Classification fundamentals



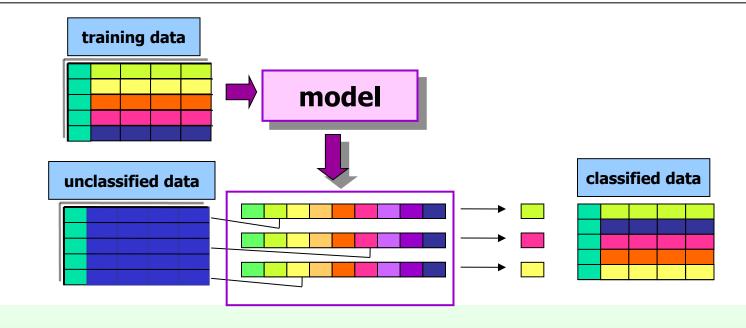
Elena Baralis

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Classification

Objectives

- prediction of a class label
- definition of an interpretable model of a given phenomenon





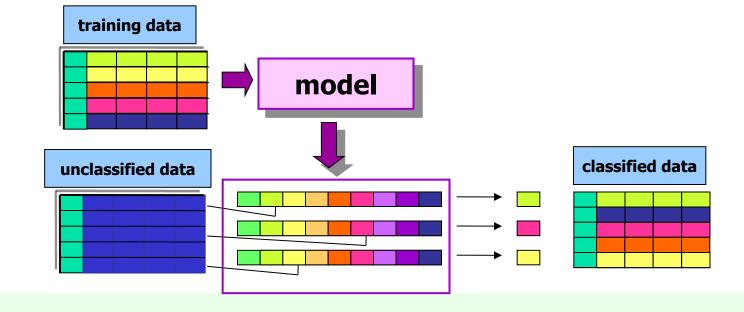


Classification

Applications

- detection of customer propension to leave a company (churn or attrition)
- fraud detection
- classification of different pathology types

• ...







Classification: definition



Given

- a collection of class labels
- a collection of data objects labelled with a class label
- Find a descriptive profile of each class, which will allow the assignment of unlabeled objects to the appropriate class



Definitions



Training set

 Collection of labeled data objects used to learn the classification model

Test set

 Collection of labeled data objects used to validate the classification model



Classification techniques

- Decision trees
- Classification rules
- Association rules
- Neural Networks
- Naïve Bayes and Bayesian Networks
- k-Nearest Neighbours (k-NN)
- Support Vector Machines (SVM)





Evaluation of classification techniques



- Accuracy
 - quality of the prediction
- Interpretability
 - model interpretability
 - model compactness
- Incrementality
 - model update in presence of newly labelled record

- Efficiency
 - model building time
 - classification time
- Scalability
 - training set size
 - attribute number
- Robustness
 - noise, missing data



Decision trees



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Example of decision tree

Marital

Status

Single

Married

Single

Married

Divorced

Married

Divorced

Single

Married

Single

Refund

Yes

No

No

Yes

No

No

Yes

No

No

No

Tid

1

2

3

4

5

6

7

8

9

10

categorical categorical continuous

Taxable

Income

125K

100K

70K

120K

95K

60K

220K

85K

75K

90K

Cheat

No

No

No

No

Yes

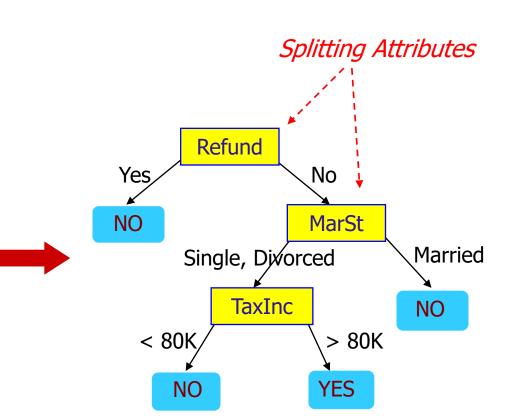
No

No

Yes

No

Yes



Training Data

Model: Decision Tree

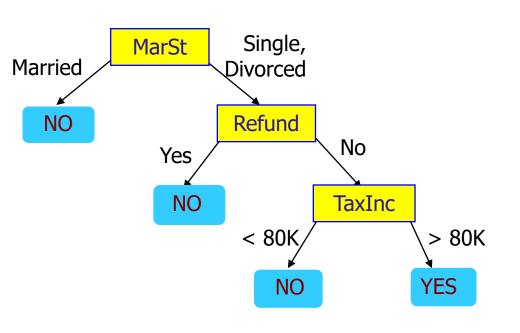




Another example of decision tree



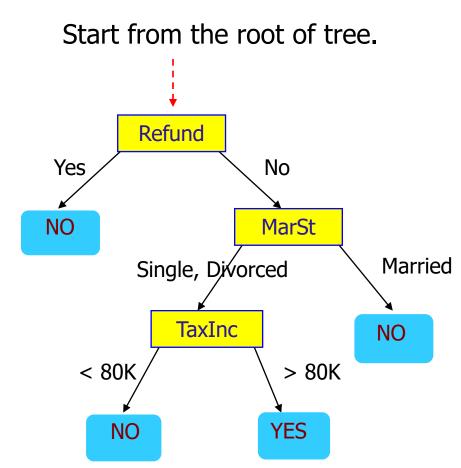
		orical	rical	NOUS
	Cate	egorical cateo	conti	Inuous class
Tid	Refund	Marital Status	Taxable Income	Cheat
1	Yes	Single	125K	No
2	No	Married	100K	No
3	No	Single	70K	No
4	Yes	Married	120K	No
5	No	Divorced	95K	Yes
6	No	Married	60K	No
7	Yes	Divorced	220K	No
8	No	Single	85K	Yes
9	No	Married	75K	No
10	No	Single	90K	Yes



There could be more than one tree that fits the same data!







Test Data

Refund	Marital Status	Taxable Income	Cheat
No	Married	80K	?





Refund Yes No NO MarSt Married Single, Divorced TaxInc NO < 80K > 80K YES NO

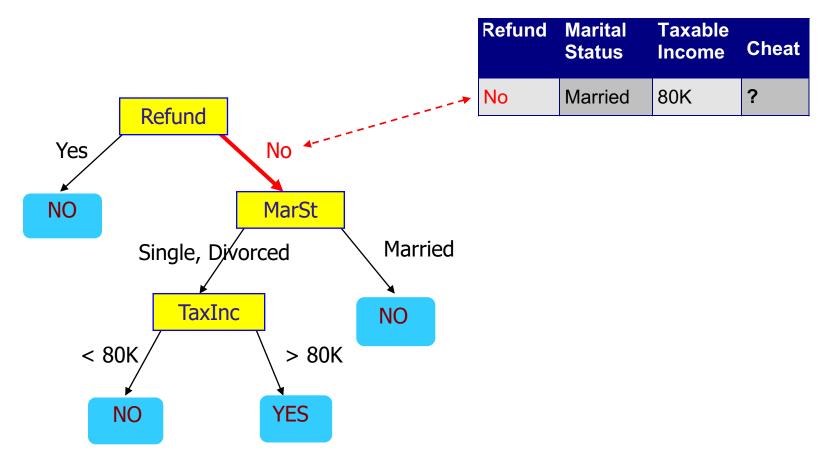
Test Data

*	Refund	Marital Status	Taxable Income	Cheat
	No	Married	80K	?





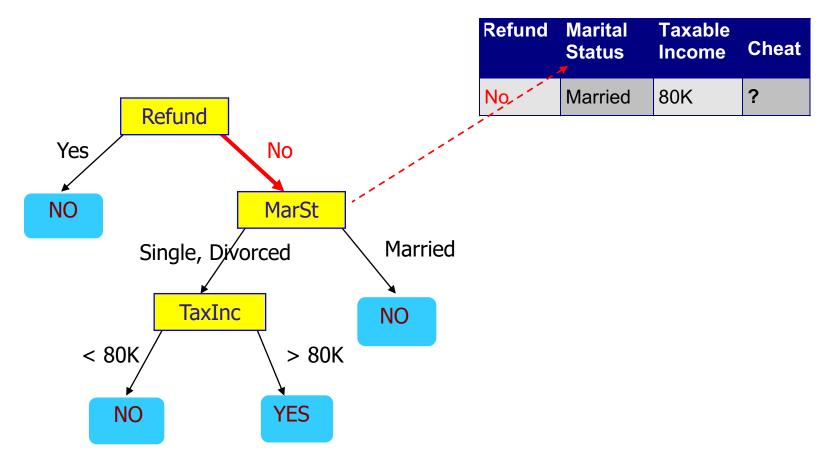
Test Data







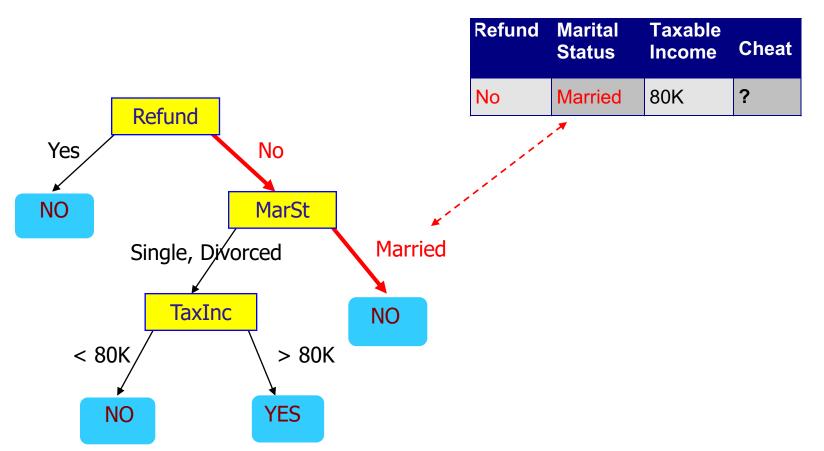
Test Data







Test Data







Test Data Refund Marital **Taxable** Cheat **Status** Income No Married 80K ? Refund Yes No NO MarSt Assign Cheat to "No" Married Single, Divorced TaxInc NO < 80K > 80K YES NO



Decision tree induction



- Many algorithms to build a decision tree
 - Hunt's Algorithm (one of the earliest)
 - CART
 - ID3, C4.5, C5.0
 - SLIQ, SPRINT



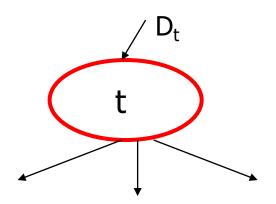
General structure of Hunt's algorithm

Basic steps

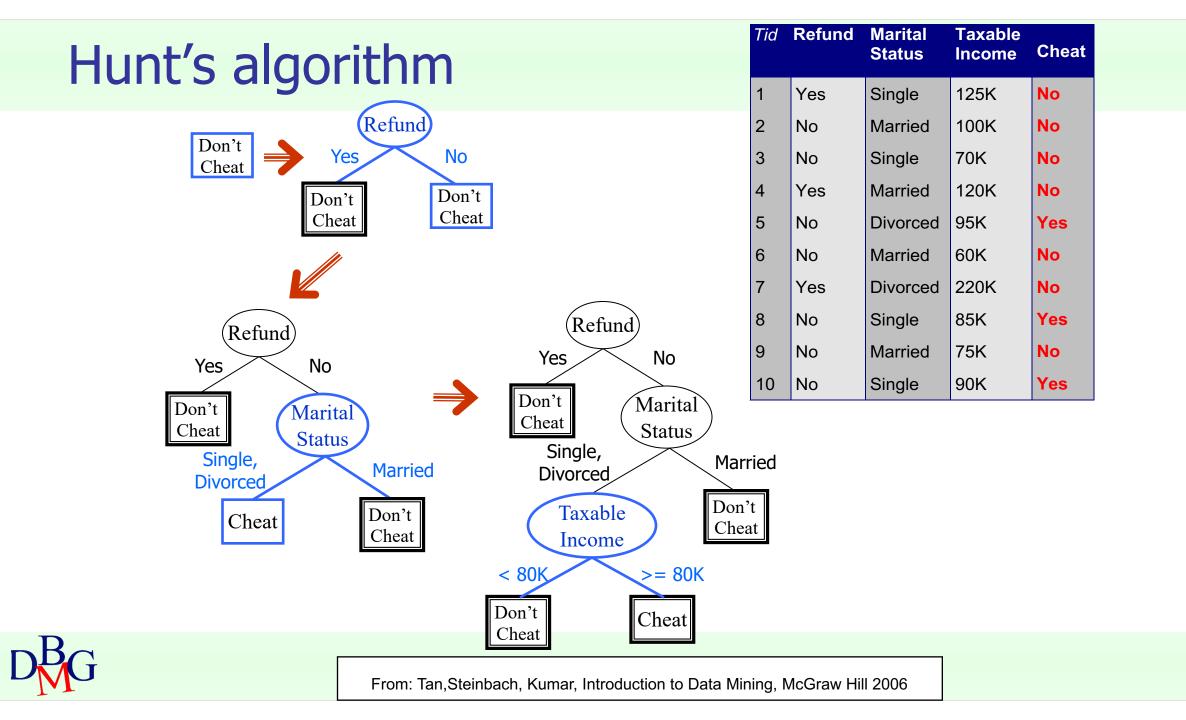
- If D_t contains records that belong to more than one class
 - select the "best" attribute A on which to split D_t and label node t as A
 - split D_t into smaller subsets and recursively apply the procedure to each subset
- If D_t contains records that belong to the same class y_t
 - then t is a leaf node labeled as \boldsymbol{y}_t
- If D_t is an empty set
 - then t is a leaf node labeled as the default (majority) class, y_d



