# Text Mining Project/Lab

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# Text Classification(2)

# Pattern Recognition and Classification

- Detecting patterns and structures is a central theme in text mining.
- We usually start with the hypothesis that certain observable patterns in text are correlated to a particular task we address in text mining.
- In the previous session we learned how to do basic classification tasks in order to detect these observable patterns:
  - Naïve Bayes
  - Decision Tree

#### Goal of this session

- We will dig more into the evaluation of a classification task;
- We will have a closer look at some learning techniques.

- Evaluation is required
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  - To guide us in the process of feature selection;
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- Evaluation is required
  - To decide whether a classifier is reliable and to quantify its quality;
  - To guide us in the process of feature selection;
  - Also, to compare classification techniques.
- Depending on the evaluation procedures, we can distinguish different types of evaluation tasks:
  - Intrinsic evaluation: an isolated classifier is evaluated with respect to a predefined gold standard result (i.e. a widely accepted dataset);
  - Extrinsic: a classifier is evaluated when it provides a precise functionality for a human user (often more complex, can say why?).

- There exists other classification of evaluation techniques:
  - Black-box vs. glass-box evaluation (similar to software engineering);
  - Automatic vs. manual evaluation (automatic is preferred).
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  - There exists an open list of performance measures, some of them task dependant:
    - Accuracy
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    - Cohen's kappa
    - Usability metrics

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- For an intrinsic evaluation we first need a test set (or evaluation set).
- We then calculate a score (performance measure) for a classifier by comparing the labels that it generates for the inputs in the test set with the correct asserted labels for them.
- Test set is often has the same structure as training set.
- However, it is very important that the test set be distinct from the training corpus to avoid over-fitting:
  - a model can simply memorized its input, without really learning how to generalize to new examples.
  - A misleading score!

- There is often a trade-off between the amount of data available for testing and the amount available for training.
- The number of instances in the test set depends on the classification task and the distribution of labels (sometimes as few as 100 instances).
- The test set must be balanced and diverse.

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  - If there are a large number of labels, or some of labels are infrequent, then the test set must be big enough to ensure that the least frequent label are occurring enough (e.g. 50 times).
  - If the test set contains many closely related instances, then the size of the test set should be increased to ensure that this lack of diversity does not skew the evaluation results.
- When large amounts of annotated data are available, it is common to use 10% of the overall data for evaluation.

- An obtained score is more reliable, if test set and train set are less similar:
  - For a PoS tagger that is trained on a new corpus, the obtained score from the evaluation of the tagger on another news corpus is less reliable than the evaluation of the tagger on a corpus of text of genres other than news.
- In short, test set must be a real representative of the actual text data that a classifier is going to deal with.

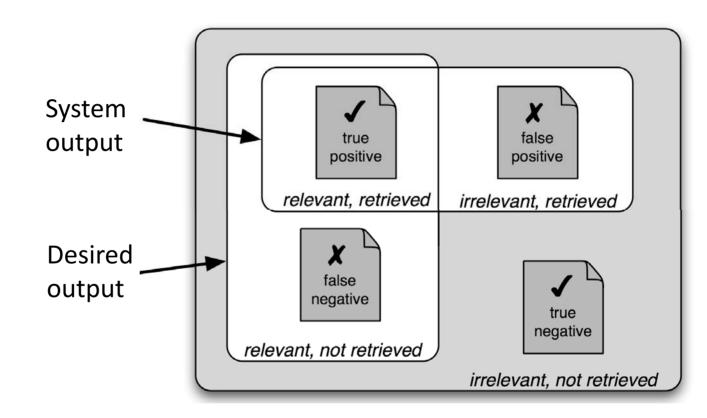
- Accuracy is one simple metric that can be used to evaluate a classifier.
- Accuracy measures the percentage of inputs in the test set that the classifier correctly labelled.
  - A classifier that predicts the correct labels 60 times in a test set containing 100 names would have an accuracy of 60/100 = 60%.
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  - The function nltk.classify.accuracy() can be used to measure accuracy.
- If the test set is not balanced, accuracy is not a very good measure :
  - The frequencies of the individual class labels in the test set are ignored.

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- Precision indicates how many of the items that are identified were relevant (or expected).
- Recall, however, indicates how many of the all relevant items we could identify.
- The goal is a high precision and high recall.
- The F-Measure (or F-Score), is thus used to combine these two measures.



