

Model Selection

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How do we measure our models' ability to predict unseen data, when we only have access to training data?

The most common method to evaluate model **generalization** performance is *cross-validation*.

It is used in two essential data analysis phases: *Model Selection* and *Model Assessment*.

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Which kind of algorithm to use, linear regression vs. K-nearest neighbors vs. SVM

Model Assessment

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The same question can be asked of a kernel parameter in an SVM.

Cross-validation is a *resampling* method to obtain estimates of **expected prediction error rate** (or any other performance measure on unseen data).

In some instances, you will have a large predefined test dataset that you should never use when training.

In the absence of access to this kind of dataset, cross validation can be used.

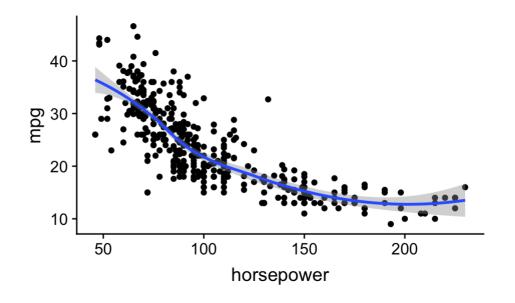
The simplest option to use cross-validation is to create a *validation* set, where our dataset is **randomly** divided into *training* and *validation* sets.

Then the *validation* is set aside, and not used at until until we are ready to compute **test error rate** (once, don't go back and check if you can improve it).



Let's look at an example using automobile data, where we want to build a regression model to predict miles per gallon given other auto attributes.

A linear regression model is not appropriate for this dataset. Use *polynomial* regression as an illustrative example.



For polynomial regression, our regression model (for a single predictor x) is given as a d degree polynomial.

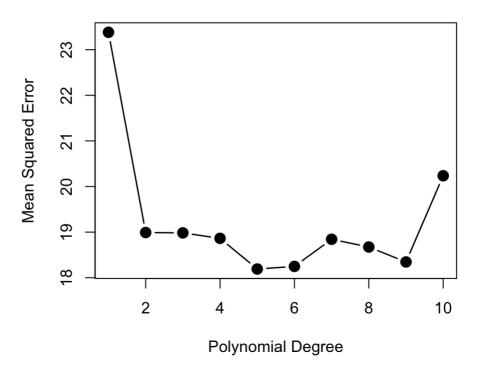
$$Y = b + w_1 x + w_2 x^2 + \dots + w_d x^d$$

For *model selection*, we want to decide what degree d we should use to model this data.

Using the *validation set* method, split our data into a training set,

fit the regression model with different polynomial degrees d on the training set,

measure test error on the validation set.



Resampled validation set

The validation set approach can be prone to sampling issues.

It can be highly variable as error rate is a random quantity and depends on observations in training and validation sets.

Resampled validation set

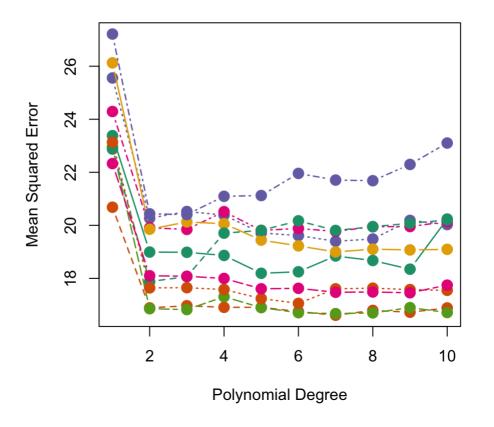
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We can improve our estimate of *test error* by averaging multiple measurements of it (remember the law of large numbers).

Resampled validation set

Resample validation set 10 times (yielding different validation and training sets) and averaging the resulting test errors.



This approach still has some issues.

Each of the training sets in our validation approach only uses 50% of data to train, which leads to models that may not perform as well as models trained with the full dataset and thus we can overestimate error.

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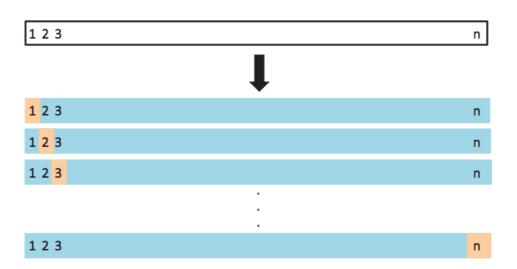
Each of the training sets in our validation approach only uses 50% of data to train, which leads to models that may not perform as well as models trained with the full dataset and thus we can overestimate error.