 $D_{t,r}$, set of training records that reach a node t









Decision tree induction



Adopts a greedy strategy

- "Best" attribute for the split is selected locally at each step
 - not a global optimum
- Issues
 - Structure of test condition
 - Binary split versus multiway split
 - Selection of the best attribute for the split
 - Stopping condition for the algorithm



Structure of test condition

Depends on attribute type

- nominal
- ordinal
- continuous

Depends on number of outgoing edges

- 2-way split
- multi-way split





From: Tan, Steinbach, Kumar, Introduction to Data Mining, McGraw Hill 2006

Splitting on nominal attributes

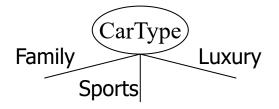
Multi-way split

use as many partitions as distinct values

Binary split

- Divides values into two subsets
- Need to find optimal partitioning







From: Tan, Steinbach, Kumar, Introduction to Data Mining, McGraw Hill 2006

OR

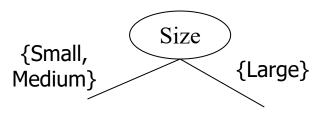
Splitting on ordinal attributes

Multi-way split

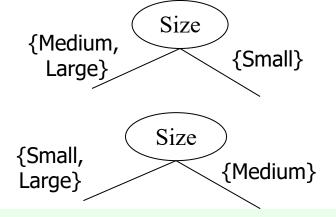
use as many partitions as distinct values

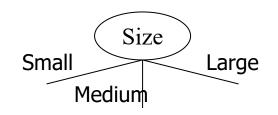
Binary split

- Divides values into two subsets
- Need to find optimal partitioning



What about this split?







Splitting on continuous attributes

Different techniques

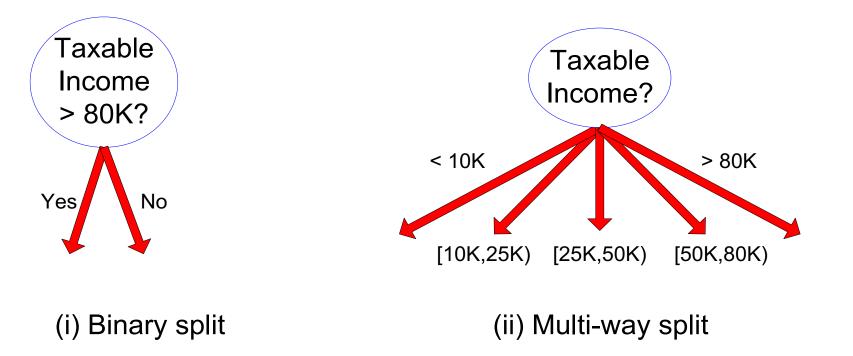
- Discretization to form an ordinal categorical attribute
 - Static discretize once at the beginning
 - Dynamic discretize during tree induction

Ranges can be found by equal interval bucketing, equal frequency bucketing (percentiles), or clustering

- Binary decision (A < v) or $(A \ge v)$
 - consider all possible splits and find the best cut
 - more computationally intensive



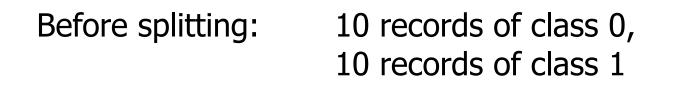
Splitting on continuous attributes

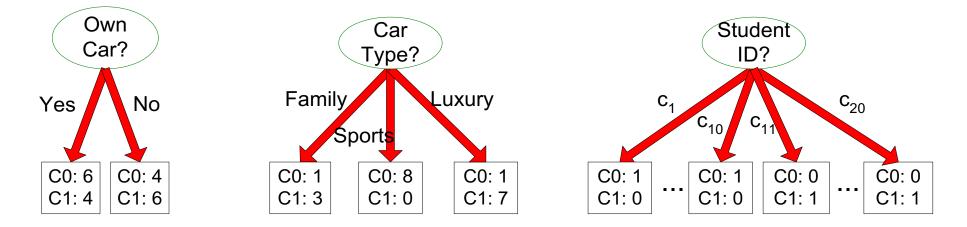




Selection of the best attribute







Which attribute (test condition) is the best?



Selection of the best attribute



- Attributes with *homogeneous* class distribution are preferred
- Need a measure of node impurity



Non-homogeneous, high degree of impurity Homogeneous, low degree of impurity



Measures of node impurity



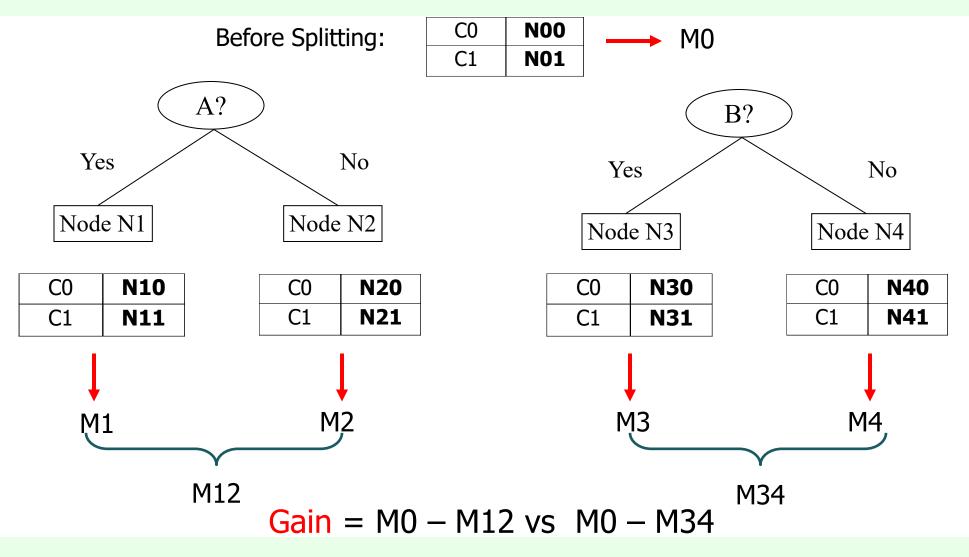
- Many different measures available
 - Gini index
 - Entropy
 - Misclassification error

Different algorithms rely on different measures



How to find the best attribute







GINI impurity measure

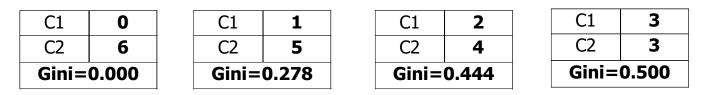


Gini Index for a given node t

$$GINI(t) = 1 - \sum_{j} [p(j | t)]^{2}$$

 $p(j \mid t)$ is the relative frequency of class j at node t

- Maximum (1 1/n_c) when records are equally distributed among all classes, implying higher impurity degree
- Minimum (0.0) when all records belong to one class, implying lower impurity degree





Examples for computing GINI



$$GINI(t) = 1 - \sum_{j} [p(j \mid t)]^2$$

C1	0
C2	6

P(C1) = 0/6 = 0	P(C2) = 6/6 = 1
Gini = $1 - P(C1)^2$ -	$-P(C2)^2 = 1 - 0 - 1 = 0$

C1	1
C2	5

$$P(C1) = 1/6$$
 $P(C2) = 5/6$
Gini = 1 - (1/6)² - (5/6)² = 0.278

C1	2
C2	4

$$P(C1) = 2/6$$
 $P(C2) = 4/6$
Gini = 1 - (2/6)² - (4/6)² = 0.444



Splitting based on GINI

- Used in CART, SLIQ, SPRINT
- When a node p is split into k partitions (children), the quality of the split is computed as

$$GINI_{split} = \sum_{i=1}^{k} \frac{n_i}{n} GINI(i)$$

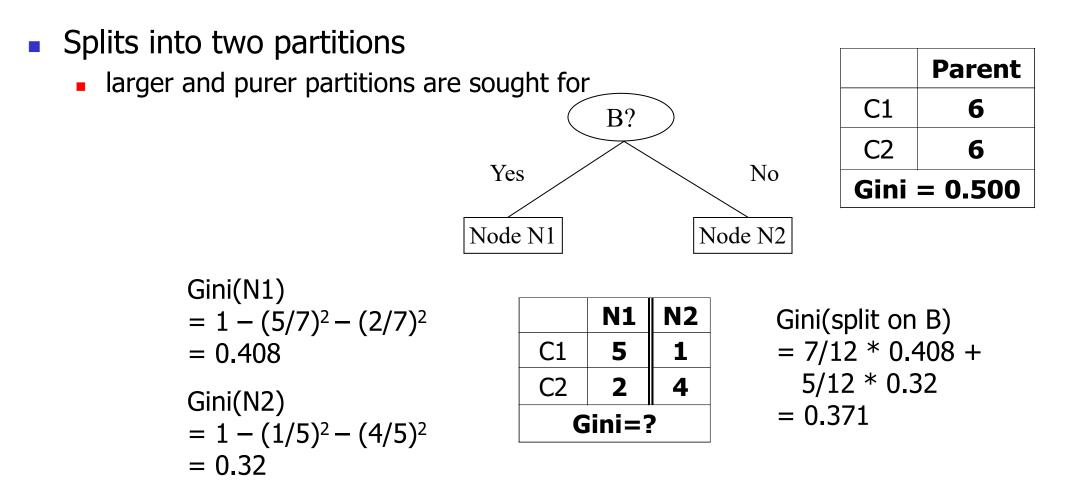
where

- n_i = number of records at child i
- n = number of records at node p





Computing GINI index: Boolean attribute





Computing GINI index: Categorical attribute

- For each distinct value, gather counts for each class in the dataset
- Use the count matrix to make decisions







Computing GINI index: Continuous attribute

- Binary decision on one splitting value
 - Number of possible splitting values
 Number of distinct values
- Each splitting value v has a count matrix
 - class counts in the two partitions
 - A < v
 - $A \ge v$

Tid	Refund	Marital Status	Taxable Income	Cheat		
1	Yes	Single	125K	No		
2	No	Married	100K	No		
3	No	Single	70K	No		
4	Yes	Married 120K		No		
5	No	Divorced	95K	Yes		
6	No	Married	60K	No		
7	Yes	Divorced	220K	No		
8	No	Single	85K	Yes		
9	No	Married	75K	No		
10	No	Single	90K	Yes		





Computing GINI index: Continuous attribute

- For each attribute
 - Sort the attribute on values
 - Linearly scan these values, each time updating the count matrix and computing gini index
 - Choose the split position that has the least gini index

	Cheat	No No		No		Ye	es Yes		S	s Yes		N	No N		0	N	lo		No				
		Taxable Income																					
Sorted Values			60		70)	7	5	85	5	9()	9	5	1()0	12	20	12	25		220	
Split Positions		5	5	6	5	7	2	8	0	8	7	9	2	9	7	11	0	12	22	17	2	23	0
opine i obliciono		<=	>	<=	>	<=	>	<=	>	<=	>	<=	>	<=	>	<=	>	<=	>	<=	>	<=	>
	Yes	0	3	0	3	0	3	0	3	1	2	2	1	3	0	3	0	3	0	3	0	3	0
	No	0	7	1	6	2	5	3	4	3	4	3	4	3	4	4	3	5	2	6	1	7	0
	Gini	0.420 0.400 0.375		0.343 0.			0.417		.417 0.400		0.400 <u>0.</u>		<u>0.300</u>		<u>.300</u> 0.3		43	3 0.37		0.4	100 0.420		20



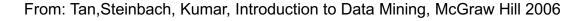


Entropy at a given node t

$$Entropy(t) = -\sum_{j} p(j | t) \log_2 p(j | t)$$

 $p(j \mid t)$ is the relative frequency of class j at node t

- Maximum (log n_c) when records are equally distributed among all classes, implying higher impurity degree
- Minimum (0.0) when all records belong to one class, implying lower impurity degree
- Entropy based computations are similar to GINI index computations



Examples for computing entropy



$$Entropy(t) = -\sum_{j} p(j \mid t) \log_{2} p(j \mid t)$$

C1	0
C2	6

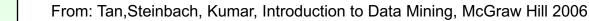
P(C1) = 0/6 = 0 P(C2) = 6/6 = 1Entropy = -0 log 0 - 1 log 1 = -0 - 0 = 0

C1	1
C2	5

P(C1) = 1/6 P(C2) = 5/6Entropy = $-(1/6) \log_2 (1/6) - (5/6) \log_2 (5/6) = 0.65$

