x_1, \ldots, x_N are drawn i.i.d from X. We want to bound $\min_i ||x - x_i||$, but since this is a bit complicated, let's instead compute

$$P(||x - x_i|| \ge \varepsilon, \forall i).$$

for some $\varepsilon > 0$.

Since the x_i are sampled independently, we can expand the probability as

$$P(\|x - x_i\| \ge \varepsilon, \forall i) = \prod_{i=1}^{N} P(\|x - x_i\| \ge \varepsilon).$$

As the x_i are also identically distributed, the product can be written as

$$\left[P(\|x - x_i\| \ge \varepsilon) \right]^N$$

which goes to 0 as $N \to \infty$ as long as the probability is not exactly 1. This shows that with infinite samples the Nearest-neighbor of x is x and so nearest neighbors yields the Bayes optimal decision boundary even with a single neighbor.

However, we often do not have enough samples to use a model-free approach to regression. The second proposal is to assume the regression function is linear in its arguments:

$$f(x) \approx x^T \beta$$

If we plug this for f into (1), we get

$$\int [y - x^T \beta]^2 p(x, y) \, \mathrm{d}x \, \mathrm{d}y.$$

We can differentiate this wrt β^1

$$\frac{\partial \text{EPE}}{\partial \beta} = 2 \int x[y - x^T \beta] p(x, y) \, dx \, dy$$
$$= 2 \left(\int xy \ p(x, y) \, dx \, dy - \int xx^T \beta \ p(x, y) \, dx \, dy \right)$$

Since β is not a random variable, we can set this to 0 to arrive at the minimizer in (2.16) in the book:

$$\beta = [\mathbf{E}(XX^T)]^{-1}\mathbf{E}(XY)$$

¹See this link for a review of matrix calculus.

Bias-Variance Decomposition

We can express the mean-squared error in terms of a squared bias term and a variance term. In equation (2.25) in the book, these vary w.r.t. the training set T. To clarify the notation a bit, x_0 is the point 0, \hat{y}_0 is the model estimate (in this case the nearest neighbor estimate), and $f(x_0)$ is the true value at 0, but the following derivation holds generally for any model approximation \hat{y} of a function $f(x)^2$:

$$MSE(x_0) = E_T [f(x_0) - \hat{y}_0]^2$$

$$= E_T [\hat{y}_0 - E_T[\hat{y}_0] + E_T[\hat{y}_0] - f(x_0)]]^2$$

$$= E_T [\hat{y}_0 - E_T[\hat{y}_0]]^2 + (f(x_0) - E_T[\hat{y}_0])^2$$

$$= Var_T(\hat{y}_0) + Bias^2(\hat{y}_0)$$

It is a somewhat instructive exercise to figure out how to go from the second line to the third. Easiest if you recall that

$$E_T[E_T[y]] = E_T[y]$$
$$E_T[f(x)] = f(x)$$

In the example in the book, the variance is consistently low, but the bias increases with dimension as the nearest point to 0 becomes increasingly distant.

We can discuss equations (2.27) and (2.28) in the book briefly. We have

$$\hat{\beta} = (X^T X)^{-1} X^T y = (X^T X)^{-1} X^T (X \beta + \varepsilon) = \beta + (X^T X)^{-1} X^T \varepsilon$$

and thus

$$\hat{y}_0 = x_0^T \hat{\beta} = x_0^T (\beta + (X^T X)^{-1} X^T \varepsilon)$$

which gives

$$\hat{y}_0 = x_0^T \beta + \sum_{i=1}^N l_i(x_0) \varepsilon_i \tag{2}$$

since $x_0^T(X^TX)^{-1}X^T\varepsilon$ is a scalar and

$$(x_0^T(X^TX)^{-1}X^T)^T = X(X^TX)^{-1}x_0$$

to give the expression in the book.

Let's write out EPE(x_0). Note that because the true data was generated from a noisy process, we have to integrate out for y_0 given a fixed x_0 :

$$EPE(x_0) = E_{y_0|x_0} E_T [y_0 - \hat{y}_0]^2$$

²See, for example, the wikipedia page.

