

# Categorization

#### • Given:

- A description of an instance,  $x \in X$ , where X is the *instance language* or *instance space*.
- A fixed set of categories:  $C = \{c_1, c_2, \dots, c_n\}$
- Determine:
  - The category of  $x: c(x) \in C$ , where c(x) is a categorization function whose domain is X and whose range is C.

# Learning for Categorization

- A training example is an instance *x*∈*X*, paired with its correct category *c*(*x*):
   <*x*, *c*(*x*)> for an unknown categorization function, *c*.
- Given a set of training examples, D.
- Find a hypothesized categorization function, *h*(*x*), such that:

 $\forall < x, c(x) > \in D : h(x) = c(x)$ Consistency

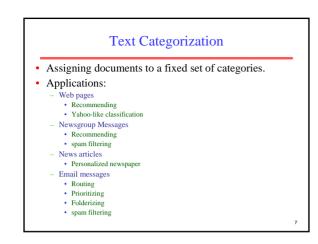
#### Sample Category Learning Problem • Instance language: <size, color, shape> - size $\in$ {small, medium, large} - color $\in$ {red, blue, green} - shape $\in$ {square, circle, triangle} $C = \{\text{positive, negative}\}$ • D: Example Size Color Shape Category small red circle positive 1 2 red circle positive large 3 small red triangle negative blue 4 large circle negative

#### General Learning Issues

- Many hypotheses are usually consistent with the training data.
- Bias
  - Any criteria other than consistency with the training data that is used to select a hypothesis.
- Classification accuracy (% of instances classified correctly).
  - Measured on independent test data.
- Training time (efficiency of training algorithm).
- Testing time (efficiency of subsequent classification).

# Generalization

- Hypotheses must generalize to correctly classify instances not in the training data.
- Simply memorizing training examples is a consistent hypothesis that does not generalize.
- Occam's razor:
  - Finding a *simple* hypothesis helps ensure generalization.



#### Learning for Text Categorization

- · Manual development of text categorization functions is difficult.
- Learning Algorithms:
  - Bayesian (naïve)
  - Neural network
  - Relevance Feedback (Rocchio)
  - Rule based (Ripper)
  - Nearest Neighbor (case based)
  - Support Vector Machines (SVM)

#### Using Relevance Feedback (Rocchio)

- Relevance feedback methods can be adapted for text categorization.
- Use standard TF/IDF weighted vectors to represent text documents (normalized by maximum term frequency).
- For each category, compute a *prototype* vector by • summing the vectors of the training documents in the category.
- Assign test documents to the category with the closest prototype vector based on cosine similarity.

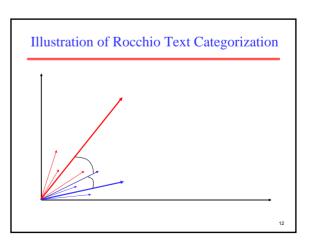
# Rocchio Text Categorization Algorithm (Training)

Assume the set of categories is  $\{c_1, c_2, \dots c_n\}$ For *i* from 1 to *n* let  $\mathbf{p}_i = \langle 0, 0, \dots, 0 \rangle$  (*init. prototype vectors*) For each training example  $\langle x, c(x) \rangle \in D$ Let **d** be the frequency normalized TF/IDF term vector for doc xLet i = j:  $(c_i = c(x))$ (sum all the document vectors in  $c_i$  to get  $\mathbf{p}_i$ ) Let  $\mathbf{p}_i = \mathbf{p}_i + \mathbf{d}$ 

#### Rocchio Text Categorization Algorithm (Test) Given test document x Let **d** be the TF/IDF weighted term vector for xLet m = -2 (*init. maximum cosSim*) For *i* from 1 to *n*: (compute similarity to prototype vector) Let $s = \cos \operatorname{Sim}(\mathbf{d}, \mathbf{p}_i)$ if s > mlet m = s



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# **Rocchio Properties**

- Does not guarantee a consistent hypothesis.
- Forms a simple generalization of the examples in each class (a *prototype*).
- Prototype vector does not need to be averaged or otherwise normalized for length since cosine similarity is insensitive to vector length.
- Classification is based on similarity to class prototypes.

