Regularized Regression Methods - Demo Examples of LASSO

STAT 32950-24620

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Shrinkage method, Ridge regression

Ridge regression shrinks the regression coefficients by imposing a penalty (also called regularization) on their 2-norm size

$$\|\boldsymbol{\beta}\|_2^2 = \boldsymbol{\beta}' \boldsymbol{\beta} = \beta_1^2 + \dots + \beta_r^2$$

Ridge regression parameters $\beta_0, \beta_1, \cdots, \beta_r$ are obtained by

$$\min_{\boldsymbol{\beta}} \left\{ \|\boldsymbol{y} - \beta_0 - \boldsymbol{Z}\boldsymbol{\beta}\|_2^2 \right\} \qquad \textit{when} \quad \|\boldsymbol{\beta}\|_2^2 \leq s_r$$

Or equivalently,

$$\min_{\boldsymbol{\beta}} \left\{ \|\boldsymbol{y} - \beta_0 - \boldsymbol{Z}\boldsymbol{\beta}\|_2^2 + \lambda_{\textit{ridge}} \|\boldsymbol{\beta}\|_2^2 \right\}$$

where $s_r \propto 1/\lambda_{ridge}$, tuning parameters aim to minimize MSE.

Review - LS regression method

Least squares (LS) linear regression model

$$\mathbf{Y} = \beta_0 + \mathbf{Z}\boldsymbol{\beta} + \boldsymbol{\varepsilon}$$

LS parameters β_0 , $\beta = (\beta_1, \dots, \beta_r)$ are obtained by minimizing

$$\|\mathbf{y} - \beta_0 - \mathbf{Z}\boldsymbol{\beta}\|_2^2$$

over all possible values of parameter vector $\boldsymbol{\beta} \in \mathbb{R}^r$.

$$\hat{oldsymbol{eta}}_{LS} = \operatorname{argmin}_{oldsymbol{eta} \in \mathbb{R}^r} \left\{ \|oldsymbol{y} - eta_0 - oldsymbol{Z}oldsymbol{eta} \|_2^2
ight\}$$

 $\hat{\beta}_{LS}$ has the smallest variance among all linear unbiased estimators of β .

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Least absolute shrinkage and selection operator (LASSO)

LASSO regression regularizes the regression coefficients by imposing a penalty or a regularization on their 1-norm size

$$\|\boldsymbol{\beta}\|_1 = |\beta_1| + \dots + |\beta_r|$$

LASSO regression parameters $\beta_0, \beta_1, \cdots, \beta_r$ are obtained by

$$\min_{\beta} \left\{ \| \boldsymbol{y} - eta_0 - \boldsymbol{Z} \boldsymbol{\beta} \|_2^2 \right\} \quad \text{when} \quad \| \boldsymbol{\beta} \|_1 \leq s_l$$

Or equivalently,

$$\min_{eta} \left\{ \| oldsymbol{y} - eta_0 - oldsymbol{Z} oldsymbol{eta} \|_2^2 + \lambda_{ extit{lasso}} \| oldsymbol{eta} \|_1
ight\}$$

where $s_l \propto 1/\lambda_{lasso}$, tuning parameters aim to minimize MSE.

Elastic Net

Elastic Net linear regression

$$\min_{\beta} \left\{ \| \boldsymbol{y} - \beta_0 - \boldsymbol{Z}\boldsymbol{\beta} \|_2^2 \right\}$$

when

$$\|\boldsymbol{\beta}\|_2^2 \leq s_r$$
 and when $\|\boldsymbol{\beta}\|_1 \leq s_l$

Or equivalently,

$$\min_{\beta} \left\{ \| \boldsymbol{y} - \beta_0 - \boldsymbol{Z}\boldsymbol{\beta} \|_2^2 + \lambda_{\textit{ridge}} \| \boldsymbol{\beta} \|_2^2 + \lambda_{\textit{lasso}} \| \boldsymbol{\beta} \|_1 \right\}$$

Another common form:

$$\min_{\boldsymbol{\beta}} \left\{ \| \boldsymbol{y} - \beta_0 - \boldsymbol{Z}\boldsymbol{\beta} \|_2^2 + \lambda \left[\alpha \| \boldsymbol{\beta} \|_2^2 + (1 - \alpha) \| \boldsymbol{\beta} \|_1 \right] \right.$$

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Example data

The covariates are

- 1 lcavol : log-cancer volume
- 2 lweight : log-prostate weight
- 3 age: age of patient
- 4 lbhp: log-amount of benign hyperplasia
- 6 svi: seminal vesicle invasion
- 6 lcp: log-capsular penetration
- 7 gleason: Gleason Score,
- 8 pgg45 : percent of Gleason scores 4 or 5
- Ipsa: is the response variable, log-psa.