Precision = TP/(TP+FP)

Recall = TP/(TP+FN)

F-Score =  $(2 \times Precision \times Recall) / (Precision + Recall).$ 

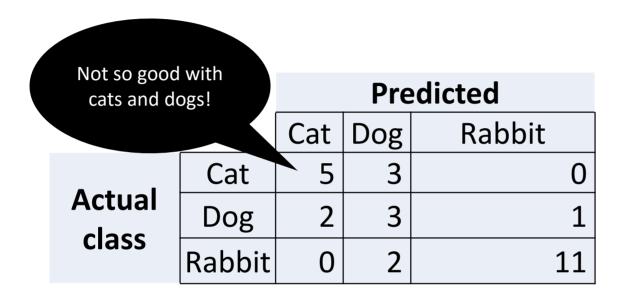
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- A confusion matrix is a table where each cell [i,j] indicates how often label j was predicted when the correct label was i.
- Thus, the diagonal entries (i.e., cells [ii]) indicate labels that were correctly predicted.
- You can use nltk.ConfusionMatrix to generate a confusion matrix.

		Predicted			
		Cat	Dog	Rabbit	
Actual class	Cat	5	3	0	
	Dog	2	3	1	
	Rabbit	0	2	11	

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Good at catching Rabbits!



Predicted

Cat Dog Rabbit

Cat 5 3 0

Dog 2 3 1

Rabbit 0 2 11

		Predicted			
		Cat	Dog	Rabbit	
Actual class	Cat	5	3	0	8
	Dog	2	3	1	6
	Rabbit	0	2	11	13
		7	8	12	

	Predicted					
		Cat	Dog	Rabbit		
Actual class	Cat	5	3	0	8	0.625
	Dog	2	3	1	6	0.5
	Rabbit	0	2	11	13	0.846154
		7	8	12		
		0.71	0.38	0.916666667		

#### **Cross-Validation**

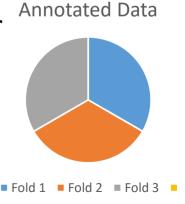
- In order to evaluate a classifier, we must reserve a portion of the annotated data for the test set.
- Test set must be big enough to have an accurate.
- However, if a limited amount of annotated data is available, making the test set larger is not plausible.
- Larger test dataset means making the training set smaller, which impact the classifier performance significantly.
- One solution to solve this problem is to perform cross-validation.

#### **Cross-Validation**

- Cross-Validation technique consists of multiple evaluations on different test sets and the combination of the obtained scores.
- We first divide the original corpus into n subsets called folds.
- For each of these folds, we train a model using all of the data except the data in that fold, and then test that model on the fold.
- The obtained scores from the individual folds (which might be too small to give accurate evaluation scores on their own) are combined to report an evaluation score.
- The final score is thus based on a large amount of data, and is therefore reliable.

#### **Cross-Validation**

- Additionally, cross-validation allows us to examine how widely the performance varies across different training sets:
  - If scores are similar for all *n* training sets, then we can be fairly confident that the score is accurate.
  - If scores vary widely across the *n* training sets, then we should probably be sceptical about the accuracy of the evaluation score.
- Remember, each fold must be balance and diverse (i.e. a fair representative of the real data).



# A closer look at learning techniques

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- It's possible to treat learning methods as black boxes:
  - Simply train and use a learning technique without understanding how it work.
- However, an understanding of a learning method can help us in several ways:
  - Selection of appropriate features;
  - Encoding feature values;
  - Setting model parameters;
  - Etc.

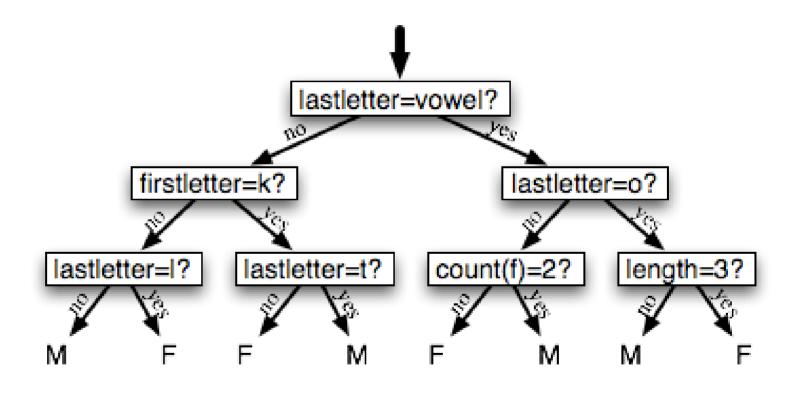
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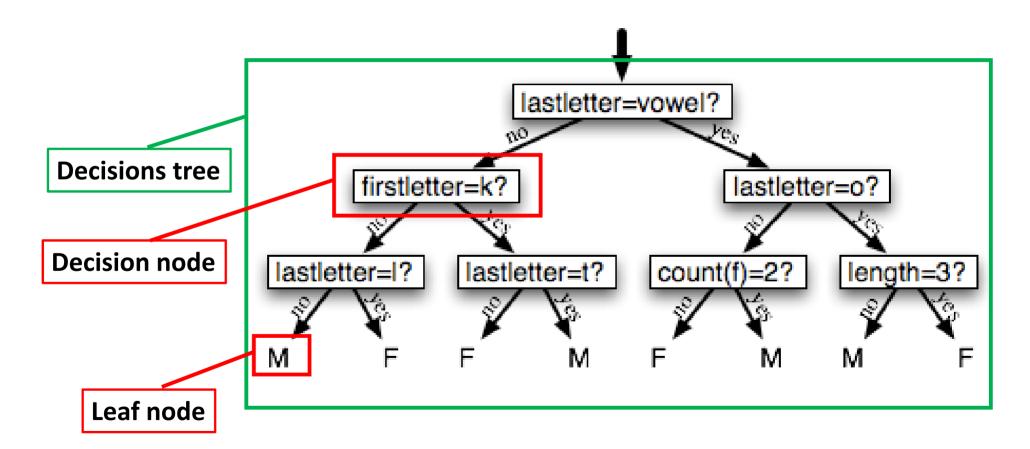
- We thus have a slightly closer look at:
  - Decisions trees
  - Naïve Bayes Classifiers
  - Maximum Entropy Classifiers

