Machine learning from scratch

Lecture 6: Non-linear models, parameter selection

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Course outline

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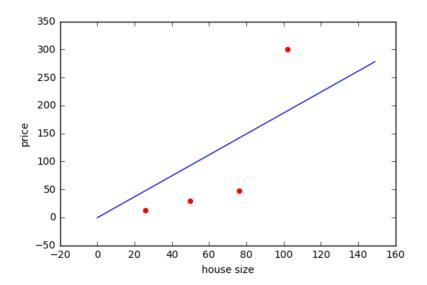
This lecture will go a bit further by introducing:

- Non linear models (polynomial kernels)
- Model evaluation
- Parameter selection

More complex models

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where λ is a **hyper-parameter** that quantifies how much we want to penalize big values of θ .

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A commonly used regularization term R is often the squared ℓ_2 norm given by

$$R(\theta) = \|\theta\|_2^2 = \sum_{i=1}^d \theta_i^2$$

Model evaluation

Parameter selection

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- Split the data into a training set and a test set
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This is often referred to as **cross-validation**.

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Another standard technique: *k*-**fold cross-validation**

- Split the data into k (equally-sized) folds
- Remove 1 fold (= test fold)
- ► Train on the other folds
- Test on the removed fold
- Do it for all the folds

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Other option: **Leave-one-out** (LOO) cross-validation:

- Remove 1 sample from the data set
- Train on all the other samples
- Test on the sample you've removed
- Evaluate the prediction
- Do it for each sample of the data set
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Alternative: Leave-p-out (LPO). LOO is LPO with p = 1.

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Most classic way: a grid-search

- ► For each parameter, define a set of possible values
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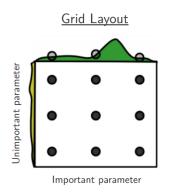
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Important note: Hyper-parameter ranges vary a lot from an application to another. It is **data-dependent**.

Random search is a more and more popular alternative to grid search, especially in this case:

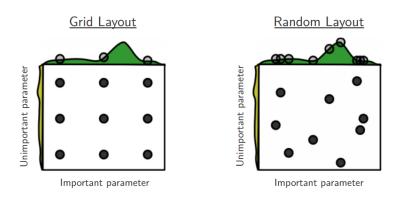
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Random Layout

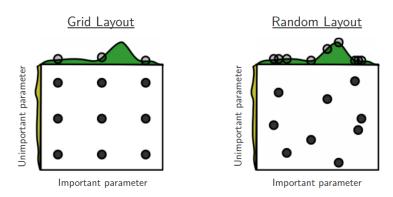
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Practical note: Each parameter combination can be trained/tested separately => possibility to distribute the tasks

Conclusion

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During the next lecture, we will work on implementing regularization to the OLS algorithm and cross-validating it and switch to classification if the time allows it.

Thank you! Questions?