C1	2
C2	4

 $P(C1) = 2/6 \qquad P(C2) = 4/6$ Entropy = - (2/6) log₂ (2/6) - (4/6) log₂ (4/6) = 0.92



Splitting Based on INFO

Information Gain

$$GAIN_{split} = Entropy(p) - \left(\sum_{i=1}^{k} \frac{n_i}{n} Entropy(i)\right)$$

Parent Node, p is split into k partitions;

n_i is number of records in partition i

- Measures reduction in entropy achieved because of the split. Choose the split that achieves most reduction (maximizes GAIN)
- Used in ID3 and C4.5
- Disadvantage: Tends to prefer splits yielding a large number of partitions, each small but pure





From: Tan, Steinbach, Kumar, Introduction to Data Mining, McGraw Hill 2006

Splitting Based on INFO

Gain Ratio

$$GainRATIO_{split} = \frac{GAIN_{split}}{SplitINFO}$$
$$SplitINFO = -\sum_{i=1}^{k} \frac{n_i}{n} \log \frac{n_i}{n}$$

Parent Node, p is split into k partitions n_i is the number of records in partition i

- Adjusts Information Gain by the entropy of the partitioning (SplitINFO). Higher entropy partitioning (large number of small partitions) is penalized
- Used in C4.5
- Designed to overcome the disadvantage of Information Gain



Comparison among splitting criteria



For a 2-class problem 0.9 Entropy 0.8 0.7 0.6 0.5 Gini 0.4 0.3 Misclassification 0.2 error 0.1 0 0.2 0.3 0.5 Ο 0.1 0.4 0.6 0.7 0.8 0.9 р



Stopping Criteria for Tree Induction

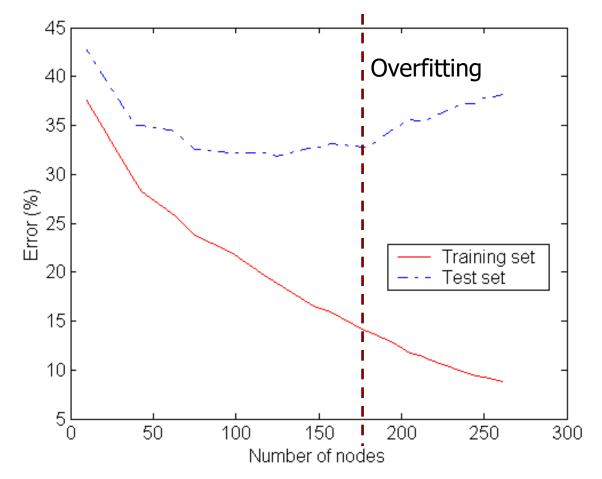


- Stop expanding a node when all the records belong to the same class
- Stop expanding a node when all the records have similar attribute values
- Early termination
 - Pre-pruning
 - Post-pruning



Underfitting and Overfitting



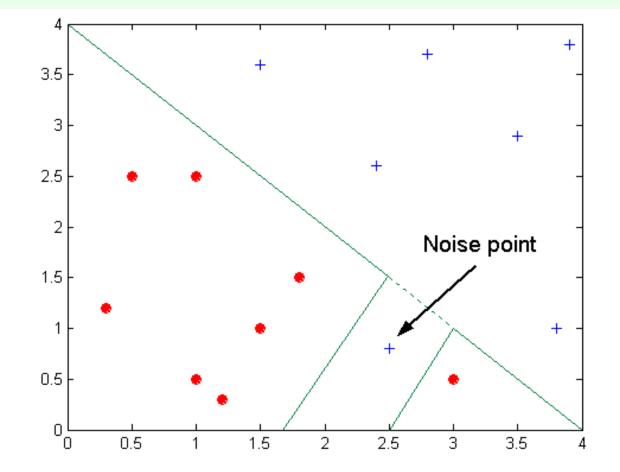


Underfitting: when model is too simple, both training and test errors are large

 $D_{M}^{B}G$

Overfitting due to Noise





Decision boundary is distorted by noise point



How to address overfitting



Pre-Pruning (Early Stopping Rule)

- Stop the algorithm before it becomes a fully-grown tree
- Typical stopping conditions for a node
 - Stop if all instances belong to the same class
 - Stop if all the attribute values are the same
- More restrictive conditions
 - Stop if number of instances is less than some user-specified threshold
 - Stop if class distribution of instances are independent of the available features (e.g., using $\chi^{\,2}$ test)
 - Stop if expanding the current node does not improve impurity measures (e.g., Gini or information gain)



How to address overfitting

11

Post-pruning

- Grow decision tree to its entirety
- Trim the nodes of the decision tree in a bottom-up fashion
- If generalization error improves after trimming, replace sub-tree by a leaf node.
- Class label of leaf node is determined from majority class of instances in the sub-tree



Data fragmentation



- Number of instances gets smaller as you traverse down the tree
- Number of instances at the leaf nodes could be too small to make any statistically significant decision



Handling missing attribute values

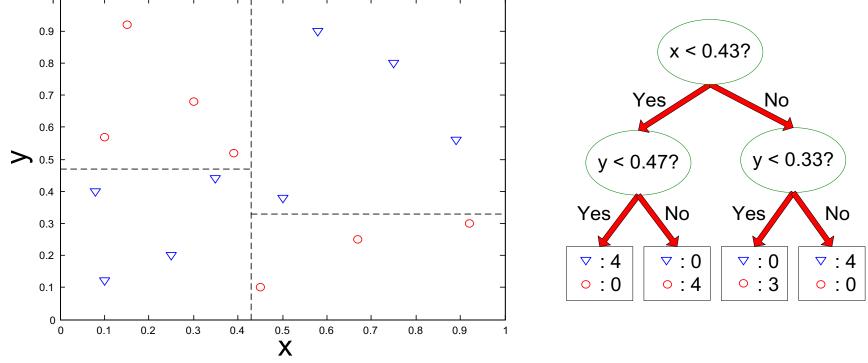


- Missing values affect decision tree construction in three different ways
 - Affect how impurity measures are computed
 - Affect how to distribute instances with missing value to child nodes
 - Affect how a test instance with missing value is classified



Decision boundary



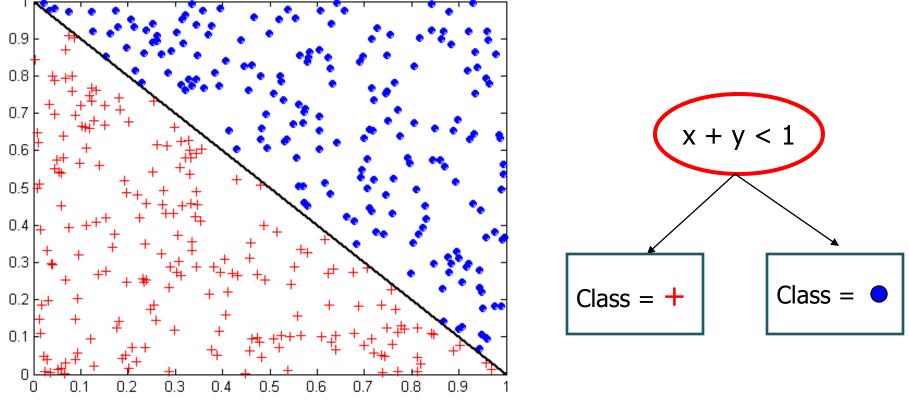


- Border line between two neighboring regions of different classes is known as decision boundary
- Decision boundary is parallel to axes because test condition involves a single attribute at-a-time

 $D_{M}^{B}G$

Oblique decision trees





- Test condition may involve multiple attributes
- More expressive representation
- Finding optimal test condition is computationally expensive

Evaluation of decision trees



Accuracy

- For simple datasets, comparable to other classification techniques
- Interpretability
 - Model is interpretable for small trees
 - Single predictions are interpretable
- Incrementality
 - Not incremental

Efficiency

- Fast model building
- Very fast classification
- Scalability
 - Scalable both in training set size and attribute number
- Robustness
 - Difficult management of missing data



Random Forest



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Random Forest



Ensemble learning technique

- multiple base models are combined
 - to improve accuracy and stability
 - to avoid overfitting

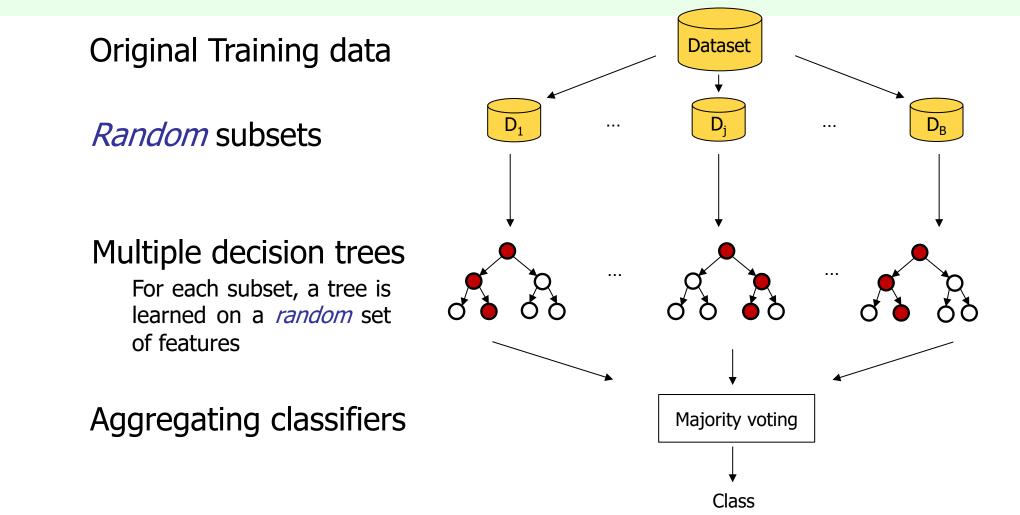
Random forest = set of decision trees

- a number of decision trees are built at training time
- the class is assigned by majority voting



Random Forest







Bootstrap aggregation



- Given a training set D of n instances, it selects B times a random sample with replacement from D and trains trees on these dataset samples
 - For b = 1, ..., B
 - Sample with replacement n' training examples, $n' \le n$
 - A dataset subset *D_b* is generated
 - Train a classification tree on D_b







- Selects, for each candidate split in the learning process, a random subset of the features
 - Being p the number of features, \sqrt{p} features are typically selected
- Trees are decorrelated
 - Feature subsets are sampled randomly, hence different features can be selected as best attributes for the split



Random Forest – Algorithm Recap



- Given a training set D of n instances with p features
- For b = 1, ..., B
 - Sample randomly with replacement n' training examples. A subset D_b is generated
 - Train a classification tree on D_b
 - During the tree construction, for each candidate split
 - $m \ll p$ random features are selected (typically $m \approx \sqrt{p}$)
 - the best split is computed among these *m* features

Class is assigned by majority voting among the B predictions



Evaluation of random forests



- Accuracy
 - Higher than decision trees
- Interpretability
 - Model and prediction are not interpretable
 - A prediction may be given by hundreds of trees
 - Provide global feature importance
 - an estimate of which features are important in the classification
- Incrementality
 - Not incremental

- Efficiency
 - Fast model building
 - Very fast classification
- Scalability
 - Scalable both in training set size and attribute number
- Robustness
 - Robust to noise and outliers



Rule-based classification



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Politecnico di Torino

Rule-based classifier



- Classify records by using a collection of "if...then..." rules
- Rule: (*Condition*) $\rightarrow \gamma$
 - where
 - *Condition* is a conjunction of simple predicates
 - y is the class label
 - LHS: rule antecedent or condition
 - *RHS*: rule consequent
- Examples of classification rules
 - (Blood Type=Warm) \land (Lay Eggs=Yes) \rightarrow Birds
 - (Taxable Income < 50K) \land (Refund=Yes) \rightarrow Cheat=No

Rule-based Classifier (Example)



			-		
Name	Blood Type	Give Birth	Can Fly	Live in Water	Class
human	warm	yes	no	no	mammals
python	cold	no	no	no	reptiles
salmon	cold	no	no	yes	fishes
whale	warm	yes	no	yes	mammals
frog	cold	no	no	sometimes	amphibians
komodo	cold	no	no	no	reptiles
bat	warm	yes	yes	no	mammals
pigeon	warm	no	yes	no	birds
cat	warm	yes	no	no	mammals
leopard shark	cold	yes	no	yes	fishes
turtle	cold	no	no	sometimes	reptiles
penguin	warm	no	no	sometimes	birds
porcupine	warm	yes	no	no	mammals
eel	cold	no	no	yes	fishes
salamander	cold	no	no	sometimes	amphibians
gila monster	cold	no	no	no	reptiles
platypus	warm	no	no	no	mammals
owl	warm	no	yes	no	birds
dolphin	warm	yes	no	yes	mammals
eagle	warm	no	yes	no	birds

R1: (Give Birth = no) \land (Can Fly = yes) \rightarrow Birds R2: (Give Birth = no) \land (Live in Water = yes) \rightarrow Fishes R3: (Give Birth = yes) \land (Blood Type = warm) \rightarrow Mammals R4: (Give Birth = no) \land (Can Fly = no) \rightarrow Reptiles R5: (Live in Water = sometimes) \rightarrow Amphibians



Rule-based classification



A rule r covers an instance x if the attributes of the instance satisfy the condition of the rule

R1: (Give Birth = no) \land (Can Fly = yes) \rightarrow Birds

R2: (Give Birth = no) \land (Live in Water = yes) \rightarrow Fishes

R3: (Give Birth = yes) \land (Blood Type = warm) \rightarrow Mammals

R4: (Give Birth = no) \land (Can Fly = no) \rightarrow Reptiles

R5: (Live in Water = sometimes) \rightarrow Amphibians

Name	Blood Type	Give Birth	Can Fly	Live in Water	Class
hawk	warm	no	yes	no	?
grizzly bear	warm	yes	no	no	?