#### **Decisions trees**

- A decision tree is a flowchart that selects labels for input values.
- This flowchart consists of **decision nodes** (check feature values), and **leaf nodes** (assign labels).
- To choose the label for an input value, we begin at the initial decision node (known as root node).

#### Decisions trees

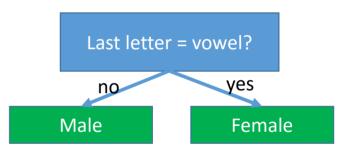




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  - What is the root node?
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# Building a decision stump

- As expected, we must first decide which feature should be used.
- Afterwards, the simplest method is to:
  - Build a decision stump for each possible feature;
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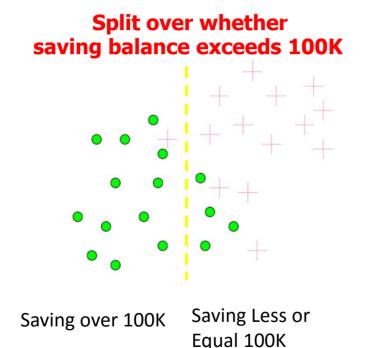


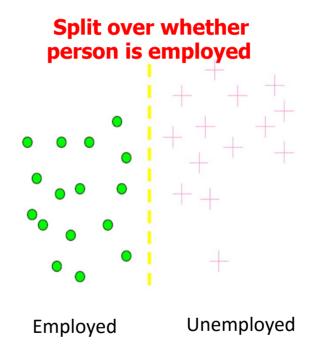
# Building a decision tree

- Now that we can build a decision stump, we can build a larger decision tree:
  - Build the decision stumps;
  - Select the overall best decision stump for the classification task;
  - Replace leaves that do not achieve sufficient accuracy with new decision stumps.
- We can use the notion of Information Gain to choose the most informative feature!

- Information gain tells us how important a feature is.
- We can use this measure to organize/order features in a decision tree.
- Vice versa, entropy measure how disorganized the original set of input values are.

Which test is more informative?



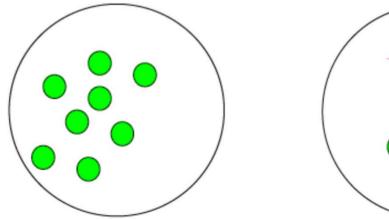


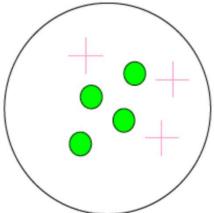
Acknowledgement: Content and Example for this introduction are taken from Shapiro and Stockman (2001)

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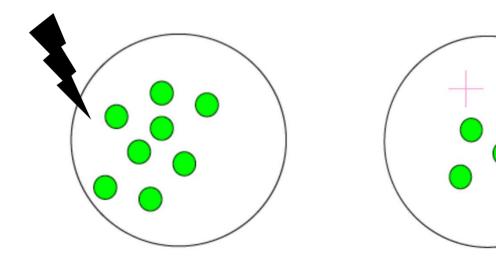
Which set is pure (not mixed)?





