Regularization Technique for Linear Regression in High Dimension:
Ridge Regression and LASSO

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Preface

Focus point of this part is still the bias-variance tradeoff by introducing the shrinkage methodology to allow the **biased** estimator. The linear regression problem here is used as an example of **model assessment and model selection**. (Chapter 7 [ESL])

Not intention to cover the followings: the optimization theory and numerical methods for the ridge regression and LASSO, which is very important in practice but more close to the optimization field and compress sensing.

For a model fit $\hat{f}_{\text{D}}(x)$ based on the data $\text{D} = (\mathbf{X}, \mathbf{y})$, where $\mathbf{X} \in \mathbb{R}^{n \times p}$ is the design matrix and $\mathbf{y} = \mathbf{X}\beta + \varepsilon \in \mathbb{R}^n$ is the response data, a good measure of the quality at a new test input $x_0 \in \mathbb{R}^p$ is the mean square error (MSE). Let $f(x) = x^{\mathsf{T}}\beta$ be the true value of the output at the point x, then[1]

$$\begin{aligned}
\text{MSE}(\hat{f}_{\text{D}}(x_0)) &:= \mathbb{E}_{\text{D}}(\hat{f}_{\text{D}}(x_0) - f(x_0))^2 \\
&= \text{Var}_{\text{D}}(\hat{f}_{\text{D}}(x_0)) + \left(\mathbb{E}_{\text{D}}\,\hat{f}_{\text{D}}(x_0) - f(x_0)\right)^2.
\end{aligned}$$

- Typically, when bias is low, variance will be high and vice-versa. Choosing estimators often involves a tradeoff between bias and variance.
- So far, OLS estimator $\hat{\beta}^{OLS} = (\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1}\mathbf{X}\mathbf{y}$ is unbiased and also has the minimal MSE if restricted to be _unbiased_ and linear in $\mathbf{y}$.

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[1] If $f(x)$ contains a measurement error ε_0, then simply add the constant $\text{Var}(\varepsilon_0)$

For the OLS fit, its MSE (written with dependency on x_0) is

$$
\begin{aligned}
\mathsf{MSE}^{OLS}(x_0) &= \mathsf{Var}_{\mathbf{y}}(x_0^{\mathsf{T}} \hat{\beta}^{OLS}) \\
&= \mathsf{Var}_{\mathbf{y}}(x_0^{\mathsf{T}} (\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1} \mathbf{X}^{\mathsf{T}} \mathbf{y}) \\
&= \mathbf{X}(\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1} x_0 \, \mathbb{V}(\mathbf{y}) x_0^{\mathsf{T}} (\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1} \mathbf{X}^{\mathsf{T}} \\
&= \sigma_{\varepsilon}^2 \mathbf{X}(\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1} x_0 x_0^{\mathsf{T}} (\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1} \mathbf{X}^{\mathsf{T}} \\
&= \sigma_{\varepsilon}^2 \left\| \mathbf{X}(\mathbf{X}^{\mathsf{T}}\mathbf{X})^{-1} x_0 \right\|_2^2
\end{aligned}
$$

Here $\mathbf{X}$ is treated as deterministic or say that we work on the condition variance of given $\mathbf{X}$.

Exercise

In the above, $x_0 \in \mathbb{R}^p$ is arbitrary. Now let x_0 be each of training sample value x_i, $1 \le i \le N$, show that the in-sample error, which is defined by

$$
\frac{1}{N} \sum_{i=1}^{N} MSE^{OLS}(x_i)
$$

equals to $\frac{p}{N} \sigma_{\varepsilon}^2$. (Equation (7.29) in [ESL])

There can be biased estimators with smaller MSE. The following property is quite general.

Exercise

Define a biased estimate of the coefficient β in the following special form

$$\widetilde{\beta} = (1 + \alpha)\hat{\beta}$$

with a scalar α where $\hat{\beta}$ is an unbiased estimate. Calculate the MSE of $\widetilde{\beta}$ and find a condition that $MSE(\widetilde{\beta}) < MSE(\hat{\beta})$.

- Generally, by regularizing the estimator in some way, its variance will be reduced; if the corresponding increase in bias is small, this will be worthwhile.

- Examples of regularization: subset selection (forward, backward, all subsets)[1]; ridge regression, Lasso.

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[1]read [ISL][ESL] by yourself; not to cover in lecturing

Ridge Regression

The problem of regression in high dimension

- High dimensional problem: the input dimension p is close or greater than the number of samples n.

- When the design matrix $\mathbf{X}$ is high-dimensional, the covariates (the columns of $\mathbf{X}$) are super-collinear. *collinearity* in regression analysis refers to the event of two (or multiple) covariates being highly linearly related.

- In OLS estimate, the dimension of the subspace X may be less than p and then the matrix inverse $(\mathbf{X}^\mathsf{T}\mathbf{X})^{-1}$ does not exist or $\mathbf{X}^\mathsf{T}\mathbf{X}$ is close to singular even if invertible[1].

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[1]Numerical linear algebra uses the **condition number** (defined as the ratio of largest to the smallest eigenvalue) to represent the illness of the problem: the larger, the worse.

