## Lecture 10: Regularised regression (cont'd)

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MSA220/MVE441 Statistical Learning for Big Data
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## Regularisation in classification

## Recall: Regularised Discriminant Analysis (RDA)

Given training samples $\left(i_{l}, \mathbf{x}_{l}\right)$, quadratic DA models

$$
p(\mathbf{x} \mid i)=N\left(\mathbf{x} \mid \mu_{i}, \boldsymbol{\Sigma}_{i}\right) \quad \text { and } \quad p(i)=\pi_{i}
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## Regularisation can help here:

- Use $\widehat{\boldsymbol{\Sigma}}_{i}=\widehat{\boldsymbol{\Sigma}}_{i}^{\mathrm{QDA}}+\lambda \widehat{\boldsymbol{\Sigma}}^{\mathrm{LDA}}$ for $\lambda>0$
- Use LDA (i.e. $\boldsymbol{\Sigma}_{i}=\boldsymbol{\Sigma}$ ) and $\widehat{\boldsymbol{\Sigma}}=\widehat{\boldsymbol{\Sigma}}^{\text {LDA }}+\lambda \boldsymbol{\Delta}$ for $\lambda>0$ and a diagonal matrix $\boldsymbol{\Delta}$


## Recall: Naive Bayes LDA

Naive Bayes LDA means that we assume that $\widehat{\Sigma}=\widehat{\Delta}$ for a diagonal matrix $\widehat{\Delta}$. The diagonal elements are estimated as

$$
\widehat{\Delta}^{(j, j)}=\frac{1}{n-K} \sum_{i=1}^{K} \sum_{i_{l}=i}\left(\mathbf{x}_{l}^{(j)}-\widehat{\mu}_{i}^{(j)}\right)^{2}
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which is the pooled within-class variance.
Classification is performed by predicting the class with the maximal discriminant function value

$$
\begin{aligned}
\delta_{i}(\mathbf{x}) & =-\frac{1}{2}\left(\mathbf{x}-\widehat{\mu}_{i}\right)^{\top} \widehat{\Delta}^{-1}\left(\mathbf{x}-\widehat{\mu}_{i}\right)+\log \left(\widehat{\pi}_{i}\right) \\
& =-\frac{1}{2}\left\|\widehat{\Delta}^{-1 / 2}\left(\mathbf{x}-\widehat{\mu}_{i}\right)\right\|_{2}^{2}+\log \left(\widehat{\pi}_{i}\right)
\end{aligned}
$$

where $\left(\widehat{\Delta}^{-1 / 2}\right)^{(i, i)}=1 / \sqrt{\widehat{\Delta}^{(i, i)}}$.

## Shrunken centroids (I)

In high-dimensional problems $(p>n)$, centroids will

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Idea: Can we perform variable selection through $\ell_{1}-/$ lasso-style regularisation? How can we account for varying variance in features and stabilise against noise?

## Shrunken centroids (II)

Nearest shrunken centroids performs variable selection and stabilises centroid estimates by solving

$$
\bar{\mu}_{i}=\underset{\mathbf{v}}{\arg \min } \frac{1}{2} \sum_{i_{l}=i}\left\|\left(\widehat{\Delta}+s_{0} \mathbf{I}_{p}\right)^{-1 / 2}\left(\mathbf{x}_{l}-\mathbf{v}\right)\right\|_{2}^{2}+\lambda n_{i} m_{i}\left\|\mathbf{v}-\widehat{\mu}_{T}\right\|_{1}
$$

where $s_{0}=\operatorname{median}\left(\widehat{\boldsymbol{\Delta}}^{(1,1)}, \ldots, \widehat{\boldsymbol{\Delta}}^{(p, p)}\right), m_{i}=\sqrt{\frac{1}{n_{i}}-\frac{1}{n}}$ and $\widehat{\mu}_{T}=\frac{1}{n} \sum_{l} \mathbf{x}_{l}$.

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- $\widehat{\boldsymbol{\Delta}}+s_{0} \mathbf{I}_{p}$ is the diagonal regularised within-class covariance matrix. Features that are highly variable across samples are scaled down (interpretability)
- $n_{i} m_{i}$ scales $\lambda$ in case of unequal class sizes


## Shrunken centroids (III)

The solution for component $j$ can be derived as

$$
\bar{\mu}_{i}^{(j)}=\widehat{\mu}_{T}^{(j)}+m_{i}\left(\widehat{\Delta}^{(j, j)}+s_{0}\right) \operatorname{ST}\left(\mathbf{t}_{i}^{(j)}, \lambda\right) \quad \text { where } \quad \mathbf{t}_{i}^{(j)}=\frac{\widehat{\mu}_{i}^{(j)}-\widehat{\mu}_{T}^{(j)}}{m_{i}\left(\widehat{\Delta}^{(j, j)}+s_{0}\right)}
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- Typically, misclassification rate improves first with increasing $\lambda$ and declines for too high values
- The larger $\lambda$ the more components will be equal to the respective component of the overall centroid.