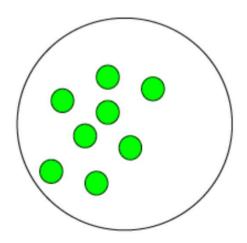
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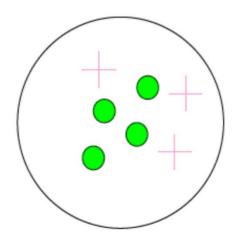


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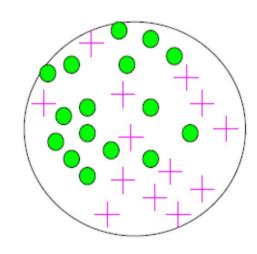
Minimum impurity



Less impure



#### Very impure

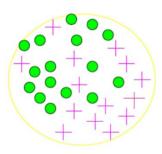


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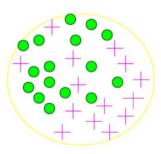
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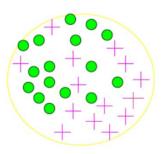
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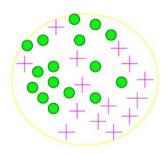


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• It comes from information theory.

Developed by Shannon for code-breaking and secure telecommunications. <a href="http://cm.bell-labs.com/cm/ms/what/shannonday/shannon1948.pdf">http://cm.bell-labs.com/cm/ms/what/shannonday/shannon1948.pdf</a>



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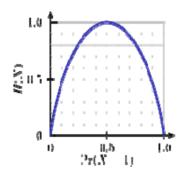
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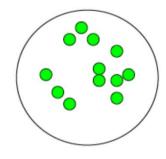


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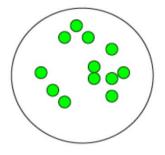
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- The higher the entropy the more the information content.
- What does entail for learning from examples?!

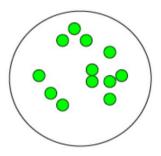
- Let's concentrate on a binary class problem:
  - What is the entropy of a group in which all examples belong to the same class?



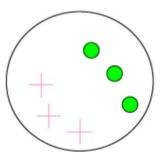
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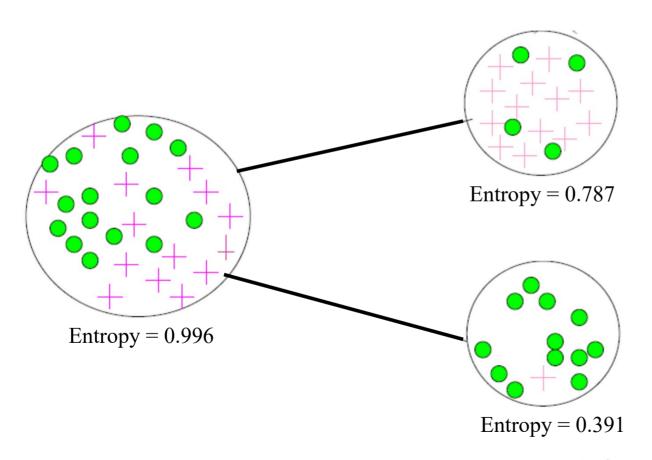
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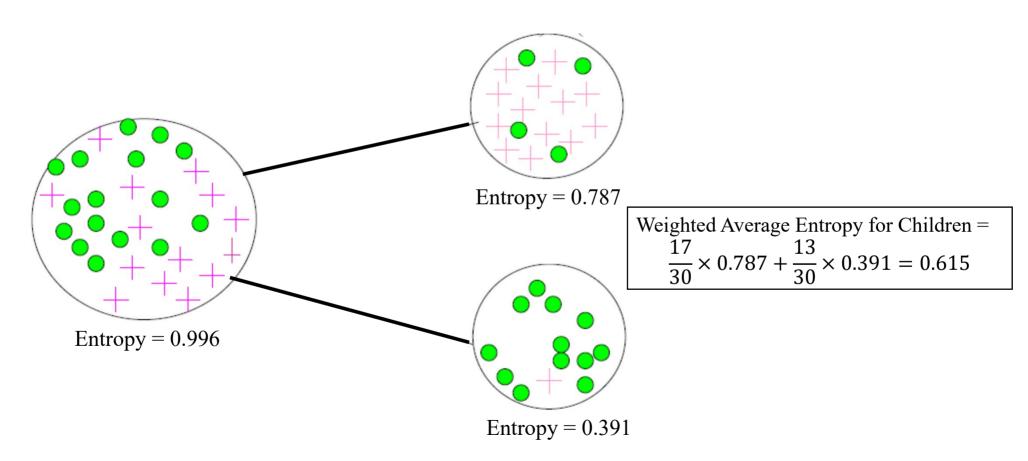