Rule R1 covers a hawk => Bird

Rule R3 covers the grizzly bear => Mammal



Rule-based classification



- R1: (Give Birth = no) \land (Can Fly = yes) \rightarrow Birds
- R2: (Give Birth = no) \land (Live in Water = yes) \rightarrow Fishes
- R3: (Give Birth = yes) \land (Blood Type = warm) \rightarrow Mammals
- R4: (Give Birth = no) \land (Can Fly = no) \rightarrow Reptiles
- R5: (Live in Water = sometimes) \rightarrow Amphibians

Name	Blood Type	Give Birth	Can Fly	Live in Water	Class
lemur	warm	yes	no	no	?
turtle	cold	no	no	sometimes	?
dogfish shark	cold	yes	no	yes	?

A lemur triggers (only) rule R3, so it is classified as a mammal

- A turtle triggers both R4 and R5
- A dogfish shark triggers none of the rules



Characteristics of rules



Mutually exclusive rules

- Two rule conditions can't be true at the same time
- Every record is covered by at most one rule

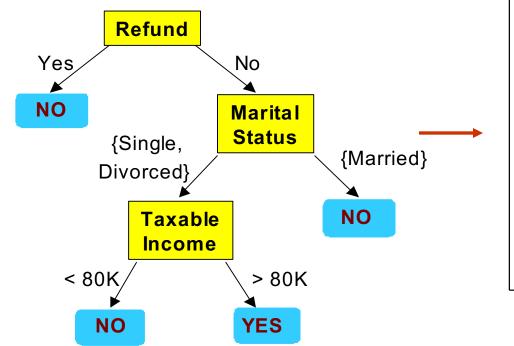
Exhaustive rules

- Classifier rules account for every possible combination of attribute values
- Each record is covered by at least one rule



From decision trees to rules





Classification Rules

(Refund=Yes) ==> No

(Refund=No, Marital Status={Single,Divorced}, Taxable Income<80K) ==> No

(Refund=No, Marital Status={Single,Divorced}, Taxable Income>80K) ==> Yes

(Refund=No, Marital Status={Married}) ==> No

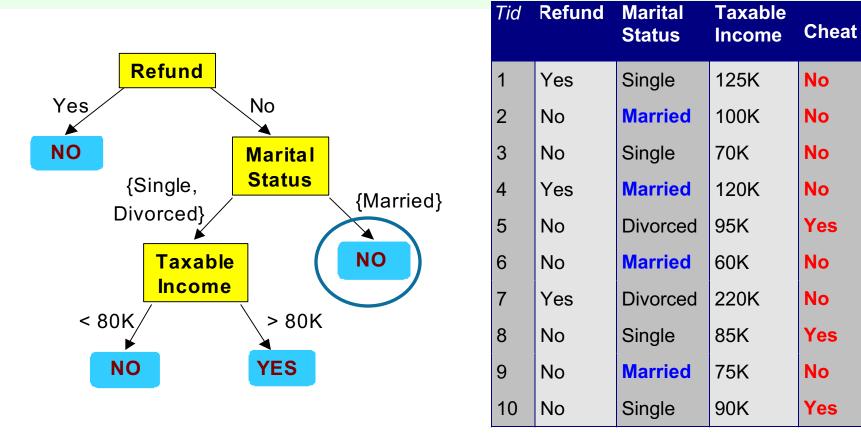
Rules are mutually exclusive and exhaustive

Rule set contains as much information as the tree



Rules can be simplified





Initial Rule:(Refund=No) \land (Status=Married) \rightarrow NoSimplified Rule:(Status=Married) \rightarrow No

Effect of rule simplification



- Rules are no longer mutually exclusive
 - A record may trigger more than one rule
 - Solution?
 - Ordered rule set
 - Unordered rule set use voting schemes
- Rules are no longer exhaustive
 - A record may not trigger any rules
 - Solution?
 - Use a default class



Ordered rule set



- Rules are rank ordered according to their priority
 - An ordered rule set is known as a decision list
- When a test record is presented to the classifier
 - It is assigned to the class label of the highest ranked rule it has triggered
 - If none of the rules fired, it is assigned to the default class

R1: (Give Birth = no) \land (Can Fly = yes) \rightarrow BirdsR2: (Give Birth = no) \land (Live in Water = yes) \rightarrow FishesR3: (Give Birth = yes) \land (Blood Type = warm) \rightarrow MammalsR4: (Give Birth = no) \land (Can Fly = no) \rightarrow ReptilesR5: (Live in Water = sometimes) \rightarrow Amphibians

¥¥					
Name	Blood Type	Give Birth	Can Fly	Live in Water	Class
turtle	cold	no	no	sometimes	?



Building classification rules

Direct Method

- Extract rules directly from data
- e.g.: RIPPER, CN2, Holte's 1R

Indirect Method

- Extract rules from other classification models (e.g. decision trees, neural networks, etc).
- e.g: C4.5rules





Evaluation of rule based classifiers



- Accuracy
 - Higher than decision trees
- Interpretability
 - Model and prediction are interpretable
- Incrementality
 - Not incremental

- Efficiency
 - Fast model building
 - Very fast classification
- Scalability
 - Scalable both in training set size and attribute number
- Robustness
 - Robust to outliers



Associative classification



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Associative classification



The classification model is defined by means of association rules

(Condition) $\rightarrow \gamma$

- rule body is an itemset
- Model generation
 - Rule selection & sorting
 - based on support, confidence and correlation thresholds
 - Rule pruning
 - Database coverage: the training set is covered by selecting topmost
 - rules according to previous sort

Evaluation of associative classifiers



Accuracy

- Higher than decision trees and rule-based classifiers
 - *correlation* among attributes is considered
- Interpretability
 - Model and prediction are interpretable
- Incrementality
 - Not incremental

Efficiency

- Rule generation may be slow
 - It depends on support threshold
- Very fast classification
- Scalability
 - Scalable in training set size
 - Reduced scalability in attribute number
 - Rule generation may become unfeasible
- Robustness
 - Unaffected by missing data
 - Robust to outliers



K-Nearest Neighbor



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Instance-Based Classifiers



Set of Stored Cases • Store the training records • Use training records to Class Atr1 AtrN predict the class label of Α unseen cases В В Unseen Case C Atr1 AtrN Α C В



Instance Based Classifiers



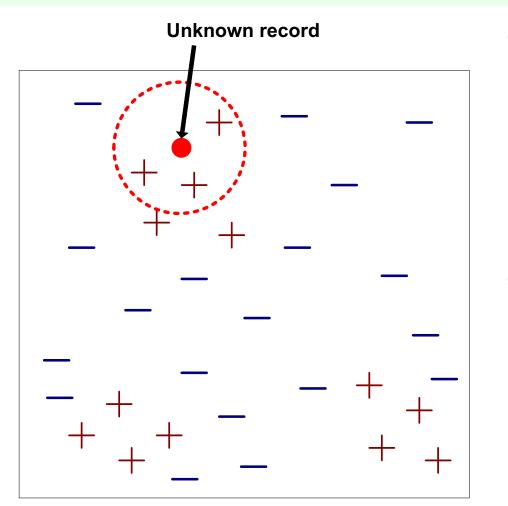
Examples

- Rote-learner
 - Memorizes entire training data and performs classification only if attributes of record match one of the training examples exactly
- Nearest neighbor
 - Uses k "closest" points (nearest neighbors) for performing classification



Nearest-Neighbor Classifiers



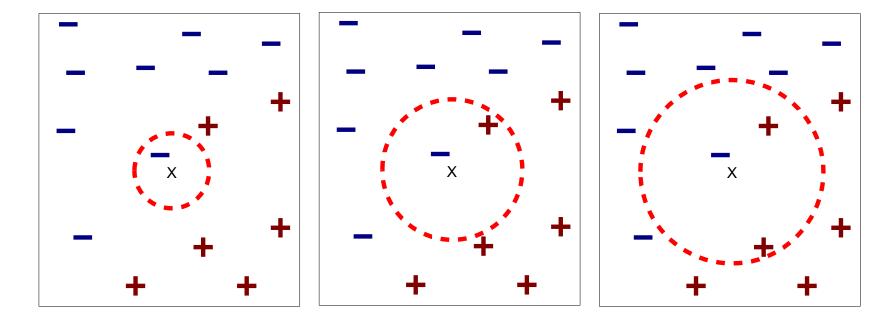


- Requires
 - The set of stored records
 - Distance Metric to compute distance between records
 - The value of k, the number of nearest neighbors to retrieve
- To classify an unknown record
 - Compute distance to other training records
 - Identify *k* nearest neighbors
 - Use class labels of nearest neighbors to determine the class label of unknown record (e.g., by taking majority vote)



Definition of Nearest Neighbor





(a) 1-nearest neighbor (b) 2-nearest neighbor (c) 3-nearest neighbor

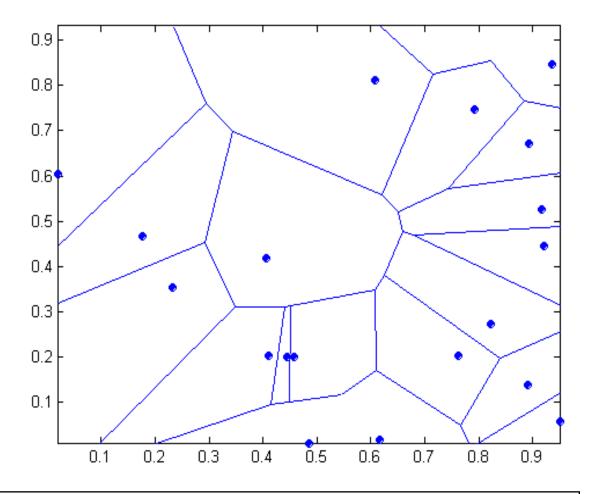
K-nearest neighbors of a record x are data points that have the k smallest distance to x



1 nearest-neighbor



Voronoi Diagram





Nearest Neighbor Classification



Euclidean distance

$$d(p,q) = \sqrt{\sum_{i} (p_i - q_i)^2}$$

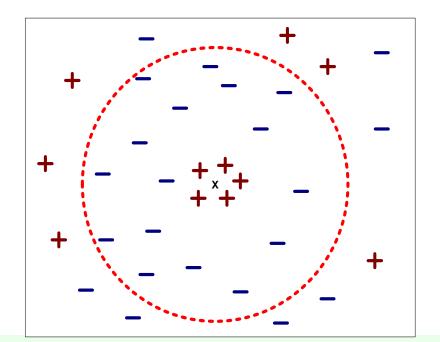
- Determine the class from nearest neighbor list
 - take the majority vote of class labels among the k-nearest neighbors
 - Weigh the vote according to distance
 - weight factor, $w = 1/d^2$



Nearest Neighbor Classification

Choosing the value of k:

- If k is too small, sensitive to noise points
- If k is too large, neighborhood may include points from other classes





Nearest Neighbor Classification

1859

Scaling issues

- Attribute domain should be normalized to prevent distance measures from being dominated by one of the attributes
- Example: height [1.5m to 2.0m] vs. income [\$10K to \$1M]
- Problem with distance measures
 - High dimensional data
 - curse of dimensionality



Evaluation of KNN



Accuracy

- Comparable to other classification techniques for simple datasets
- Interpretability
 - Model is not interpretable
 - Single predictions can be "described" by neighbors

Incrementality

- Incremental
- Training set *must* be available

Efficiency

- (Almost) no model building
- Slower classification, requires computing distances
- Scalability
 - Weakly scalable in training set size
 - Curse of dimensionality for increasing attribute number
- Robustness
 - Depends on distance computation



Bayesian Classification



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P(C|X) = P(X|C) P(C) / P(X)

and also

Hence

P(C|X) P(X) = P(X|C) P(C)

P(C,X) = P(C|X) P(X)P(C,X) = P(X|C) P(C)





Bayesian classification



- Let the class attribute and all data attributes be random variables
 - C = any class label
 - $X = \langle x_1, ..., x_k \rangle$ record to be classified
- Bayesian classification
 - compute P(C|X) for all classes
 - probability that record X belongs to C
 - assign X to the class with *maximal* P(C|X)
- Applying Bayes theorem

