

MATH 829: Introduction to Data Mining and
Analysis
Subset selection

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Testing multiple coefficients

We saw before how to use the t -statistic to test

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$$H_1 : \beta_{i_1} \neq 0 \text{ or } \beta_{i_2} \neq 0 \text{ or } \dots \text{ or } \beta_{i_k} \neq 0.$$

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$$F = \frac{(\text{RSS}_0 - \text{RSS}_1)/(p - p_0)}{\text{RSS}_1/(n - p)},$$

where

RSS_1 = residual sum of squares for full model,

RSS_0 = residual sum of squares for the nested smaller model.

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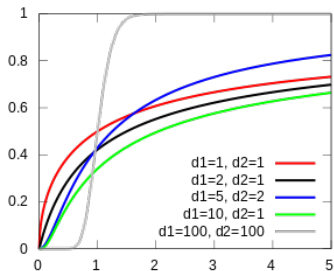
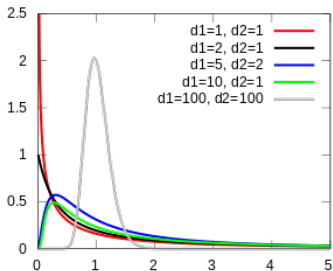
RSS_0 = residual sum of squares for the nested smaller model.

Can be seen as a measure of the *change in residual sum-of-squares per additional parameter in the bigger model*.

Testing multiple coefficients (cont.)

Under the H_0 assumption that the smaller model is correct, the F statistic has an F -distribution

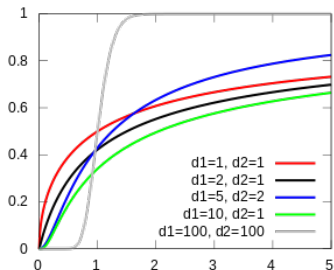
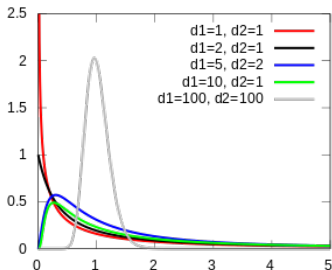
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To test if a group of coefficients are 0:

- 1 Compute the F -statistic.
- 2 Reject H_0 for large values of the F -statistic.

A simple illustration of the previous ideas.

```
import numpy as np
import statsmodels.api as sm
# Generate random data
n = 50
epsilon = np.random.randn(n,1) # Try varying the sample size
X = np.random.randn(n,5)
y = 3*X[:,0] + 4*X[:,1] + epsilon # Try changing coefficients
results = sm.OLS(y,X).fit()
print(results.summary())
R = [[0,0,1,0,0],
      [0,0,0,1,0],
      [0,0,0,0,1]]
print(results.f_test(R))
R = [[1,0,0,0,0], [0,1,0,0,0]]
print(results.f_test(R))
```


Python (cont.)

```
=====
                        OLS Regression Results
=====
Dep. Variable:          y      R-squared:                0.954
Model:                  OLS    Adj. R-squared:           0.949
Method:                 Least Squares    F-statistic:              187.2
Date:                   Tue, 19 Jan 2016  Prob (F-statistic):      6.23e-29
Time:                   12:40:31    Log-Likelihood:          -78.513
No. Observations:      50      AIC:                     167.0
Df Residuals:          45      BIC:                     176.6
Df Model:               5
=====
                coef    std err          t      P>|t|      [95.0% Conf. Int.]
-----
x1                3.3360    0.208     16.071    0.000     2.918    3.754
x2                4.0380    0.167     24.139    0.000     3.701    4.375
x3               -0.1904    0.167     -1.143    0.259    -0.526    0.145
x4                0.1282    0.186     0.689    0.495    -0.247    0.503
x5                0.1163    0.155     0.751    0.456    -0.195    0.428
=====
Omnibus:                0.748    Durbin-Watson:           2.074
Prob(Omnibus):          0.688    Jarque-Bera (JB):        0.755
Skew:                   -0.002    Prob(JB):                0.686
Kurtosis:               2.398    Cond. No.:               1.91
=====
<F test: F=array([[ 0.76049081]]), p=[[ 0.52218257]], df_denom=45, df_num=3>
<F test: F=array([[ 390.38886666]]), p=[[ 3.69709216e-29]], df_denom=45, df_num
=2>
```