Solution of ridge regression

Ridge Regression solves

$$\min_{\beta \in \mathbb{R}^p} \quad \|\mathbf{y} - \mathbf{X}\beta\|_2^2 + \lambda \|\beta\|_2^2$$

The solution is

$$\hat{\beta}^\lambda = (\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}\mathbf{X}^\mathsf{T}\mathbf{y}$$

- Inclusion of $\lambda > 0$ makes problem non-singular even if $\mathbf{X}^\mathsf{T}\mathbf{X}$ is not invertible: This was the original motivation for ridge regression (Hoerl and Kennard, 1970)
- Note $\lambda = 0$ gives the ordinary least squares estimator, and if $\lambda \to \infty$ then $\hat{\beta}^\lambda \to 0$. As λ increases, $\left\|\hat{\beta}^\lambda\right\|_2$ decreases.
- In general, with a good choice of λ, $\hat{\beta}^\lambda$ is a biased estimator that may have smaller mean squared error than the least squares estimator

Let $\mathbf{X} = \mathbf{U}\mathbf{D}\mathbf{V}^\mathsf{T}$ is a singular value decomposition of $\mathbf{X}$: $\mathbf{D}$ is the $p \times p$ diagonal matrix consisting of singular values $d_1 \geq d_2 \geq \ldots \geq d_p$. $\mathbf{U}$ and $\mathbf{V}$ are $n \times p$ and $p \times p$ matrices, respectively [1]. Then $\mathbf{X}^\mathsf{T}\mathbf{X} = \mathbf{V}\mathbf{D}^2\mathbf{V}^\mathsf{T}$ and the ridge regression solution

$$\begin{aligned}
\hat{\beta}^\lambda &= (\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}\mathbf{X}^\mathsf{T}\mathbf{y}, \\
&= (\mathbf{V}\mathbf{D}^2\mathbf{V}^\mathsf{T} + \lambda\mathbf{V}\mathbf{V}^\mathsf{T})^{-1}\mathbf{V}\mathbf{D}\mathbf{U}^\mathsf{T}\mathbf{y} \\
&= \mathbf{V}(\mathbf{D}^2 + \lambda\mathbf{I})^{-1}\mathbf{D}\mathbf{U}^\mathsf{T}\mathbf{y} \\
&= \mathbf{V}\,\mathsf{diag}\left\{\frac{d_j}{d_j^2 + \lambda}\right\}\mathbf{U}^\mathsf{T}\mathbf{y},
\end{aligned}$$

which is well defined for any d_j (some of d_j may be zero) when λ is strictly positive.

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[1] The column space of $\mathbf{U}$ is the column space of $\mathbf{X}$ in $\mathbb{R}^n$ and the column space of $\mathbf{V}$ is the row space of $\mathbf{X}$ in $\mathbb{R}^p$. The columns of $\mathbf{V}$ are the orthonormal eigenvectors of the symmetric positive semidefinite matrix $\mathbf{X}^\mathsf{T}\mathbf{X}$. $\mathbf{U}^\mathsf{T}\mathbf{U} = \mathbf{I}_p$ and $\mathbf{V}\mathbf{V}^\mathsf{T} = \mathbf{V}^\mathsf{T}\mathbf{V} = \mathbf{I}_p$. The number of nonzero $\{d_j\}$ is the rank of $\mathbf{X}$ (which is $\leq \min(n, p)$).

Let $\widetilde{\beta} = \mathbf{V}^\mathsf{T} \hat{\beta}$ and $\widetilde{\mathbf{y}} = \mathbf{U}^\mathsf{T} \mathbf{y}$. Then

$$\widetilde{\beta}^\lambda = \mathbf{V}^\mathsf{T} \hat{\beta}^\lambda = \mathsf{diag} \left\{ \frac{d_j}{d_j^2 + \lambda} \right\} \widetilde{\mathbf{y}}$$

If $\mathbf{D}$ is nonsingular, then $\widetilde{\beta}^{OLS} = \widetilde{\beta}^0 = \mathbf{D}^{-1} \widetilde{\mathbf{y}}$ exists. So

$$\widetilde{\beta}^\lambda = \mathbf{D}_\lambda \mathbf{D}^{-1} \widetilde{\mathbf{y}} = \mathbf{D}_\lambda \widetilde{\beta}^{OLS} \quad \text{where} \quad \mathbf{D}_\lambda := \mathsf{diag} \left\{ \frac{d_j^2}{d_j^2 + \lambda} \right\} \leq \mathbf{I}$$

$$\boxed{\widetilde{\beta}_j^{ridge} = \frac{d_j^2}{d_j^2 + \lambda} \, \widetilde{\beta}_j^{OLS}}$$

Exercise

(1) Find the bias, the variance and the MSE for the transformed ridge coefficient $\widetilde{\beta}^\lambda$ in terms of $\beta, \mathbf{X}, \mathbf{y}, \sigma^2$. Find the optimal λ in theory that minimize the MSE.

(2) Ex 18.7 in [ESL].

Smoother matrix and effective degree of freedom

- A smoother matrix S is a linear operator transforming the observations into predictions:

$$\hat{\mathbf{y}} = \mathbf{S}\mathbf{y}$$

where $\mathbf{S}$ may depend on $\mathbf{X}$, but not $\mathbf{y}$.

- define the **effective degrees of freedom** (or effective number of parameters) for a smoother $\mathbf{S}$:

$$\mathrm{df}(\mathbf{S}) = \mathrm{Trace}(\mathbf{S}) := \sum_j S_{jj}$$

- The smoother matrix of the ridge regression is

$$\mathbf{S} = \mathbf{X}(\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}\mathbf{X}^\mathsf{T}$$

Exercise

Ex. 7.4 and 7.5 in [ESL]

Effective degree of freedom

The ridge solution can be rewritten as the linear mapping from the observations $\mathbf{y}$ to the predictions $\hat{\mathbf{y}}^{ridge}$:

$$
\begin{aligned}
\hat{\mathbf{y}}^{ridge} &= \mathbf{X}\hat{\beta}^{ridge} = \mathbf{U}\mathbf{D}\mathbf{V}^{\mathsf{T}}\,\mathbf{V}\mathbf{D}_\lambda\mathbf{D}^{-1}\mathbf{U}^{\mathsf{T}}\mathbf{y} \\
&= \mathbf{U}\mathbf{D}_\lambda\mathbf{U}^{\mathsf{T}}\mathbf{y} \\
&= \sum_{j=1}^{p}\Big(\frac{d_j^2}{d_j^2+\lambda}\Big)\mathbf{U}_j\mathbf{U}_j^{\mathsf{T}}\mathbf{y},
\end{aligned}
$$

