## The Linear Model

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The *linear model* is a theoretical framework that unifies a number of statistical concepts, like ANOVA and regression.

# 1 Derivatives and linear algebra

We will need a few results concerning derivatives of linear algebra expressions. Consider a linear function:

$$f: \mathbb{R}^n \to \mathbb{R}, f(\beta) = a^t \beta = \begin{pmatrix} a_1 & a_2 & \cdots & a_n \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \vdots \\ \beta_n \end{pmatrix}$$
 (1)

Here,  $a \in \mathbb{R}^{n \times 1}$  and  $\beta \in \mathbb{R}^{n \times 1}$ , so in other words:

$$f(\beta) = a_1 \beta_1 + a_2 \beta_2 + \dots + a_n \beta_n \tag{2}$$

The (multidimensional) derivative is therefore:

$$\frac{\partial f}{\partial \beta} = \nabla_{\beta} f = \begin{pmatrix} a_1 \\ a_2 \\ \vdots \\ a_n \end{pmatrix} = a \tag{3}$$

If  $A \in \mathbb{R}^{n \times m}$  instead, each column in a will map according to equation 3 under differentiation, and so generalizes to:

$$\frac{\partial}{\partial \beta} A^t \beta = A \tag{4}$$

Since  $a^t\beta = \beta^t a$ , by analogy this means that we also have:

$$\frac{\partial}{\partial \beta} \beta^t A = A \tag{5}$$

Similarly, consider a quadratic form in  $\beta$ :

$$g: \mathbb{R}^n \to \mathbb{R}, g(\beta) = \beta^t A \beta = \begin{pmatrix} \beta_1 & \beta_2 & \cdots & \beta_n \end{pmatrix} \begin{pmatrix} a_{11} & a_{12} & \cdots & a_{1n} \\ a_{21} & a_{22} & \cdots & a_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{n1} & a_{n2} & \cdots & a_{nn} \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \vdots \\ \beta_n \end{pmatrix}$$
(6)

Here,  $A \in \mathbb{R}^{n \times n}$  and  $\beta \in \mathbb{R}^{n \times 1}$ . Furthermore, A is assumed to be symmetric, such that  $a_{ij} = a_{ji}$ . Multiplying out, this means that:

$$g(\beta) = \sum_{i=1}^{n} \sum_{j=1}^{n} \beta_i a_{ij} \beta_j \tag{7}$$

Differentiating with respect to  $\beta_k$  only terms where i = k or i = j will contribute. However, the case i = j = k is distinct. So, when i = k we get the contribution  $a_{kj}\beta_j$ . When j = k we get  $\beta_i a_{ij}$ . And when i = j = k we get  $2a_{kk}\beta_k$ . All in all, when summing up, we get two of each a- $\beta$  set (because of the symmetry of A). So:

$$\frac{\partial g}{\partial \beta_k} = 2\sum_{i=1}^n a_{ik}\beta_i \tag{8}$$

Or more compactly:

$$\frac{\partial g}{\partial \beta} = \nabla_{\beta} g = 2A\beta \tag{9}$$

## 2 Ordinary least squares estimation (OLS)

## 2.1 Statement of the problem

The general problem is this: We wish to model a linear relationship between a response variables Y and p predictor variables  $X_1, X_2, \dots X_p$ . In other words:

$$Y = \beta_1 X_1 + \beta_2 X_2 + \dots + \beta_p X_p + \epsilon \tag{10}$$

Here, the  $\beta$ 's are the coefficients corresponding to the X's. The random term  $\epsilon$  is known as the *error term* and represents the deviations from the exact model. Since it is a random variable, so is Y. Now, assume that we have n realizations (data points), so that  $y_i$  corresponds to  $x_{i1}, x_{i2}, \dots x_{ip}$ . In matrix form equation (10) now becomes:

$$y = X\beta + \epsilon \tag{11}$$

Here,  $y \in \mathbb{R}^{n \times 1}$ ,  $X \in \mathbb{R}^{n \times p}$  and  $\beta \in \mathbb{R}^{p \times 1}$ :

$$y = \begin{pmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{pmatrix}, \quad X = \begin{pmatrix} x_{11} & x_{12} & \dots & x_{1p} \\ x_{21} & x_{22} & \dots & x_{2p} \\ \vdots & \vdots & \ddots & \vdots \\ x_{n1} & x_{n2} & \dots & x_{nn} \end{pmatrix}, \quad \beta = \begin{pmatrix} \beta_1 \\ \beta_2 \\ \vdots \\ \beta_n \end{pmatrix}, \quad \epsilon = \begin{pmatrix} \epsilon_1 \\ \epsilon_2 \\ \vdots \\ \epsilon_n \end{pmatrix}$$
(12)

X is known as the design matrix. Given y and X, we seek the best fit for  $\beta$ .

### 2.2 Least squares

There's a number of criteria one could use to pick the best fitting  $\beta$ . Here, we will search for the one that minimizes the square of the differences in predicted and actual y values. When predicting, we can't include the error term, and so the predicted values for a set of parameters  $\hat{\beta}$  are simply:

$$y = X\hat{\beta} \tag{13}$$

The squared difference is:

$$||y - X\hat{\beta}||^{2} = (y - X\hat{\beta})^{t}(y - X\hat{\beta}) = (y^{t} - \hat{\beta}^{t}X^{t})(y - X\hat{\beta})$$
$$= y^{t}y - y^{t}X\hat{\beta} - \hat{\beta}^{t}X^{t}y + \hat{\beta}^{t}X^{t}X\hat{\beta}$$

Taking the derivative with respect to  $\beta$  we can now use equations 4, 5, and 9 to yield:

$$-2X^t y + 2X^t X \hat{\beta} \tag{14}$$

Since we're looking for a minimum, this vector should be equal to zero:

$$-2X^{t}y + 2X^{t}X\hat{\beta} = 0 \Leftrightarrow \hat{\beta} = (X^{t}X)^{-1}X^{t}y \tag{15}$$

Here it has been assumed that  $X^tX$  is invertible. If  $X^tX$  is not invertible, we have a case of *perfect* (multi)collinearity. The equations in 15 are known as the normal equations for the model. Inserting into equation (11) we get the corresponding predicted y-values, also denoted by a hat:

$$\hat{y} = X\hat{\beta} = \underbrace{X(X^t X)^{-1} X^t}_{H} y \tag{16}$$

The matrix  $H = X(X^tX)^{-1}X^t$  is often called the *hat matrix*, since it puts the hat on the y's. The hat matrix can also be used to find *residuals*, i.e. the difference between actual and predicted y-values:

$$e = y - \hat{y} = y - Hy = \underbrace{(I - H)}_{M} y \tag{17}$$

### 2.3 Properties of the OLS estimator

First of all we note, that the estimation is a linear function of the y values. The estimated value of  $\beta$  according to OLS is:

$$\hat{\beta} = (X^t X)^{-1} X^t y = (X^t X)^{-1} X^t (X\beta + \epsilon)$$
(18)

Here, we have re-inserted equation 11. We may now consider the expected value:

$$E[\hat{\beta}] = \beta + (X^t X) X^t E[\epsilon] \tag{19}$$

# 3 Geometric picture

It is useful to adapt the picture of the columns of X spanning a p-dimensional hyperplane in n-dimensional space. y is then a vector, and  $X\hat{\beta}$  is found by projecting y onto the hyperplane; The corresponding point is exactly the one that minimizes the distance between y (as a point) and the hyperplane.

### 3.1 Projection operators

A linear map that is symmetric and idempotent is called a *projection*. A matrix corresponding to such a mapping is a projection matrix.

