Statistical Natural Language Processing Machine learning: evaluation

Çağrı Çöltekin

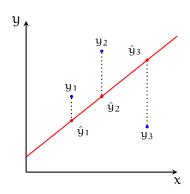
University of Tübingen Seminar für Sprachwissenschaft

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Measuring success/failure in regression

Root mean squared error (RMSE)

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

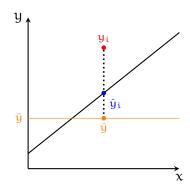


Measures average error in the units compatible with the outcome variable

Measuring success/failure in regression

Coefficient of determination

$$R^{2} = \frac{\sum_{i}^{n} (\hat{y}_{i} - \overline{y})^{2}}{\sum_{i}^{n} (y_{i} - \overline{y})^{2}}$$
$$= 1 - \frac{MSE}{\sigma_{y}^{2}}$$



- r^2 is a standardized measure in range [0, 1]
- Indicates the ratio of variance of y explained by x
- For single predictor it is the square of the correlation coefficient r

Measuring success in classification

Accuracy

- In classification, we do not care (much) about the average of the error function
- We are interested in how many of our predictions are correct
- Accuracy measures this directly

$$accuracy = \frac{number\ of\ correct\ predictions}{total\ number\ of\ predictions}$$

Accuracy may go wrong

- Think about a 'dummy' search engine that always returns an empty document set (no results found)
- If we have
 - 1000 000 documents
 - 1000 relevant documents (including the term in the query)
 the accuracy is:

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- If we have
 - 1000000 documents
 - 1000 relevant documents (including the term in the query)
 the accuracy is:

$$\frac{999\,000}{1\,000\,000} = 99.90\,\%$$

 In general, if our class distribution is skewed accuracy will be a bad indicator of success

Measuring success in classification

Precision, recall, F-score

$$precision = \frac{TP}{TP + FP}$$

$$recall = \frac{TP}{TP + FN}$$

$$F_{1}\text{-score} = \frac{2 \times precision \times recall}{precision + recall}$$

		true value				
þ		positive	negative			
predicted	pos.	TP	FP			
	neg.	FN	TN			

Example: back to the search engine

- We had a 'dummy' search engine that returned false for all queries
- For a query
 - 1000000 documents
 - 1000 relevant documents

accuracy =
$$\frac{999\,000}{1\,000\,000}$$
 = 99.90 %
precision = $\frac{0}{1\,000\,000}$ = 0 %
recall = $\frac{0}{1\,000\,000}$ = 0 %

Precision and recall are asymmetric, the choice of the 'positive' class is important.

Classifier evaluation: another example

Consider the following two classifiers:

		true	value	true value		value
predicted		positive	negative		positive	negative
	pos.	7	9		1	3
	neg.	3	1		9	7

Classifier evaluation: another example

Consider the following two classifiers:

		true value			true value		
ğ		positive	negative		positive	negative	
predicted	pos.	7	9		1	3	
	neg.	3	1		9	7	

Accuracy both
$$8/20 = 0.4$$

Precision $7/16 = 0.44$ and $1/4 = 0.25$
Recall $7/10 = 0.7$ and $1/10 = 0.1$
F-score 0.54 and 0.14

Multi-class evaluation

- For multi-class problems, it is common to report average precision/recall/f-score
- For C classes, averaging can be done two ways:

$$precision_{M} = \frac{\sum_{i}^{C} \frac{TP_{i}}{TP_{i} + FP_{i}}}{C} \qquad recall_{M} = \frac{\sum_{i}^{C} \frac{TP_{i}}{TP_{i} + FN_{i}}}{C}$$

$$precision_{\mu} = \frac{\sum_{i}^{C} TP_{i}}{\sum_{i}^{C} TP_{i} + FP_{i}} \qquad recall_{\mu} = \frac{\sum_{i}^{C} TP_{i}}{\sum_{i}^{C} TP_{i} + FN_{i}}$$

$$(M = macro, \mu = micro)$$

• The averaging can also be useful for binary classification, if there is no natural positive class

Confusion matrix

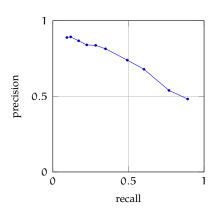
 A confusion matrix is often useful for multi-class classification tasks

		true		
		a	b	c
redicted	a	10	3	4
dic	b	2	12	8
pre	c	0	7	7

- Are the classes balanced?
- What is the accuracy?
- What is per-class, and averaged precision/recall?

Precision-recall trade-off

- Increasing precision (e.g., by changing a hyperparameter) results in decreasing recall
- Precision–recall graphs are useful for picking the correct models
- Area under the curve (AUC) is another indication of success of a classifier

