# **MACHINE LEARNING IN BIOINFORMATICS** FEATURE SELECTION

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#### Feature selection problem

$$\hat{\theta} = \begin{cases} \underset{\theta}{\arg\min} & \|y - X\theta\|_2^2 \\ & \\ \text{subject to} & \|\theta\|_0 = m \end{cases} \quad \text{with } \binom{p}{m} \text{ possible subsets}$$

- Required are computationally efficient methods to approximate the feature selection problem
- Offline methods: Select features before estimating parameters
- Online methods: Features are selected during parameter estimation

#### FEATURE SELECTION METHODS

#### Offline methods:

- Safe and Strong rules
- Sure independence screening (SIS)
- Estimation of mutual information
- Online methods:
  - (Orthogonal) matching pursuit
  - Least angle regression (LARS) / Homotopy algorithm
  - Penalty methods

# LINEAR REGRESSION - RECAP

$$\begin{array}{c} y = X\theta + \epsilon \\ \begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{bmatrix} = \begin{bmatrix} x_1^{(1)} & x_1^{(2)} & \dots & x_1^{(p)} \\ x_2^{(1)} & x_2^{(2)} & \dots & x_2^{(p)} \\ \vdots & \vdots & \ddots & \vdots \\ x_n^{(1)} & x_n^{(2)} & \dots & x_n^{(p)} \end{bmatrix} \begin{bmatrix} \theta_1 \\ \theta_2 \\ \vdots \\ \theta_p \end{bmatrix} + \begin{bmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \vdots \\ \varepsilon_n \end{bmatrix}$$

response :	$y \in \mathbb{R}^n$
covariates :	$X \in \mathbb{R}^{n \times p}$
coefficients :	$\theta \in \mathbb{R}^p$
residuals :	$\epsilon \in \mathbb{R}^n$

Geometric interpretation of ordinary least squares [Hastie et al., 2009]:

$$\hat{\theta} = \arg\min_{\theta} \|\epsilon\|_{2}^{2}$$
$$= \arg\min_{\theta} \|y - X\theta\|_{2}^{2}$$



- Consider the case of ultrahigh-dimensional data, where the number of features p is much larger than the number of observations n
- Specifically, we assume that p is so large that we cannot compute an estimate of  $\theta$
- Assuming  $\theta$  is sparse, we can first select a promising subset of q features  $M_q$  (called feature screening)
- **The coefficients**  $\theta$  are estimated based on the subset  $M_q$

• Consider the solution of rigde regression:

$$\hat{\theta}(\lambda) = (X^{\top}X + \lambda I)^{-1}X^{\top}y$$

- $\blacksquare$  For  $\lambda \rightarrow$  0 we obtain the OLS solution
- For  $\lambda \to \infty$  it follows that  $\lambda \hat{\theta}(\lambda)$  converges to the componentwise regression estimator

$$\hat{\theta}_k(\lambda) = \tilde{X}^\top y$$

where  $\tilde{X}$  is the data matrix X with normalized columns  $f_j$  such that  $f_j^\top f_j = \mathbf{1}$ 

■ Traditionally, for very large *p* we would select  $\lambda$  large in order to decrease the variance of  $\hat{\theta}$ 

- $\tilde{X}^{\top}y = (f_1^{\top}y, \dots, f_p^{\top}y)$  can be interpreted as the correlation of features  $f_j$  with y
- Sure independence screening (SIS) [Fan and Lv, 2008] selects a subset of features

$$\Omega = \left\{ j \mid |f_j^\top y| > t \right\}$$
(1)

based on their correlation with y, where t is a threshold such that  $|\Omega| = q < p$ 

- The OLS estimate  $\hat{\theta}$  is computed using only the selected features  $\Omega$
- **All remaining components of**  $\hat{\theta}$  are set to zero

The same idea can be applied to more complex models [Fan and Song, 2010], such as logistic regression, where

$$\hat{\theta} = \arg\max_{\theta} \operatorname{pr}_{\theta}(y \,|\, X)$$

Select a subset of features

$$\Omega = \left\{ j \mid \text{score}(f_j, y) > t \right\}$$
(2)

