

COMS 4771-2 Fall 2018 Homework 2

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Problem 1

@inproceedingsseabold2010statsmodels, title=Statsmodels: Econometric and statistical modeling with python, author=Seabold, Skipper and Perktold, Josef, booktitle=9th Python in Science Conference, year=2010,

- (a) Test risks of the ordinary least squares estimator is 0.53978372.
Test risks of the sparse linear predictor is 0.5565456.
- (b) Names of the variables with non-zero coefficients in the sparse linear predictor, along with the coefficient values are:
 $\hat{\beta}_0$: 5.82610034
volatile acidity : -0.24474264
sulphates : 0.10847122
alcohol : 0.37362602
- (c) volatile acidity :
total sulfur dioxide : -0.4220072736318842
citric acid : -0.3755630209692122

sulphates :
chlorides : 0.4071152150059981
fixed acidity : 0.29686322762159284

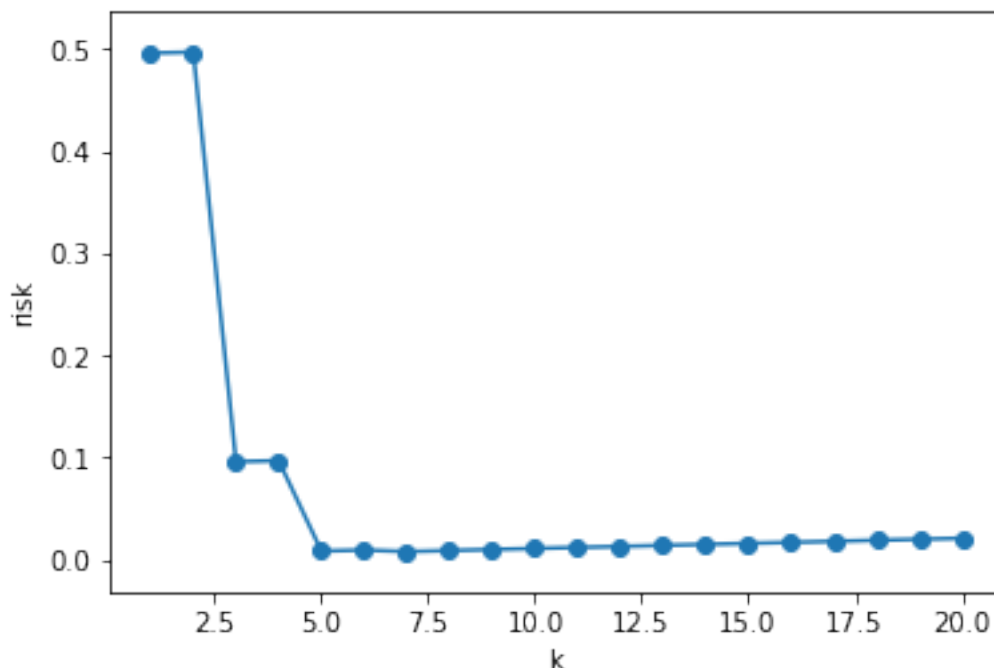
alcohol :
density : -0.6795934959123543
residual sugar : -0.34626243022889536

Problem 2

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- (a) As the conclusion in handout, we know that $E[R(\hat{\beta})] \leq R(\beta^*) + \frac{\bar{\sigma}^2 r}{n}$
 so, $E \left[\frac{1}{n} \sum_{i=1}^n (x_i^T \hat{\beta} - E(Y_i))^2 \right] \leq R \left[\frac{1}{n} \sum_{i=1}^n (x_i^T \beta^* - E(Y_i))^2 \right] + \frac{\bar{\sigma}^2 r}{n}$
 $A^T A \hat{\beta}^* = A^T b$
 $E(Y)$ using the Taylor's (remainder) theorem, so only the $(k+1)$ -th will remain.
 $\text{var}(Y)=1$
 $\text{rank of } A \text{ is } (k+1)$
 so, $E \left[\frac{1}{n} \sum_{i=1}^n (x_i^T \hat{\beta} - E(Y_i))^2 \right] \leq \left(\frac{\pi^{k+1}}{(k+1)!} \right)^2 + \frac{k+1}{n}$
- (b)
- (c) The average fixed design risks for each k are shown in the table and plot as below.

k	risk	k	risk	k	risk	k	risk
1	0.496164	6	0.009518	11	0.012068	16	0.017014
2	0.497105	7	0.008000	12	0.012997	17	0.018002
3	0.096005	8	0.009021	13	0.014052	18	0.019102
4	0.097012	9	0.010011	14	0.015061	19	0.020058
5	0.008510	10	0.011038	15	0.016037	20	0.021035



Problem 3

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(a) Features remain after the screening is 73

(b) The empirical risk of $\hat{\beta}(\hat{J})$ is 0.5184631137998037

(c) The empirical risk of $\hat{\beta}(I)$ is 0.838553272038375

(d) I randomly generated 10 pairs of data and labels, computed their $\hat{\beta}(\hat{J})$ and $\hat{\beta}(I)$. Below are the outcomes:

n is 236 ; d is 753 ; random risk is 0.410838458083448 ; Irisk is 0.7323342590472934

n is 217 ; d is 447 ; random risk is 0.5105407622750471 ; Irisk is 0.8575392192518303

n is 269 ; d is 772 ; random risk is 0.4170041318711085 ; Irisk is 0.7552813920061883

n is 235 ; d is 560 ; random risk is 0.5092864102832999 ; Irisk is 0.7832188161174825

n is 217 ; d is 812 ; random risk is 0.2794990858610098 ; Irisk is 0.6756825703025061

n is 765 ; d is 903 ; random risk is 0.7074898586441014 ; Irisk is 0.9530342963972447

n is 224 ; d is 637 ; random risk is 0.45179746254645453 ; Irisk is 0.7863053297575778

n is 339 ; d is 342 ; random risk is 0.750291984275502 ; Irisk is 0.9355944079878

n is 485 ; d is 469 ; random risk is 0.6744905836310303 ; Irisk is 0.9382904951357046

n is 332 ; d is 311 ; random risk is 0.7060903365984399 ; Irisk is 0.9407575809911612

I find that when n is much larger than k, the difference between two risks will decrease, while when n is equal or smaller than k, the difference will grow larger.

This may be caused by the screening procedure. We add the relationship to x and y by counting their correlations. So the risk we compute by that way is smaller than other methods. But with the number of n grows, this relationship can be reduced, so the difference between two risks gets smaller.

(e) I think the $\hat{\beta}(\hat{J})$ will become smaller but the empirical risk will be larger.

Because when we choose the same $\hat{\beta}$ on a different data-label set, the $\hat{\beta}$ cannot fully reflect the new relationships between data and labels. So the ordinary least squares estimate $\hat{\beta}(\hat{J})$ will be smaller as the correlation decreases. At the same time, square loss risk will become larger because of the inaccurate parameters.