**Theorem 1.** The hat matrix H is a projection matrix.

*Proof.* We need to show that H is symmetric and idempotent. Symmetry:

$$X(X^{t}X)^{-1}X^{t})^{t} = X\left[(X^{t}X)^{-1}\right]^{t}X^{t}$$
(20)

But the transpose of an inverse is the same as the inverse of the transpose, so:

$$[(X^{t}X)^{-1}]^{t} = [(X^{t}X)^{t}]^{-1} = (X^{t}X)^{-1}$$
(21)

This proves the symmetry of H. Idempotency:

$$H^{2} = \left[X(X^{t}X)^{-1}X^{t}\right]^{2} = X(X^{t}X)^{-1}X^{t}X(X^{t}X)^{-1}X^{t} = X(X^{t}X)^{-1}X^{t} = H$$
(22)

This also turns out to be true for the matrix used to find residuals:

**Theorem 2.** The matrix M = I - H is a projection matrix.

*Proof.* Symmetry follows from the symmetry of H. Idempotency:

$$M^{2} = (I - H)^{2} = I^{2} + H^{2} - 2H = I + H - 2H = I - H = M$$
 (23)

### 4 The Gauss-Markov theorem

So far, we have considered only the ordinary least squares (OLS) estimator of the vector  $\beta$ . But clearly it is not the only possibility. What is the justification for picking this particular estimator? The answer lies in the Gauss-Markov theorem. According to this theorem, under certain conditions, the OLS is the best linear unbiased estimator. This is often abbreviated to BLUE. Let's examine the meaning of this.

#### 4.1 Linear and unbiased

We already know that the OLS estimator is linear in terms of the y's.

Recall, that an estimator is *unbiased* if its expectation value is the true value. Here it means:

$$E[\hat{\beta}] = \beta \tag{24}$$

From equation 19 we know, that this is true exactly when the expectation value of  $\epsilon$  is zero.

#### 4.2 'Best'

In this context, "best" means having the smallest possible variance. We could express this by requiring every estimator element of the  $\hat{\beta}$  vector to have a minimal variance. But we will go further than this: Let  $\hat{\gamma} = \sum_{i=1}^p c_i \hat{\beta}_i = C \hat{\beta}$  be an arbitrary linear combination of the predictors. Then the variance of every such expression should be minimal. According to the usual rules of calculating variance:

$$\operatorname{var}(\hat{\gamma}) = \operatorname{var}(C\hat{\beta}) = C\operatorname{var}(\hat{\beta})C^{t}$$
(25)

Consider another estimator  $\tilde{\beta}$  which has a covariance matrix of:

$$var(\tilde{\beta}) = var(\hat{\beta}) + \Delta \tag{26}$$

Here  $\Delta$  is the deviation from our proposed best estimator. The variance of a linear combination of the tilde estimator is:

$$\operatorname{var}(\tilde{\gamma}) = C\operatorname{var}(\tilde{\beta})C^{t} = C(\operatorname{var}(\hat{\beta}) + \Delta)C^{t} = \operatorname{var}(\tilde{\beta}) + \underline{C\Delta C^{t}}$$
 (27)

Hence  $\hat{\beta}$  is the best estimator if and only if the underbraced quantity is always positive, except when C=0. In other words, exactly when  $\Delta$  is positive definite.

#### 4.3 The theorem

**Theorem 3.** (Gauss-Markov) Given a linear model with design matrix X, responses y, and true parameters  $\beta$ . Assume the following three conditions for the error terms  $\epsilon_i$  are met:

- The expected value is zero:  $E[\epsilon_i] = 0$ .
- The variance of the error terms are finite and constant:  $var(\epsilon_i) = \sigma^2 < \infty$ . This is known as homoscedasticity.
- The error terms are pairwise uncorrelated:  $cov(\epsilon_i, \epsilon_j) = 0, i \neq j$ .

Then, the OLS estimator  $\hat{\beta} = (X^t X)^{-1} X^t y$  is BLUE.

