Machine Learning

Regularization and Feature Selection

Fabio Vandin

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Ridge Regression: Matrix Form

Linear regression: pick

$$\arg\min_{\mathbf{w}} (\mathbf{y} - \mathbf{X}\mathbf{w})^T (\mathbf{y} - \mathbf{X}\mathbf{w})$$

Ridge regression: pick

$$\arg\min_{\mathbf{w}} \left(\lambda ||\mathbf{w}||^2 + \left(\mathbf{y} - \mathbf{X} \mathbf{w} \right)^T \left(\mathbf{y} - \mathbf{X} \mathbf{w} \right) \right)$$

Want to find \mathbf{w} which minimizes

$$f(\mathbf{w}) = \lambda ||\mathbf{w}||^2 + (\mathbf{y} - \mathbf{X}\mathbf{w})^T (\mathbf{y} - \mathbf{X}\mathbf{w}).$$

$$\text{How?} \quad 2\lambda \vec{w} \quad -2 \times^T (\vec{y} - \vec{x} \vec{w})$$

Compute gradient $\frac{\partial f(\mathbf{w})}{\partial \mathbf{w}}$ of objective function w.r.t \mathbf{w} and compare it to 0.

Want to find \mathbf{w} which minimizes $f(\mathbf{w}) = \lambda ||\mathbf{w}||^2 + (\mathbf{y} - \mathbf{X}\mathbf{w})^T (\mathbf{y} - \mathbf{X}\mathbf{w}).$

How?

Compute gradient $\frac{\partial f(\mathbf{w})}{\partial \mathbf{w}}$ of objective function w.r.t \mathbf{w} and compare it to 0.

$$\frac{\partial f(\mathbf{w})}{\partial \mathbf{w}} = 2\lambda \mathbf{w} - 2\mathbf{X}^T (\mathbf{y} - \mathbf{X}\mathbf{w})$$

Then we need to find w such that

$$2\lambda \mathbf{w} - 2\mathbf{X}^{T}(\mathbf{y} - \mathbf{X}\mathbf{w}) = 0$$

$$\lambda \overrightarrow{w} - \lambda^{T}(\overrightarrow{y} - \lambda \overrightarrow{w}) = 0$$

$$\lambda \overrightarrow{w} + \lambda^{T} \lambda \overrightarrow{w} = \lambda^{T}$$

$$(\lambda T + \lambda^{T} \lambda) \overrightarrow{w} = \lambda^{T} \lambda^{T}$$

$$\Rightarrow \overrightarrow{w} = (\lambda T + \lambda^{T} \lambda) \lambda^{T} \lambda^{$$

$$2\lambda \mathbf{w} - 2\mathbf{X}^T(\mathbf{y} - \mathbf{X}\mathbf{w}) = 0$$

is equivalent to

$$\left(\lambda \mathbf{I} + \mathbf{X}^{T} \mathbf{X}\right) \mathbf{w} = \mathbf{X}^{T} \mathbf{y}$$

Note:

- X^TX is positive semidefinite
- **\lambda** is positive definite
- $\Rightarrow \lambda \mathbf{I} + \mathbf{X}^T \mathbf{X}$ is positive definite
- $\Rightarrow \lambda \mathbf{I} + \mathbf{X}^T \mathbf{X}$ is invertible

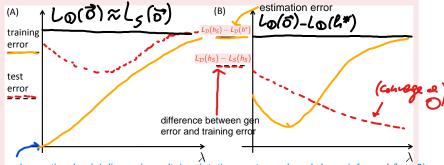
Ridge regression solution:

$$\mathbf{w} = \left(\lambda \mathbf{I} + \mathbf{X}^T \mathbf{X}\right)^{-1} \mathbf{X}^T \mathbf{y}$$

Exercise 5

Consider the ridge regression problem $\arg\min_{\mathbf{w}} \lambda ||\mathbf{w}||^2 + \sum_{i=1}^m (\langle \mathbf{w}, \mathbf{x_i} \rangle - y_i)^2$. Let: h_S be the hypothesis obtained by ridge regression on with training set S; h^* be the hypothesis of minimum generalization error among all linear models.

- (A) Draw, in the plot below, a *typical* behaviour of (i) the training error and (ii) the test/generalization error of h_S as a function of λ .
- (B) Draw, in the plot below, a *typical* behaviour of (i) $L_{\mathcal{D}}(h_S) L_{\mathcal{D}}(h^*)$ and (ii) $L_{\mathcal{D}}(h_S) L_S(h_S)$ as a function of λ .



puo' anche partire da piu' di zero in realta', asintoticamente per lamabda a +inf va a L(\vec0)

Feature Selection

In general, in machine learning one has to decide what to use as features (= input) for learning.