Example for LASSO Regression

Example — Prostate cancer dataset

- Number of explanatory variables p = 9, number of observations n = 96.
- measuring the correlation between the level of a prostate-specific antigen and some covariates.
- Used in Hastie, Tibshirani and Friedman (2009)
- from a study by Stamey et al. (1989) of prostate cancer

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Partition the data into Training and Calibration parts

```
pcancer=read.table("pcancer.dat",header=T)
str(pcancer)
'data.frame': 97 obs. of 10 variables:
$lcavol : num -0.58 -0.994 -0.511 -1.204 0.751 ...
$lweight: num 2.77 3.32 2.69 3.28 3.43 ...
$age : int 50 58 74 58 62 50 64 58 47 63 ...
$lbph : num -1.39 -1.39 -1.39 -1.39 -1.39 ...
$svi : int 0 0 0 0 0 0 0 0 0 0 ...
$lcp : num -1.39 -1.39 -1.39 -1.39 -1.39 ...
$gleason: int 6 6 7 6 6 6 6 6 6 6 6 ...
$pgg45 : int 0 0 20 0 0 0 0 0 0 0 ...
$lpsa : num -0.431 -0.163 -0.163 -0.163 0.372 ...
$train : logi TRUE TRUE TRUE TRUE TRUE TRUE ...

train <- pcancer[which(pcancer$train),1:9]
calibrate <- pcancer[which(pcancer$train),1:9]
str(train) # 67 obs. of 9 variables
```

Fit LASSO regression model on the training set

Training data: 70% Calibrating data: 30%

Note: Calibration data is used for validation, just not "cross validation".

Using glmnet to fit a LASSO regression model on the training set.

library(MASS) library(glmnet) y <- as.numeric(train[,9]) x <- as.matrix(train[,1:8]) trainfit = glmnet(x,y) plot(trainfit,label=T)

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Set of models from a LASSO regression fit

For a set of possible λ values, model fit is evaluated on the calibration set.

Different λ values apply different levels of regularization on the L1 norm $\|\beta\|_1$, resulting in different numbers of explanatory variables kept in the model.

In the following plots:

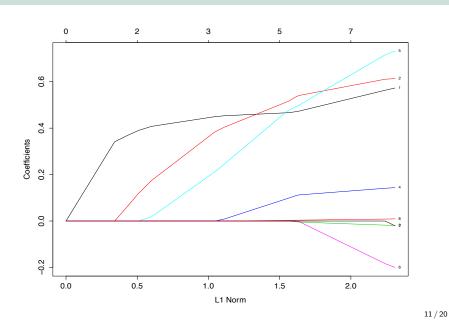
plot(trainfit,label=T,xvar="lambda")

Top axis: DF = number of explanatory variables kept in the model Bottom axis: L1 norm of β (can ask for xvar="lambda" to get log(λ)) Vertical axis: values of coefficients β_i

Visualization of the coefficients in the model (on $log(\lambda)$)

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Visualization of the coefficients in the model (on L1 norm)

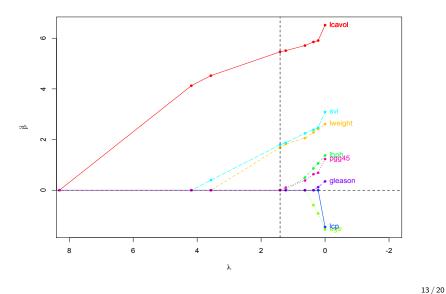


Coefficients

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Log Lambda

A fancier visualization of the coefficients, x-axis in λ



Model selection: Many choices of tuning parameters $(s, or \lambda)$

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```
> coef(trainfit,s=0.1)
                            # 6 term (including intercept)
9 x 1 sparse Matrix of class "dgCMatrix"
(Intercept) -0.064140435
lcavol
             0.462712201
             0.483361985
lweight
1bph
             0.072282456
svi
             0.410200201
1cp
gleason
             0.002245818
> coef(trainfit,s=0.05)
                           # 7 terms (including intercept)
9 x 1 sparse Matrix of class "dgCMatrix"
(Intercept) -0.112643640
             0.470217454
lcavol
lweight
             0.532127691
            -0.002943140
1bph
             0.107622441
             0.489955346
svi
lcp
gleason
pgg45
             0.003463712
```

Model fitting process

```
> print(trainfit) # no of nonzero coeff(DF), %dev explained, lambda value
Call: glmnet(x = x, y = y)
     Df %Dev Lambda
 [1,] 0 0.00000 0.87890
     1 0.09126 0.80080
 [3.] 1 0.16700 0.72970
 [4,] 1 0.22990 0.66480
      1 0.28220 0.60580
     1 0.32550 0.55200
     1 0.36150 0.50290
      1 0.39140 0.45820
     2 0.42810 0.41750
[10,] 2 0.45980 0.38040
      3 0.48770 0.34660
     3 0.51310 0.31590
     3 0.53420 0.28780
      3 0.55180 0.26220
      3 0.56630 0.23890
     3 0.57840 0.21770
      5 0.59170 0.19840
      5 0.60450 0.18070
      5 0.64650 0.10340
      5 0.64990 0.09424
      6 0.66320 0.05393
      6 0.66530 0.04914
[69,] 8 0.69430 0.00157
[70,] 8 0.69430 0.00143
[71,] 8 0.69430 0.00130
```

Parameter-dependent model prediction on validation set