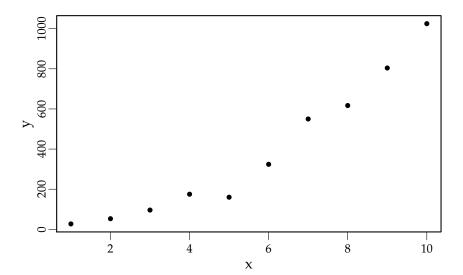


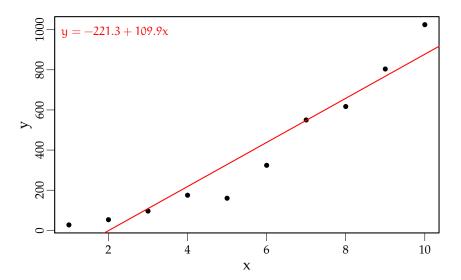
Performance metrics a summary

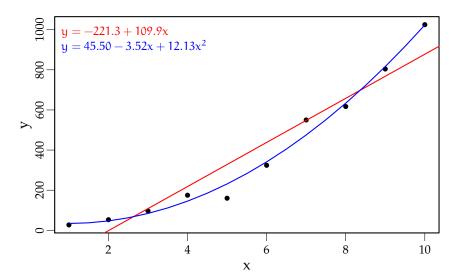
- Accuracy does not reflect the classifier performance when class distribution is skewed
- Precision and recall are binary and asymmetric
- For multi-class problems, calculating accuracy is straightforward, but others measures need averaging
- These are just the most common measures: there are more
- You should understand what these metrics measure, and use/report the metric that is useful for the purpose

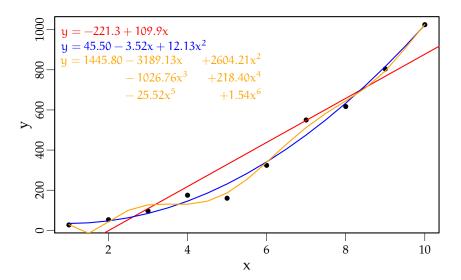
Model selection/evaluation

- Our aim is to fit models that are (also) useful outside the training data
- Evaluating a model on the training data is wrong: complex models tend to fit to the noise in the training data
- The results should always be tested on a test set that does not overlap with the training data
- Test set is ideally used only once to evaluate the final model
- Often, we also need to tune the model, find best hyperparameters (e.g., regularization constant)
- Tuning has to be done on a separate development set

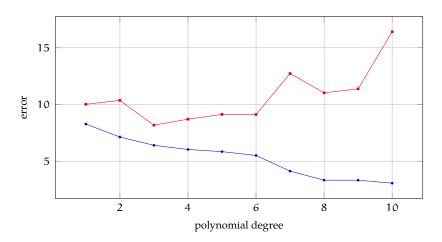








Training/test error



Bias and variance (revisited)

Bias of an estimate is the difference between the value being estimated, and the expected value of the estimate

$$B(\hat{\boldsymbol{w}}) = E[\hat{\boldsymbol{w}}] - \boldsymbol{w}$$

An unbiased estimator has 0 bias

Variance of an estimate is, simply its variance, the value of the squared deviations from the mean estimate

$$var(\hat{\boldsymbol{w}}) = E\left[(\hat{\boldsymbol{w}} - E[\hat{\boldsymbol{w}}])^2\right]$$

w is the parameters that define the model

Bias-variance relationship is a trade-off: models with low bias result in high variance.

Some issues with bias and variance

- *Overfitting* occurs when the model learns the idiosyncrasies of the training data
- Underfitting occurs when the model is not flexible enough for the data at hand

Some issues with bias and variance

- Overfitting occurs when the model learns the idiosyncrasies of the training data
- Underfitting occurs when the model is not flexible enough for the data at hand
- Complex models tend to overfit and exhibit high variance
- Simple models tend to show low variance, but likely to have (high) bias

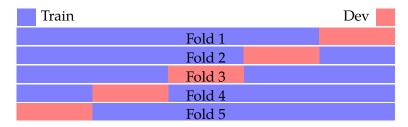
Model selection & hyperparamater tuning

- Our aim is to reduce the test error
- We can estimate the test error on a development set, or held-out data:
 - Split the data at hand as training and development set
 - Train alternative models (different hyperparameters) on the training set
 - Choose the model with best development set performance

Cross validation

- To avoid overfitting, we want to tune our models on a development set
- But (labeled) data is valuable
- Cross validation is a technique that uses all the data, for both training and tuning with some additional effort
- Besides tuning hyper-parameters, we may also want to get 'average' parameter estimates over multiple folds

K-fold Cross validation



- At each fold, we hold part of the data for testing, train the model with the remaining data
- Typical values for k is 5 and 10
- In *stratified* cross validation each fold contains (approximately) the same proportions of class labels.
- A special case, when k is equal to n (the number of data points is called *leave-one-out cross validation*

The choice of k in k-fold CV

- Increasing k
 - reduces the bias: the estimates converge to true value of the measure (e.g., accuracy) in the limit
 - increases the variance: smaller held-out sets produce more varied parameter estimates
 - is generally computationally expensive
- 5- or 10-fold cross validation is common practice (and found to have a good balance between bias and variance)

Comparing with a basline

- The performance measures are only meaningfull if we have something to compare against
- random does the model do anything useful at all?
- majority class does the classifier better than predicting the majority class all the time?
- state-of-the-art how does your model compare against known (non-trivial) models?
 - Diferences between models are reliable only if the same data set is used
 - Differences are stable if your test set size is large enough
 - Use statistical tests when comparing different models/methods

Summary

The first principle is that you must not fool yourself and you are the easiest person to fool. – Richard P. Feynman

- The measures of success in ML systems include
 - RMSE / r²
 Accuracy
 Precision / recall / F-score
- We want models with low bias and low variance
- Evaluating ML system requires special care:
 - Never use your test set during training / development
 - Tuning your system on a development set
 - Cross-validation allows efficient use of labeled data

Next:

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Next:

• Have good holiday! We'll start with sequence learning after the break.