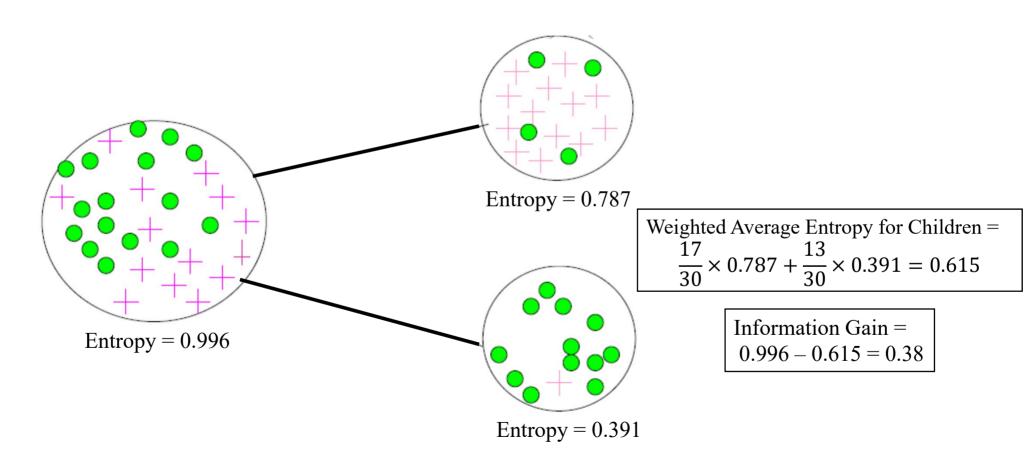
- What is the entropy of a group in which with 50% in either class?
  - The answer is  $E = -0.50 \log_2 0.50 0.50 \log_2 0.50 = 1$



- We would like to find which feature in a given set of features is most discriminative between classes to be learned.
- Information gain tell can tell us how important a given feature is!
- We use that to decide the order of features/decision nodes in a decision tree!





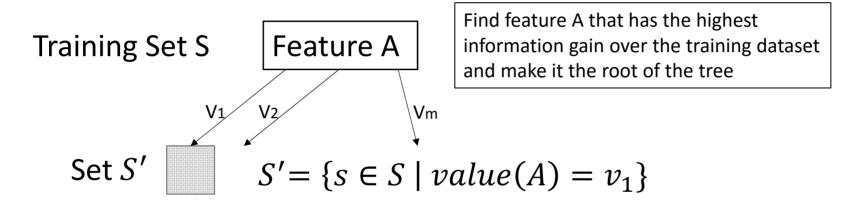


- To construct a decision tree we must answer two questions (reminder):
  - What features and what values must be used?
- ID3 (Iterative Dichotomiser 3) by Quinlan exploits information gain to construct a decision tree!
- Use entropy to calculate information gain and then answer the above questions.

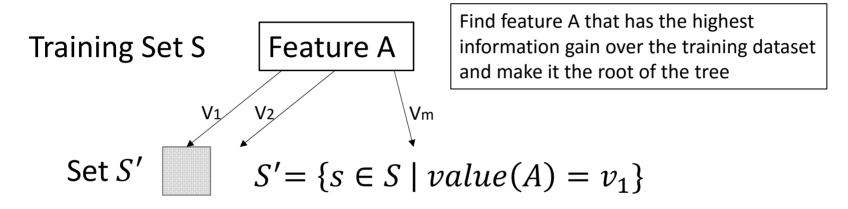
Training Set S

Feature A

Find feature A that has the highest information gain over the training dataset and make it the root of the tree

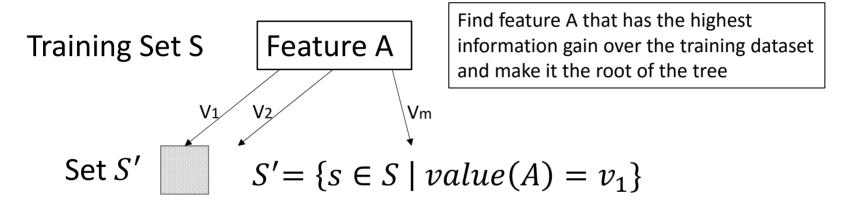


Construct child nodes for each value of A, i.e. find a subset of training samples of which A has certain value.



Construct child nodes for each value of A, i.e. find a subset of training samples of which A has certain value.

Repeat recursively until certain conditions are met!



Construct child nodes for each value of A, i.e. find a subset of training samples of which A has certain value.

Until all the elements of the S' are the same, or there is no more feature to be used!