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#### **Rocchio Time Complexity**

- Note: The time to add two sparse vectors is proportional to minimum number of non-zero entries in the two vectors.
- Training Time:  $O(|D|(L_d + |V_d|)) = O(|D||L_d)$ where  $L_d$  is the average length of a document in D and  $V_d$ is the average vocabulary size for a document in D.
- Test Time: O(L<sub>t</sub> + /C/|V<sub>t</sub>/) where L<sub>t</sub> is the average length of a test document and |V<sub>t</sub>/ is the average vocabulary size for a test document.
   Assumes lengths of p, vectors are computed and stored during
  - training, allowing  $\cos Sim(\mathbf{d}, \mathbf{p}_i)$  to be computed in time proportional to the number of non-zero entries in  $\mathbf{d}$  (i.e.  $|V_i|$ )

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#### Nearest-Neighbor Learning Algorithm

- Learning is just storing the representations of the training examples in *D*.
- Testing instance x:
  - Compute similarity between *x* and all examples in *D*.
  - Assign x the category of the most similar example in D.
- Does not explicitly compute a generalization or category prototypes.
- Also called:
  - Case-based
  - Memory-based
  - Lazy learning

# K Nearest-Neighbor Using only the closest example to determine categorization is subject to errors due to: A single atypical example. Noise (i.e. error) in the category label of a single training example. More robust alternative is to find the *k* most-similar examples and return the majority category of these *k* examples. Value of *k* is typically odd to avoid ties, 3 and 5 are most common.

# Similarity Metrics

- Nearest neighbor method depends on a similarity (or distance) metric.
- Simplest for continuous *m*-dimensional instance space is *Euclidian distance*.
- Simplest for *m*-dimensional binary instance space is *Hamming distance* (number of feature values that differ).
- For text, cosine similarity of TF-IDF weighted vectors is typically most effective.

3 Nearest Neighbor Illustration (Euclidian Distance)

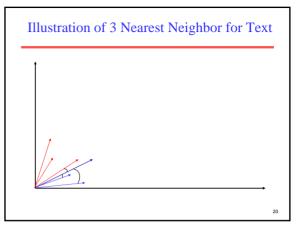
# K Nearest Neighbor for Text

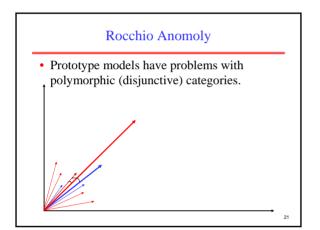
#### Training:

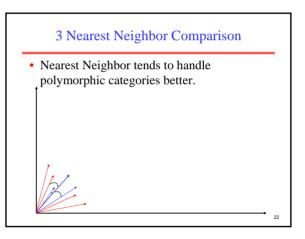
For each training example  $\langle x, c(x) \rangle \in D$ Compute the corresponding TF-IDF vector,  $\mathbf{d}_x$ , for document *x* 

#### Test instance y:

Compute TF-IDF vector **d** for document y For each  $\langle x, c(x) \rangle \in D$ Let  $s_x = \cos Sim(\mathbf{d}, \mathbf{d}_x)$ Sort examples, x, in D by decreasing value of  $s_x$ Let N be the first k examples in D. (get most similar neighbors) Return the majority class of examples in N







#### Nearest Neighbor Time Complexity

- Training Time:  $O(|D| L_d)$  to compose TF-IDF vectors.
- Testing Time:  $O(L_t + /D//V_t/)$  to compare to all training vectors.
  - Assumes lengths of d<sub>x</sub> vectors are computed and stored during training, allowing cosSim(d, d<sub>x</sub>) to be computed in time proportional to the number of non-zero entries in d (i.e. /V<sub>t</sub>)
- Testing time can be high for large training sets.

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# Nearest Neighbor with Inverted Index

- Determining *k* nearest neighbors is the same as determining the *k* best retrievals using the test document as a query to a database of training documents.
- Testing Time: O(B/V<sub>t</sub>/) where *B* is the average number of training documents in which a test-document word appears.
- Therefore, overall classification is O(L<sub>1</sub> + B/V<sub>1</sub>/)
   Typically B << |D|</li>

#### **Bayesian Methods**

- Learning and classification methods based on probability theory.
- Bayes theorem plays a critical role in probabilistic learning and classification.
- Uses *prior* probability of each category given no information about an item.
- Categorization produces a *posterior* probability distribution over the possible categories given a description of an item.