The score is given by the independent estimate

$$\operatorname{score}(f_j, y) = \arg \max_{\theta_j} \operatorname{pr}_{\theta_j}(y \mid f_j)$$

for all j = 1, ..., p

#### Feature selection problem

$$\hat{\theta} = \begin{cases} \arg\min_{\theta} & \|y - X\theta\|_2^2 \\ \text{subject to} & \|\theta\|_0 = m \end{cases} \quad \text{with } \begin{pmatrix} p \\ m \end{pmatrix} \text{ possible subsets}$$

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#### **Matching Pursuit**

Greedy approximation to feature selection problem.

#### Feature selection problem

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If we must represent *y* with only one feature, which one should we take?

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#### **Matching Pursuit**

Greedy approximation to feature selection problem.

If we must represent *y* with only one feature, which one should we take?

$$j_1 = rgmin_j \left\| y - f_j \hat{ heta}_j \right\|_2^2$$
, where  $\hat{ heta}_j = rgmin_{ heta_j} \left\| y - f_j heta_j \right\|_2^2$ 

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$$\begin{split} & j_1 = \arg\min_j \left\| y - f_j \hat{\theta}_j \right\|_2^2, \quad \text{where} \quad \hat{\theta}_j = \arg\min_{\theta_j} \left\| y - f_j \theta_j \right\|_2^2 \\ & = \arg\max_j \frac{(f_j^\top y)^2}{f_j^\top f_j} \\ & = \arg\max_j \left| f_j^\top y \right| \end{split}$$

[assuming normalized data, i.e.  $f_j^{\top} f_j = 1$ ]

 $\Rightarrow$  select feature *j* with maximal scalar projection of *y* onto  $f_i$ 

$$\epsilon = y - X\theta$$

$$= \underbrace{y}_{r_0} - f_{j_1}\theta_{j_1} - f_{j_2}\theta_{j_2} - \dots - f_{j_p}\theta_{j_p}$$

$$\underbrace{y}_{r_1} - f_{r_2} - f_{r_1}\theta_{r_2} - \dots - f_{r_p}\theta_{r_p}$$

$$\begin{aligned} j_1 &= \arg\min_{j} \left\| y - f_j \hat{\theta}_j \right\|_2^2 &= \arg\min_{j} \left\| r_0 - f_j \hat{\theta}_j \right\|_2^2 \\ &= \arg\max_{j} \left| f_j^\top r_0 \right| \end{aligned}$$

$$\epsilon = \mathbf{y} - X\theta$$

$$= \underbrace{\mathbf{y} - f_{j_1}\theta_{j_1} - f_{j_2}\theta_{j_2} - \ldots - f_{j_p}\theta_{j_p}}_{r_1}$$

$$j_{1} = \arg\min_{j} \left\| y - f_{j}\hat{\theta}_{j} \right\|_{2}^{2} = \arg\min_{j} \left\| r_{0} - f_{j}\hat{\theta}_{j} \right\|_{2}^{2}$$

$$= \arg\max_{j} \left| f_{j}^{\top}r_{0} \right|$$

$$j_{2} = \arg\min_{j} \left\| y - f_{j_{1}}\hat{\theta}_{j_{1}} - f_{j}\hat{\theta}_{j} \right\|_{2}^{2} = \arg\min_{j} \left\| r_{1} - f_{j}\hat{\theta}_{j} \right\|_{2}^{2}$$

$$= \arg\max_{j} \left| f_{j}^{\top}r_{1} \right|$$

# Matching pusuit (MP) [Tropp et al., 2007]

The MP feature selection rule is given by

$$j_k = \arg \max_j \left| f_j^\top r_{k-1} \right| \qquad k = 1, \dots, m$$

where  $r_k$  are the residuals at step k:

$$\epsilon = \mathbf{y} - X\theta$$

$$= \underbrace{\mathbf{y}_{r_0}}_{r_0} - f_{j_1}\theta_{j_1} - f_{j_2}\theta_{j_2} - \dots - f_{j_p}\theta_{j_p}$$

$$\underbrace{\mathbf{y}_{r_0}}_{r_1} - f_{r_2}\theta_{r_2} - \dots - f_{r_p}\theta_{r_p}$$

#### Orthogonal Matching Pursuit

Orthogonal Matching Pursuit: Re-estimate parameters after every iteration.

