

Regression Models (Part IV)

A. Notation for data

We write X_1, X_2, \dots, X_n to describe n data points. As an example, consider the data set $\{1, 2, 5\}$ then $X_1 = 1, X_2 = 2, X_3 = 5$ and $n = 3$.

Of course, there's nothing in particular about the variable X . We often use a different letter, such as Y_1, \dots, Y_n to describe a data set. We will typically use Greek letters for things we don't know. Such as, μ being a population mean that we'd like to estimate.

B. The empirical mean

The empirical mean is a measure of center of our data. Under sampling assumptions, it estimates a population mean of interest. Define the empirical mean as

$$\bar{X} = \frac{1}{n} \sum_{i=1}^n X_i$$

Notice if we subtract the mean from data points, we get data that has mean 0. That is, if we define

$$\tilde{X}_i = X_i - \bar{X}$$

then the mean of the \tilde{X}_i is 0. This process is called centering the random variables. Recall from the previous lecture that the empirical mean is the least squares solution for minimizing

$$\sum_{i=1}^n (X_i - \mu)^2$$

C. The empirical standard deviation and variance

The variance and standard deviation are measures of how spread out are data is. Under sampling assumptions, they estimate variability in the population. We define the empirical variance as

$$S^2 = \frac{1}{n-1} \sum_{i=1}^n (X_i - \bar{X})^2 = \frac{1}{n-1} \left(\sum_{i=1}^n X_i^2 - n\bar{X}^2 \right)$$

The empirical standard deviation is defined as $S = \sqrt{S^2}$

Notice that the standard deviation has the same units as the data. The data defined by X_i / s have empirical standard deviation 1. This is called **scaling** the data.

D. Normalization

We can combine centering and scaling of data as follows to get normalized data. In particular, the data defined by:

$$Z_i = \frac{X_i - \bar{X}}{s}$$

have empirical mean zero and empirical standard deviation 1. The process of centering then scaling the data is called **normalizing** the data. Normalized data are centered at 0 and have units equal to standard deviations of the original data. Example, a value of 2 from normalized data means that data point was two standard deviations larger than the mean.

Normalization is very useful for creating data that comparable across experiments by getting rid of any shifting or scaling effects.

E. The empirical covariance

This class is largely considering how variables covary. This is estimated by the empirical covariance. Consider now when we have pairs of data, (X_i, Y_i) . Their empirical covariance is defined as:

$$Cov(X, Y) = \frac{1}{n-1} \sum_{i=1}^n (X_i - \bar{X})(Y_i - \bar{Y}) = \frac{1}{n-1} \left(\sum_{i=1}^n X_i Y_i - n \bar{X} \bar{Y} \right)$$

This measure is of limited utility, since its units are the product of the units of the two variables. A more useful definition normalizes the two variables first.

The **correlation** is defined as:

$$Cor(X, Y) = \frac{Cov(X, Y)}{S_x S_y}$$

where S_x and S_y are the estimates of standard deviations for the X observations and Y observations, respectively. The correlation is simply the covariance of the separately normalized X and Y data. Because the the data have been normalized, the correlation is a unit free quantity and thus has more of a hope of being interpretable across settings.

F. Some facts about correlation

First, the order of the arguments is irrelevant $Cor(X, Y) = Cor(Y, X)$ Secondly, it has to be between -1 and 1 , $-1 \leq Cor(X, Y) \leq 1$. Thirdly, the correlation is exactly -1 or 1 only

when the observations fall perfectly on a negatively or positively sloped, line, respectively.

Fourthly, $Cor(X, Y)$ measures the strength of the linear relationship between the two variables, with stronger relationships as $Cor(X, Y)$ heads towards -1 or 1 . Finally, $Cor(X, Y) = 0$ implies no linear relationship.