 $P(C|X) = P(X|C) \cdot P(C) / P(X)$

- P(X) constant for all C, disregarded for maximum computation
- P(C) a priori probability of C

 $P(C) = N_c/N$



Bayesian classification

1859

- How to estimate P(X|C), i.e. $P(x_1,...,x_k|C)$?
- Naïve hypothesis

 $P(x_1,...,x_k|C) = P(x_1|C) P(x_2|C) ... P(x_k|C)$

- statistical independence of attributes x₁,...,x_k
- not always true
 - model quality may be affected
- Computing P(x_k|C)
 - for discrete attributes

 $P(x_k|C) = |x_{kC}| / N_c$

- where $|x_{kC}|$ is number of instances having value x_k for attribute k and belonging to class C
- for continuous attributes, use probability distribution
- Bayesian networks
 - allow specifying a subset of dependencies among attributes

Bayesian classification: Example



Outlook	Temperature	Humidity	Windy	Class
sunny	hot	high	false	N
sunny	hot	high	true	N
overcast		high	false	Ρ
rain	mild	high	false	Ρ
rain	cool	normal	false	Ρ
rain	cool	normal	true	Ν
overcast	cool	normal	true	Ρ
sunny	mild	high	false	Ν
sunny	cool	normal	false	Ρ
rain	mild	normal	false	Ρ
sunny	mild	normal	true	Ρ
overcast	mild	high	true	Р
overcast	hot	normal	false	Ρ
rain	mild	high	true	Ν



From: Han, Kamber,"Data mining; Concepts and Techniques", Morgan Kaufmann 2006

Bayesian classification: Example



outlook		
P(sunny p) = 2/9	P(sunny n) = 3/5	Р
P(overcast p) = 4/9	P(overcast n) = 0	
P(rain p) = 3/9	P(rain n) = 2/5	[P
temperature		
P(hot p) = 2/9	P(hot n) = 2/5	
P(mild p) = 4/9	P(mild n) = 2/5	
P(cool p) = 3/9	P(cool n) = 1/5	
humidity		
P(high p) = 3/9	P(high n) = 4/5	
P(normal p) = 6/9	P(normal n) = 2/5	
windy		
P(true p) = 3/9	P(true n) = 3/5	
P(false p) = 6/9	P(false n) = 2/5	
7		

P(p)		9/14
P(n)	=	5/14

From: Han, Kamber,"Data mining; Concepts and Techniques", Morgan Kaufmann 2006

Bayesian classification: Example

Data to be labeled

X = <rain, hot, high, false>

- For class p
 - $P(X|p) \cdot P(p) =$ = P(rain|p) \cdot P(hot|p) \cdot P(high|p) \cdot P(false|p) \cdot P(p) = 3/9 \cdot 2/9 \cdot 3/9 \cdot 6/9 \cdot 9/14 = 0.010582
- For class n

 $P(X|n) \cdot P(n) =$ = P(rain|n) \cdot P(hot|n) \cdot P(high|n) \cdot P(false|n) \cdot P(n) = 2/5 \cdot 2/5 \cdot 4/5 \cdot 2/5 \cdot 5/14 = 0.018286



Evaluation of Naïve Bayes Classifiers



- Accuracy
 - Similar or lower than decision trees
 - Naïve hypothesis simplifies model
- Interpretability
 - Model and prediction are not interpretable
 - The weights of contributions in a single prediction may be used to explain
- Incrementality
 - Fully incremental
 - Does *not* require availability of training data

- Efficiency
 - Fast model building
 - Very fast classification
- Scalability
 - Scalable both in training set size and attribute number
- Robustness
 - Affected by attribute correlation

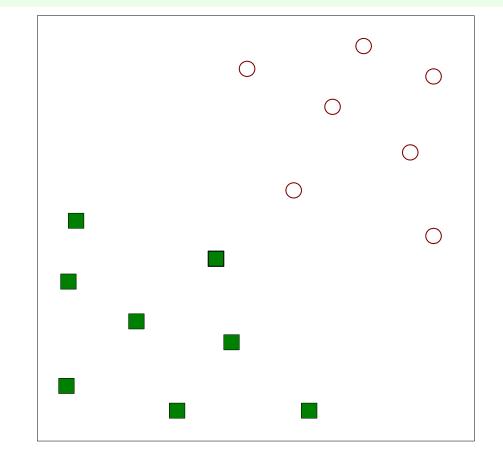




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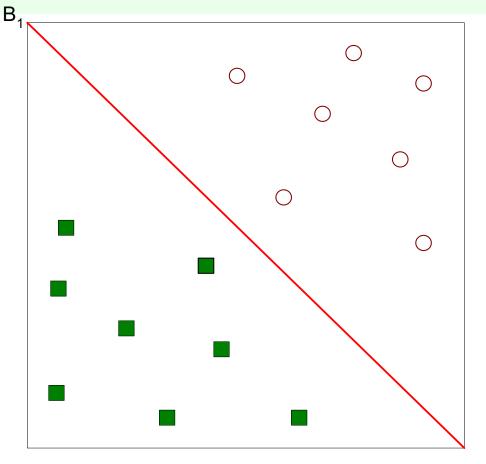




• Find a linear hyperplane (decision boundary) that will separate the data



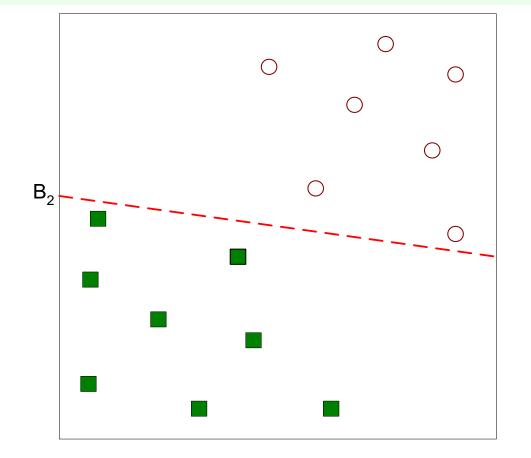




One Possible Solution



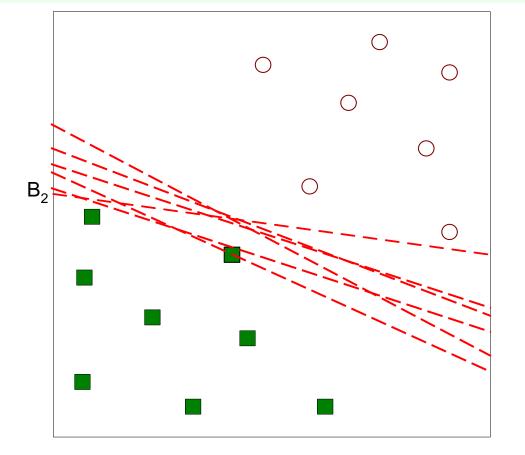




Another possible solution



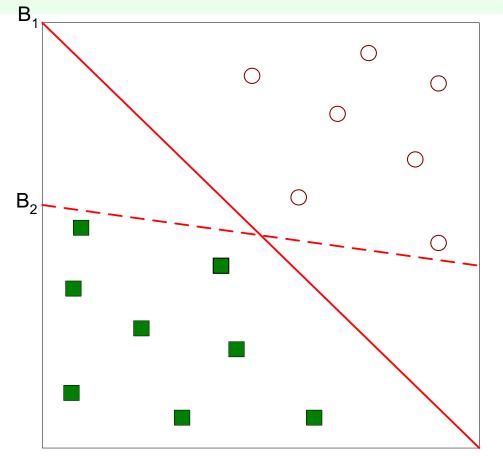




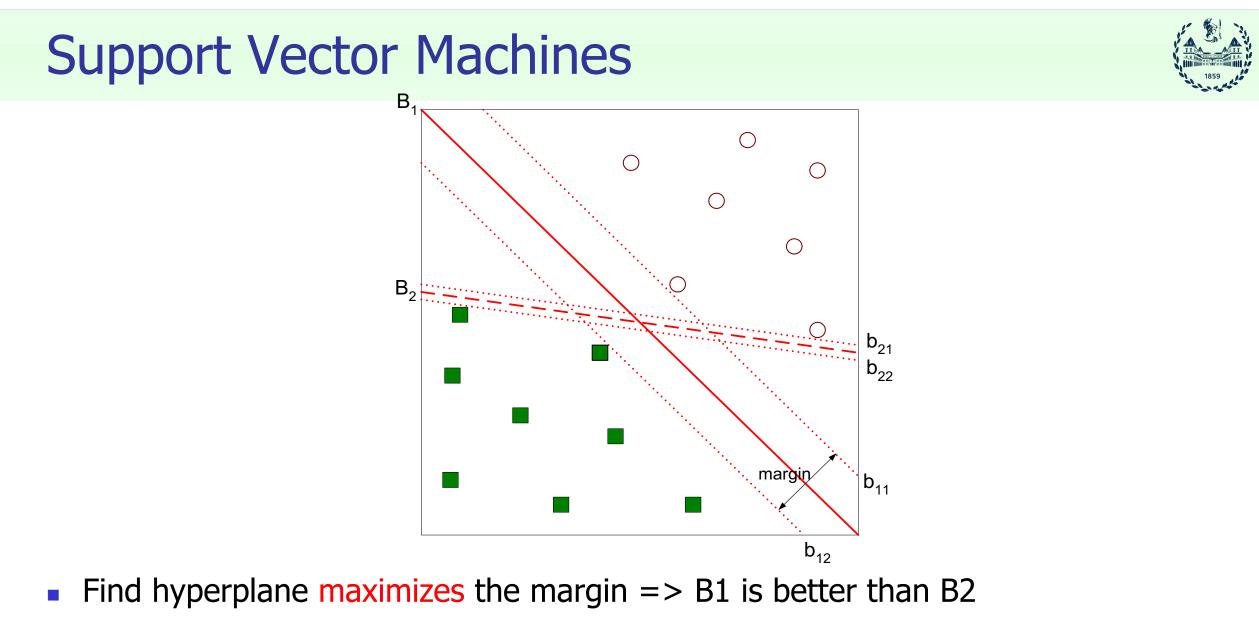
Other possible solutions







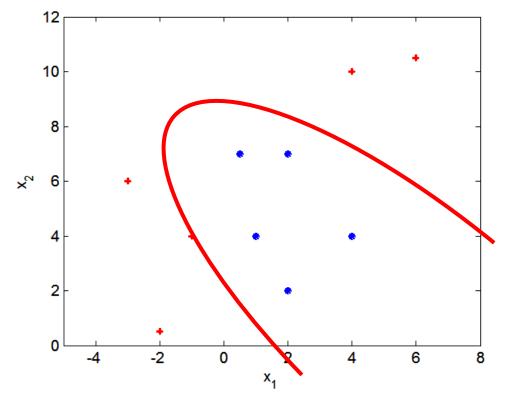
- Which one is better? B1 or B2?
- How do you define better?







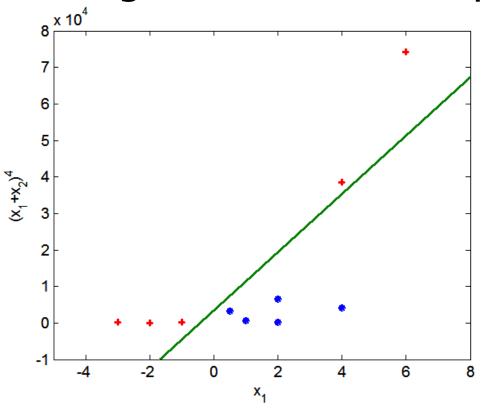
What if decision boundary is not linear?