After every iteration *t*, update all  $\theta_{\Omega_t}$  entries, where  $\Omega_t = \{j_1, j_2, \dots, j_t\}$ , i.e. compute

$$heta_{\Omega_t} = rgmin_{ heta} \|y_{\Omega_t} - X_{\Omega_t}\theta\|_2^2 \;.$$

This update changes the residuals

$$r_t = y - f_{j_1}\theta_{j_1} - f_{j_2}\theta_{j_2} - \cdots - f_{j_t}\theta_{j_t}$$

used in the next iteration of the algorithm.

#### LOGISTIC REGRESSION

$$\begin{bmatrix} \operatorname{pr}_{\theta}(y_{1} = 1) \\ \operatorname{pr}_{\theta}(y_{2} = 1) \\ \vdots \\ \operatorname{pr}_{\theta}(y_{n} = 1) \end{bmatrix} = \sigma \begin{pmatrix} \begin{bmatrix} x_{1}^{(1)} & x_{1}^{(2)} & \dots & x_{1}^{(p)} \\ x_{2}^{(1)} & x_{2}^{(2)} & \dots & x_{2}^{(p)} \\ \vdots & \vdots & \ddots & \vdots \\ x_{n}^{(1)} & x_{n}^{(2)} & \dots & x_{n}^{(p)} \end{bmatrix} \begin{bmatrix} \theta_{1} \\ \theta_{2} \\ \vdots \\ \theta_{p} \end{bmatrix} \end{pmatrix}$$

class labels : $y \in \{0,1\}^n$ covariates : $X \in \mathbb{R}^{n \times p}$ coefficients : $\theta \in \mathbb{R}^p$ 

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Parameter estimation for logistic regression:

$$\begin{split} \hat{\theta} &= \arg\max_{\theta} \operatorname{pr}_{\theta}(\mathbf{y}) \approx \arg\min_{\theta} \|\mathbf{y} - \sigma(\mathbf{X}\theta)\|_{2}^{2} \quad [\text{but not convex}] \\ &= \arg\max_{\theta} \sum_{i=1}^{n} \log \operatorname{pr}_{\theta}(\mathbf{y}_{i}) \\ &= \arg\max_{\theta} \sum_{i=1}^{n} \{\mathbf{y}_{i} \log \sigma(\mathbf{x}_{i}\theta) + (1 - \mathbf{y}_{i}) \log(-\mathbf{x}_{i}\theta)\} \\ &= \arg\max_{\theta} \sum_{i=1}^{n} \log \sigma(\tilde{\mathbf{y}}_{i}\mathbf{x}_{i}\theta) \,, \end{split}$$

where  $\tilde{y}_i = 2y_i - 1 \in \{-1, 1\}$ 

#### Pseudo-residuals

$$\begin{aligned} \mathbf{r}_k &= \mathbf{y} - \sigma(f_{j_1}\theta_{j_1} + f_{j_2}\theta_{j_2} + \dots + f_{j_k}\theta_{j_k}) \\ \mathbf{X}^\top \mathbf{r}_p &= \nabla \log \mathrm{pr}_{\theta}(\mathbf{y}) \end{aligned}$$

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$$\begin{split} r_k &= y - \sigma(f_{j_1}\theta_{j_1} + f_{j_2}\theta_{j_2} + \dots + f_{j_k}\theta_{j_k}) \\ X^\top r_p &= \nabla \log \operatorname{pr}_{\theta}(y) \end{split}$$

$$\begin{split} j_1 &= \arg\min_{j} \left\| \mathbf{y} - \sigma(f_j \hat{\theta}_j) \right\|_2^2 \\ &\approx \arg\max_{j} \left| f_j^\top \mathbf{r}_0 \right| \end{split}$$

