Problem 1 Penalized likelihood and soft thresholding

(A) --Prove that the quadratic term in the objective $S_{\lambda}(y) = \arg\min_{\theta} \frac{1}{2} (y - \theta)^2 + \lambda |\theta|$ is the negative

likelihood of a Gaussian distribution with mean θ and variance 1.

Since the Gaussian $(\theta, 1)$ distribution had the following probability density function:

$$f(y \mid \theta, 1) = (2\pi\sigma^2)^{-1/2} \exp[-\frac{(x-\theta)^2}{2\sigma^2}] = (2\pi)^{-1/2} \exp[-\frac{(x-\theta)^2}{2}]$$

The likelihood function is $L(\theta \mid y, 1) = (2\pi)^{-1/2} \exp[-\frac{(y-\theta)^2}{2}]$ (for a single y)

Then the log likelihood function is $l(\theta \mid y, 1) = \log[L(\theta \mid y, 1)] = -\frac{1}{2}\log(2\pi) - \frac{1}{2}(y - \theta)^2$, and the first term is not dependent on θ , so we drop it.

Therefore we have the negative log likelihood function $\frac{1}{2}(y-\theta)^2$ of the Gaussian $(\theta, 1)$,

which is the quadratic term of the objective minimization framework $S_{\lambda}(y)$.

--Prove that $S_{\lambda}(y) = sign(y) \cdot (|y| - \lambda)_{+}$, where $a_{+} = \max(a, 0)$ is the positive part of a.

Take derivative of $S_{\lambda}(y)$, we have

$$\frac{\partial S_{\lambda}(y)}{\partial \theta} = \frac{\partial}{\partial \theta} \left[\frac{1}{2} (y - \theta)^2 + \lambda |\theta| \right] = -(y - \theta) + \lambda \frac{|\theta|}{\theta} = -(y - \theta) + \lambda \cdot sign(\theta) \tag{*}$$

We will separate this problem into three parts: (1) $\theta > 0$; (2) $\theta < 0$; (3) $\theta = 0$.

(1) If $\theta > 0$, set function (*) equal to zero, which can be written as $-(y - \theta) + \lambda = 0$

$$\theta = v - \lambda > 0 \Rightarrow v > \lambda$$

(2) If $\theta < 0$, set function (*) equal to zero, which can be written as $-(y - \theta) - \lambda = 0$

$$\Rightarrow \theta = v + \lambda < 0 \Rightarrow v < -\lambda$$

(3) If $\theta = 0$, we need to employ the subdifferential or subgradient:

A vector g is a subgradient of a convex function f at $x \in dom f$ if

$$f(y) \ge f(x) + g^{T}(y - x) \quad \forall y \in dom f$$

The subdifferential $\partial f(x)$ (always a closed convex set) of f at x is the set of all subgradients:

$$\partial f(x) = \{ g \mid g^T(y - x) \le f(y) - f(x), \forall y \in dom f \}$$

Based on the above, we have $|\theta| \ge |0| + \tau \cdot (\theta - 0) \Longrightarrow |\theta| \ge \tau \theta \Longrightarrow -1 \le \tau \le 1$

So for
$$\theta = 0$$
, $\frac{\partial S_{\lambda}(y)}{\partial \theta} = -y + \lambda \tau = 0 \Rightarrow \tau = y / \lambda \in [-1,1]$

 \Rightarrow $y \in [-\lambda, \lambda]$, which means that $\theta = 0$ is optimal for $y \in [-\lambda, \lambda]$.