Transform data into higher dimensional space





Evaluation of Support Vector Machines



Accuracy

- Among best performers
- Interpretability
 - Model and prediction are not interpretable
 - Black box model
- Incrementality
 - Not incremental

- Efficiency
 - Model building requires significant parameter tuning
 - Very fast classification
- Scalability
 - Medium scalable both in training set size and attribute number
- Robustness
 - Robust to noise and outliers



Artificial Neural Networks



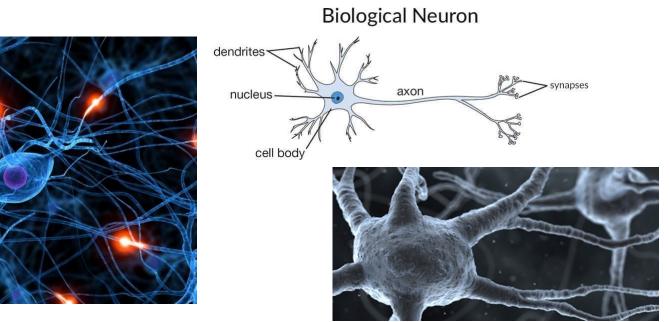
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Artificial Neural Networks

Inspired to the structure of the human brain

- Neurons as elaboration units
- Synapses as connection network

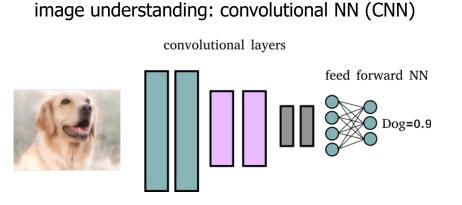




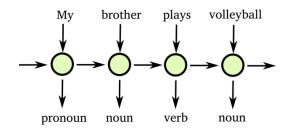


Artificial Neural Networks

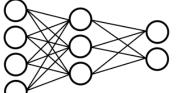
Different tasks, different architectures



time series analysis: recurrent NN (RNN)



numerical vectors classification: feed forward NN (FFNN)

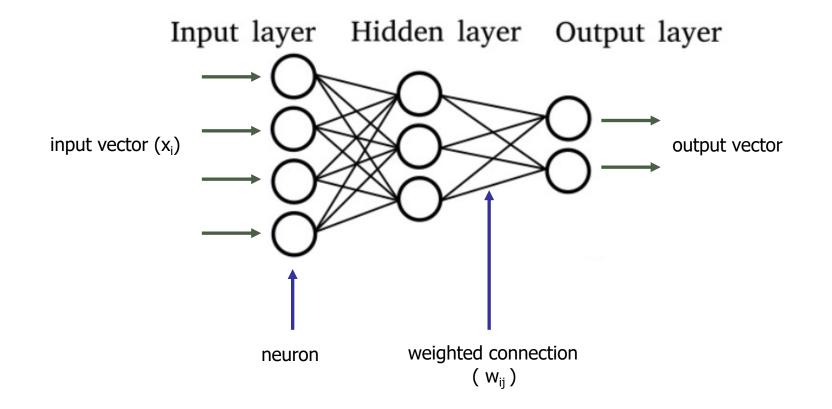


denoising: auto-encoders





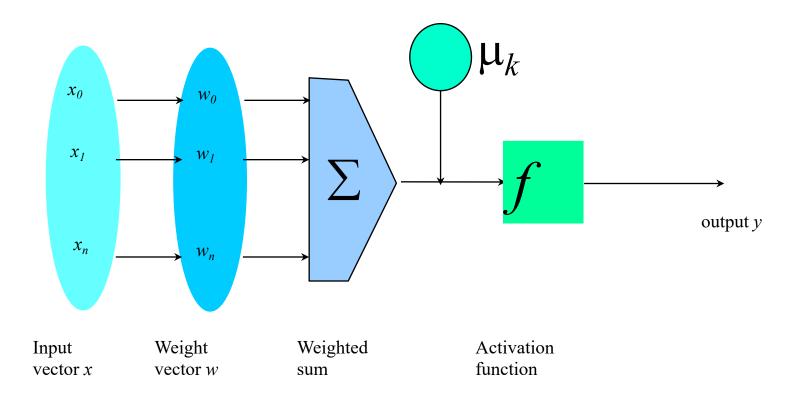
Feed Forward Neural Network





Structure of a neuron







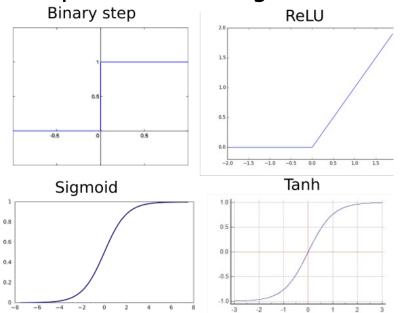
From: Han, Kamber,"Data mining; Concepts and Techniques", Morgan Kaufmann 2006

Activation Functions



Activation

- simulates biological activation to input stymula
- provides non-linearity to the computation
- may help to saturate neuron outputs in fixed ranges

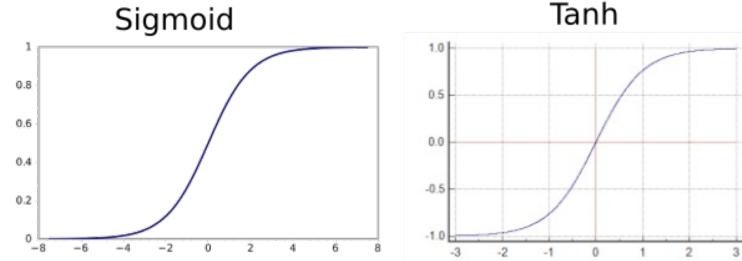




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Activation Functions

- Sigmoid, tanh
 - saturate input value in a fixed range
 - non linear for all the input scale
 - typically used by FFNNs for both hidden and output layers
 - E.g. *sigmoid* in output layers allows generating values between 0 and 1 (useful when output must be interpreted as likelihood)





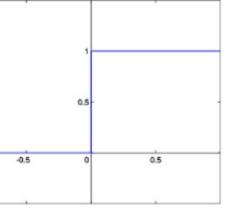
Activation Functions

Binary Step

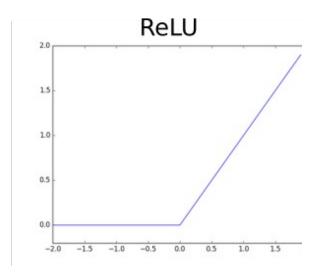
- outputs 1 when input is non-zero
- useful for binary outputs
- **issues**: not appropriate for gradient descent
 - derivative not defined in x=0
 - derivative equal to 0 in every other position

ReLU (Rectified Linear Unit)

- used in deep networks (e.g. CNNs)
 - avoids vanishing gradient
 - does not saturate
- neurons activate linearly only for positive input



Binary step





Activation Functions



Softmax

- differently to other activation functions
 - it is applied only to the **output** layer
 - works by considering all the neurons in the layer
- after softmax, the output vector can be interpreted as a discrete distribution of probabilities
 - e.g. the probabilities for the input pattern of belonging to each class

$$softmax(z_j) = \frac{e^{z_j}}{\sum_{i=0}^{N-1} e^{z_i}} \qquad \begin{array}{c} \text{softmax} \\ z_0 \\ z_1 \\ z_2 \\ z$$



Building a FFNN



For each node, definition of

- set of weights
- offset value

providing the highest accuracy on the training data

Iterative approach on training data instances



Building a FFNN



Base algorithm

- Initially assign random values to weights and offsets
- Process instances in the training set one at a time
 - For each neuron, compute the result when applying weights, offset and activation function for the instance
 - Forward propagation until the output is computed
 - Compare the computed output with the expected output, and evaluate error
 - Backpropagation of the error, by updating weights and offset for each neuron
- The process ends when
 - % of accuracy above a given threshold
 - % of parameter variation (error) below a given threshold
 - The maximum number of epochs is reached



Evaluation of Feed Forward NN



- Accuracy
 - Among best performers
- Interpretability
 - Model and prediction are not interpretable
 - Black box model
- Incrementality
 - Not incremental

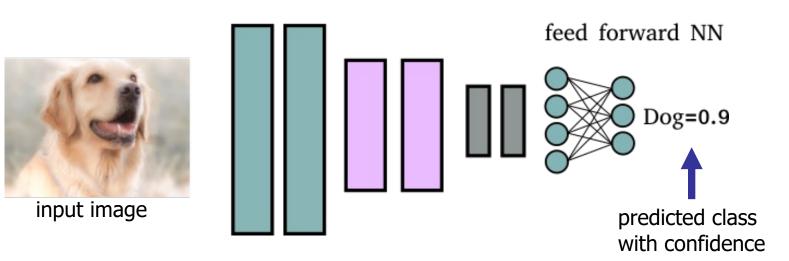
- Efficiency
 - Model building requires very complex parameter tuning
 - It requires significant time
 - Very fast classification
- Scalability
 - Medium scalable both in training set size and attribute number
- Robustness
 - Robust to noise and outliers
 - Requires large training set
 - Otherwise unstable when tuning parameters





 Allow automatically extracting features from images and performing classification

convolutional layers

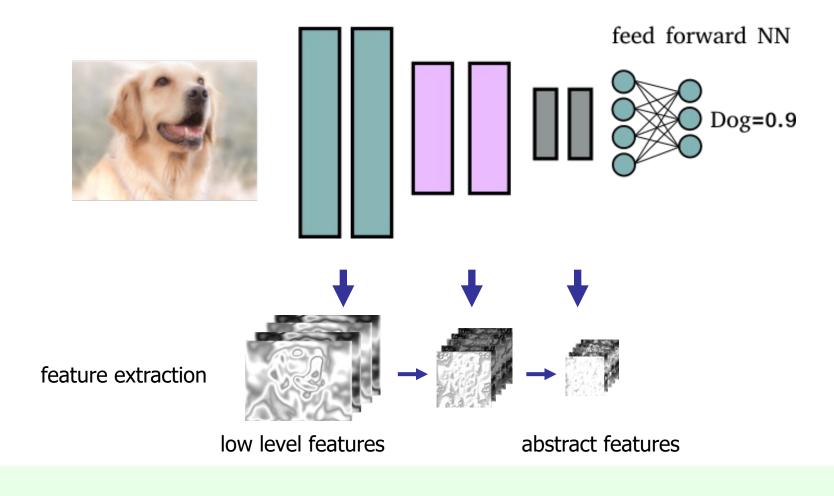


Convolutional Neural Network (CNN) Architecture





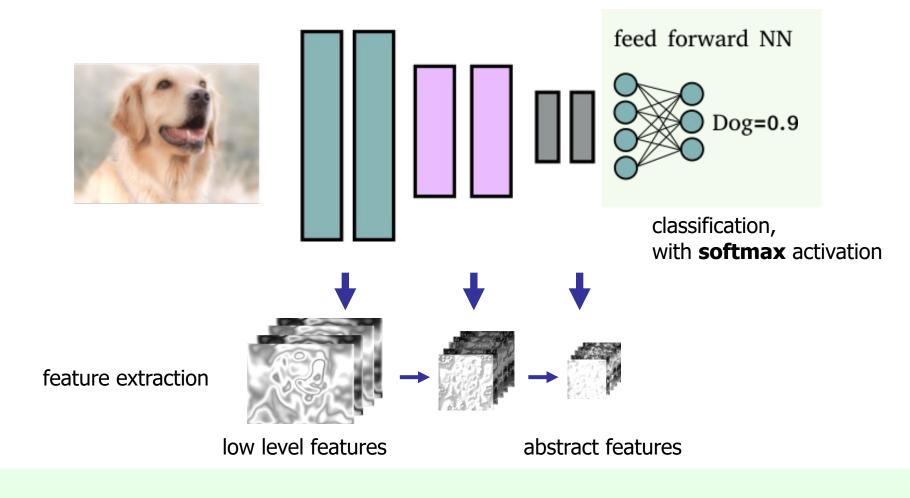
convolutional layers







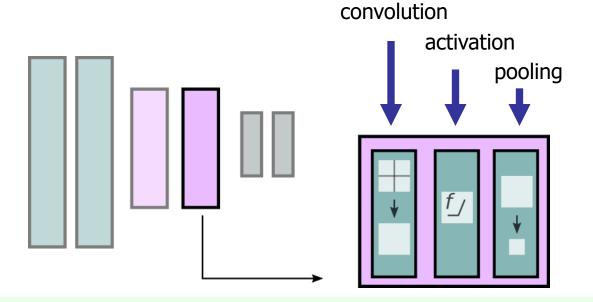
convolutional layers







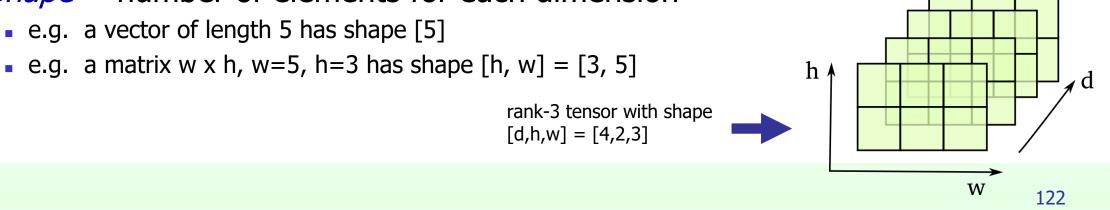
- Typical convolutional layer
 - *convolution* stage: feature extraction by means of (hundreds to thousands) sliding filters
 - sliding filters *activation*: apply activation functions to input tensor
 - *pooling*: tensor downsampling





Tensors

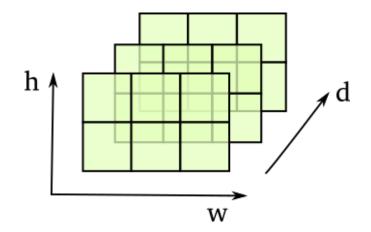
- data flowing through CNN layers is represented in the form of *tensors*
- Tensor = N-dimensional vector
- Rank = number of dimensions
 - scalar: rank 0
 - 1-D vector: rank 1
 - 2-D matrix: rank 2
- Shape = number of elements for each dimension