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- Note: there are $\binom{p}{k}$ subsets of size k and 2^k possible subsets. So the procedure is only computationally feasible for small values of p .

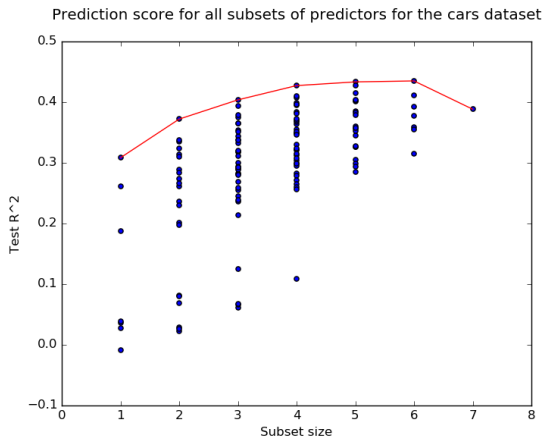
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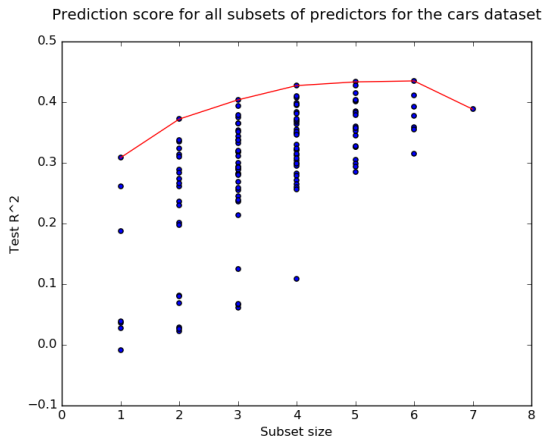
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- The leaps and bounds procedure (Furnival and Wilson, 1974) makes this feasible for p as large as 30 or 40.

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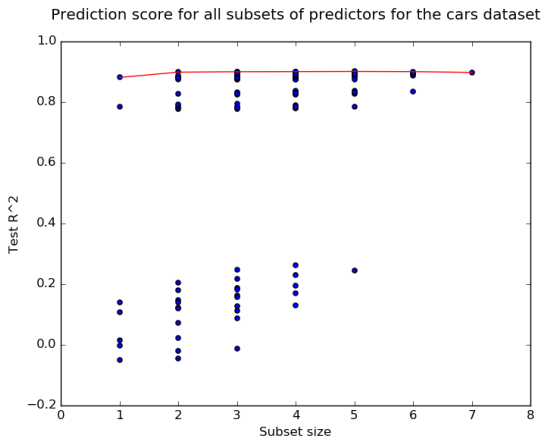
Best subset = ['Mileage', 'Liter', 'Doors', 'Cruise', 'Sound', 'Leather'].

Not included = ['Cylinder']

Best subset of 4 elements: ['Mileage', 'Liter', 'Cruise', 'Leather']

Best subset selection: cars dataset, Chevrolet

Restricting to Chevrolet only:



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Nevertheless, the stepwise approaches often return predictors similar to the predictors obtained from more complex methods with better theory.