$\mathbf{U}_j$ is the column vector of $\mathbf{U}$. The effective degree of freedom of the ridge regression is

$$
df(\lambda) := \mathrm{Trace}\left(\mathbf{X}(\mathbf{X}^{\mathsf{T}}\mathbf{X}+\lambda\mathbf{I})^{-1}\mathbf{X}^{\mathsf{T}}\right) = \mathrm{Trace}(\mathbf{U}\mathbf{D}_\lambda\mathbf{U}^{\mathsf{T}})
$$

$$
= \mathrm{Trace}(\mathbf{D}_\lambda\mathbf{U}^{\mathsf{T}}\mathbf{U}) = \mathrm{Trace}(\mathbf{D}_\lambda) = \sum_{j=1}^{p}\frac{d_j^2}{d_j^2+\lambda}
$$

decreases from p to 0 as λ increases from 0 to ∞.

Ridge regression equivalent form

$$\hat{\beta}^\lambda = \underset{\beta}{\operatorname{argmin}} \ \|\mathbf{y} - \mathbf{X}\beta\|_2^2 + \lambda \|\beta\|_2^2,$$

$$\hat{\beta}^t = \underset{\|\beta\|_2 \leq t}{\operatorname{argmin}} \ \|\mathbf{y} - \mathbf{X}\beta\|_2^2.$$

One can prove that there exists a bijection between λ and t.

Exercise

Find the bijection between two positive scalars λ and t if two vectors β^λ and β^t are the same.

The sum of squared errors can be rewritten as the $\mathbf{D}$-weighted L_2 distance to $\hat{\beta}^{OLS}$ in the transformed coordinate system

$$\|\mathbf{y} - \mathbf{X}\beta\|_2^2$$
$$= \left\|\mathbf{U}\mathbf{U}^{\mathsf{T}}\mathbf{y} - \mathbf{U}\mathbf{D}\mathbf{V}^{\mathsf{T}}\beta\right\|_2^2 = \left\|\mathbf{U}^{\mathsf{T}}\mathbf{y} - \mathbf{D}\mathbf{V}^{\mathsf{T}}\beta\right\|_2^2$$
$$= \|\widetilde{\mathbf{y}} - \mathbf{D}\widetilde{\beta}\|_2^2 = \|\mathbf{D}(\widetilde{\beta}^{OLS} - \widetilde{\beta})\|_2^2$$

For the constraint $\|\beta\|_2 = \|\mathbf{V}^{\mathsf{T}}\beta\| = \|\widetilde{\beta}\| \leq t$, one can show the inequality problem $\min_{\|\widetilde{\beta}\|_2^2 \leq t} \|\widetilde{\mathbf{y}} - \mathbf{D}\widetilde{\beta}\|_2^2$ actually attains the equality constraint for this case. The Lagrangian function is $\|\widetilde{\mathbf{y}} - \mathbf{D}\widetilde{\beta}\|_2^2 + \mu(\|\widetilde{\beta}\|_2^2 - t)$ So, KKT gives $\mathbf{D}^2\widetilde{\beta} - \mathbf{D}\widetilde{\mathbf{y}} + \mu\widetilde{\beta} = 0$, i.e., $\widetilde{\beta} = (\mathbf{D}^2 + \mu\mathbf{I})^{-1}\mathbf{D}\widetilde{\mathbf{y}}$. So, μ is the same as λ. The equality constratin $\|\widetilde{\beta}\| = t$ determines uniquely $\lambda = \mu = \mu(t)$.

The solution path of ridge regression
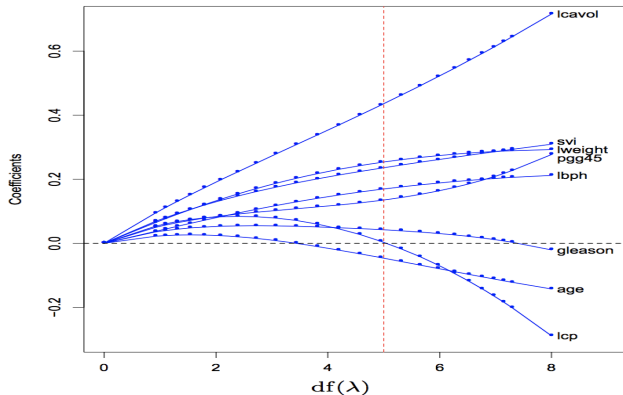
FIGURE 3.8. *Profiles of ridge coefficients for the prostate cancer example, as the tuning parameter λ is varied. Coefficients are plotted versus* $df(\lambda)$*, the effective degrees of freedom. A vertical line is drawn at* $df = 5.0$*, the value chosen by cross-validation.*

[ESL].

LASSO

(Frank and Friedman, 1993)
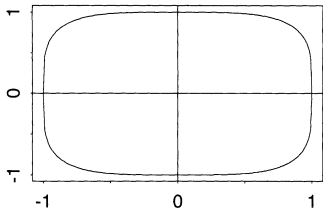With $L_r(\beta) = \sum_{j=0}^{p} |\beta_j|^r$,

$$\hat{\beta}^{bridge} = \underset{\beta}{\operatorname{argmin}} \|\mathbf{y} - \mathbf{X}\beta\|^2 + \lambda L_r(\beta)$$

- $L_0(\beta) = \sum_{j=0}^{p} I(\beta_j \neq 0)$; (Hard thresholding) [1]
- $L_1(\beta) = \sum_{j=0}^{p} |\beta_j|$; (Lasso)
- $L_2(\beta) = \sum_{j=0}^{p} \beta_j^2$; (Ridge regression)
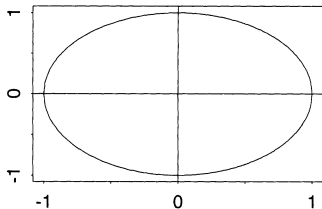- $L_\infty(\beta) = \max_j \ |\beta_j|$.