To alleviate this situation, we can extend our approach to the extreme: Make each single training point it's own validation set.

Procedure:

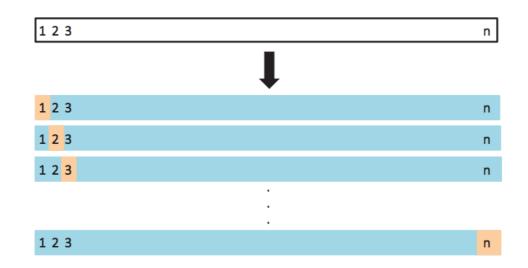
For each observation i in data set:

- a. Train model on all but i-th observation
- b. Predict response for i-th observation
- c. Calculate prediction error



This gives us the following *cross-validation* estimate of error.

$$CV_{(n)} = \frac{1}{n} \sum_{i} (y_i - \hat{y}_i)^2$$



Advantages:

- use n-1 observations to train each model
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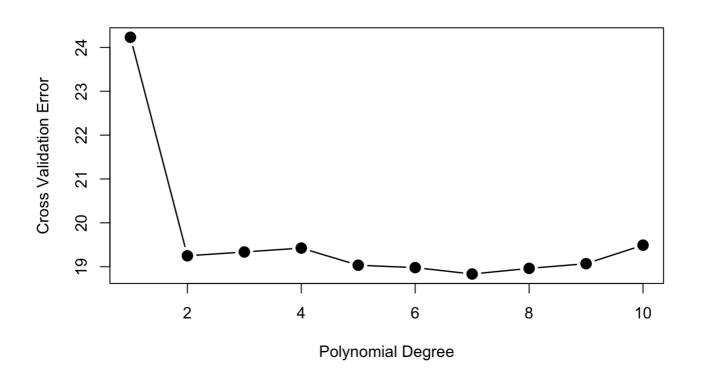
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Disadvantages:

- Depending on the models we are trying to fit, it can be very costly to train n-1 models.
- Error estimate for each model is highly variable (since it comes from a single datapoint).

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On our running example



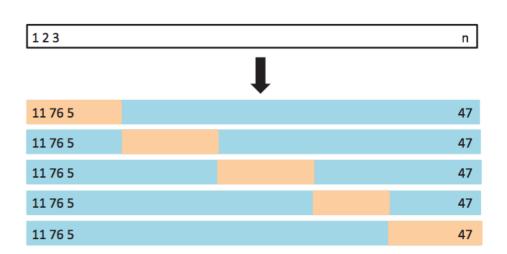
This discussion leads us to the most commonly used cross-validation approach *k-fold Cross-Validation*.

Procedure:

Partition observations randomly into k groups (folds).

For each of the k groups of observations:

- Train model on observations in the other k-1 folds
- Estimate test-set error (e.g.,
 Mean Squared Error) on this fold

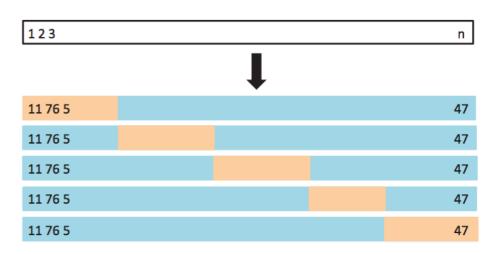


Procedure:

Compute average error across **k** folds

$$CV_{(k)} = \frac{1}{k} \sum_{i} MSE_{i}$$

where MSEi is mean squared error estimated on the i-th fold

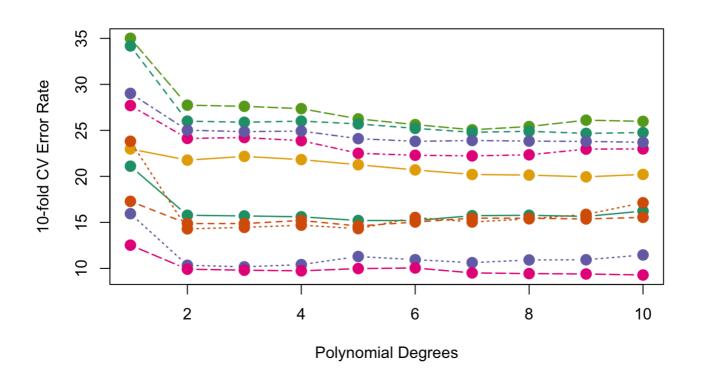


- Fewer models to fit (only k of them)
- Less variance in each of the computed test error estimates in each fold.

