Fitting Generalized Lasso Models and Post-Selection Inference for the Lasso

Taeyoung Chang

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Fitting Generalized Lasso Models

Post-selection Inference for the Lasso

Taeyoung Chang

Notation

- x_i : p-dimensional vector for i-th observation of predictor variables
- \mathbf{x}_j : *n*-dimensional vector for *j*-th predictor of *n* observations
- $s_j \in \text{sign}(\beta_j)$: notation for sign function given as

$$s_j = egin{cases} 1 & eta_j > 0 \ -1 & eta_j < 0 \ c_j \in [-1,1] & eta_j = 0 \end{cases}$$

• $\mathbf{s} = (s_1, \cdots, s_p)$



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Motivation

- So far we have focused on the Lasso for squared-error loss, and exloited the piecewise-lilnearity of its coefficient profile to efficiently compute the entire path.
- Unfortunately this is not the case for most other loss functions.
 - Obtaining the coefficient path is potentially more costly.

Logistic regression example

- We will use logistic regression as an example.
- Use loss function L which is the negative log-likelihood.
- The problem is given as

$$\operatorname{minimize}_{\beta \in \mathbb{R}^p, \, \beta_0 \in \mathbb{R}} - \left\{ \frac{1}{n} \sum_{i=1}^n y_i \log \mu_i + (1 - y_i) \log (1 - \mu_i) \right\} + \lambda \|\beta\|_1$$

where $y_i \overset{indep}{\sim} \mathsf{Bern}(\mu_i)$ and $\mathsf{logit}(\mu_i) = \beta_0 + x_i^T \beta \quad \forall \ i = 1, \cdots, n$

The solution satisfies the subgradient condition

 As in the case of the lasso for squared-error loss, the solution satisfies the subgradient condition (KKT condition).

$$\frac{\partial}{\partial \beta} f(\beta, \beta_0) = \mathbf{0}$$
 and $\frac{\partial}{\partial \beta_0} f(\beta, \beta_0) = 0$

where $f(\beta, \beta_0)$ is the given objective function.

We shall taking advantage of

$$\frac{\partial}{\partial \beta}\mu_i = \mu_i(1-\mu_i)x_i$$
 and $\frac{\partial}{\partial \beta_0}\mu_i = \mu_i(1-\mu_i)$

Derivation of the subgradient condition

First condition

$$\frac{\partial}{\partial \beta} f(\beta, \beta_0) = \mathbf{0}$$

$$\Leftrightarrow \frac{\partial}{\partial \beta} - \frac{1}{n} \sum_{i=1}^n y_i \log \mu_i + (1 - y_i) \log(1 - \mu_i) + \lambda \|\beta\|_1 = \mathbf{0}$$

$$\Leftrightarrow -\frac{1}{n} \sum_{i=1}^n y_i (1 - \mu_i) x_i - (1 - y_i) \mu_i x_i + \lambda \mathbf{s} = \mathbf{0}$$

$$\Leftrightarrow -\frac{1}{n} \sum_{i=1}^n (y_i - \mu_i) x_i + \lambda \mathbf{s} = \mathbf{0}$$

$$\Leftrightarrow -\frac{1}{n} \langle \mathbf{x}_j, \mathbf{y} - \mu \rangle + \lambda s_j = \mathbf{0} \quad \forall j = 1, \dots, p$$

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Derivation of the subgradient condition

Second condition

$$\frac{\partial}{\partial \beta_0} f(\beta, \beta_0) = 0$$

$$\Leftrightarrow \frac{\partial}{\partial \beta_0} - \frac{1}{n} \sum_{i=1}^n y_i \log \mu_i + (1 - y_i) \log(1 - \mu_i) + \lambda \|\beta\|_1 = 0$$

$$\Leftrightarrow -\frac{1}{n} \sum_{i=1}^n y_i (1 - \mu_i) - (1 - y_i) \mu_i = 0$$

$$\Leftrightarrow -\frac{1}{n} \sum_{i=1}^n (y_i - \mu_i) = 0$$

$$\Leftrightarrow \frac{1}{n} \sum_{i=1}^n y_i = \sum_{i=1}^n \mu_i$$

Solution path on λ grid

- The nonlinearity of μ_i in β_j results in piecewise nonlinear coefficient profiles.
- \bullet Instead, we settle for a solution path on a sufficiently fine grid of values for λ
- ullet The largest value of λ we need to consider is

$$\lambda_{max} = \max_{j=1,\cdots p} |\langle \mathbf{x}_j, \mathbf{y} - \overline{y}\mathbf{1}\rangle|$$

• This is because it is the smallest value of λ for which $\hat{\beta}=0$ and $\hat{\beta}_0=\operatorname{logit}(\overline{y})$

Solution path on λ grid

- A reasonable sequence is 100 values $\lambda_1 > \lambda_2 > \cdots > \lambda_{100}$ equally spaced on the log-scale from λ_{max} down to $\varepsilon \lambda_{max}$ where ε is some small fraction such as 0.001
- An approach that has proven to be surprisingly efficient is path-wise coordinate descent.

Coordinate descent

For the problem

minimize
$$f(\mathbf{x})$$

with convex and differentiable function $f: \mathbb{R}^m \to \mathbb{R}$, coordinatewise minimization can yield a global minimization.

$$f(\mathbf{x}^* + \delta e_i) \ge f(\mathbf{x}^*) \quad \forall \ \delta > 0 \ , \ i = 1, \cdots, m \ \Rightarrow f(\mathbf{x}^*) = \min f(\mathbf{x})$$

where e_i is the *i*-th standard basis vector of \mathbb{R}^m

- Coordinate descent method is proceeded as the following :
 - **1** Take initial value $\mathbf{x}^{(0)} \in \mathbb{R}^m$
 - Iterate

$$x_i^{(k)} = \operatorname{argmin}_{x_i} f(x_1^k, \dots, x_{i-1}^{(k)}, x_i, x_{i+1}^{(k-1)}, \dots, x_m^{(k-1)}) \quad \forall \ i = 1, \dots, m$$

for step $k = 1, 2, \cdots$ and so on until convergence.