## Application of nearest shrunken centroids (I)

A gene expression data set with $n=63$ and $p=2308$. There are four classes (cancer subtypes) with $n_{\mathrm{BL}}=8, n_{\mathrm{EWS}}=23, n_{\mathrm{NB}}=12$, and $n_{\mathrm{RMS}}=20$.


5-fold cross-validation curve and largest $\lambda$ that leads to minimal misclassification rate

## Application of nearest shrunken centroids (II)



Grey lines show the original centroids and red lines show the shrunken centroids

## Extensions of the lasso

## The lasso and groups of highly correlated variables

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- Example: Two groups of highly correlated variables, e.g.

$$
\mathbf{X} \sim N(\mathbf{0}, \boldsymbol{\Sigma}) \quad \text { where } \quad \boldsymbol{\Sigma}=\left(\begin{array}{cc}
\boldsymbol{\Sigma}_{1} & \mathbf{0} \\
\mathbf{0} & \boldsymbol{\Sigma}_{1}
\end{array}\right) \in \mathbb{R}^{200 \times 200}
$$

where

$$
\Sigma_{1} \in \mathbb{R}^{100 \times 100}, \quad \Sigma_{1}^{(i, i)}=1.04 \quad \text { and } \quad \Sigma_{1}^{(i, j)}=1, \quad i \neq j
$$

The response is generated for $n=100$ samples as

$$
\mathbf{y}=\mathbf{x}_{1}-\mathbf{x}_{102}+\varepsilon \quad \text { where } \quad \varepsilon \sim N\left(\mathbf{0}, 4 \mathbf{I}_{p}\right)
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- Expectation: Since the predictors in each group are strongly correlated, all could be considered equally as predictors.
- Possible caveat: The lasso makes a sparsity assumption and tries to set as many coefficients to zero as possible.


## The lasso and groups of highly correlated variables in practice

Lasso


Elastic net ( $\alpha=0.2$ )

$$
\log (\lambda)
$$

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- At optimal $\lambda$ the lasso selects 8 non-zero coefficients 0 of which were in the true coefficient vector.
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## The lasso and groups of highly correlated variables in practice



- At optimal $\lambda$ the lasso selects 8 non-zero coefficients 0 of which were in the true coefficient vector.
- Very precise but 'wrong' estimates.
- An alternative algorithm, the elastic net estimates 95 non-zero coefficients. (44 in the $1^{\text {st }}$ group and 51 in the $2^{\text {nd }}$ group, group-wise close coefficients)
- 'Shares' responsibility among correlated variables


## The elastic net (I)

The elastic net solves the problem

$$
\underset{\beta}{\arg \min } \frac{1}{2}\|\mathbf{y}-\mathbf{X} \boldsymbol{\beta}\|_{2}^{2}+\lambda\left(\frac{1-\alpha}{2}\|\boldsymbol{\beta}\|_{2}^{2}+\alpha\|\boldsymbol{\beta}\|_{1}\right)
$$

striking a balance between lasso (variable selection) and ridge regression (grouping of variables)

Lasso


Elastic net ( $\alpha=0.7$ )


## Notes on the elastic net (II)

- The solution can be found through cyclic coordinate descent
- $\alpha$ is an additional tuning parameter that should be determined by cross-validation
- The lasso and ridge regression are special cases of the elastic net ( $\alpha=1$ and $\alpha=0$, respectively).


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- Categorical variables in dummy encoding
- Domain-knowledge (e.g. genes in the same signal pathway, signals that only appear in groups in a compressed sensing problem,...)


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- Domain-knowledge (e.g. genes in the same signal pathway, signals that only appear in groups in a compressed sensing problem,...)
- Ideally the whole group is either present or not
- The elastic net can find groups, but only does so for highly correlated variables and without external influence. Sometimes more control is necessary.


