

Linear Regression for Quantitative Finance

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1 Introduction

We begin by defining terms.

Consider some variable you are trying to understand, y . You have a dataset that contains n data points, and each data point, i has its own value of y , y_i . You describe this collection of values for y in a vector representation, as follows:

$$y = \langle y_1, y_2, \dots, y_n \rangle \quad (1)$$

You decide to try to predict the value of this variable as a linear combination of a series of k predictor variables, X . X is best described as an $n \times k$ matrix, where each of the n rows corresponds to a single data point, and each of the k columns corresponds to a single predictor variable. You can refer to a single row (i.e. a single data point with k variables) as X_i . These rows can also be described in a vector representation, as follows:

$$X_i = \langle X_{i1}, X_{i2}, \dots, X_{ik} \rangle \quad (2)$$

A linear combination of k predictor variables requires k coefficients. You call these k coefficients β , and you also describe them using a vector representation:

$$\beta = \langle \beta_1, \beta_2, \dots, \beta_k \rangle \quad (3)$$

Call your final prediction \hat{y} . The linear combination of your X predictor variables with β coefficients is defined in vector notation as:

$$\hat{y} = X\beta \quad (4)$$

Equivalently, you can write this using subscripts:

$$\hat{y}_i = \sum_{j=1}^k X_{ij} \cdot \beta_j \quad (5)$$

Your estimate \hat{y} might not be perfectly equal to y , so you should define some residual variable, ϵ to describe the difference:

$$\epsilon_i = y_i - \sum_{j=1}^k X_{ij} \cdot \beta_j \quad (6)$$

This allows you to finally describe the true variable, y , in terms of your predictor variables:

$$y_i = \sum_{j=1}^k X_{ij} \cdot \beta_j + \epsilon_i \quad (7)$$

Equivalently, in the vector representation:

$$y = X\beta + \epsilon \quad (8)$$

The goal of the regression is to pick some vector of coefficients β (length k) such that the residuals of our prediction, ϵ (length n), are minimized.

2 Regression

There are many metrics to use to determine whether a given vector of coefficients, β minimizes ϵ . By far the most common is the sum of the difference of squares, S , which is defined as follows:

$$S(\beta) = \sum_{i=1}^n \left| y_i - \sum_{j=1}^k X_{ij} \cdot \beta_j \right|^2 \quad (9)$$

Equivalently, in the vector representation:

$$S(\beta) = \|y - X\beta\|^2 \quad (10)$$

One nice feature of the sum of the difference of squares is that the minimization problem has a unique solution, as long as the k columns of matrix X are linearly independent. This solution is derived below.

We begin by redefining S in terms of the residuals, ϵ :

$$S(\beta) = \sum_{i=1}^n \epsilon_i^2 \quad (11)$$

Equivalently, in the vector representation:

$$S(\beta) = \|\epsilon\|^2 \quad (12)$$

If we think of β as a surface, then $S(\beta)$ is minimized when the gradient vector of $S(\beta)$ has magnitude zero. There is a geometric argument for this: if the gradient is not zero, then there is a direction that we can move such that the gradient may be reduced. The gradient, $\nabla S(\beta)$, is a vector of length k , where each element of the vector is defined as a partial derivative:

$$\frac{\partial S}{\partial \beta_j} = 2 \sum_{i=1}^n \epsilon_i \frac{\partial \epsilon_i}{\partial \beta_j} \quad (13)$$

Recalling the definition of ϵ from Equation 6, the partial derivate is easy to determine:

$$\frac{\partial \epsilon_i}{\partial \beta_j} = -X_{ij} \quad (14)$$

This allows us to completely determine each element of the gradient of $S(\beta)$:

$$\frac{\partial S}{\partial \beta_j} = -2 \sum_{i=1}^n \epsilon_i \cdot X_{ij} \quad (15)$$

$$\frac{\partial S}{\partial \beta_j} = -2 \sum_{i=1}^n \left(y_i - \sum_{l=1}^k X_{il} \cdot \beta_l \right) \cdot X_{ij} \quad (16)$$

We define $\hat{\beta}$ as the value of β that minimizes $S(\beta)$ such that $\frac{\partial S}{\partial \beta_j} = 0$ for all j between 1 and k :

$$-2 \sum_{i=1}^n \left(y_i - \sum_{l=1}^k X_{il} \cdot \hat{\beta}_l \right) \cdot X_{ij} = 0 \quad (17)$$

We can divide out a factor of 2 and rearrange this. For all j between 1 and k :

$$\sum_{i=1}^n \sum_{l=1}^k X_{ij} X_{il} \hat{\beta}_l = \sum_{i=1}^n X_{ij} y_i \quad (18)$$

Equivalently, in vector notation:

$$(X^T X) \hat{\beta} = X^T y \quad (19)$$

Solving for $\hat{\beta}$:

$$\hat{\beta} = (X^T X)^{-1} X^T y \quad (20)$$

3 Explained variance

We define the number of degrees of freedom, ndof , to be the difference between the number of entries in the dataset, n , and the number of discriminating variables, k :

$$\text{ndof} = n - k \quad (21)$$

Note that in classical linear regression n must be greater than k , or else the dataset could be fit perfectly. With this in mind, we can define the residual standard deviation, $\hat{\sigma}$, as follows:

$$\hat{\sigma} = \sqrt{\sum_{i=1}^n \frac{\epsilon_i^2}{(n - k)}} \quad (22)$$

Here, $\hat{\sigma}$ summarizes the scale of the residuals. You can think of this as a measure of the average distance that the linear model output falls from the true value.

If we call the true standard deviation of the data s_y , then we can define R^2 such that:

$$R^2 = 1 - \frac{\hat{\sigma}^2}{s_y^2} \quad (23)$$