Images

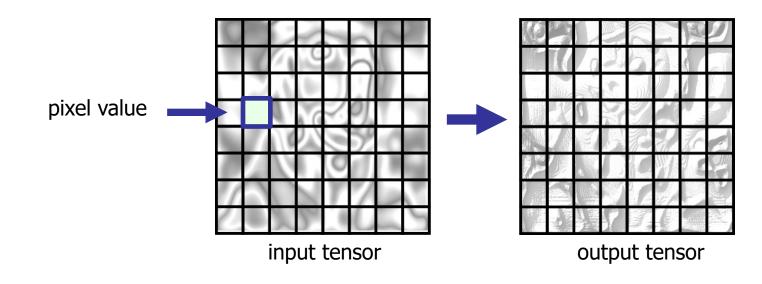
- rank-3 tensors with shape [d,h,w]
- where h=height, w=width, d=image depth (1 for grayscale, 3 for RGB colors)





Convolution

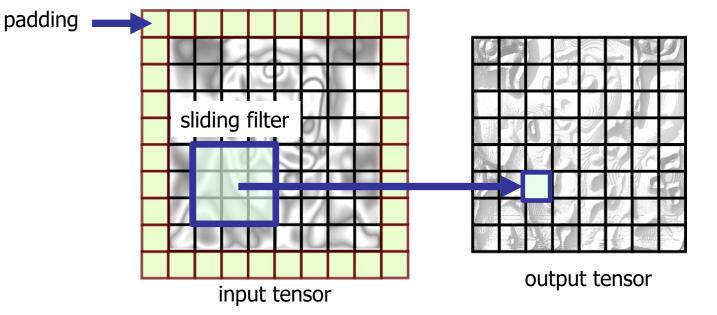
- processes data in form of *tensors* (multi-dimensional matrices)
- input: input image or intermediate features (tensor)
- **output**: a tensor with the extracted features





Convolution

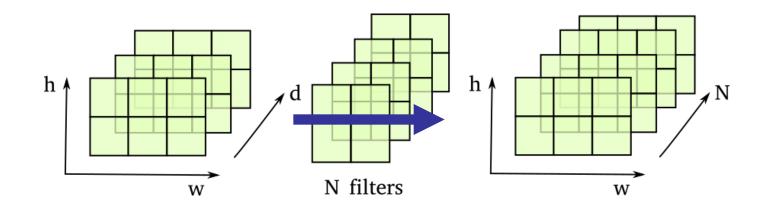
- a *sliding filter* produces the values of the output tensor
- sliding filters contain the trainable weights of the neural network
- each convolutional layer contains many (hundreds) filters





Convolution

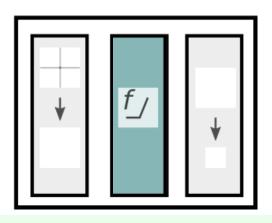
- images are transformed into features by convolutional filters
- after convolving a tensor [d,h,w] with *N filters* we obtain
 - a rank-3 tensor with shape [N,h,w]
 - hence, each filter generates a layer in the depth of the output tensor

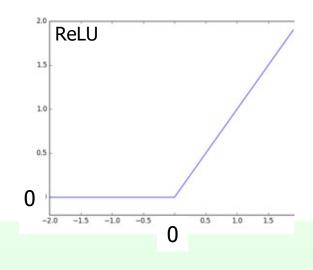




Activation

- symulates biological activation to input stymula
- provides non-linearity to the computation
- ReLU is typically used for CNNs
 - faster training (no vanishing gradients)
 - does not saturate
 - faster computation of derivatives for backpropagation





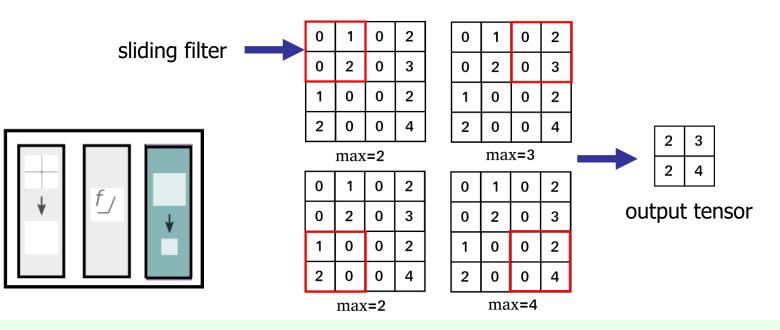




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Pooling

- performs tensor *downsampling*
- sliding filter which replaces tensor values with a summary statistic of the nearby outputs
- *maxpool* is the most common: computes the maximum value as statistic

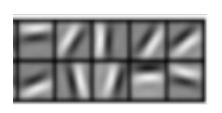




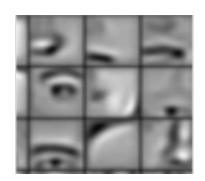


Convolutional layers training

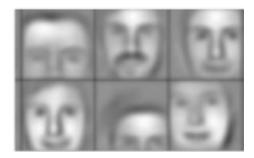
- during training each sliding filter learns to recognize a particular *pattern* in the input tensor
- filters in *shallow layers* recognize textures and edges
- filters in *deeper layers* can recognize objects and parts (e.g. eye, ear or even faces)



shallow filters



deeper filters

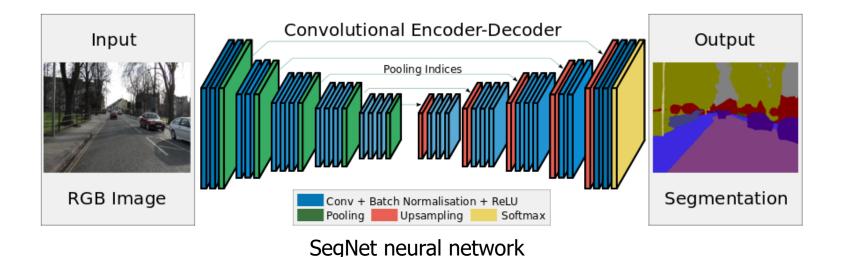






Semantic segmentation CNNs

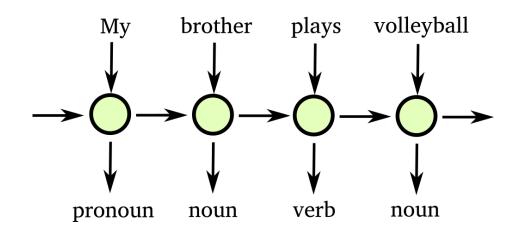
- allow assigning a class to each pixel of the input image
- composed of 2 parts
 - encoder network: convolutional layers to extract abstract features
 - **decoder network**: deconvolutional layers to obtain the output image from the extracted features



 $D_{M}^{B}G$

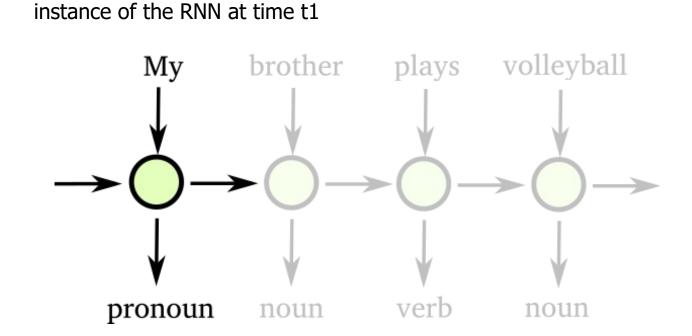


- Allow processing *sequential* data x(t)
- Differently from normal FFNN they are able to keep a *state* which evolves during time
- Applications
 - machine translation
 - time series prediction
 - speech recognition
 - part of speech (POS) tagging



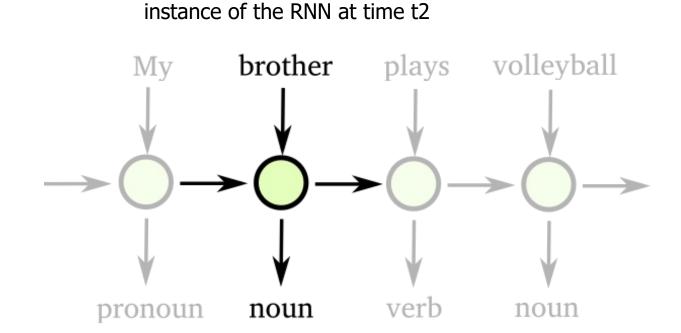






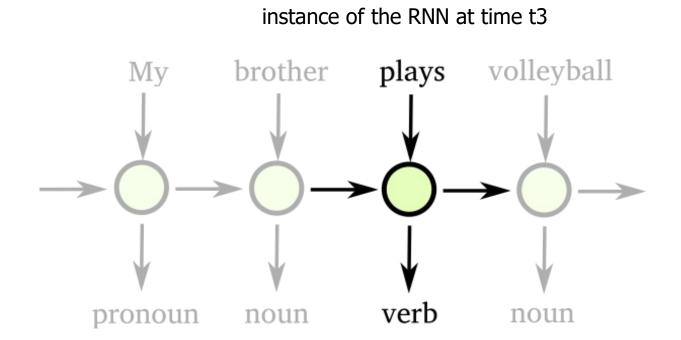






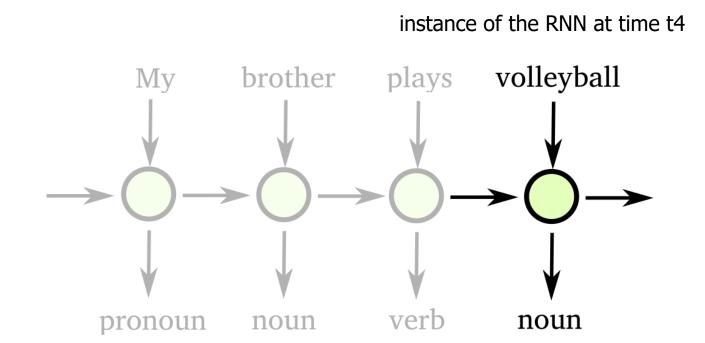








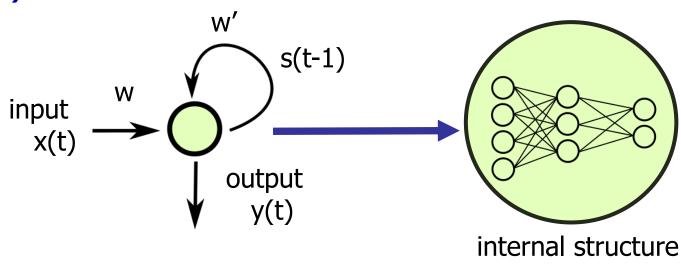








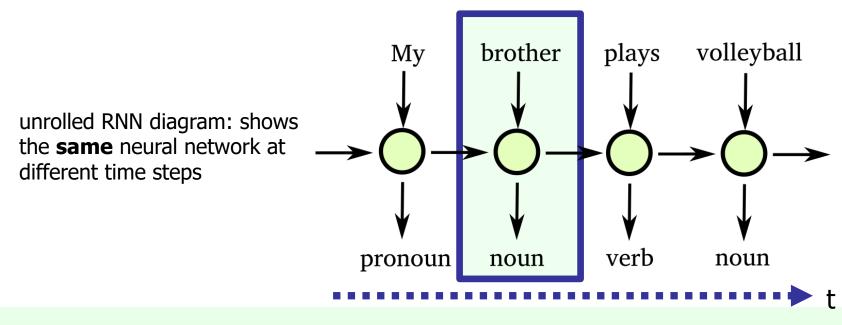
- A RNN receives as input a vector x(t) and the state at previous time step s(t-1)
- A RNN typically contains many *neurons organized in different layers* state retroaction







- Training is performed with *Backpropagation Through Time*
- Given a pair training sequence x(t) and expected output y(t)
 - error is propagated through time
 - weights are updated to minimize the error across all the time steps

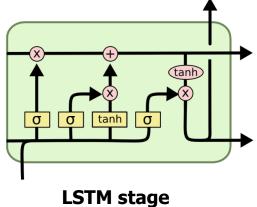


instance of the RNN at time t



Issues

- vanishing gradient: error gradient decreases rapidly over time, weights are not properly updated
- this makes harder having RNN with *long-term* memories
- Solution: *LSTM* (Long Short Term Memories)
 - RNN with "gates" which encourage the state information to flow through long time intervals

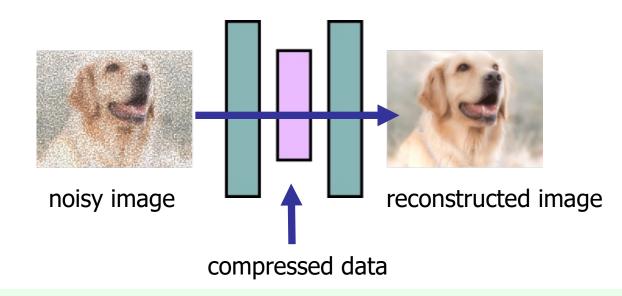




Autoencoders



- Autoencoders allow *compressing* input data by means of compact representations and from them *reconstructing* the initial input
 - for feature extraction: the compressed representation can be used as significant set of features representing input data
 - for image (or signal) *denoising*: the image reconstructed from the abstract representation is denoised with respect to the original one



