#### POIR 613: Computational Social Science

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# Overview of QTA (Grimmer and Stewart, 2013)

- 1. Acquire textual data:
  - Existing corpora; scraped data; digitized text
- 2. Preprocess the data:
  - Bag-of-words vs word embeddings
- 3. Apply method appropriate to research goal:
  - Describe and compare documents
    - Readability; similarity; keyness metrics
  - Classify documents into known categories
    - Dictionary methods
    - Supervised machine learning
  - Classify documents into unknown categories
    - Document clustering
    - Topic models
  - Scale documents on latent dimension
    - Known dimension: wordscores
    - Unknown dimensions: wordfish

# Supervised machine learning

#### Outline

#### Supervised learning overview

- Creating a labeled set and evaluating its reliability
- Classifier performance metrics
- One classifier for text
  - Regularized regression

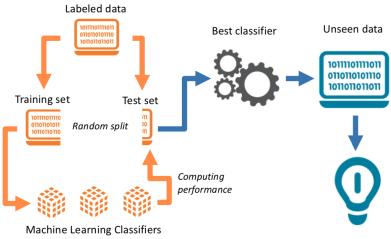
# Supervised machine learning

Goal: classify documents into pre existing categories.

e.g. authors of documents, sentiment of tweets, ideological position of parties based on manifestos, tone of movie reviews...

#### What we need:

- Hand-coded dataset (labeled), to be split into:
  - Training set: used to train the classifier
  - Validation/Test set: used to validate the classifier
- Method to extrapolate from hand coding to unlabeled documents (classifier):
  - Naive Bayes, regularized regression, SVM, K-nearest neighbors, BART, ensemble methods...
- Performance metric to choose best classifier and avoid overfitting: confusion matrix, accuracy, precision, recall...



Predictions

Supervised learning v. dictionary methods

- Dictionary methods:
  - Advantage: not corpus-specific, cost to apply to a new corpus is trivial
  - Disadvantage: not corpus-specific, so performance on a new corpus is unknown (domain shift)
- Supervised learning can be conceptualized as a generalization of dictionary methods, where features associated with each categories (and their relative weight) are learned from the data
- By construction, they will outperform dictionary methods in classification tasks, as long as training sample is large enough

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#### Creating a labeled set

How do we obtain a labeled set?

- External sources of annotation
  - Disputed authorship of Federalist papers estimated based on known authors of other documents
  - Party labels for election manifestos
  - Legislative proposals by think tanks (text reuse)
- Expert annotation
  - "Canonical" dataset in Comparative Manifesto Project
  - In most projects, undergraduate students (expertise comes from training)
- Crowd-sourced coding
  - Wisdom of crowds: aggregated judgments of non-experts converge to judgments of experts at much lower cost (Benoit et al, 2016)
  - Easy to implement with FigureEight or MTurk

#### Code the Content of a Sample of Tweets

Instructions -

In this job, you will be presented with tweets about the recent protests related to race and law enforcement in the U.S.

You will have to read the tweet and answer a set of questions about its content.

Read the tweet below paying close attention to detail:

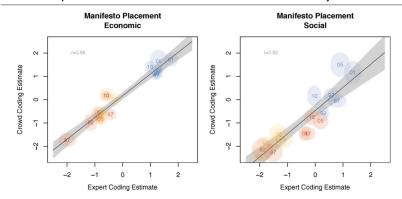
Tweet ID: 447



Is this tweet related to the ongoing debate about law enforcement and race in the United States?

- Yes
- O No
- O Don't Know

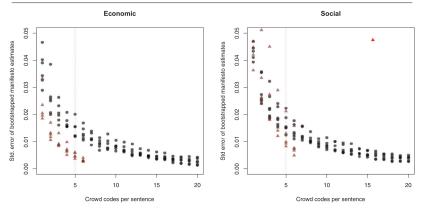
#### Crowd-sourced text analysis (Benoit et al, 2016 APSR)



#### FIGURE 3. Expert and Crowd-sourced Estimates of Economic and Social Policy Positions

#### Crowd-sourced text analysis (Benoit et al, 2016 APSR)





Note: Each point is the bootstrapped standard deviation of the mean of means aggregate manifesto scores, computed from sentencelevel random n subsamples from the codes.

## Evaluating the quality of a labeled set

Measures of agreement:

Percent agreement Very simple:

(number of agreeing ratings) / (total ratings) \* 100%

- Correlation
  - (usually) Pearson's r
  - May also be ordinal, such as Spearman's rho or Kendall's tau-b
  - Range is [-1,1]
- Agreement measures
  - Take into account not only observed agreement, but also agreement that would have occurred by chance
  - Cohen's κ is most common
  - Krippendorf's α is a generalization of Cohen's κ
  - Both range from [0,1]

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# Computing performance

Binary outcome variables:

|                | Actual value   |                |  |
|----------------|----------------|----------------|--|
| Classification | Ham            | Spam           |  |
| Ham            | True negative  | False negative |  |
| Spam           | False positive | True positive  |  |

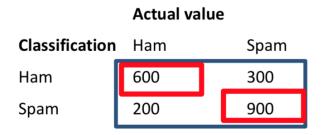
Confusion matrix:

- True negatives and true positives are correct predictions (to maximize)
- False positives and false negatives are incorrect predictions (to minimize)

• Performance metrics

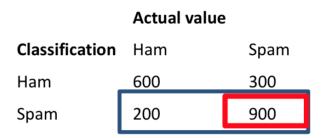
|                | Actual value   |                |  |
|----------------|----------------|----------------|--|
| Classification | Ham            | Spam           |  |
| Ham            | True negative  | False negative |  |
| Spam           | False positive | True positive  |  |