```
nx=as.matrix(calibrate[,1:8])
                                              # saved validation data
cbind(predict(trainfit,newx=nx, s=c(0.1,0.05)),calibrate[,9]) # predict of s=.1, .05, true y
         1
7 2.000388 1.960186 0.7654678
9 1.187210 1.118166 1.0473190
10 1.507195 1.384180 1.0473190
15 2.068642 1.986382 1.3987169
22 2.732481 2.770052 1.6582281
25 2.002415 1.989092 1.7316555
26 2.137355 2.062555 1.7664417
28 1.795435 1.823814 1.8164521
32 1.982366 1.989699 2.0082140
34 1.419724 1.322725 2.0215476
36 2.700799 2.757486 2.0856721
42 2.305471 2.210431 2.3075726
44 2.552468 2.486081 2.3749058
48 2.638200 2.704350 2.5687881
49 2.334230 2.293865 2.5915164
50 2.324743 2.301932 2.5915164
53 2.194626 2.276820 2.6844403
54 3.129746 3.209202 2.6912431
55 3.005596 2.935674 2.7047113
57 1.680437 1.596517 2.7880929
62 3.275113 3.423275 2.8535925
64 3.460785 3.648120 2.8820035
65 2.546161 2.452543 2.8820035
66 2.605518 2.646575 2.8875901
73 2.701967 2.701174 3.0563569
74 2.995277 3.146681 3.0750055
80 3.082012 3.067295 3.5130369
84 3.267951 3.360083 3.5709402
95 3 255211 3 284046 5 1431245
97 3.950538 4.041404 5.5829322
```

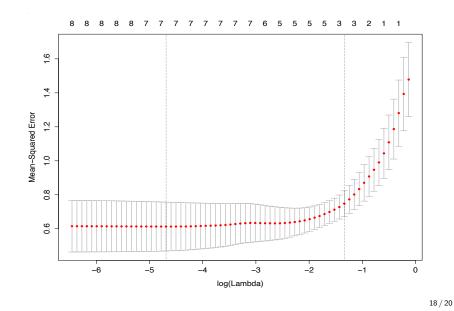
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Alternative model selection: Cross validation

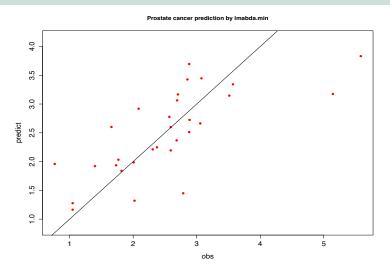
```
> tcvfit = cv.glmnet(x,y) # cross-validation, on training data; nfold=10 default
> tcvfit$lambda.1se
                    # 0.26, log=-1.338545, for example
> coef(tcvfit, s="lambda.min") #8 x's
9 x 1 sparse Matrix of class "dgCMatrix"
(Intercept) 0.193668298
            0.553290578
lweight
           -0.016384627
1bph
           0.137824087
           0.691699046
svi
           -0.163512605
lcp
gleason
pgg45
           0.007860603
> coef(tcvfit, s="lambda.1se") #4 x's
9 x 1 sparse Matrix of class "dgCMatrix"
(Intercept) 0.7185225
          0.4350437
          0.3115982
lweight
          0.1452013
svi
lcp
gleason
pgg45
```

Choose tuning parameters by Mean Squared Error



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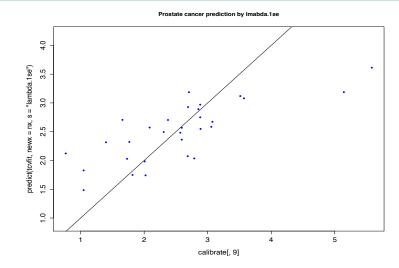
lambda.min: The λ minimizing cross-validated mean squared error



plot(calibrate[,9],predict(tcvfit,newx=nx,s="lambda.min"),asp=1,xlab="obs",ylab="predict",col=2,pch=16)
abline(0,1)

title(main="Prostate cancer prediction by lmabda.min",cex.main=.8)

lambda.1se: The λ of cross-validated MSE within one s.e. of the min



plot(calibrate[,9],predict(tcvfit,newx=nx,s="lambda.1se"),asp=1,,col=4,pch=16,cex=.8)
abline(0,1)
title(main="Prostate cancer prediction by lmabda.1se",cex.main=.8)

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