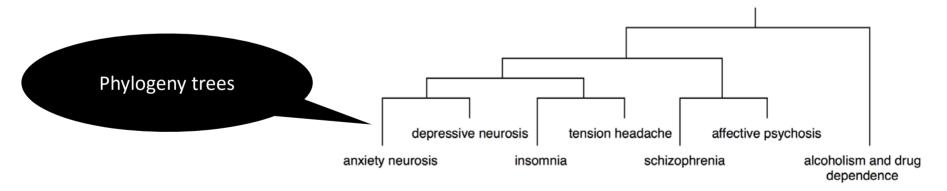
```
>>> import math def entropy(labels):
    freqdist = nltk.FreqDist(labels)
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>>> print(entropy(['male', 'male', 'male', 'male']))
0.0
>>> print(entropy(['male', 'female', 'male', 'male']))
0.811
```

#### Decision Trees: a summary

- Some of Advantages:
  - Decision trees are easy to understand and interpret.
  - Decision trees are very well suited for hierarchical classification task. distinctions can be made.



#### Decision Trees: a summary

- Some of Disadvantages:
  - Decision trees are prone to overfitting, specifically at lower decision nodes (can you explain why?); although some solutions are available:
    - Stop dividing nodes once the amount of training data becomes too small.
    - Using **pruning**, i.e. reduces the size of decision trees by removing nodes that do not improve performance on a dev-test.

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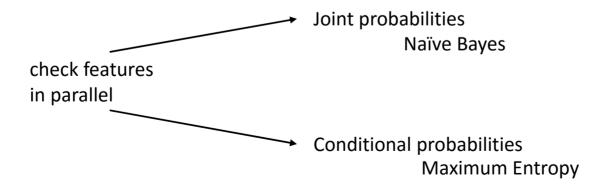
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    - Stop dividing nodes once the amount of training data becomes too small.
    - Using **pruning**, i.e. reduces the size of decision trees by removing nodes that do not improve performance on a dev-test.
  - They force features to be checked in a particular order, which often results in huge decision trees:
    - E.g. in document classification some words are strong indicative of labels independently. These words are pushed to the bottom of the tree and repeated in different branch (an exponential growth!).

#### Alternatives to Decision Trees

- As discussed, decision trees check features in a specific order, one at a time and that is not desirable!
- We can however check features in parallel!
- To do so, we can use joint probabilities as well as conditional probabilities!

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 To review these concepts, we start with the famous "Sex, Math and English" example

	Math	English	Total
Female	1	17	18
Male	37	20	57
Total	38	37	75

Acknowledgement: Slides for this section are based on materials from Mark Liberman and Stephen Isard (see <a href="http://www.ling.upenn.edu/courses/cogs501">http://www.ling.upenn.edu/courses/cogs501</a>)

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Number of enrolled students by their gender

P	fema	عا	Math)	_	013
$\Gamma$	IEIIIa	IE,	ivia (i i <i>)</i>		.UIS

P(female, English) = .227

P(male, Math) = .493

P(male, English) = .267

which are symmetric!

"joint probabilities",

probability of picking a female math professor!

	Math	English	Total
Female	.013	.227	.240
Male	.493	.267	.760
Total	.506	.494	1.00

"joint distribution" of sex and department: probabilities of picking a female or male for each subject in the whole set

Text Mining Project --- Behrang QasemiZadeh ©

• Let's start with a little game, we want to guess the sex:

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P(male) = 57/75 = .760
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P(male | math) = 37/38 = .974
P(math | male) = 37/57 = .649
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• Let's start with a little game, we want to guess the sex:

	Math	English	Total
Female	1	17	18
Male	37	20	57
Total	38	37	75

 $P(male \mid math) = 37/38 = .974$ 

The probability of bumping into a male professor in the math department!

• Let's start with a little game, we want to

	Math	English	Total
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Male	37	20	57
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The probability of bumping into a math professor in the male employees of the department! Formally, read it as "the probability of *male* given *math*"

$$P(math \mid male) = 37/57 = .649$$

• Let's start with a little game, we want to guess the sex:

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Male	37	20	57
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These probabilities are different because they represent different assumptions!

 Please think about these numbers, write down the table and calculate the numbers again!

 Would it be possible to calculate conditional probabilities from the joint distributions?!