Therefore,
$$S_{\lambda}(y) = \arg\min_{\theta} \frac{1}{2} (y - \theta)^{2} + \lambda |\theta| = \begin{cases} y - \lambda, & y > \lambda \\ y + \lambda, & y < -\lambda \\ 0, & -\lambda \leq y \leq \lambda \end{cases}$$
$$= sign(y) \cdot (|y| - \lambda)_{+}$$

--Compare soft-thresholding with hard-thresholding:

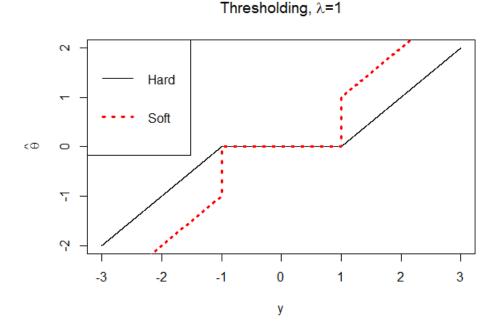


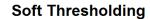
Figure 1: Comparison of soft thresholding and hard thresholding

(B) Toy examples:

- i) Plot $\hat{\theta}(y_i)$ versus θ_i across a discrete grid of different λ values (0, 1, 2, 3, 4, 5) and use 80% sparsity across n = 100 data points, where the remaining points are given $\theta = 1, \ldots, 20$. We observe that soft-thresholding function both selects certain θ_i 's by sparsifying the estimate, as well as shrinks the nonzero estimates towards 0 (Figure 2).
- ii) Plot the mean-squared error of the estimate as a function of λ :

$$MSE(\lambda) = \frac{1}{n} \sum_{i=1}^{n} [\hat{\theta}(y_i) - \theta_i]^2$$

Since each curve comes from a different randomly chosen vector θ and generated data, I've scaled each MSE to minimize to 1. The plot in Figure 3 shows how MSE changes with λ for different levels of sparsity in θ . Clearly, the λ that minimizes MSE increase as θ becomes more sparse. The bigger λ , the more aggressive the shrinkage affect.



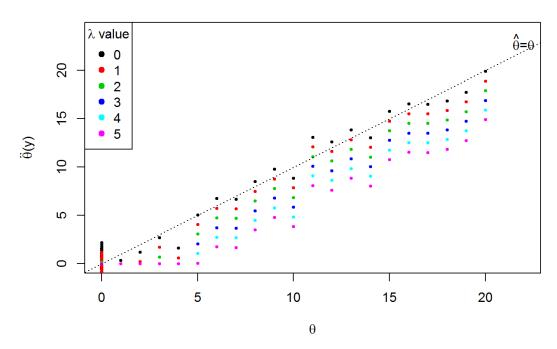


Figure 2: soft-thresholding function selects certain θ_i 's by sparsifying the estimate and shrinks the nonzero estimates towards 0

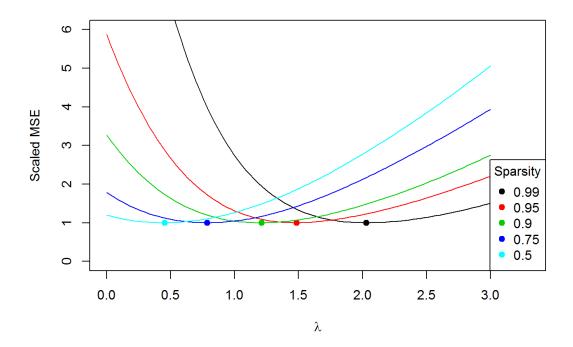


Figure 3: MSE changes with λ for different levels of sparsity in θ

Problem 2 The Lasso

Consider the standard linear regression model

$$y = X\beta + e$$

where y is an *n*-vector of responses, X is an $n \times p$ features matrix whose *i*th row x_i is the vector of features for observation *i*, and *e* is a vector of errors/residuals.

The Lasso involves estimating β as the solution to the penalized least squares problem

$$\hat{\beta} = \arg\min_{\beta} \frac{1}{2} \| y - X\beta \|_{2}^{2} + \lambda \| \beta \|_{1}$$

where $\|\beta\|_1$ is the L_1 norm of coefficient vector: $\|\beta\|_1 = \sum_{j=1}^p |\beta_j|$. When λ is very small, the

Lasso solution should be very close to the OLS solution, and all of the coefficients are in the model. As λ grows, the regularization term has greater effect and we will see fewer variables in your model (because more and more coefficients will be zero valued).