In this particular case, because Y depends on X stochastically, $E_T = E_X E_{Y|X}$ We will write out $y_0 - \hat{y}_0$ as:

$$y_0 - \hat{y}_0 = (y_0 - x_0^T \beta) + (x_0^T \beta - E_T[\hat{y}_0]) + (E_T[\hat{y}_0] - \hat{y}_0)$$

Let's square this, and keep in mind that $E[\varepsilon] = 0$, and $Var[\varepsilon] = \sigma^2$:

$$E_{y_0|x_0}E_T[y_0 - \hat{y}_0]^2 = E_{y_0|x_0}[y_0 - x_0^T \beta]^2 + (x_0^T \beta - E_T[\hat{y}_0])^2 + E_T[\hat{y}_0 - E_T[\hat{y}_0]]^2 +$$
cross terms

For the cross terms, we notice the following:

$$E_{y_0|x_0}(y_0 - x_0^T \beta) = 0$$

$$E_T(E_T[\hat{y}_0] - \hat{y}_0) = 0$$

and

$$(x_0^T \beta - \mathcal{E}_T[\hat{y}_0]) = \mathcal{E}_T \left[\sum_{i=1}^N l_i(x_0) \varepsilon_i \right]$$

$$= \mathcal{E}_X \left[\sum_{i=1}^N l_i(x) \mathcal{E}_{Y|X}(\varepsilon_i) \right]$$

$$= 0$$
(3)

where we have used (2).

This gives

$$E_{y_0|x_0}E_T[y_0 - \hat{y}_0]^2 = E_{y_0|x_0}[y_0 - x_0^T \beta]^2 + (x_0^T \beta - E_T[\hat{y}_0])^2 + E_T[\hat{y}_0 - E_T[\hat{y}_0]]^2 +$$

$$= Var[y_0|x_0] + Bias^2(\hat{y}_0) + Var_T(\hat{y}_0)$$

but the bias is 0 by (3), and $Var[y_0|x_0] = \sigma^2$, we have:

$$E_{y_0|x_0}E_T[y_0 - \hat{y}_0]^2 = \sigma^2 + Var_T(\hat{y}_0).$$

To finish the derivation, let's write out $\operatorname{Var}_T(\hat{y}_0)$. We have just proved that $\operatorname{E}_T(\hat{y}_0) = x_0^T \beta$, and so

$$\operatorname{Var}_{T}(\hat{y}_{0}) = \operatorname{E}_{T} \left[x_{0}^{T} (X^{T} X)^{-1} X^{T} \varepsilon \right]^{2}$$
$$= \operatorname{E}_{T} \left[x_{0}^{T} (X^{T} X)^{-1} X^{T} \varepsilon \varepsilon^{T} X (X^{T} X)^{-1} x_{0} \right]$$

Since $\varepsilon \sim N(0, \sigma^2)$, $\varepsilon \varepsilon^T = \sigma^2 I_N$, and we can replace above:

$$Var_{T}(\hat{y}_{0}) = E_{T} \left[x_{0}^{T} (X^{T} X)^{-1} X^{T} X (X^{T} X)^{-1} x_{0} \right] \sigma^{2}$$
$$= E_{T} \left[x_{0}^{T} (X^{T} X)^{-1} x_{0} \right] \sigma^{2}$$

which is the value in the book.

To derive (2.28), we assume large N and that $X^TX \to N\text{Cov}(X)$, hence:

$$E_{x_0} EPE(x_0) = \sigma^2 + E_{x_0} \left[x_0^T (X^T X)^{-1} x_0 \right] \sigma^2$$

$$\sim \sigma^2 + E_{x_0} \left[x_0^T Cov(X)^{-1} x_0 \right] \sigma^2 / N$$
(4)

Now $x_0^T \text{Cov}(X)^{-1} x_0$ is a scalar, and can be written as $\text{trace}[x_0^T \text{Cov}(X)^{-1} x_0]$. Exploiting the cyclic properties of the trace operator:

$$\begin{aligned} \mathbf{E}_{x_0} \mathrm{trace}[x_0^T \mathrm{Cov}(X)^{-1} x_0] &= \mathbf{E}_{x_0} \mathrm{trace}[\mathrm{Cov}(X)^{-1} x_0 x_0^T] \\ &= \mathrm{trace}[\mathrm{Cov}(X)^{-1} \mathbf{E}_{x_0}(x_0 x_0^T)] \\ &= \mathrm{trace}[\mathrm{Cov}(X)^{-1} \mathrm{Cov}(x_0)] \end{aligned}$$

Since $Cov(X) = Cov(x_0)$, and since each training set point is p-dimensional:

$$\operatorname{trace}[\operatorname{Cov}(X)^{-1}\operatorname{Cov}(x_0)] = \operatorname{trace}[I_p] = p$$

which when replaced in (4) gives equation (2.28) in the book.