- Accuracy: correct predictions / total of predictions
  - % of units that are correctly predicted
- Precision for positive labels: (true positive) / (false positive + true positive)
  - % of units predicted to be positive that are indeed positive
- Recall for positive labels: (true positive) / (true positive + false negative)
  - % of units that are positive and are predicted as such



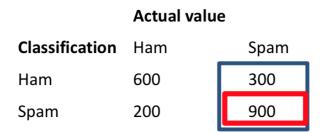
Accuracy = (600 + 900) / (600 + 900 + 200 + 300) = 0.75

75% of all emails are correctly classified



Precision = (900) / (900 + 200) = 0.82

83% of all emails predicted to be spam are indeed **spam** 



**Recall = (900) / (900 + 300)** = 0.75

75% of all spam emails are correctly classified

|                | Actual value |      |                          |
|----------------|--------------|------|--------------------------|
| Classification | Ham          | Spam | Total emails:            |
| Ham            | 1700         | 50   | Ham = 1850<br>Spam = 150 |
| Spam           | 150          | 100  |                          |

Accuracy = (1700+100) / (1700+50+150+100) = 0.90 Precision (spam) = (100) / (150+100) = 0.40 Recall (spam) = (100) / (50+100) = 0.67

Accuracy can be misleadingly high!

Imagine extreme scenario: we classify everything as ham – accuracy would be 92.5%

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#### The trade-off between precision and recall

Two extreme scenarios (but same underlying data):

1) Model predicts always spam

 Model predicts (almost) always ham (e.g. only emails with 10+ links as spam)
Actual value

#### Actual value

| Classif. | Ham | Spam |
|----------|-----|------|
| Ham      | 0   | 0    |
| Spam     | 800 | 1200 |

Accuracy = 1200 / 2000 = 0.60 Precision (spam) = 1200 / (800 + 1200) = 0.60 Recall (spam) = 1200 / 1200 = 1.00

High recall but low precision.

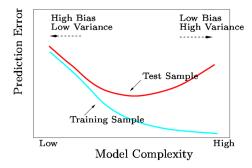
#### Classif. Ham Spam Ham 800 1190 Spam 0 10

Accuracy = (800 + 10) / 2000 = 0.40 Precision (spam) = 10 / 10 = 1.0 Recall (spam) = 10 / 1200 = 0.01

High precision but low recall.

# Measuring performance

- Classifier is trained to maximize in-sample performance
- But generally we want to apply method to new data
- Danger overfitting: In-sample performance better than out-of-sample performance (low generalizability)

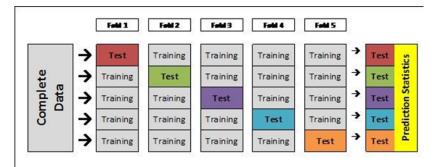


- Solutions?
  - Randomly split dataset into training and test set
  - Cross-validation

## **Cross-validation**

Intuition:

- Create K training and test sets ("folds") within training set.
- For each k in K, run classifier and estimate performance in test set within fold.
- Choose best classifier based on cross-validated performance



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# Types of classifiers

General thoughts:

- Trade-off between accuracy and interpretability
- Parameters need to be cross-validated

Frequently used classifiers:

- Naive Bayes
- Regularized regression
- SVM
- Others: k-nearest neighbors, tree-based methods, etc.
- Ensemble methods

#### Regularized regression

Assume we have:

- $i = 1, 2, \dots, N$  documents
- Each document *i* is in class  $y_i = 0$  or  $y_i = 1$
- $j = 1, 2, \ldots, J$  unique features
- And x<sub>ij</sub> as the count of feature j in document i

We could build a linear regression model as a classifier, using the values of  $\beta_0$ ,  $\beta_1$ , ...,  $\beta_J$  that minimize:

$$RSS = \sum_{i=1}^{N} \left( y_i - \beta_0 - \sum_{j=1}^{J} \beta_j x_{ij} \right)^2$$

But can we?

- If J > N, OLS does not have a unique solution
- Even with N > J, OLS has low bias/high variance (overfitting)

#### Regularized regression

What can we do? Add a penalty for model complexity, such that we now minimize:

$$\sum_{i=1}^{N} \left( y_i - \beta_0 - \sum_{j=1}^{J} \beta_j x_{ij} \right)^2 + \lambda \sum_{j=1}^{J} \beta_j^2 \rightarrow \text{ridge regression}$$

or

$$\sum_{i=1}^{N} \left( y_i - \beta_0 - \sum_{j=1}^{J} \beta_j x_{ij} \right)^2 + \lambda \sum_{j=1}^{J} |\beta_j| \rightarrow \text{lasso regression}$$

where  $\lambda$  is the **penalty parameter** (to be estimated)

#### Regularized regression

Why the penalty (shrinkage)?

- Reduces the variance
- Identifies the model if J > N
- Some coefficients become zero (feature selection)

The penalty can take different forms:

- Ridge regression: λ Σ<sup>J</sup><sub>j=1</sub> β<sup>2</sup><sub>j</sub> with λ > 0; and when λ = 0 becomes OLS
- ► Lasso  $\lambda \sum_{j=1}^{J} |\beta_j|$  where some coefficients become zero.
- ► Elastic Net:  $\lambda_1 \sum_{j=1}^J \beta_j^2 + \lambda_2 \sum_{j=1}^J |\beta_j|$  (best of both worlds?)

How to find best value of  $\lambda$ ? Cross-validation.

Evaluation: regularized regression is easy to interpret, but often outperformed by more complex methods.

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