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$$j_{1} = \arg\min_{j} \left\| y - \sigma(f_{j}\hat{\theta}_{j}) \right\|_{2}^{2}$$

$$\approx \arg\max_{j} \left| f_{j}^{\top} r_{0} \right|$$

$$j_{2} = \arg\min_{j} \left\| y - \sigma(f_{j_{1}}\hat{\theta}_{j_{1}} - f_{j}\hat{\theta}_{j}) \right\|_{2}^{2}$$

$$\approx \arg\max_{j} \left| f_{j}^{\top} r_{1} \right|$$

#### Matching pursuit feature selection rule [Lozano et al., 2011]

Assuming normalized data, i.e.  $f_j^ op f_j =$  1, the OMP rule is given by

$$j_k = \arg\max_j \left| f_j^\top r_{k-1} \right|$$

where  $r_k$  are the kth pseudo-residuals

$$\begin{aligned} \mathbf{r}_k &= \mathbf{y} - \sigma(f_{j_1}\theta_{j_1} + f_{j_2}\theta_{j_2} + \dots + f_{j_k}\theta_{j_k}) \\ \mathbf{X}^\top \mathbf{r}_p &= \nabla \log \mathrm{pr}_{\theta}(\mathbf{y}) \end{aligned}$$

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$$r_k = y - \sigma(f_{j_1}\theta_{j_1} + f_{j_2}\theta_{j_2} + \dots + f_{j_k}\theta_{j_k})$$
$$X^{\top}r_p = \nabla \log \operatorname{pr}_{\theta}(y)$$

#### **OMP** Performance

Greedy strategy causes poor performance of Orthogonal Matching Pursuit in practice

 $\blacksquare$  Consider  $\ell_1\text{-}\mathsf{penalized}$  linear regression (LASSO) where

$$\hat{\theta}(\lambda) = \arg\min_{\theta} \|y - X\theta\|_{2}^{2} + \lambda \|\theta\|_{1}$$

- There exists a regularization strength  $\lambda = \lambda_{max}$  for which all estimated coefficients are zero
- Least Angle Regression (LARS) [Efron et al., 2004] is a method to efficiently compute  $\hat{\theta}(\lambda)$  for all  $0 \le \lambda \le \lambda_{max}$
- LARS computes breakpoints  $\lambda_k$  at which individual coefficients  $\hat{\theta}_j(\lambda_k) \in \mathbb{R}$  change its value from
  - zero to non-zero, or from
  - non-zero to zero
- Between breakpoints the values of coefficients can be linearly interpolated

**Remember that the OLS solution**  $\hat{\theta}(\mathbf{0})$  for  $\lambda = \mathbf{0}$  requires that

$$abla_{ heta} \| \mathbf{y} - \mathbf{X} \mathbf{\theta} \|_2^2 = 2\mathbf{X}^{ op}(\mathbf{y} - \mathbf{X} \mathbf{\theta}) = \mathbf{0}$$

• For  $\lambda > 0$  the solution requires

$$X^{\top}(y - X\theta) \in \frac{\lambda}{2} \partial \|\theta\|_{1}$$

where  $\partial \|\theta\|_1$  is the subgradient with respect to  $\theta$ 

#### We define

$$\mathsf{c}(\theta) = \mathsf{X}^\top (\mathsf{y} - \mathsf{X}\theta)$$

which is interpreted as the correlation of features  $X = [f_1, f_2, \dots, f_p]$  with the residuals  $\epsilon = y - X\theta$ 

The correlation  $\hat{c}(\lambda) = c(\hat{\theta}(\lambda))$  varies with  $\lambda$  as follows:

• 
$$\hat{c}(\lambda) = c_{\max}$$
 for  $\lambda = \lambda_{\max}$ 

• 
$$\hat{c}(\lambda) = 0$$
 for  $\lambda = 0$ 



- LARS maintains a set of active features  $\Omega \subset \{1, ..., p\}$  all equally correlated with the residuals  $y X\hat{\theta}(\lambda)$  for the current estimate  $\hat{\theta}(\lambda)$
- Let  $X_{\Omega} = (f_j)_{j \in \Omega}$  denote the covariate matrix and  $\theta_{\Omega} = (\theta_j)_{j \in \Omega}$ the coefficients restricted to the features in the active set  $\Omega$
- $\blacksquare$  In each iteration, the coefficients  $\theta$  are updated

$$\theta \leftarrow \theta + \gamma^* \mathbf{V} \,,$$

where  $\gamma^*$  is the amount by which the correlation  $c_{\Omega}(\theta)$  is reduced and  $v \in \mathbb{R}^p$  defines the direction and relative size of the update

The vector  $\mathbf{v}$  is selected so that for features in  $\Omega$  the difference in correlation  $c_{\Omega}(\theta) - c_{\Omega}(\theta + \gamma \mathbf{v})$  shrinks uniformly towards zero with rate  $\gamma$ , i.e.

$$\begin{split} & c_{\Omega} \left( \theta \right) - c_{\Omega} \left( \theta + \gamma \mathbf{v} \right) = \gamma \operatorname{sign} c_{\Omega}(\theta) \,, \quad \text{while} \\ & c_{\Omega^{c}}(\theta) - c_{\Omega^{c}}(\theta + \gamma \mathbf{v}) = \mathrm{O} \,. \end{split}$$

Both conditions can be combined into

$$\mathbf{C}(\theta) - \mathbf{C}(\theta + \gamma \mathbf{V}) = \gamma \operatorname{sign} \mathbf{C}(\theta),$$

since sign  $c_{\Omega^c}(\theta) = 0$ 

It follows that

$$V_{\Omega} = [X_{\Omega}^{\top}X_{\Omega}]^{-1}\operatorname{sign} C_{\Omega}(\theta)$$

and  $v_{\Omega^c} = 0$ 

#### ■ LARS stop shrinking the correlations whenever:

- Case 1: A non-active feature becomes equally correlated with the residuals
- Case 2: A coefficient of an active feature becomes zero<sup>1</sup>
- Case 1: More formally,  $\gamma$  is increased until some feature  $j' \in \Omega^c$  outside the active group satisfies

$$\begin{aligned} |\mathbf{c}_{j'}(\theta + \gamma \mathbf{v})| &= |\mathbf{c}_{j}(\theta + \gamma \mathbf{v})| \\ &= \lambda - \gamma \,, \end{aligned}$$

where  $j \in \Omega$ , and  $\lambda = |c_j(\theta)|$  is the absolute correlation of the active features

<sup>1</sup>This case was not part of the initial LARS algorithm but was later on added in order to ensure equivalence with the LASSO (see also Homotopy algorithm [Osborne et al., 2000])

■ The solution is given by

$$\gamma^{+} = \min_{j \in \Omega^{c}}^{+} \left\{ \frac{\lambda - c_{j}(\theta)}{1 - f_{j}^{\top} X \mathbf{v}}, \frac{\lambda + c_{j}(\theta)}{1 + f_{j}^{\top} X \mathbf{v}} \right\} ,$$

where min<sup>+</sup> is the minimum over positive elements and note that  $f_j^\top X v = f_j^\top X_\Omega v_\Omega$ 

Case 2: The algorithm also removes a feature j from the active set when for some  $\gamma$ 

$$\theta_j + \gamma \mathbf{V}_j = \mathbf{O}$$

so that  $\gamma^- = \min_{j \in \Omega} \{-\theta_j / \mathbf{v}_j\}$ 

■ The subsequent breakpoint is given by  $\gamma^* = \min\{\gamma^+, \gamma^-\}$ 

# SAFE AND STRONG RULES

# Penalized regression

$$\begin{split} \omega(\theta) &= -\log \operatorname{pr}_{\theta}(\mathbf{y}) \\ \omega(\theta) &= \|\mathbf{y} - \mathbf{X}\theta\|_2^2 \end{split}$$

(logistic regression), or (linear regression)

$$\hat{\theta} = \begin{cases} \arg\min & \omega(\theta) \\ \\ \theta \\ \text{subject to} & \|\theta\|_{1} = N \end{cases}$$