	Math	English	Total
Female	.013	.227	.240
Male	.493	.267	.760
Total	.506	.494	1.00

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P(A | B) = P(A, B) / P(B)
```

## Bayes' Theorem

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Using simple math [1] and [2] can be written as:

$$P(A \mid B) P(B) = P(A, B)$$

$$P(B \mid A) P(A) = P(B, A)$$

#### Bayes' Theorem

• So far we know that:

[1] 
$$P(A \mid B) = P(A, B) / P(B)$$

[2] 
$$P(B \mid A) = P(B, A) / P(A)$$

[3] 
$$P(A, B) = P(B, A)$$

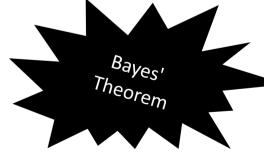
Using simple math [1] and [2] can be written as:

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Using [3] we arrive to:

$$P(A \mid B) P(B) = P(B \mid A) P(A)$$
 and thus  $P(A \mid B) = P(B \mid A) P(A) / P(B)$ 



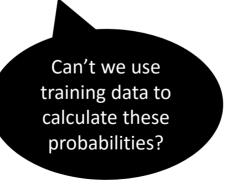
- Think about the relationship between evidence and theory, or feature and class label!
- Suppose we have a set of class labels C<sub>1</sub>, C<sub>2</sub>, ..., and we've observed some features.
- We'd like to pick the class label that is more likely to be true given our observations.

- Think about the relationship between evidence and theory, or feature and class label!
- Suppose we have a set of class labels C<sub>1</sub>, C<sub>2</sub>, ..., and we've observed some features.
- We'd like to pick the class label that is more likely to be true given our observations.
- This can be formulated by conditional probability **P(C | E)**, i.e. the probability of *class label* given *observed features*.

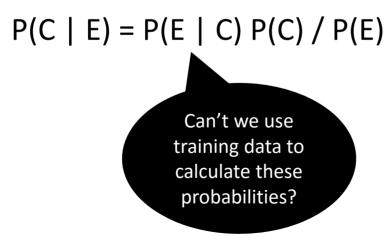
- The classification problem can be solved by calculating  $P(C \mid E)$  for each class label and picking the maximum  $P(C \mid E)$ .
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$$P(C \mid E) = P(E \mid C) P(C) / P(E)$$



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$$P(C \mid E) = P(E \mid T) P(T) / P(E)$$
  
**ARGMAX<sub>i</sub> P(E | C<sub>i</sub>) P(C<sub>i</sub>)**

The best class label accordingly!

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ARGMAX<sub>i</sub>  $P(E \mid C_i) P(C_i)$ 

The best cla accordir

Bayes rule is a mathematical formulation of Hermann von Helmholtz statement: what we perceive is our "best guess" given both sensory data and our prior experience.

- What if we have a set of features instead of one?
  - i.e. if we want to decide class labels by observation made over a feature set?
  - Or, how to calculate  $P(C_i \mid E_1 ... E_n)$ ?!

- What if we have a set of features instead of one?
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  - Or, how to calculate  $P(C_i \mid E_1 \dots E_n)$ ?!
- Let's use the Bayes theorem and the chain rule:

```
\begin{split} P(C_{i} \mid E_{1} \dots E_{n}) &= P(C_{i}) \ P(E_{1} \dots E_{n} \mid C_{i}) \\ &= P(C_{i}) \ P(E_{1} \mid C_{i}) \ P(E_{2} \dots E_{n} \mid C_{i,} E_{1}) \\ &= P(C_{i}) \ P(E_{1} \mid C_{i}) \ P(E_{2} \mid C_{i,} E_{1}) \ P(E_{3} \dots E_{n} \mid C_{i,} E_{1,} E_{2}) \\ &= P(C_{i}) \ P(E_{1} \mid C_{i}) \ P(E_{2} \mid C_{i,} E_{1}) \dots \ P(E_{n} \mid C_{i,} E_{1,} E_{2, \dots,} E_{n-1}) \end{split}
```

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• Let's use the Bayes theorem and the chain rule:

 $P(C_{i} | E_{1} ... E_{n}) = P(C_{i}) P(E_{1} ... E_{n} | C_{i})$   $= P(C_{i}) P(E_{1} | C_{i}) P(E_{2} ... E_{n} | C_{i}, E_{1})$   $= P(C_{i}) P(E_{1} | C_{i}) P(E_{2} | C_{i}, E_{1}) P(E_{3} ... E_{n} | C_{i}, E_{1})$   $= P(C_{i}) P(E_{1} | C_{i}) P(E_{2} | C_{i}, E_{1}) ... P(E_{n} | C_{i}, E_{1}, E_{2}, ..., E_{n-1})$ 

Here naivety can

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• Or, how to calculate  $P(C_i \mid E_1 ... \mid E_i \mid$ • Let's use the Bayes theorem and the conditionally independent of every other **E**; for i≠j