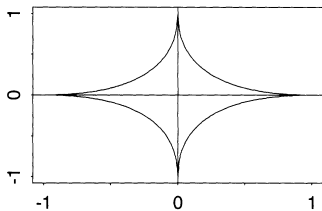[1] This gives subset selection with $L_0(\beta)$ number of variables selected.
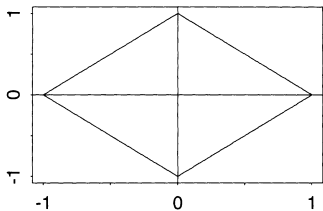
Figure 1.   Constrained Areas of Bridge Regressions with t = 1.

Least Absolute Shrinkage and Selection Operator (Lasso)

Tibshirani (Journal of the Royal Statistical Society 1996) introduced the LASSO.

Lasso estimator: let $r = 1$ in bridge estimator

$$\hat{\beta}^\lambda = \operatorname*{argmin}_{\beta} \left\{ \|\mathbf{y} - \mathbf{X}\beta\|_2^2 + \lambda\|\beta\|_1 \right\},$$

$$\hat{\beta}^s = \operatorname*{argmin}_{\|\beta\|_1 \leq s} \|\mathbf{y} - \mathbf{X}\beta\|_2^2,$$

where $\|\beta\|_1 = \sum_{j=1}^{p} |\beta_j|$.
Again there exists a bijection between λ and s.

Sparse solution

- Due to the nature of the l_1-norm constraint, if s is small enough some coefficients of the lasso solution become *exactly* zero.

- The elliptical contour is likely to hit the corner of the polytope, corresponding to sparse $\hat{\beta}$. The l_1 penalty can perform *variable selection*: drop the features with $\hat{\beta}_j = 0$.

- The l_r regularization results in sparsity when $0 \leq r \leq 1$, and is convex when $1 \leq r < \infty$.

- **Lasso is sparse and convex**:[1] the original implementation involves quadratic programming techniques from convex optimization

- Efron et al. (Annals of Statistics 2004) proposed **LARS (least angle regression)**, which computes the LASSO path efficiently
  - Interesting modification called is called forward stagewise
  - In many cases it is the same as the LASSO solution
  - Forward stagewise is easy to implement

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[1] not strictly convex. At a fixed λ, the coefficient $\hat{\beta}$ exists but may not be unique but the prediction $\hat{\mathbf{y}} = \mathbf{X}\hat{\beta}$ is unique (Tibshirani, R. J. (2013). The lasso problem and uniqueness. Electronic Journal of Statistics, 7, 1456– 1490.).
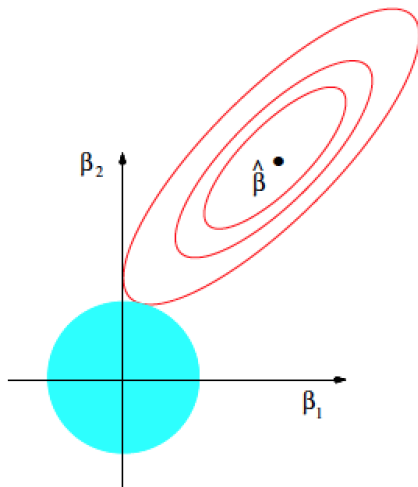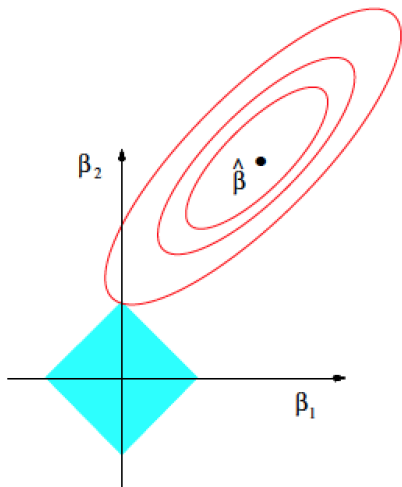
$\hat{\beta}$ in the center is the $\hat{\beta}^{OLS}$

Ridge vs Lasso

Consider a simple but illuminating example:

Exercise

Show that the solutions

$$\hat{\beta}^{ridge} = \underset{\beta}{\operatorname{argmin}} \left\{ \|\beta - \hat{\beta}\|_2^2/2 + \lambda\|\beta\|_2^2 \right\},$$

$$\hat{\beta}^{lasso} = \underset{\beta}{\operatorname{argmin}} \left\{ \|\beta - \hat{\beta}\|_2^2/2 + \lambda\|\beta\|_1 \right\}$$

are

- $\hat{\beta}_j^{ridge} = \hat{\beta}_j/(1 + 2\lambda)$
- $\hat{\beta}_j^{lasso} = \operatorname{sign}(\hat{\beta}_j) \left(|\hat{\beta}_j| - \lambda \right)_+ = \left(1 - \frac{\lambda}{|\hat{\beta}_j|} \right)_+ \hat{\beta}_j$

Visualization of $\hat{\beta}^{ridge}$ and $\hat{\beta}^{lasso}$ from desmos.com webpage
Ridge regression shrinks the $\hat{\beta}$ in all components/directions. Lasso translates them towards zero by a constant, truncating at λ.