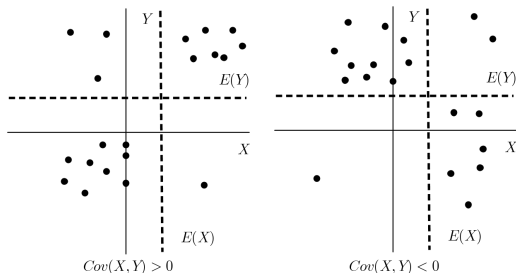
Recall: **Covariance** is a measure of linear dependence between random variables:

$$\text{Cov}(X, Y) = E((X - E(X))(Y - E(Y))).$$

Correlation

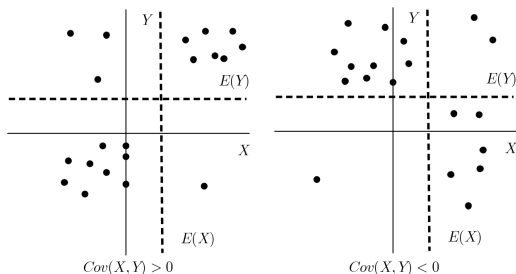
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Properties:

- 1 $\text{Cov}(\cdot, \cdot)$ is bilinear and symmetric.
- 2 $\text{Cov}(X, X) = \text{Var}(X)$.
- 3 $\text{Cov}(X, Y) = E(XY) - E(X)E(Y)$.
- 4 X, Y independent $\Rightarrow \text{Cov}(X, Y) = 0$.

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Theorem: Assume $\text{Var}(X), \text{Var}(Y) < \infty$. The correlation coefficient $\rho(X, Y)$ satisfies

$$-1 \leq \rho(X, Y) \leq 1.$$

Moreover, $\rho(X, Y) = \pm 1$ if and only if $\mathbb{P}(Y = aX + b) = 1$ for some constants a, b . In this case, $a > 0$ if $\rho(X, Y) = 1$ and $a < 0$ if $\rho(X, Y) = -1$.

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- Continued till none of the variables have correlation with the residuals.

In other words:

- $C = \emptyset, \hat{y}_1 = \bar{y}, \beta_1 = \dots = \beta_p = 0.$
- Suppose X_{i_1} is most correlated to $y.$

$$C \rightarrow C \cup \{X_{i_1}\}.$$

- Solve $y - \hat{y}_1 = \alpha_{i_1} X_{i_1} + \epsilon.$

$$\beta_{i_1} \rightarrow \beta_{i_1} + \alpha_{i_1}.$$

- etc.

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- 1 Unlike forward-stepwise regression, none of the other variables are adjusted when a term is added to the model.
- 2 The process can take **more than p** steps to reach the least squares fit.
- 3 Historically, forward stagewise regression has been dismissed as being inefficient.
- 4 However, it can be quite competitive, especially in very high-dimensional problems.

Forward stagewise regression (cont.)

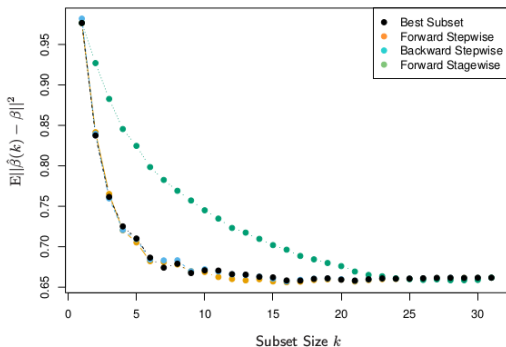


FIGURE 3.6. Comparison of four subset-selection techniques on a simulated linear regression problem $Y = X^T\beta + \varepsilon$. There are $N = 300$ observations on $p = 31$ standard Gaussian variables, with pairwise correlations all equal to 0.85. For 10 of the variables, the coefficients are drawn at random from a $N(0, 0.4)$ distribution; the rest are zero. The noise $\varepsilon \sim N(0, 6.25)$, resulting in a signal-to-noise ratio of 0.64. Results are averaged over 50 simulations. Shown is the mean-squared error of the estimated coefficient $\hat{\beta}(k)$ at each step from the true β .

ESL, Fig. 3.6