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Coordinate descent

- Coordinate descent example : linear regression
- minimize $\frac{1}{2}||y X\beta||_2^2$ over β_i with all $\beta_j \quad \forall j \neq i$ are fixed.
- Using $\frac{\partial \beta}{\partial \beta_i} = e_i$ where e_i is *i*-th standard basis of \mathbb{R}^p

$$\begin{split} \hat{\beta}_i \text{ minimizes } \frac{1}{2} \|y - X\beta\|_2^2 \text{ over } \beta_i \text{ with all } \beta_j \quad \forall \, j \neq i \text{ are fixed} \\ \Leftrightarrow \frac{\partial}{\partial \beta_i} \frac{1}{2} \|y - X\beta\|_2^2 = 0 \quad \text{at } \beta_i = \hat{\beta}_i \\ \Leftrightarrow \frac{\partial \beta}{\partial \beta_i} \frac{\partial}{\partial \beta} \frac{1}{2} \|y - X\beta\|_2^2 = 0 \quad \text{at } \beta_i = \hat{\beta}_i \\ \Leftrightarrow e_i^T (X^T X\beta - X^T y) = 0 \quad \text{at } \beta_i = \hat{\beta}_i \\ \Leftrightarrow \mathbf{x}_i^T (X\beta - y) = \mathbf{x}_i^T (X_i \beta_i + X_{-i} \beta_{-i} - y) = 0 \quad \text{at } \beta_i = \hat{\beta}_i \\ \Leftrightarrow \hat{\beta}_i = \frac{\mathbf{x}_i^T (y - X_{-i} \beta_{-i})}{\mathbf{x}_i^T \mathbf{x}_i} \end{split}$$

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Coordinate descent

- Coordinate descent example : the Lasso problem for squared-error loss
- minimize $\frac{1}{2}||y X\beta||_2^2 + \lambda ||\beta||_1$ over β_i with all $\beta_j \quad \forall j \neq i$ are fixed.
- ullet By similar logic we used for the linear regression case , solution \hat{eta}_i should satisfy

$$\hat{\beta}_i + \frac{\lambda}{\|\mathbf{x}_i\|_2^2} s_i = \frac{\mathbf{x}_i^T (y - X_{-i}\beta_{-i})}{\mathbf{x}_i^T \mathbf{x}_i}$$

• We have the solution $\hat{\beta}_i$ given as

$$\hat{\beta}_i = S_{\lambda/\|\mathbf{x}_i\|_2^2} \left(\frac{\mathbf{x}_i^T (y - X_{-i}\beta_{-i})}{\mathbf{x}_i^T \mathbf{x}_i} \right)$$

where $S_{\lambda}(x)$ is soft-thresholding defined as

$$S_{\lambda}(x) = \begin{cases} x - \lambda & \text{if } x > \lambda \\ 0 & \text{if } -\lambda \le x \le \lambda \\ x + \lambda & \text{if } x < \lambda \end{cases}$$

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Pathwise coordinate descent

- Outer loop
 - Find optimal value β for each λ_k in the order of $\lambda_1 > \lambda_2 > \cdots > \lambda_{100}$
 - By starting at λ_1 , where all parameters are zero, we use warm starts in computing the solutions at the decreasing sequence of λ values.
 - resulting β for λ_k is used as an initial value of coordinate descent algorithm for λ_{k+1}

Pathwise coordinate descent

Inner loop

- For each value λ_k , solve the lasso problem for one β_j only, holding the others fixed. This is done by coordinate descent. One or several coordinate cycles are implemented until the estimates stabilize.
- Store the nonzero coefficients in the active set \mathcal{A} . (The active set grows slowly as λ decreases.)
- Iterates coordinate descent using only those variables until convergence.
- One more sweep through all the variables to check optimality conditions. If there is a variable not satisfying the condition, then add it in active set $\mathcal A$ and go back to the first step of inner loop.

Comments

- The R package glmnet employs a 'proximal-Newton' strategy at each value λ_k , which takes advantage of a weighted least-squares and coordinate descent.
- We can consider another penalty term called as 'elastic net' penalty which bridges the gap between the lasso and ridge regression. It is defined as

$$P_{\alpha}(\beta) = \frac{1}{2} \{ (1 - \alpha) \|\beta\|_{2}^{2} + \alpha \|\beta\|_{1} \}$$

for some $\alpha \in [0,1]$



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Post-selection Inference

- Inference is generally difficult for adaptively selected models.
- Suppose we have fit a lasso regression model with a particular value for λ , which ends up selecting a subset \mathcal{A} of size $|\mathcal{A}|=k$ of p available variables.
- Question : interest in the population regression parameters using the full set of p predictors VS interest is restricted to the population regression parameters using only the subset $\mathcal A$

Post-selection Inference

- Focus on the second case
- ullet The idea is to condition on the selected set ${\mathcal A}$ itself, and then perform conditional inference on the unrestricted (not lasso-shrunk) regression coefficients of the response on only the variables in ${\mathcal A}$
- For the case of the lasso with squared-error loss, using the fact about convexity along with delicate Gaussian conditioning arguments, it leads to truncated Gaussian and t-distributions for parameters of interest.

Reference

Bradley Efron and Trevor Hastie. *Computer Age Statistical Inference:* Algorithms, Evidence, and Data Science. Institute of Mathematical Statistics Monographs. Cambridge University Press.

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- Convex optimization for All