Even if somebody gives us a representation as a feature vector, maybe there is a "better" representation?

What is "better"?

Example

- features x₁, x₂, output y
- $x_1 \sim Uniform(-1,1)$
- $y = x_1^2$
- $x_2 \sim y + Uniform(-0.01, 0.01)$

If we want to predict y, which feature is better: x1 or x2?

-X1: because with (x1) we perfectly predict y.

But what if you ase x3 ds feature of a linear model.

Prediction is not great!

- x2, because with linear nodels we predict y exactly up to hoise (Unitorn (-0.01, 0.01)). But we could do better by booking at x1.

Example

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- $x_2 \sim y + Uniform(-0.01, 0.01)$

If we want to predict y, which feature is better: x_1 or x_2 ?

No-free lunch...

Feature Selection: Scenario

We have a large pool of features

Goal: select a small number of features that will be used by our (final) predictor

Assume $\mathcal{X} = \mathbb{R}^d$.

Goal: learn (final) predictor using $k \ll d$ predictors

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Motivation?

- prevent overfitting: less predictors ⇒ hypotheses of lower complexity!
- predictions can be done faster
- useful in many applications!

Feature Selection: Computational Problem

Assume that we use the Empirical Risk Minimization (ERM) procedure.

Assumption: an hypothesis $h \in \mathcal{H}$ corresponds to a vector of The problem of selecting k features that minimize the empirical well risk can be written as:

$$\min_{\mathbf{w}} L_S(\mathbf{w})$$
 subject to $||\mathbf{w}||_0 \le k$

where
$$||\mathbf{w}||_0 = |\{i : w_i \neq 0\}|$$

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How can we solve it?

Subset Selection

How do we find the solution to the problem below?

$$\min_{\mathbf{w}} L_{\mathcal{S}}(\mathbf{w})$$
 subject to $||\mathbf{w}||_0 \le k$

Note: the solution will always include *k* features

Let:

- $\mathcal{I} = \{1, \ldots, d\};$
- given $p = \{i_1, \dots, i_k\} \subseteq \mathcal{I}$: $\mathcal{H}_p = \text{hypotheses/models where}$ only features $w_{i_1}, w_{i_2}, \dots, w_{i_k}$ are used

Complexity? Learn $\Theta\left(\binom{d}{k}\right) \in \Theta\left(d^k\right)$ models \Rightarrow exponential algorithm!

2 it is unlikely that there is a poly-ting olg. to solve it.

Can we do better?

Proposition

The optimization problem of feature selection NP-hard.

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The optimization problem of feature selection NP-hard.

What can we do?

Heuristic solution ⇒ greedy algorithms

Greedy Algorithms for Feature Selection

Forward Selection: start from the empty solution, add one feature at the time, until solution has cardinality k

Complexity? Learns $\Theta(kd)$ models

Backward Selection: start from the solution which includes all features, remove one feature at the time, until solution has cardinality k

Pseudocode: analogous to forward selection [Exercize!]

Complexity? Learns $\Theta((d-k)d)$ models

Notes

We have used only training set to select the best hypothesis...

⇒ we may overfit!

Solution? Use validation! (or cross-validation)

Split data into training data and validation data, learn models on training, evaluate (= pick among different hypothesis models) on validation data. Algorithms are similar.

Note: now the best model (in terms of validation error) may include less than k features!

Subset Selection with Validation Data

```
S = \text{training data (from data split)}
V = validation data (from data split)
Using training and validation:
for \ell \leftarrow 0 to k do
     P^{(\ell)} \leftarrow \{J \subseteq \mathcal{I} : |J| = \ell\};
foreach p \in P^{(\ell)} do
     h_p \leftarrow \arg\min_{h \in \mathcal{H}_p} L_S(h);
     p_{\ell} \leftarrow \arg\min_{p \in P^{(\ell)}} L_V(h_p);
return arg \min_{p \in \{p_0, p_1, \dots, p_{\ell}\}} L_V(h_p)
```

Forward Selection with Validation Data

```
Using training and validation: sol \leftarrow \emptyset;
while |sol| < k \text{ do}
| foreach | i \in \mathcal{I} \setminus sol | foreach | i \in \mathcal{I} \setminus sol | foreach | f
```

Backward Selection with validation: similar [Exercize]

Similar approach for all algorithms with cross-validation [Exercize]

Bibliography [UML]

Regularization and Ridge Regression: Chapter 12

- no Section 13.3;
- Section 13.4 only up to Corollary 13.8 (excluded)

Feature Selection and LASSO: Chapter 25

• only Section 25.1.2 (introduction and "Backward Elimination") and 25.1.3