Find Lasso solution in the above simple case.

Let $L(\beta) = (\beta - \hat{\beta})^2/2 + \lambda |\beta|$. The derivative w.r.t. β is

$L'(\beta) = \beta - \hat{\beta} + \lambda \operatorname{sign} \beta$.

1. $\beta > 0$, then $L'(\beta) = \beta - (\hat{\beta} - \lambda)$.

   1. If $\hat{\beta} - \lambda > 0$, the optimal β^* is indeed $\hat{\beta} - \lambda > 0$ where $L'(\beta^*) = 0$;
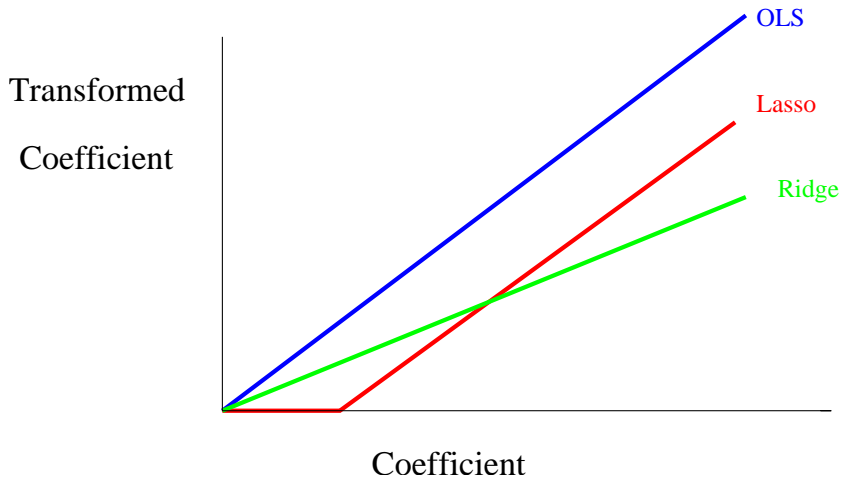   2. If $\hat{\beta} - \lambda < 0$, then $L'(\beta) > \beta > 0$; L is the increasing function in β. The minimal point of L in $[0, +\infty)$ is $\beta^* = 0$.

2. $\beta < 0$, then $L'(\beta) = \beta - (\hat{\beta} + \lambda)$.

   1. If $\hat{\beta} + \lambda < 0$, the optimal β^* is indeed $\hat{\beta} + \lambda > 0$ where $L'(\beta^*) = 0$;
   2. If $\hat{\beta} + \lambda > 0$, then $L'(\beta) < \beta < 0$; L is the decreasing function in β. The minimal point of L in $(-\infty, 0]$ is $\beta^* = 0$.

Transformed Coefficient (vertical axis)

Coefficient (horizontal axis)

OLS

Lasso

Ridge

Maximum number of selected covariates

The number of parameter/covariates selected by the lasso regression estimator is bounded non-trivially. The cardinality (i.e. the number of included covariates) of every lasso estimated linear regression model is smaller than or equal to $\min\{n, p\}$.

For any $\lambda > 0$, $\hat{\beta}^{lasso}(\lambda)$ has at most $\min(n, p)$ non-zeros entries. Here p is the number of parameters, and n is the number of training samples. [1]

The number of all possible sub-models from variable selection (such as subset selection) is 2^p. However, the typical number of different Lasso-estimated sub-models is at the order of

$$O(\min(n, p))$$

which is significantly smaller than 2^p if $p \gg n$.

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[1]Osborne, M. R., Presnell, B., and Turlach, B. A. (2000). On the lasso and its dual. Journal of Computational and Graphical Statistics, 9(2), 319–337.

sparsity of the lasso solution

- A large λ helps increase the sparsity [1].
- However, $\left\| \hat{\beta}^{lasso}(\lambda) \right\|_1$ decreases monotonically as λ increases, but the number of non-zero coefficients does not.

source of figure

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[1]Theory: blog article of Terence Tao
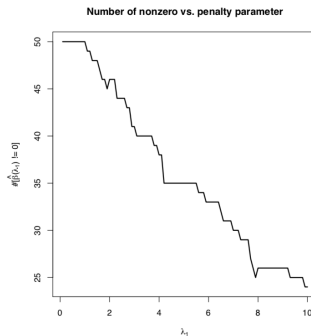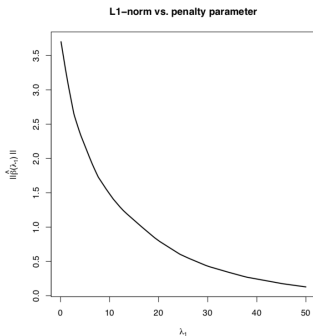
Solution path of LASSO

- The solution $\hat{\beta}^{lasso}(\lambda)$ is a piecewise linear function of λ. [1] See Figure 3.10 in [ESL].
- The path algorithm starts at $\lambda = \infty$ or $s = 0$, and traces the solution path by continuously changing λ [2]. Each new λ-solution is computed successively by solving the KKT conditions with a good initial guess set as the precise solution for a neighboring old λ. The key is to find the turning knots $\lambda_1, \ldots, \lambda_T$. An interesting reading: Efron et al. (AOS; 2003)
- Read the reference Regularization Paths for Generalized Linear Models via Coordinate Descent
- Path algorithms are available for many methods, such as fused lasso, trend filtering, locally adaptive regression splines, SVMs, 1-norm SVMs, relaxed maximum entropy method ...
- One can compute all possible Lasso sub-models $\cup_{\lambda > 0} \left\{ j : \hat{\beta}^{lasso}(\lambda) \neq 0 \right\}$ with $O(np\min\{n, p\})$ operation counts.