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It can be shown that there is a slight bias (over estimating usually) in error estimate obtained from this procedure.

Running Example



Cross-Validation in Classification

Each of these procedures can be used for classification as well.

In this case we would substitute MSE with performance metric of choice. E.g., error rate, accuracy, TPR, FPR, AUROC.

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In this case we would substitute MSE with performance metric of choice. E.g., error rate, accuracy, TPR, FPR, AUROC.

Note however that not all of these work with LOOCV (e.g. AUROC since it can't be defined over single data points).

Evaluating Classification

The AUROC statistic

The AUROC statistic is related to the Mann-Whitney non-parametric statistical test for distributional differences.

Null hypothesis: for randomly drawn pair of samples from two populations, it is equally likely that sample from first population is greater than sample from second population.

Evaluating Classification

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The AUROC statistic is related to the Mann-Whitney non-parametric statistical test for distributional differences.

Null hypothesis: for randomly drawn pair of samples from two populations, it is equally likely that sample from first population is greater than sample from second population.

Specifically, if x_A and x_B are drawn randomly from populations A and B respectively, $P(x_A < x_B) = P(x_A > x_B)$.

The AUROC statistic

Consider a classifier c trained to distinguish between two classes, using a training set containing n_A and n_B instances for each of the two classes respectively.

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Denote as c_i the score given by classifier i with higher c_i indicating predictions for class A.

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Null hypothesis: $P(C_i < C_j) = P(C_j < C_i)$ for randomly drawn pairs C_i from class A and C_i from class B.

The AUROC statistic

The Mann-Whitney test uses the U statistic to perform this test:

$$U = \sum_{i=1}^{n_A} \sum_{j=1}^{n_B} \frac{I\{C_i > C_j\}}{n_A n_B}$$

This is an empirical estimate of $P(C_i > C_j)$, which under the null hypothesis of the Mann-Whitney test is 0.5.

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It can be shown that ${\tt U}$ is exactly the AUCROC.

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Note that the $\mbox{$\mbox{$\mbox{$$}$}$}$ statistic, and thus AUROC, is only dependent on the rank of scores c_i *not on their magnitude*.

This implies that we can compare AUCROC for classifiers that produce scores in different scales, e.g., probabilities or not.

The AUROC statistic

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See http://papers.nips.cc/paper/2645-confidence-intervals-for-the-area-under-the-roc-curve.pdf

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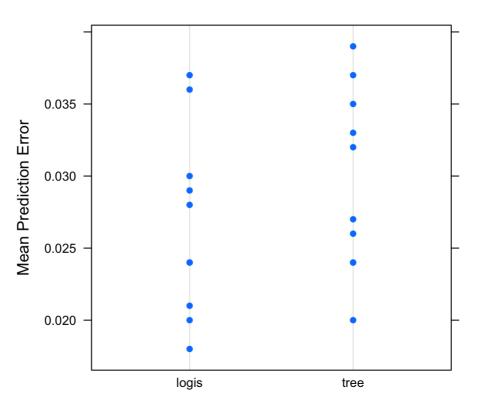
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There are methods to compare AUCROC statistics from multiple classifiers. See http://ieeexplore.ieee.org/document/6851192/ for the most practical.

Comparing models using cross-validation

Suppose you want to compare two classification models (logistic regression vs. a decision tree) on the Default dataset. We can use Cross-Validation to determine if one model is better than the other, using a t-test for example.



Comparing models using cross-validation

Using hypothesis testing:

| term | estimate | std.error | statistic | p.value |
|-------------|----------|-----------|-----------|-----------|
| (Intercept) | 0.0267 | 0.0020306 | 13.148828 | 0.0000000 |
| methodtree | 0.0030 | 0.0028717 | 1.044677 | 0.3099998 |

In this case, we do not observe any significant difference between these two classification methods.

Summary

Model selection and assessment are critical steps of data analysis.

Resampling methods are general tools used for this purpose.

k-fold cross-validation can be used to provide larger training sets to algorithms while stabilizing empirical estimates of expected prediction error