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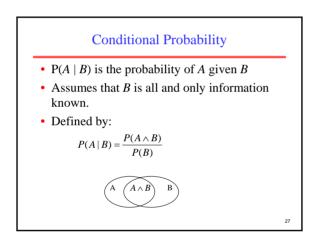
#### Axioms of Probability Theory

- All probabilities between 0 and 1  $0 \le P(A) \le 1$
- True proposition has probability 1, false has probability 0.

P(true) = 1 P(false) = 0.

• The probability of disjunction is:  $P(A \lor B) = P(A) + P(B) - P(A \land B)$ 

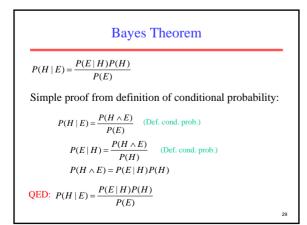
 $(A \land B)$ 

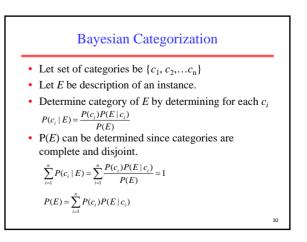


# Independence

- A and B are *independent* iff: P(A | B) = P(A) P(B | A) = P(B)These two constraints are logically equivalent
- Therefore, if *A* and *B* are independent:

$$P(A | B) = \frac{P(A \land B)}{P(B)} = P(A)$$
$$P(A \land B) = P(A)P(B)$$





#### Bayesian Categorization (cont.)

- Need to know:
  - Priors:  $P(c_i)$
  - Conditionals:  $P(E | c_i)$
- P(c<sub>i</sub>) are easily estimated from data.
   If n<sub>i</sub> of the examples in D are in c<sub>i</sub>, then P(c<sub>i</sub>) = n<sub>i</sub>/|D/
- Assume instance is a conjunction of binary features:

 $E = e_1 \wedge e_2 \wedge \cdots \wedge e_m$ 

• Too many possible instances (exponential in *m*) to estimate all P(*E* | *c<sub>i</sub>*)

#### Naïve Bayesian Categorization

• If we assume features of an instance are independent given the category (*c<sub>i</sub>*) (*conditionally independent*).

 $P(E \mid c_i) = P(e_1 \land e_2 \land \dots \land e_m \mid c_i) = \prod_{i=1}^m P(e_i \mid c_i)$ 

• Therefore, we then only need to know  $P(e_i | c_i)$  for each feature and category.

= {allergy	cold	woll)	
- {anergy	, colu,	wenj	
sneeze;	$e_2 = co$	ugh; $e_3$	= feve
{sneeze,	cough	feve	r]
- Usineeze,	cougn	, 1000	IJ
Prob	Well	Cold	Allergy
<b>Prob</b> P(c <sub>i</sub> )	<b>Well</b> 0.9	<b>Cold</b> 0.05	Allergy 0.05
$P(c_i)$	0.9	0.05	0.05

Probability	Well	Cold	Allergy	
$P(c_i)$	0.9	0.05	0.05	7
$P(\text{sneeze} \mid c_i)$	0.1	0.9	0.9	E={sneeze, cough, ¬fever}
$P(\text{cough}   c_i)$	0.1	0.8	0.7	-
P(fever   c <sub>i</sub> )	0.01	0.7	0.4	1
$P(\text{well} \mid \text{E}) = (0.9)$		, , ,		
P(cold   E) = (0.0) P(allergy   E) = (0.0)	0.05)(0.9)(0.8) 0.05)(0.9)(0.	(0.3)/P(E)=0 7)(0.6)/P(E)	01/P(E)	
$P(\text{cold} \mid \text{E}) = (0.0)$	05)(0.9)(0.8) 0.05)(0.9)(0. ategory: aller	(0.3)/P(E)=0 7)(0.6)/P(E)	01/P(E)	