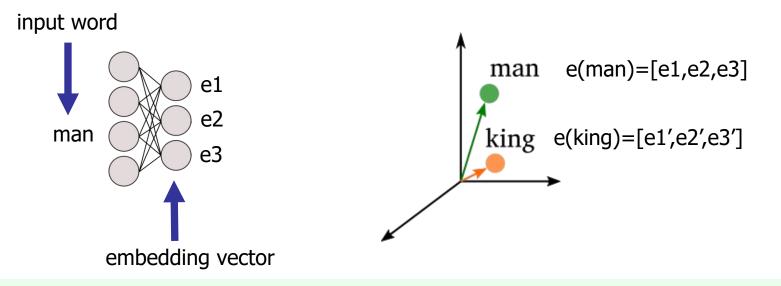


Word Embeddings (Word2Vec)



Word *embeddings* associate words to n-dimensional vectors

- trained on big text collections to model the word distributions in different sentences and contexts
- able to capture the semantic information of each word
- words with similar *meaning* share vectors with similar characteristics

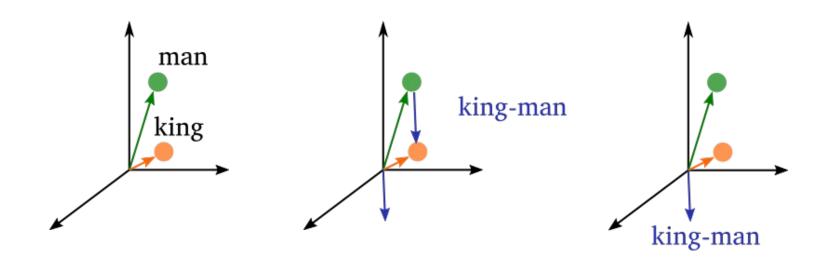




Word Embeddings (Word2Vec)



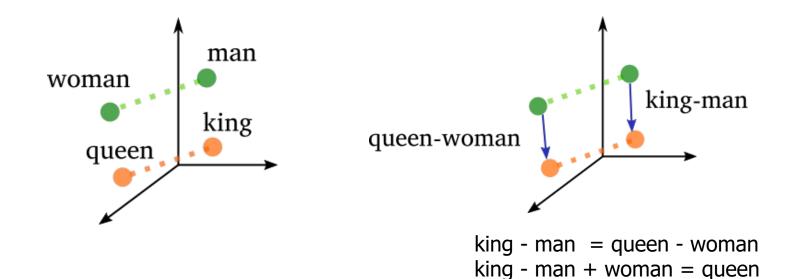
 Since each word is represented with a vector, operations among words (e.g. difference, addition) are allowed





Word Embeddings (Word2Vec)

 Semantic relationships among words are captured by vector positions





Model evaluation



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Model evaluation



- Methods for performance evaluation
 - Partitioning techniques for training and test sets
- Metrics for performance evaluation
 - Accuracy, other measures
- Techniques for model comparison
 - ROC curve



Methods for performance evaluation

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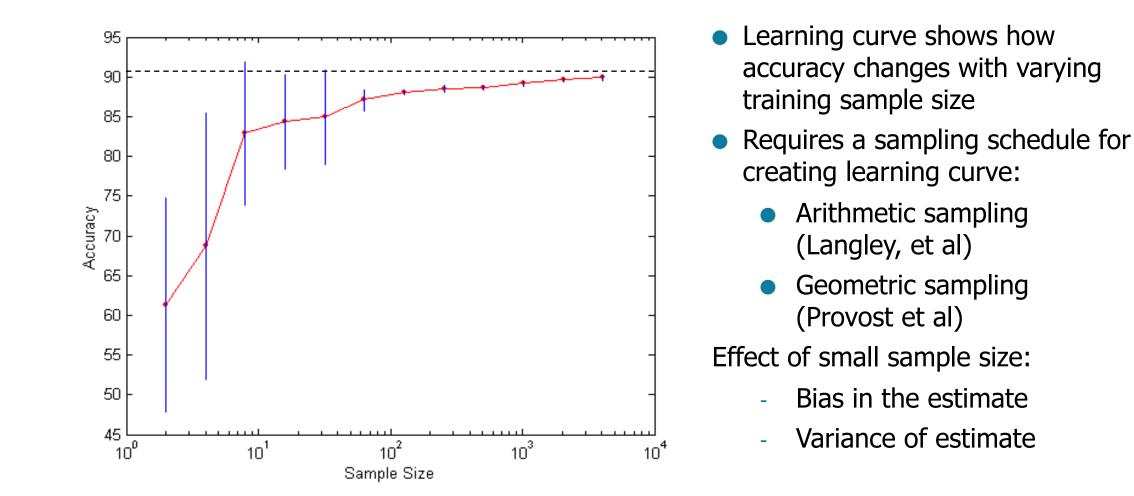
Objective

- reliable estimate of performance
- Performance of a model may depend on other factors besides the learning algorithm
 - Class distribution
 - Cost of misclassification
 - Size of training and test sets



Learning curve







From: Tan, Steinbach, Kumar, Introduction to Data Mining, McGraw Hill 2006

Partitioning data



- Several partitioning techniques
 - holdout
 - cross validation
- Stratified sampling to generate partitions
 - without replacement
- Bootstrap
 - Sampling with replacement



Holdout



Fixed partitioning

- Typically, may reserve 80% for training, 20% for test
- Other proportions may be appropriate, depending on the dataset size

Appropriate for large datasets

- may be repeated several times
 - repeated holdout



Cross validation



Cross validation

- partition data into k disjoint subsets (i.e., folds)
- k-fold: train on k-1 partitions, test on the remaining one
 - repeat for all folds
- reliable accuracy estimation, not appropriate for very large datasets
- Leave-one-out
 - cross validation for k=n
 - only appropriate for very small datasets



Model performance estimation

- Model training step
 - Building a new model
- Model validation step
 - Hyperparameter tuning
 - Algorithm selection
- Model test step
 - Estimation of model performance





Model performance estimation

Typical dataset size

- Training set 60% of labeled data
- Validation set 20% of labeled data
- Test set 20% of labeled data

Splitting labeled data

- Use hold-out to split in
 - training+validation
 - test

Use cross validation to split in

- training
- validation



Metrics for model evaluation



- Evaluate the predictive accuracy of a model
- Confusion matrix
 - binary classifier

	PREDICTED CLASS			
		Class=Yes	Class=No	
ACTUAL CLASS	Class=Yes	а	b	a: TP (true positive)b: FN (false negative)c: FP (false positive)d: TN (true negative)
	Class=No	С	d	







Most widely-used metric for model evaluation

 $Accuracy = \frac{Number of correctly classified objects}{Number of classified objects}$

Not always a reliable metric







For a binary classifier

ACTUAL CLASSClass=YesClass=NoCLASSClass=YesabClass=Nocd		PREDICTED CLASS		
ACTUAL (TP) (FN)			Class=Yes	Class=No
CLASS Class=No c d		Class=Yes		b (FN)
(FP) (TN)	CLASS	Class=No	c (FP)	d (TN)

Accuracy =
$$\frac{a+d}{a+b+c+d} = \frac{TP+TN}{TP+TN+FP+FN}$$



Limitations of accuracy

Consider a binary problem

- Cardinality of Class 0 = 9900
- Cardinality of Class 1 = 100
- Model

() \rightarrow class 0

Model predicts everything to be class 0

accuracy is 9900/10000 = 99.0 %

Accuracy is misleading because the model does not detect any class 1 object





Limitations of accuracy



Classes may have different importance

- Misclassification of objects of a given class is more important
- e.g., ill patients erroneously assigned to the healthy patients class
- Accuracy is not appropriate for
 - unbalanced class label distribution
 - different class relevance



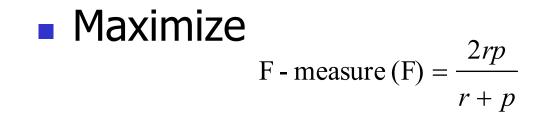
Class specific measures



Evaluate separately for each class C

 $Recall (r) = \frac{Number of objects correctly assigned to C}{Number of objects belonging to C}$

 $Precision (p) = \frac{Number of objects correctly assigned to C}{Number of objects assigned to C}$



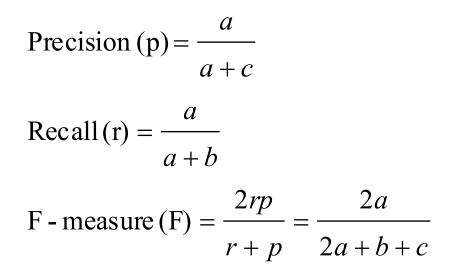


Class specific measures



For a binary classification problem

on the confusion matrix, for the positive class





ROC (Receiver Operating Characteristic)



- Developed in 1950s for signal detection theory to analyze noisy signals
 - characterizes the trade-off between positive hits and false alarms
- ROC curve plots
 - TPR, True Positive Rate (on the y-axis) TPR = TP/(TP+FN)

against

FPR, False Positive Rate (on the x-axis)

FPR = FP/(FP + TN)



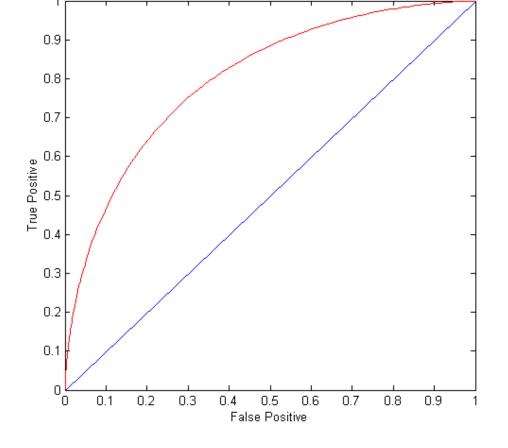
From: Tan, Steinbach, Kumar, Introduction to Data Mining, McGraw Hill 2006

(FPR, TPR)

ROC curve

- (0,0): declare everything to be negative class
- (1,1): declare everything to be positive class
- (0,1): ideal
- Diagonal line
 - Random guessing
 - Below diagonal line
 - prediction is opposite of the true class







How to build a ROC curve

Instance	P(+ A)	True Class
1	0.95	+
2	0.93	+
3	0.87	-
4	0.85	-
5	0.85	-
6	0.85	+
7	0.76	-
8	0.53	+
9	0.43	-
10	0.25	+

- Use classifier that produces posterior probability for each test instance P(+|A)
- Sort the instances according to P(+|A) in decreasing order
- Apply threshold at each unique value of P(+|A)
- Count the number of TP, FP, TN, FN at each threshold

TP rate

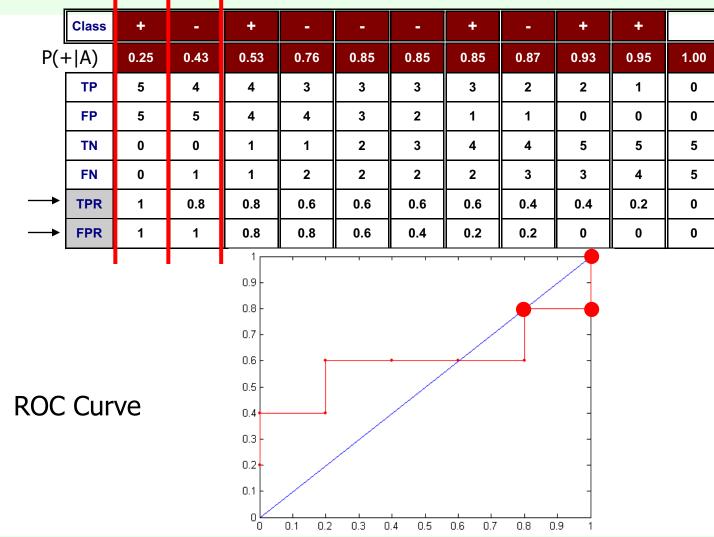
TPR = TP/(TP+FN)

FP rate

FPR = FP/(FP + TN)

How to build a ROC curve

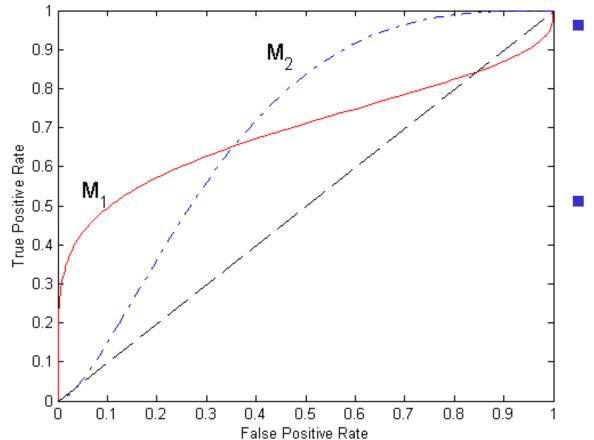






Using ROC for Model Comparison





- No model consistently outperforms the other
 - M₁ is better for small FPR
 - M₂ is better for large FPR
- Area under ROC curve
 - Ideal
 - Area = 1.0
 - Random guess
 - Area = 0.5