Basic idea: Select 
$$\wedge$$
 such that  $\|\theta\|_{o} = m$ 

## Numerical solution of penalized regression

Identify saddle points of Lagrangian

$$\mathcal{L}(\theta, \lambda) = \omega(\theta) + \lambda(\|\theta\|_{1} - \Lambda)$$

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Identify saddle points of Lagrangian

$$\mathcal{L}(\theta, \lambda) = \omega(\theta) + \lambda(\|\theta\|_{1} - \Lambda)$$

In practice the constraint  $\|\theta\|_1 = \Lambda$  is ignored, but  $\lambda$  is chosen such that classification performance is optimal:

#### Penalized regression in practice

$$\hat{ heta}(\lambda) = \mathop{\arg\min}\limits_{ heta} \omega( heta) + \lambda \left\| heta 
ight\|_{1}$$

# SAFE RULE FOR LINEAR REGRESSION

SAFE rule: What features can we neglect for a fixed  $\lambda$ ?

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SAFE rule [Ghaoui et al., 2010, Kim et al., 2007] for  $l_1$ -penalized linear regression

*j*th component of  $\hat{\theta}$  must be zero if

$$\begin{split} \|f_j^\top \mathbf{y}\| &< \lambda - \left\|f_j\right\|_2 \|\mathbf{y}\|_2 \frac{\lambda_{\max} - \lambda}{\lambda_{\max}} \\ \lambda_{\max} &= \max_j |f_j^\top \mathbf{y}| \end{split}$$

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$$\begin{split} \|\boldsymbol{f}_{j}^{\top}\boldsymbol{y}\| &< \lambda - \left\|\boldsymbol{f}_{j}\right\|_{2} \|\boldsymbol{y}\|_{2} \, \frac{\lambda_{\max} - \lambda}{\lambda_{\max}} \\ \lambda_{\max} &= \max_{j} |\boldsymbol{f}_{j}^{\top}\boldsymbol{y}| \end{split}$$

$$|f_j^{\top}(y - \underbrace{X\theta}_{\theta = 0})| < \lambda - \left\|f_j\right\|_2 \left\|y\right\|_2 \frac{\lambda_{\max} - \lambda}{\lambda_{\max}}$$

#### STRONG RULE FOR LINEAR REGRESSION

SAFE rule for linear regression: *j*th component of  $\hat{\theta}$  must be zero if

$$\begin{split} |f_j^{\top} \mathbf{y}| < \lambda - \left\| f_j \right\|_2 \| \mathbf{y} \|_2 \, \frac{\lambda_{\max} - \lambda}{\lambda_{\max}} \\ \lambda_{\max} = \max_j |f_j^{\top} \mathbf{y}| \end{split}$$

Strong rule for  $\ell_1$ -penalized linear regression [Tibshirani et al., 2012]

Discard jth component if

$$egin{aligned} &|f_j^ op \mathbf{y}| < \lambda - (\lambda_{max} - \lambda) = \mathbf{2}\lambda - \lambda_{max} \ &\lambda_{max} = \max_j |f_j^ op \mathbf{y}| \end{aligned}$$

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#### Remark

Strong rule may drop features that should not be discarded  $\Rightarrow$  KKT conditions must be checked, i.e.

$$\boldsymbol{X}^{\top}(\boldsymbol{y} - \boldsymbol{X}\hat{\theta}) \in \lambda \partial_{\theta = \hat{\theta}} \left\| \boldsymbol{\theta} \right\|_{1}$$

#### STRONG SEQUENTIAL RULE FOR LINEAR REGRESSION

# Strong rule for $\ell_1$ -penalized linear regression [Tibshirani et al., 2012]

Discard jth component if

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# Strong sequential rule for $\ell_1$ -penalized linear regression[Tibshirani et al., 2012]

Discard *j*th feature if

$$|f_j^{ op}\{\mathbf{y} - \sigma(\mathbf{X}\hat{ heta}(\lambda_{k-1}))\}| < 2\lambda_k - \lambda_{k-1}$$