## The group lasso (I)

The group lasso solves the problem

$$
\underset{\beta}{\arg \min } \frac{1}{2}\|\mathbf{y}-\mathbf{X} \boldsymbol{\beta}\|_{2}^{2}+\lambda \sum_{k=1}^{K}\left\|\mathbf{B}_{k}\right\|_{2}
$$

where $\mathbf{B}_{k}$ is a vector of coefficients $\beta_{i}$ for the $k$-th group. Note that $\left\|\beta_{i}\right\|_{2}=\left|\beta_{i}\right|$ for singleton groups.

Lasso


Group lasso (\{ $\left.\left.\beta_{1}, \beta_{3}\right\},\left\{\beta_{2}\right\}\right)$


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- The elastic net similarly sets variables exactly to zero on a corner or along an edge. The curved edges encourage remaining coefficients to be closer together.


## Comparison: Lasso, elastic net and group lasso

Elastic net $(\alpha=0.7)$


Group lasso $\left(\left\{\beta_{1}, \beta_{3}\right\},\left\{\beta_{2}\right\}\right)$


- The lasso sets variables exactly to zero either on a corner or along an edge.
- The elastic net similarly sets variables exactly to zero on a corner or along an edge. The curved edges encourage remaining coefficients to be closer together.
- The group lasso has actual information about groups of variables. It encourages whole groups to be zero or non-zero with similar coefficients.


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Given $p(y \mid \beta, \mathbf{x})$ the log-likelihood of the model is

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Instead of penalising the minimisation of the residual sum of squares (RSS), the minimisation of the negative log-likelihood is penalized, i.e.

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\underset{\beta}{\arg \min }-\mathcal{L}(\beta \mid \mathbf{y}, \mathbf{X})+\lambda\|\beta\|_{1}
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Note: If $p(y \mid \boldsymbol{\beta}, \mathbf{x})$ is Gaussian and the linear model $\mathbf{y}=\mathbf{X} \boldsymbol{\beta}+\boldsymbol{\varepsilon}$ is assumed, this is equivalent to the lasso.

## Sparse logistic regression

Recall: For logistic regression with $i_{l} \in\{0,1\}$ it holds that

$$
p(1 \mid \boldsymbol{\beta}, \mathbf{x})=\frac{\exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}\right)}{1+\exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}\right)} \quad \text { and } \quad p(0 \mid \boldsymbol{\beta}, \mathbf{x})=\frac{1}{1+\exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}\right)}
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- The minimisation problem is still convex, but non-linear in $\beta$. Iterative quadratic approximations combined with coordinate descent can be used to solve this problem.
- Another way to perform sparse classification (like e.g. nearest shrunken centroids)


## Sparse multi-class logistic regression

In multi-class logistic regression with $i_{l} \in\{1, \ldots, K\}$, there is a matrix of coefficients $\mathbf{B} \in \mathbb{R}^{p \times(K-1)}$ and it holds for $i=1, \ldots, K-1$ that

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p(i \mid \mathbf{B}, \mathbf{x})=\frac{\exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}_{i}\right)}{1+\sum_{j=1}^{K-1} \exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}_{j}\right)} \quad \text { and } \quad p(K \mid \mathbf{B}, \mathbf{x})=\frac{1}{1+\sum_{j=1}^{K-1} \exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}_{j}\right)}
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- As in two-class case, the absolute value of each entry in $\mathbf{B}$ can be penalised.


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p(i \mid \mathbf{B}, \mathbf{x})=\frac{\exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}_{i}\right)}{1+\sum_{j=1}^{K-1} \exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}_{j}\right)} \quad \text { and } \quad p(K \mid \mathbf{B}, \mathbf{x})=\frac{1}{1+\sum_{j=1}^{K-1} \exp \left(\mathbf{x}^{\top} \boldsymbol{\beta}_{j}\right)}
$$

- As in two-class case, the absolute value of each entry in $\mathbf{B}$ can be penalised.
- Another possibility is to use the group lasso on all coefficients for one variable, i.e. penalise with $\left\|\mathbf{B}_{j} .\right\|_{2}$ for $j=1, \ldots, p$.


## Example for sparse multi-class logistic regression

## MNIST-derived zip code digits ( $n=7291, p=256$ )

Sparse multi-class logistic regression was applied to the whole data set and the penalisation parameter was selected by $10-$ fold CV.


Orange tiles show positive coefficients and blue tiles show negative coefficients. Class averages are shown in the background.

## Take-home message

- Penalisation methods are not only restricted to regression, also applicable to classification
- Sparsity is a very important concept when interpretability of models is important
- Many extensions to the lasso exist, which make it more suitable for a variety of different situations