#### **Estimating Probabilities**

- Normally, probabilities are estimated based on observed frequencies in the training data.
- If *D* contains *n<sub>i</sub>* examples in category *c<sub>i</sub>*, and *n<sub>ij</sub>* of these *n<sub>i</sub>* examples contains feature *e<sub>j</sub>*, then:

$$P(e_j \mid c_i) = \frac{n}{n}$$

- However, estimating such probabilities from small training sets is error-prone.
- If due only to chance, a rare feature,  $e_k$ , is always false in the training data,  $\forall c_i : P(e_k | c_i) = 0$ .
- If  $e_k$  then occurs in a test example, E, the result is that  $\forall c_i$ :  $P(E \mid c_i) = 0$  and  $\forall c_i$ :  $P(c_i \mid E) = 0$

Smoothing
To account for estimation from small samples, probability estimates are adjusted or *smoothed*.
Laplace smoothing using an *m*-estimate assumes that each feature is given a prior probability, *p*, that is assumed to have been previously observed in a "virtual" sample of size *m*.

$$P(e_j \mid c_i) = \frac{n_{ij} + mp}{n_i + m}$$

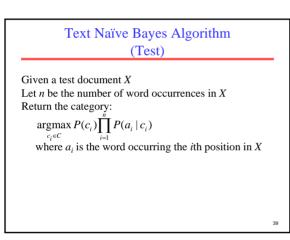
• For binary features, p is simply assumed to be 0.5.

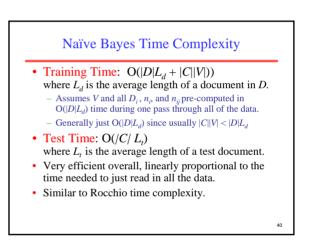
# Naïve Bayes for Text

- Modeled as generating a bag of words for a document in a given category by repeatedly sampling with replacement from a vocabulary  $V = \{w_1, w_2, \dots, w_m\}$  based on the probabilities  $P(w_i | c_i)$ .
- Smooth probability estimates with Laplace *m*-estimates assuming a uniform distribution over all words (p = 1/|V|) and m = |V|
  - Equivalent to a virtual sample of seeing each word in each category exactly once.

# Text Naïve Bayes Algorithm (Train)

Let *V* be the vocabulary of all words in the documents in *D* For each category  $c_i \in C$ Let  $D_i$  be the subset of documents in *D* in category  $c_i$  $P(c_i) = |D_i| / |D|$ Let  $T_i$  be the concatenation of all the documents in  $D_i$ Let  $n_i$  be the total number of word occurrences in  $T_i$ For each word  $w_j \in V$ Let  $n_{ij}$  be the number of occurrences of  $w_j$  in  $T_i$ Let  $P(w_i | c_i) = (n_{ij} + 1) / (n_i + |V|)$ 





# **Underflow Prevention**

- Multiplying lots of probabilities, which are between 0 and 1 by definition, can result in floating-point underflow.
- Since log(xy) = log(x) + log(y), it is better to perform all computations by summing logs of probabilities rather than multiplying probabilities.
- Class with highest final un-normalized log probability score is still the most probable.

Naïve Bayes Posterior Probabilities

- Classification results of naïve Bayes (the class with maximum posterior probability) are usually fairly accurate.
- However, due to the inadequacy of the conditional independence assumption, the actual posterior-probability numerical estimates are not.
  - Output probabilities are generally very close to 0 or 1.

# **Evaluating Categorization**

- Evaluation must be done on test data that are independent of the training data (usually a disjoint set of instances).
- Classification accuracy: c/n where n is the total number of test instances and c is the number of test instances correctly classified by the system.
- Results can vary based on sampling error due to different training and test sets.
- Average results over multiple training and test sets (splits of the overall data) for the best results.

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#### N-Fold Cross-Validation

- Ideally, test and training sets are independent on each trial.
  - But this would require too much labeled data.
- Partition data into *N* equal-sized disjoint segments.
- Run *N* trials, each time using a different segment of the data for testing, and training on the remaining *N*-1 segments.
- This way, at least test-sets are independent.
- Report average classification accuracy over the *N* trials.
- Typically, N = 10.

#### Learning Curves

- In practice, labeled data is usually rare and expensive.
- Would like to know how performance varies with the number of training instances.
- *Learning curves* plot classification accuracy on independent test data (*Y* axis) versus number of training examples (*X* axis).

#### *N*-Fold Learning Curves

- Want learning curves averaged over multiple trials.
- Use *N*-fold cross validation to generate *N* full training and test sets.
- For each trial, train on increasing fractions of the training set, measuring accuracy on the test data for each point on the desired learning curve.