### STRONG SEQUENTIAL RULE FOR LINEAR REGRESSION

Compute  $\hat{\theta}(\lambda_k)$  for all  $\lambda_1 > \cdots > \lambda_k > \cdots > \lambda_K$ 



Assumption :  $|f_j^{\top}(y - X\hat{\theta}(\lambda_{k-1} - \epsilon)) - f_j^{\top}(y - X\hat{\theta}(\lambda_{k-1}))| \le \epsilon$  $\Rightarrow |f_j^{\top}(y - X\hat{\theta}(\lambda_{k-1}))| < 2\lambda_k - \lambda_{k-1}$ 

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# DERIVATION OF THE SAFE RULE FOR LINEAR REGRESSION

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

Define

$$\beta = \mathbf{y} - \mathbf{X}\boldsymbol{\theta}$$

Equivalent optimization problem

$$\hat{\theta} = \begin{cases} \arg\min_{\theta} & \beta^{\top}\beta + \lambda \left\|\theta\right\|_{1} \\ \text{subject to} & \beta = y - X\theta \end{cases}$$

# DERIVATION OF THE SAFE RULE FOR LINEAR REGRESSION

Lagrangian

$$\mathcal{L}(\theta, \beta, \nu) = \beta^{\top}\beta + \lambda \left\|\theta\right\|_{1} + \nu^{\top}(\mathbf{y} - \mathbf{X}\theta - \beta)$$

Dual function

$$\inf_{\theta,\beta} \mathcal{L}(\theta,\beta,\nu) = \begin{cases} \mathsf{G}(\nu) & \text{if } |f_j^\top \nu| \leq \lambda \,, \, j = 1,\dots,p \\ -\infty & \text{otherwise} \end{cases}$$

where  $G(\nu) = -\frac{1}{4}\nu^{\top}\nu + \nu^{\top}y$ . Lagrange dual

$$\hat{\theta}^* = \begin{cases} \underset{\nu}{\arg\max} & \mathcal{G}(\nu) \\ \\ \text{subject to} & |f_j^\top \nu| \le \lambda \,, \, j = 1, \dots, p \end{cases}$$

# DERIVATION OF THE SAFE RULE FOR LINEAR REGRESSION

Side note: Since the primal problem satisfies Slater's condition, we know that the duality gap  $\gamma = \hat{\theta} - \hat{\theta}^*$  is zero, i.e.

$$\hat{\theta} = \hat{\theta}^*$$

For a dual feasible point  $\nu_0$ , we solve for each  $j = 1, \ldots, p$ 

$$\xi_j(\nu_0) = \begin{cases} \arg \max_{\nu} & |f_j^\top \nu| \\ \text{subject to} & G(\nu) \ge G(\nu_0) \\ & = |f_j^\top y| + \sqrt{(y^\top y - 2G(\nu_0))f_j^\top f_j} \end{cases}$$

If  $\xi_j(\nu_0) < \lambda$  we know that  $\hat{\theta}_j = 0$ . A simple dual feasible point is  $\nu_0 = y\lambda/\lambda_{max}$ . The SAFE rule is obtained from

 $\xi_j(y\lambda/\lambda_{max}) < \lambda$ 

# SAGA algorithm [Defazio et al., 2014]: select $j \in \{1, \dots, n\}$ at random

$$\begin{split} \vartheta_{j,t+1} &= \theta_t \\ \vartheta_{i,t+1} &= \vartheta_{i,t+1} \text{ for all } i \neq j \\ \theta_{t+1}^* &= \theta_t - \gamma \left[ \nabla \ell_j(\vartheta_{j,t+1}) - \nabla \ell_j(\vartheta_{j,t}) + \frac{1}{n} \sum_{i=1}^n \nabla \ell_i(\vartheta_{i,t}) \right] \\ \theta_{t+1} &= \arg\min_{\theta} \left\{ \lambda \|\theta\|_1 + \frac{1}{2\gamma} \|\theta - \theta_{t+1}^*\|_2^2 \right\} \end{split}$$