Consider a simple regression model

$$y_i = \beta_0 + \beta_1 x_i + \varepsilon_i \tag{1}$$

for i = 1, 2, ..., n. Stacking the observations, we can write the model for the entire sample as

$$\mathbf{y} = \beta_0 \mathbf{l}_n + \beta_1 \mathbf{x} + \boldsymbol{\varepsilon} \tag{2}$$

where l_n is the n by 1 vector of 1's; and $\boldsymbol{y}, \boldsymbol{x}, \boldsymbol{\varepsilon}$ are all n by 1 vectors. Here $\{y_i, x_i\}$'s are random observations (i.i.d.), and β_0 and β_1 are unknown coefficients. The least-squares methodology tries to infer about the unknown coefficients by finding a best fit to sample data taken from the population. The best fit refers to values of β_0 and β_1 so that the sum of the squared residuals (or prediction errors) is the least possible, i.e.

$$(\widehat{\beta}_0, \widehat{\beta}_1)' = \underset{\beta_0, \beta_1}{\operatorname{arg \, min}} \sum_{i=1}^n (y_i - \beta_0 - \beta_1 x_i)^2.$$

It can be shown that the least squares estimator for the slope coefficient β_1 is given by

$$\widehat{\beta}_{1} = \frac{\sum_{i=1}^{n} (x_{i} - \bar{x})(y_{i} - \bar{y})}{\sum_{i=1}^{n} (x_{i} - \bar{x})^{2}}$$

and the least-squares estimator for the intercept is given by

$$\widehat{\beta}_0 = \bar{y} - \widehat{\beta}_1 \bar{x},$$

where \bar{y} and \bar{x} denote the sample means of y and x, respectively.

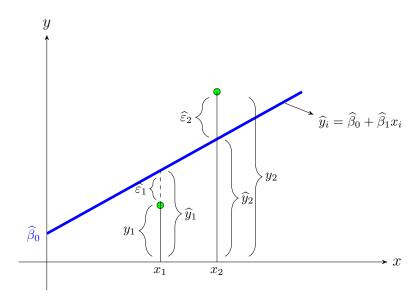


Figure 1. The least-squares regression line.

- (0) Write a function which will take \boldsymbol{y} and \boldsymbol{x} as inputs and return estimates of β_0 and β_1 using the least-squares methodology.
- (1) Set the seed to 37 for the random number generator in numpy, i.e., np.random.seed(37).

- (2) Generate an array of zeros of size 5000, and name it tmp1. For 5000 times, do the following in a loop:
 - (a) generate 1000 observations on \boldsymbol{x} by drawing randomly from the standard normal distribution, (i.e., \boldsymbol{x} is a 1000 by 1 vector of standard Gaussian random variables);
 - (b) generate 1000 observations on ε by drawing randomly from the standard normal distribution, (i.e., ε is a 1000 by 1 vector of standard Gaussian random variables);
 - (c) generate the \boldsymbol{y} vector using

$$\mathbf{y} = 0.5\,\mathbf{l}_n + 1.8\,\mathbf{x} + \boldsymbol{\varepsilon}$$

using the draws from (a) and (b);

- (d) estimate the simple linear regression model using your function;
- (e) save β_1 estimate to tmp1.
- (3) Calculate the mean of tmp1. Is it 1.8? Is it close to 1.8?
- (4) Redo (2) and (3), but this time in 2(b), generate 1000 observations on ε as

$$\varepsilon_i = -0.5x_i + v_i$$

for i = 1, 2, ..., 1000, where v_i is drawn from the standard normal distribution.

(5) Calculate the mean of tmp1 again. Compared to your finding in (3), is it further away from 1.8? Why?