Exercises

2.1. (Classifying to max)

The wording here is a bit confusing, but I am assuming that \hat{y} is constant, and we want to show

$$\arg\min_{k} \|\hat{y} - t_k\| = \arg\max_{k} \hat{y}_k.$$

To prove this, note that $||t_k|| = 1, \forall k$, and the length of \hat{y} is constant with $0 \leq \hat{y}_k \leq 1$. For this setup, we can square the cost function without affecting the optimal value. Then

$$\min_{k} \|\hat{y} - t_k\|^2 = \min_{k} \langle \hat{y} - t_k, \hat{y} - t_k \rangle = \|t_k\|^2 + \|\hat{y}\|^2 + 2\min_{k} (-\langle \hat{y}, t_k \rangle)$$

Since $\min_k(-\langle \hat{y}, t_k \rangle) = \max_k \langle \hat{y}, t_k \rangle$ and $\langle \hat{y}, t_k \rangle = \hat{y}_k$ we have

$$\arg\min_{k} \|\hat{y} - t_k\|^2 = \arg\max_{k} \langle \hat{y}, t_k \rangle = \arg\max_{k} \hat{y}_k.$$

2.2. (Bayes decision boundary)

Computing the Bayes decision boundary only makes sense after the means of the two classes have been sampled. If we integrate out that step, the decision boundary is the diagonal through the origin that separates $[0, 1]^T$ and $[1, 0]^T$.

The procedure for sampling each data point looks something like this (for the blue means):

$$(\mu_1, \dots, \mu_{10}) \sim \mathcal{N}([1, 0]^T, I_n)$$
 for $i = 1..100$
$$k \sim \text{Cat}(10, [1/10, \dots, 1/10])$$

$$x \sim \mathcal{N}(\mu_k, I_n/5)$$
 end for

The decision boundary is where

$$P(\text{blue}) \sum_{i=1}^{10} \exp\left\{-5\|x - \mu_k\|^2 / 2\right\} = P(\text{orange}) \sum_{i=1}^{10} \exp\left\{-5\|x - \nu_k\|^2 / 2\right\}$$

which can be computed easily.

2.3. (Median distance in high dimensions)

Let x be a point sampled from the unit hypersphere. The probability that $||x|| \le r$ is equal to the ratio between the volume of the radius r hypersphere and the volume of the unit sphere. That is, the cdf of ||x|| in d-dimensions is

$$F(r) = r^d.$$

Now let's say we sample n points x_1, x_2, \ldots, x_n . The probability that the closest of these is distance r away can be written as

$$F(\min_{i} ||x||_{i} \le r) = 1 - (1 - F(r))^{n} = 1 - (1 - r^{d})^{n}$$
(5)

since each of the x_i is independent, and the event that the closest is distance r away is the same as the event that all of them are at least distance r away.

The median is defined as the distance r for which $F(\min_i ||x||_i \le r) = 1/2$. If we substitute in (5), we have

$$1 - (1 - r^{d})^{n} = \frac{1}{2}$$

$$1 - r^{d} = \left(\frac{1}{2}\right)^{1/n}$$

$$r = \left(1 - \left(\frac{1}{2}\right)^{1/n}\right)^{1/d}$$

To compute the distance to the mean is a bit more of a challenge. Recall that the pdf of a random variable is the derivative of the cdf, and so, from (5) we have

$$p(\min_{i} ||x||_{i} \le r) = nd(1 - r^{d})^{n-1}r^{d-1}$$

The mean distance then integrates this over $r \in [0, 1]$:

$$d_{\text{mean}} = nd \int_0^1 (1 - r^d)^{n-1} r^{d-1} dr$$

Integrating this explicitly is impossible, but we can relate it to the Beta function:

$$B(z_1, z_2) = \int_0^1 t^{z_1 - 1} (1 - t)^{z_2 - 1} dt$$

which can be integrated numerically and is available in most programming languages (e.g. scipy.special.beta).