Consider the fitted values that result from performing linear regression without an intercept. In this setting, the ith fitted value takes the form

$$\hat{y}_i = x_i \hat{\beta},$$

where

$$\hat{\beta} = \left(\sum_{i=1}^{n} x_i y_i\right) / \left(\sum_{i'=1}^{n} x_{i'}^2\right).$$
 (3.38)

Show that we can write

$$\hat{y}_i = \sum_{i'=1}^n a_{i'} y_{i'}.$$

What is $a_{i'}$?

$$\hat{\mathbf{y}}_{\hat{i}} = \chi_{\hat{i}} \hat{\boldsymbol{\beta}}$$

$$= \chi_{\hat{i}} \frac{\sum_{\hat{i}=1}^{n} \chi_{\hat{i}} \, \mathbf{y}_{\hat{i}}}{\sum_{\hat{i}'=1}^{n} \chi_{\hat{i}'} \, \mathbf{y}_{\hat{i}'}}$$

$$= \chi_{\hat{i}} \cdot \frac{\sum_{\hat{i}'=1}^{n} \chi_{\hat{i}'} \, \mathbf{y}_{\hat{i}'}}{\sum_{\hat{j}=1}^{n} \chi_{\hat{j}}^{2}}$$

$$= \sum_{\hat{i}'=1}^{n} \frac{\chi_{\hat{i}} \cdot \chi_{\hat{i}'} \, \mathbf{y}_{\hat{i}'}}{\sum_{\hat{j}=1}^{n} \chi_{\hat{j}}^{2}}$$

$$= \sum_{\hat{i}'=1}^{n} \alpha_{\hat{i}'} \, \mathbf{y}_{\hat{i}'}$$

$$= \sum_{\hat{i}'=1}^{n} \alpha_{\hat{i}'} \, \mathbf{y}_{\hat{i}'}$$

$$= \sum_{\hat{i}'=1}^{n} \alpha_{\hat{i}'} \, \mathbf{y}_{\hat{i}'}$$

ESL 3,4

Ex. 3.4 Show how the vector of least squares coefficients can be obtained from a single pass of the Gram–Schmidt procedure (Algorithm 3.1). Represent your solution in terms of the QR decomposition of X.