[1] Rosset, S. and Zhu, J. (2007). The Annals of Statistics, pp 1012–1030.
[2] homotopy method

https://scikit-learn.org/stable/auto_examples/linear_model/
plot_lasso_coordinate_descent_path.html

Consistence of variable selection

- Empirical fact: The set of variables selected by Lasso contains the true (non-zero) variables in the sparse model, with high chance:

$$\left\{ j : \hat{\beta}_j^{lasso} \neq 0 \right\} \supseteq \left\{ j : \hat{\beta}_j^{true} \neq 0 \right\}$$

- The Theory(Meinshausen and Buhlmann 2006): Under some (restrictive and crucial) assumptions, if $\lambda = \lambda_n \gg \sqrt{\log(p)/n}$, then

$$\left\{ j : \hat{\beta}_j^{lasso} \neq 0 \right\} \rightarrow \left\{ j : \hat{\beta}_j^{true} \neq 0 \right\}, \quad p \gg n \rightarrow \infty$$

with prob. 1.

- For high-dimensional model selection, with strongly correlated design $\mathbf{X}$, the Lasso can perform very poorly for variable selection.

The motivation for the Lasso came from an interesting proposal of Breiman (1993). Breiman's non-negative Garotte minimizes

$$\min_{c_j \geq 0, \forall j} \frac{1}{2} \sum_{i=1}^{n} \left(y_i - \sum_{j=1}^{p} c_j \hat{\beta}_j x_{ij} \right)^2 + \lambda \sum_{j=1}^{p} c_j$$

and then $\hat{\beta}_j^{ng} = \hat{c}_j \hat{\beta}_j$.

- It shrinks small $|\hat{\beta}_j|$ to zero. It is almost unbiased for large $|\hat{\beta}_j|$.
- Garotte starts with the OLS estimates and shrinks them by non-negative factors whose sum is constrained.
- In contrast, Lasso avoids the explicit use of the OLS estimates.
- Lasso is also closely related to the wavelet soft-thresholding method by Donoho and Johnstone (1994), and boosting method.

Exercise

Show when $\mathbf{X}^T\mathbf{X} = n\mathbf{I}$, the non-negative Garotte estimator is

$$\hat{\beta}_j^{ng} = \left(1 - \frac{\lambda}{2\hat{\beta}_j^2}\right)_+ \hat{\beta}_j.$$

Compare this with

$$\hat{\beta}_j^{lasso} = \text{sign}(\hat{\beta}_j)(|\hat{\beta}_j| - \lambda)_+ = \left(1 - \frac{\lambda}{\left|\hat{\beta}_j\right|}\right)_+ \hat{\beta}_j.$$

One can just focus on the component-wise formulation

$$\min_{\beta_j} \frac{1}{2}(z_j - \beta_j)^2 + J(|\beta_j|)$$

- Hard thresholding: $J(|\beta_j|) = \lambda^2 - (|\beta_j| - \lambda)^2 I(|\beta_j| < \lambda)$, then the optimal solution [1]

$$\hat{\beta}_j = z_j I(|z_j| > \lambda).$$

- Soft thresholding (Lasso): $J(|\beta_j|) = \lambda|\beta_j|$, then

$$\hat{\beta}_j = \text{sign}(z_j)(|z_j| - \lambda)_+.$$

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[1] the proof is exercise. Note J here is not continuous

Ridge regression vs LASSO

- Both can yield a reduction in variance at the expense of a small increase in bias.
- Bayesian Interpretation for Ridge Regression[1] and the Lasso: Figure 6.11, [ISL]. The prior distribution for β in ridge regression is Gaussian distribution; the prior distribution for β in LASSO is Laplace distribution $p(x) = \exp(-|x|/\lambda)/2\lambda$.
- Unlike ridge regression, the lasso performs variable selection when estimating the coefficients, and hence results in models easier to interpret.
- Ridge regression tends to give similar coefficient values to *correlated variables*, whereas the lasso may give quite different coefficient values to correlated variables.

Exercise
Ex. 7, Ex. 5, in Section 6.8, [ISL]

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[1]Ex. 3.6, Ex.3.7 in [ESL]

The Grouped Lasso

- In some problems, the predictors belong to pre-defined groups: genes that belong to the same biological pathway;
- Want to shrink and select the members of a group together: an entire group of predictors may drop out of the model.
- Suppose that the p predictors are divided into L groups[1]. Group ℓ has p_ℓ number of predictors, with design matrix $\mathbf{X}_\ell$ and the coefficient $\beta_\ell \in \mathbb{R}^{p_\ell}$. The grouped-lasso minimizes the convex criterion:

$$\min_{\beta \in \mathbb{R}^p} \left(\left\| \mathbf{y} - \beta_0 \mathbf{1} - \sum_{\ell=1}^{L} \mathbf{X}_\ell \beta_\ell \right\|_2^2 + \lambda \sum_{\ell=1}^{L} \sqrt{p_\ell} \left\| \beta_\ell \right\|_2 \right).$$

  The l_2 norm is applied since l_1-norm can not distinguish groups: $\|\beta\|_1 = \sum_l \|\beta_l\|_1$.
- This procedure encourages sparsity at both the group and individual levels.

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[1]This grouping is the known prior knowledge, required to apply the group lasso.

elastic-net penalty

Zou and Hastie (2005) introduced the elastic-net penalty:

$$\lambda \sum_{j=1}^{p} \left(\alpha \beta_j^2 + (1 - \alpha) \left| \beta_j \right| \right)$$

a different compromise between ridge and lasso from using $\left\| \cdot \right\|_r$ norm.

The elastic-net selects variables like the lasso, and shrinks together the coefficients of correlated predictors like ridge. The elastic-net regression is good at dealing with situations when there are correlations between parameters while Lasso regression tends to pick up just one of the correlated terms and eliminates the others.