```
P(C_i | E_1 ... E_n) = P(C_i) P(E_1 ... E_n | C_i)
                    = P(C_i) P(E_1 | C_i) P(E_2 ... E_n | C_i E_1)
                    = P(C_i) P(E_1 | C_i) P(E_2 | C_i, E_1) P(E_3 ... E_n | C_i, E_1, E_2)
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• Or, how to calculate  $P(C_i \mid E_1 ... \mid E_i$  Let's assume that each feature  $E_i$  is • Let's use the Bayes theorem and the conditionally independent of every other **E**; for i≠j

$$P(C_{i} | E_{1} ... E_{n}) = P(C_{i}) P(E_{1} ... E_{n} | C_{i})$$

$$= P(C_{i}) P(E_{1} | C_{i}) P(E_{2} ... E_{n} | C_{i}, E_{1})$$

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$$= P(E_{2} | C_{i})$$

$$= P(E_{n} | C_{i})$$

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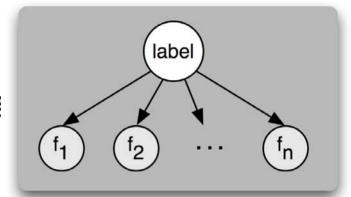
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                           = P(C_i) P(E_1 | C_i) P(E_2 | C_{i, E_1}) ... P(E_n | C_{i, E_{1, E_{2, ..., E_{n-1}}})
under our naïve
  assumption
                            \approx P(C_i) \prod_{i=1}^n P(E_i|C)
   (independence
    assumption)
```

 Now that we have a clear picture of what is going on (hopefully), the classification task can be easily formalized by:

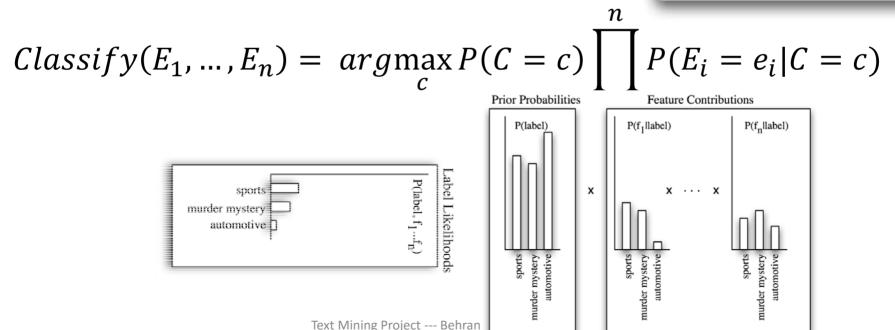
Classify
$$(E_1, ..., E_n) = arg \max_{c} P(C = c) \prod_{i=1}^{n} P(E_i = e_i | C = c)$$

# Naive Bayes Classifiers

 Now that we have a clear picture of what is going classification task can be easily formalized by:



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# Problems with Naive Bayes Classifiers (1)

- What if in a training set a certain feature does not appear with a certain class label, i.e. P(E|C) = 0?!
  - The class label likelihood in this situation is 0 and thus regardless of other features, an input will never be assigned this label!
  - Therefore, does a zero count shows an impossible event?!

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  - The class label likelihood in this situation is 0 and thus regardless of other features, an input will never be assigned this label!
  - Therefore, does a zero count shows an impossible event?!
  - For several reasons, the answer can be "NO"!
- Use smoothing to address the limitation stated above:
  - Estimate the value of P(E|C) using techniques other than simple counting:
    - Additive counting (generally a horrible choice!!!), Good-Tring smoothing, ...
  - nltk.probability implements a number of smoothing methods.
  - Also, see Bill MacCartney's tutorial slides on smoothing (goo.gl/9LCfHE)

# Problems with Naive Bayes Classifiers (2)

- What if we want to decide about labels based on multiple features?
  - Put it simply, we question the naïve independence assumption!
  - In other words, using the Naïve Bayes Classifiers we cannot incorporate complex features when making decisions about class labels.

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- For example:
  - Is it reasonable to assume independence between Bavaria and Weißbier?
  - Or, is it reasonable to assume that the features *ends-with(a)* and *ends-with(vowel)* are independent from each other?.

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- Bird et. al. describe this as the double counting-counting problem!

## **Double-Counting Problem**

- During training, the contribution of dependant features are computed separately, i.e. the calculation of P(feature, label);
- But when using the classifier to choose labels for new inputs, feature contributions are combined, i.e. we use  $\prod P(feature, label)$ .

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#### Solution:

- Do not assume the conditional independence between features!
- Put it simply, this statement means "do not use Naïve Bayes Classifiers"!

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- Solution:
  - Do not assume the conditional independence between features!
  - Put it simply, this statement means "do not use Naïve Bayes Classifiers"!
- Use Maximum Entropy Classifiers (a logistic regression technique).

#### Maximum Entropy Models

 What is the discarding of the assumption of independence in Naïve Bayes Approach means?

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```

Lets continue with Mark Johnson's slides (goo.gl/k4TYnd)

Next Session: syntax and parsing!