$$= R^{-1}(R^{T})^{-1}R^{T}Q^{T}y$$

```
In [1]: # Data Import
                       import ssl
                       import pandas as pd
                        ssl._create_default_https_context = ssl._create_unverified_context #Github에서 데이터를 바로 불러오도록 하는 세팅입니다. 해당 코드 무시하고 데이
                        터 받아서 쓰셔도 됩니다!
                       data = pd.read_csv('https://github.com/YonseiESC/ESC-21SUMMER/blob/main/week1/HW/week1_data.csv?raw=True')
                       y = data['mpg']
                        x = data.drop(['mpg'],axis=1)
In [2]:
                      idx = x['horsepower']!='?'
                       x = x[idx]
                       y = y[idx]
                       x = x.astype({'horsepower': 'float'})
                       import numpy as np
                        # numpy 모듈만을 이용해주세요.
                       def YourOwnRegression(x,y):
                                  x = np.array(x)
                                   y = np.array(y)
                                   XTX = np.dot(np.transpose(x), x)
                                  betahat = np.dot(np.dot(np.linalq.inv(XTX), np.transpose(x)), y)
                                  yhat = np.dot(x, betahat)
                                   return betahat, yhat
                        # 결과물 반화
                        bhat, yhat = YourOwnRegression(x,y)
                       print("beta hat: ", bhat,'\n', "vhat: ", vhat)
                     beta hat: [-0.5226089 0.01022108 -0.020873 -0.00639456 -0.05202195 0.61025869]
                        yhat: [15.93081361 14.45720405 16.11263762 15.93670419 16.10071197 10.51021782
10.27543153 10.53128391 9.67525181 13.49634208 15.59946068 15.17857818
                        14.95056695 18.23756934 23.85149163 20.70116218 21.04691637 22.47738545
                       25.4075601 27.88849004 21.93938914 23.54965655 23.61026333 24.5700506 22.02475307 7.48992449 8.73758029 8.68094775 6.39448743 26.01781879 25.50668622 25.43458909 22.95714525 17.50355598 18.56684981 18.98997897
                       18.64504732 11.74185816 10.43958016 12.2762232 12.39844199 7.0217271 8.71466792 6.09084578 20.89073634 24.7795066 18.89340516 20.0843144
                      8.71466792 6.09084578 20.89073634 24.7795066 18.89340516 20.0843144 25.76559234 26.24104628 26.30753902 26.59195154 28.28090927 29.28279009 28.2609542 27.13912286 25.6470868 26.69569542 26.07663641 24.9880403 26.2074288 11.93647037 11.52899822 12.73330182 13.0723569 15.65493241 9.60277136 10.60920749 10.79899231 10.95329347 25.45996983 14.19610698 13.24891771 11.63170439 13.0781754 21.23759877 24.50545484 21.19494316 26.4550875 25.15275892 25.40418994 24.27095036 26.56185652 26.71587043 13.3986307 16.2651911 14.74101165 13.99404444 15.68483539 8.35588404 12.1559116 12.08192325 12.63929331 9.52710571 8.09046672 15.38896808 20.7055219 19.98343495 22.03288824 21.95534612 22.05191343 28.9278806 8.64356655 8.94866353 10.06261966 10.76884679 23.08246175 26.05144779
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                       12.92654367 27.29002111 29.11881818 26.97024946 31.3571203 28.84708125 28.08036911 28.1370547 27.59838633 25.57416431 26.13565935 28.85620844 21.24635801 20.04621563 20.64756809 22.47771079 11.68048944 13.0191654 12.19083021 11.60515866 16.62277687 17.11534294 18.13462735 17.75872575
                       22.48538772 20.67124626 21.09405268 28.38999264 25.6178386 23.45163703 25.94471506 25.06438254 28.04340891 25.7120626 22.52118358 30.02318129 21.7091604 24.86417853 23.27558145 23.28386212 24.73315284 31.11419428 27.02526368 28.70906066 26.57792197 28.30139996 28.71233562 14.76463858
                       14.84864669 16.74288177 14.99473679 21.98169626 21.41233074 23.42150596 23.09392064 29.90480163 29.05099549 30.6230356 31.72965517 19.18911918
                       23.04352004 23.59480183 29.0393549 30.0230530 31.72565317 19.18518918 20.22456108 19.35231654 22.55428679 31.49394493 30.08735698 29.08175554 27.10777245 22.52930717 16.43688261 21.62943493 23.04180738 17.16039851 13.38138443 16.19537407 17.04478509 17.57825889 30.44250197 29.8139795 11.85863268 28.62881666 30.99577532 17.43872127 16.22016388 15.84658308 14.94255744 20.751491 21.24312777 19.93366366 20.83473714 15.57784903 15.61161917 14.64046595 14.81182982 31.10306759 26.25259415 28.89497299
                       26.0337752 30.58653253 30.07282718 30.94607365 29.52523606 24.56685474 26.54846444 25.84824622 31.61605985 32.87446415 31.51916094 30.71307435
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                        28.75362625 31.67091587 32.82504694 31.90800579 32.48874873 28.41207605
                       28.75362625 31.67091587 32.82504694 31.90800579 32.48874873 28.41207605 27.03039926 26.14656464 23.51339355 31.28036395 28.07931906 29.22595587 29.68156125 30.99424682 31.82881074 27.47919606 31.75096112 32.18671672 30.48403252 26.14671839 24.9076443 33.71404606 31.67048481 33.75190878 25.44572433 29.75593179 29.36847327 30.7931412 30.22783049 29.42774108 29.5118248 27.68686195 31.04492087 34.83572405 34.04069724 34.82406496 32.80271279 33.2140812 32.94723852 33.3257177 32.15965292 33.06580622 30.68952944 32.1263293 31.98760285 30.97773269 28.97782474 29.22077581 25.39617051 24.96456589 26.39147031 25.8288513 23.70096728 21.82763916 6.04576044 23.80012019 29.58125731 29.4046443 31.00734511 30.01258259
                       33.35582651 33.00291086 32.77391248 32.62463215 32.35858312 34.13676256 34.1003472 33.88249939 26.77440244 27.64530275 30.34069094 28.04577073 29.6541942 31.74589579 27.85165521 28.93428241 32.95667079 32.29831861
                       29.77500853 29.05306756]
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