Talk of Hui Zou on elastic-net

Exercise

Prove that the minimal point of

$$\min_\beta (\beta - \hat{\beta})^2/2 + \lambda_2 \beta^2 + \lambda_1 |\beta|$$

is $sign(\hat{\beta}) \left(\left| \hat{\beta} \right| - \lambda_1 \right)_+ \frac{1}{1+2\lambda_2} = \frac{1}{1+2\lambda_2} \hat{\beta}^{lasso}$ which is effectively the ridge coefficient applied to $\hat{\beta}^{lasso}$.

Exercise

Consider only two (highly correlated) samples $(x_{11}, x_{12}, y_1) = (1, 1, 1)$ and $(x_{21}, x_{22}, y_2) = (-1, -1, -1)$. Find the solutions of ridge, lasso, elastic-net regressions, respectively.
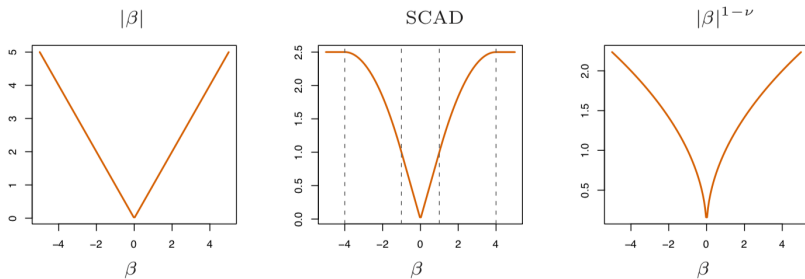
SCAD

One can modify the lasso penalty function so that larger coefficients are shrunken less severely (this may help reduce the unnecessary bias); the smoothly clipped absolute deviation (SCAD) penalty of Fan and Li (2005):

$$\min_{\beta} \ \frac{1}{2}\|\mathbf{y} - \mathbf{X}\beta\|^2 + \sum_{j=1}^{p} q_\lambda(|\beta_j|),$$

where for some $a \geq 2$, such that the derivative $q_\lambda(|\beta|)$ is

$$\frac{dq_\lambda(\beta)}{d\beta} = \lambda \cdot \text{sign}(\beta) \left[I(|\beta| \leq \lambda) + \frac{(a\lambda - |\beta|)_+}{(a-1)\lambda} I(|\beta| > \lambda) \right]$$

- The derivatives decreases from λ to 0 as $|\beta|$ increases.
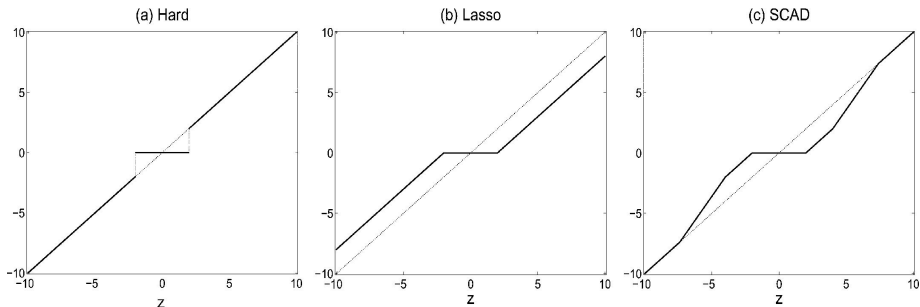- $q_\lambda(\beta)$ is not convex, a drawback for computation.

FIGURE 3.20. *The lasso and two alternative non-convex penalties designed to penalize large coefficients less. For SCAD we use $\lambda = 1$ and $a = 4$, and $\nu = \frac{1}{2}$ in the last panel.*

Figure source : [ESL]

Reference: Antoniadis and Fan (2001); Fan and Li (2001)

$\beta \to \hat{\beta}$

- All of them can select variables by setting some small coefficient to exactly zero. Hard thresholding is NP hard in computation.
- Lasso decreases the large coefficient uniformly by a gap λ which may induce unnecessary bias.
- SCAD leaves those coefficients with large value unchanged (thus the bias reduced) while still keeping the sparsity of the coefficients by thresholding for $|\beta| < \lambda$: a simple interpolation between hard and soft thresholding.

Adaptive Lasso (Zou, 2006; JASA)

Adaptive Lasso is an important two-stage procedure to address some bias problems of the Lasso.

$$\min_{\beta} \ \frac{1}{2}\|\mathbf{y} - \mathbf{X}\beta\|^2 + \lambda \sum_{j=1}^{p} w_j|\beta_j|,$$

where $w_j > 0$ and adjusts the penalty on each β_j.

- The resulting estimator is

$$\beta_j = \text{sign}(\hat{\beta}_j)(|\hat{\beta}_j| - \lambda w_j)_+.$$

- A reasonable choice for w_j is obtained from the first stage:

$$w_j = |\hat{\beta}_j^{OLS}|^{-\nu}, \text{ or } |\hat{\beta}_j^{lasso}|^{-\nu}$$

  with $\nu > 0$

- If $|w_j|$ is small (i.e., $\hat{\beta}_j$ from the first stage is large), the adaptive Lasso employs a small regularization (small adjustment) for the coefficient β_j, which implies less bias.

What not covered here

- Optimization theory and computational method for LASSO: Least Angle Regression, Dantzig Selector
- Principal Components Regression, partial linear regression (Section 3.5 in [ESL])
- Further reading on LASSO: *Statistics for High-Dimensional Data: Methods, Theory and Applications*, by Peter Buhlmann and Sara van de Geer, Springer Series in Statistics (2011)

Exercise

Ex. 3.28, 3.29, 3.30 in [ESL].