Leave the intercept in a Lasso fit unpenalized and write the objective as

$$\frac{1}{2n} \parallel y - (\alpha 1 + X\beta) \parallel_2^2 + \lambda \parallel \beta \parallel_1$$

where α is a scalar and 1 is a vector of all 1's.

(A) Fit the Lasso model across a range of λ values and plot the solution path $\hat{\beta}_{\lambda}$ as a function of λ . Also, track the in-sample mean-squared prediction error of the fit across the solution path.

$$MSE(\hat{\beta}_{\lambda}) = \frac{1}{n} \sum_{i=1}^{n} (y_i - x_i^T \hat{\beta}_{\lambda})^2 = \frac{1}{n} ||y - X \hat{\beta}_{\lambda}||_2^2$$

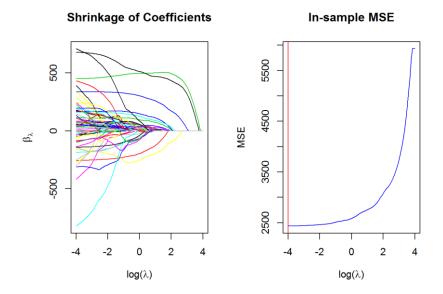


Figure 4: $\log(\lambda)$ v.s. $\hat{\beta}_{\lambda}$ and in-sample mean-squared prediction errors (MSE)

- The in-sample MSE decreases as λ goes to zero. In other words, in sample fit does not want shrinkage, it wants all 64 coefficients.
- (B) Randomly split the datasets into two equal parts (training and test datasets) and use 10-fold cross-validation across training data to find the best λ . *glmnet* packages have a *cv.glmnet* to choose best tuning parameters (in a chosen sequence of λ) which gives the minimum MSE. From the R output, the best $\lambda = 0.935$ with 35 parameters in the models (as shown in Figure 5). To avoid simply choosing part of λ (90 out of 100) in *cv.glmnet*, code up the 10-folds cross-validation across all λ 's. The resulting best $\lambda = 0.960$ with 34 parameters in the models and other parameters shrink to zero.

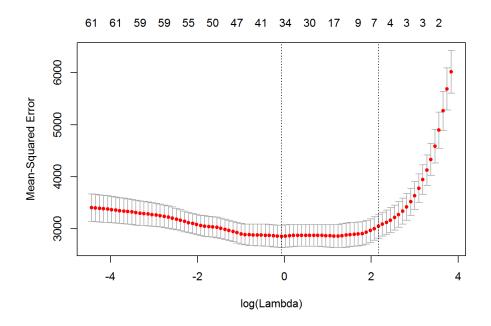


Figure 5: *cv.glmnet* chooses the best $\lambda = 0.935$ with 35 parameters (MSE =2808.36)

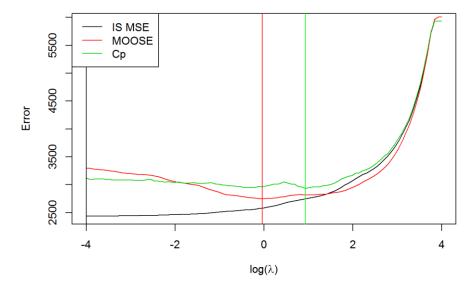


Figure 6: Comparison of MSE, MOOSE and C_p with respective best λ

(C) The C_p statistic (Mallow's C_p is defined as

$$C_p(\hat{\beta}_{\lambda}) = MSE(\hat{\beta}_{\lambda}) + 2 \cdot \frac{s_{\lambda}}{n} \hat{\sigma}^2$$

Where s_{λ} is the degrees of freedom of the fit (i.e. the number of nonzero parameters selected at that particular value of λ), and $\hat{\sigma}^2$ is an estimate of the residual variance. Figure 6 shows the comparison between MSE, MOOSE, and C_p . The C_p statistic closely follows the cross-validated MOOSE when λ increases. The minimum from C_p with best $\lambda = 2.533$, which is different from that of MOOSE with best $\lambda = 0.960$, implies 16 non-zero coefficients. The in-sample MSE clearly prefers over-fitting.