*Proof.* Let  $\tilde{\beta} = Cy$  be another unbiased linear estimator of  $\beta$ . We may then write the matrix C as  $(X^tX)^{-1}X^t + D$ , where D is the deviation from the OLS estimator. Then we may calculate the expected value:

$$E[\tilde{\beta}] = E[Cy] = E[(X^t X)^{-1} X^t + D)(X\beta + \epsilon)] = (X^t X)^{-1} X^t + D)(X\beta + E[\epsilon])$$
(28)

By the first assumption this is:

$$((X^{t}X)^{-1}X^{t} + D)X\beta = (X^{t}X)^{-1}X^{t}X\beta + DX\beta = \beta + DX\beta$$
 (29)

Since  $\tilde{\beta}$  is an unbiased estimator, we must have DX = 0. Now, let's compute the variance:

$$\operatorname{var}(\tilde{\beta}) = \operatorname{var}(Cy) = C\operatorname{var}(y)C^{t}$$
 (30)

Here, we've used a property of variances. By the homoscedasticity assumptions, this is simply:

$$\sigma^2 C C^t = \sigma^2 ((X^t X)^{-1} X^t + D) ((X^t X)^{-1} X^t + D)^t$$
 (31)

Since  $X^tX$  is symmetric, so is the inverse, so  $((X^tX)^{-1}X^t+D)^t = X(X^tX)^{-1}+D^t$ . So we get:

$$\sigma^{2}((X^{t}X)^{-1}X^{t} + D)(X(X^{t}X)^{-1} + D^{t})$$
(32)

Ignoring the  $\sigma^2$  factor for a while, this is:

$$(X^{t}X)^{-1}X^{t}X(X^{t}X)^{-1} + (X^{t}X)^{-1}X^{t}D^{t} + DX(X^{t}X)^{-1} + DD^{t}$$
 (33)

But since we just concluded DX = 0 the two middle terms vanish (since  $X^tD^t = (DX)^t = 0$ ). So, reinstating the  $\sigma^2$ , the variance is

$$\operatorname{var}(\tilde{\beta}) = \sigma^2 (X^t X)^{-1} + \sigma^2 D D^t$$
(34)

The first term is what we would get without the D term, and is therefore the variance of the OLS estimator.  $DD^t$  is a positive definite matrix, and hence according to the section above,  $\hat{\beta}$  is the least variance estimator.

Note that no assumptions of independence, identical distribution or normality is assumed of the error terms.

#### 4.4 Omitted variable bias

Assume we have forgotten, missed or simply not had access to a (set of) important predictor variables  $z_1, z_2, \ldots, z_n$ . The parameters corresponding to these are called  $\gamma$ , and we may now rewrite the model:

$$y = X\beta + \underbrace{Z\gamma + \delta}_{\epsilon} \tag{35}$$

Here,  $\delta$  is the error terms associated with the new variables. We will assume that since the missing variables are important/good, the expectation value of this error is zero:  $E[\delta] = 0$ . The error term of the original model is now the underbraced part of the equation.

## 5 Abstract definition of a linear model

This section will deal with the linear model in its most abstract form.

Let V be a vector space of finite dimension N. To specify a linear model we need two ingredients:

- A subspace  $L \subset V$ . Do note that we require L to be a proper subset of V, i.e.  $\dim L < N$ . This subspace is known as the *mean value subspace*.
- An inner product  $\langle \cdot, \cdot \rangle$  on V.

The inner product induces a family of inner products  $\langle \cdot, \cdot \rangle_{\sigma^2}$  parametrized by  $\sigma^2 > 0$ :

$$\langle\!\langle \cdot, \cdot \rangle\!\rangle_{\sigma^2} = \frac{\langle \cdot, \cdot \rangle}{\sigma^2}$$
 (36)

These inner products are known as *precisions*. While they do not agree on distances, the precisions do agree on orthogonality.

The linear model