Revisit Bias-Variance

Revisit Bias-Variance

- In the beginning, We calculated the MSE for ordinary least square where the bias vanishes.
- We now focus on the ridge regression, $\hat{\beta}^\lambda = (\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}\mathbf{X}\mathbf{y}$, where $\lambda > 0$ is the penalty parameter. The corresponding MSE for $\hat{\beta}^\lambda$ at the new testing point x_0 then is written as [1]

$$
\begin{aligned}
\mathsf{MSE}^\lambda(x_0) &= \mathsf{Var}_{\mathbf{y}}(x_0^\mathsf{T}\hat{\beta}^\lambda) + \left(\mathbb{E}_{\mathbf{y}}(x_0^\mathsf{T}\hat{\beta}^\lambda) - f(x_0)\right)^2 \\
&= \mathsf{Var}_{\mathbf{y}}(x_0^\mathsf{T}(\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}\mathbf{X}^\mathsf{T}\mathbf{y}) + \left(x_0^\mathsf{T}\,\mathbb{E}_{\mathbf{y}}(\hat{\beta}^\lambda) - f(x_0)\right)^2 \\
&= \sigma_\varepsilon^2\,\|\mathbf{h}_\lambda x_0\|_2^2 + \left(x_0^\mathsf{T}(\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}\mathbf{X}^\mathsf{T}\mathbf{X}\beta - x_0^\mathsf{T}\beta\right)^2
\end{aligned}
$$

where $\mathbf{h}_\lambda = \mathbf{X}(\mathbf{X}^\mathsf{T}\mathbf{X} + \lambda\mathbf{I})^{-1}$. Exercise: with the aid of SVD, find the minimum point λ for MSE^λ.

---

[1] again, the training data's $\mathbf{X}$ part is fixed or conditioned.

Decomposition of average squared bias

- Note that we have <u>assumed</u> that the ground truth is a linear model $f(x) = x^{\mathsf{T}}\beta$.
- From now we do not assume f is linear and it could be a nonlinear function for a general consideration. This is a more realistic setting.
- The additive error model $Y = f(X) + \varepsilon$ is still assumed for data.
- The best-fitting approximation in the *linear model class* [1] is given by

$$\beta^* = \underset{\beta}{\operatorname{argmin}} \, \mathbb{E}_X(f(X) - X^{\mathsf{T}}\beta)^2$$

- Note that β_* satisfies the normal equation:

$$\mathbb{E}(X^{\mathsf{T}}X)\beta_* = \mathbb{E}[f(X)X]$$

---

[1] In other words, $x \to x^{\mathsf{T}}\beta^*$ is $f_{\mathcal{H}}(x)$ defined in Topic 1 ($\mathcal{H}$ is hypothesis space)

Decomposition of average (squared) bias

- We still consider the MSE$^\lambda$ for the ridge regression. The variance part Var$(x_0^\mathsf{T} \hat{\beta}^\lambda)$ is unchanged. Now the squared bias becomes
$\left(x_0^\mathsf{T} \mathbb{E}_{\mathbf{y}}(\hat{\beta}^\lambda) - f(x_0)\right)^2 = \left(x_0^\mathsf{T} \mathbb{E}_{\mathbf{y}}(\hat{\beta}^\lambda) - x_0^\mathsf{T} \beta_* + x_0^\mathsf{T} \beta_* - f(x_0)\right)^2 =$
$\left(x_0^\mathsf{T} \mathbb{E}_{\mathbf{y}}(\hat{\beta}^\lambda) - x_0^\mathsf{T} \beta_*\right)^2 + \left(x_0^\mathsf{T} \beta_* - f(x_0)\right)^2 +$
$2\left(x_0^\mathsf{T} \mathbb{E}_{\mathbf{y}}(\hat{\beta}^\lambda) - x_0^\mathsf{T} \beta_*\right)\left(x_0^\mathsf{T} \beta_* - f(x_0)\right)$

- Taking expectation for $x_0 \sim X$ and using the normal equation, we have the average squared bias is

$$\mathbb{E}_{x_0} \left(x_0^\mathsf{T} \mathbb{E}_{\mathbf{y}}(\hat{\beta}^\lambda) - f(x_0)\right)^2$$
$$= \underbrace{\mathbb{E}_{x_0} \left(x_0^\mathsf{T} (\mathbb{E}_{\mathbf{y}} \hat{\beta}^\lambda - \beta_*)\right)^2}_{\text{Ave (Estimation Bias}^2)} + \underbrace{\mathbb{E}_{x_0} \left(x_0^\mathsf{T} \beta_* - f(x_0)\right)^2}_{\text{Ave (Model Bias}^2)}$$

This is equation (7.14) in [ELS].

Remarks

- Model Bias, by definition, involves the ground truth, which is not accessible; Estimation Bias, by definition, involves β^*, which is not accessible either.

- However, this decomposition is conceptually inspiriting to have Model Bias.

- The Model Bias corresponds to the approximation error in Topic 1 ($f_{\mathcal{H}}$ vs f^*) and the Estimation Bias is similar to the sampling error in Topic 1 ($\hat{f}_{D}$ vs $f_{\mathcal{H}}$). In approximation theory viewpoint, the variance $\mathrm{Var}_{D} \, \hat{f}_{D}$ is not considered.

- Ridge method and LASSO further restricts the linear model $\mathcal{H}_{linear}$ to a smaller set in the form $\mathcal{H}_{\lambda} := \left\{ f : f(x) = \beta^{\mathsf{T}}x, \|\beta\| \leq \lambda \right\}$; they affect the estimation bias (and the variance of the prediction). The improvement of Model Bias needs go from linear to nonlinear if the ground truth is far away from being linear.