

Lecture 10: 12 February, 2026

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Data Mining and Machine Learning
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Bayesian classifiers

- As before
 - Attributes $\{A_1, A_2, \dots, A_k\}$ and
 - Classes $C = \{c_1, c_2, \dots, c_\ell\}$
- Each class c_i defines a probabilistic model for attributes
 - $Pr(A_1 = a_1, \dots, A_k = a_k \mid C = c_i)$
- Given a data item $d = (a_1, a_2, \dots, a_k)$, identify the best class c for d
- Maximize $Pr(C = c_i \mid A_1 = a_1, \dots, A_k = a_k)$

Generative models

- To use probabilities, need to describe how data is randomly generated
 - Generative model
- Typically, assume a random instance is created as follows
 - Choose a class c_j with probability $Pr(c_j)$
 - Choose attributes a_1, \dots, a_k with probability $Pr(a_1, \dots, a_k | c_j)$
- Generative model has associated parameters $\theta = (\theta_1, \dots, \theta_m)$
 - Each class probability $Pr(c_j)$ is a parameter
 - Each conditional probability $Pr(a_1, \dots, a_k | c_j)$ is a parameter
- We need to estimate these parameters

Bayesian classification

- Maximize $Pr(C = c_i | A_1 = a_1, \dots, A_k = a_k)$
- By Bayes' rule,

$$\begin{aligned} & Pr(C = c_i | A_1 = a_1, \dots, A_k = a_k) \\ &= \frac{Pr(A_1 = a_1, \dots, A_k = a_k | C = c_i) \cdot Pr(C = c_i)}{Pr(A_1 = a_1, \dots, A_k = a_k)} \\ &= \frac{Pr(A_1 = a_1, \dots, A_k = a_k | C = c_i) \cdot Pr(C = c_i)}{\sum_{j=1}^{\ell} Pr(A_1 = a_1, \dots, A_k = a_k | C = c_j) \cdot Pr(C = c_j)} \end{aligned}$$

- Denominator is the same for all c_i , so sufficient to maximize

$$Pr(A_1 = a_1, \dots, A_k = a_k | C = c_i) \cdot Pr(C = c_i)$$

Naïve Bayes classifier

- Strong simplifying assumption: attributes are pairwise independent

$$Pr(A_1 = a_1, \dots, A_k = a_k \mid C = c_i) = \prod_{j=1}^k Pr(A_j = a_j \mid C = c_i)$$

- $Pr(C = c_i)$ is fraction of training data with class c_i
 - $Pr(A_j = a_j \mid C = c_i)$ is fraction of training data labelled c_i for which $A_j = a_j$
- Final classification is

$$\arg \max_{c_i} Pr(C = c_i) \prod_{j=1}^k Pr(A_j = a_j \mid C = c_i)$$

Naïve Bayes classifier . . .

- Conditional independence is not theoretically justified
- For instance, text classification
 - Items are documents, attributes are words (absent or present)
 - Classes are topics
 - Conditional independence says that a document is a set of words: ignores sequence of words
 - Meaning of words is clearly affected by relative position, ordering
- However, naive Bayes classifiers work well in practice, even for text classification!
 - Many spam filters are built using this model

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- “Pad” training data with one sample for each value a_j — m_i extra data items
- Adjust $Pr(A_i = a_i | C = c_j)$ to $\frac{n_{ij} + 1}{n_j + m_i}$ where
 - n_{ij} is number of samples with $A_i = a_i, C = c_j$
 - n_j is number of samples with $C = c_j$

- Laplace's law of succession

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- $\lambda = 1$ is Laplace's law of succession

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- Want to use a naïve Bayes classifier
- Need to define a generative model
- How do we represent documents?

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No reference to length of d

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- $Pr(d) = \sum_{c \in C} Pr(d | c)$

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- Recall $Pr(d | c) = \prod_{w_i \in d} Pr(w_i | c) \prod_{w_i \notin d} (1 - Pr(w_i | c))$

Bag of words model

- Each document is a **multiset** or **bag** of words over a vocabulary

$$V = \{w_1, w_2, \dots, w_m\}$$

- Count multiplicities of each word

Universe U

Set $S \subseteq U$

$$f_S : U \rightarrow \{0, 1\}$$

char. fn

$$f_S : U \rightarrow \mathbb{N}_0$$

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- As before
 - Each topic c has probability $Pr(c)$
 - Each word $w_i \in V$ has conditional probability $Pr(w_i | c_j)$ with respect to each $c_j \in C$ (but we will estimate these differently)
 - Note that $\sum_{i=1}^m Pr(w_i | c_j) = 1$
 - Assume document length is independent of the class

Bag of words model

- Generating a random document d
 - Choose a document length ℓ with $Pr(\ell)$ *Will disappear*
 - Choose a topic c with probability $Pr(c)$
 - Recall $|V| = m$.
 - To generate a single word, throw an m -sided die that displays w with probability $Pr(w | c)$
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- $Pr(d | c) = Pr(\ell) \ell! \prod_{j=1}^m \frac{Pr(w_j | c)^{n_j}}{n_j!}$

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$$\blacksquare Pr(w_i | c_j) = \frac{\sum_{d \in D_j} n_{id}}{\sum_{t=1}^m \sum_{d \in D_j} n_{td}}$$

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only if $d \in D_j$

$$\text{since } Pr(c_j | d) = \begin{cases} 1 & \text{if } d \in D_j, \\ 0 & \text{otherwise} \end{cases}$$

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Classification

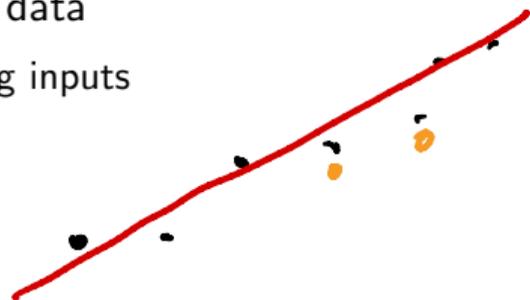
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Limitations of classification models

- **Bias** : Expressiveness of model limits classification
 - For instance, linear logistic regression
- **Variance**: Variation in model based on sample of training data
 - Shape of a decision tree varies with distribution of training inputs



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*Bias - Variance
tradeoff*

Models with high variance are expressive but **unstable**

- In principle, a decision tree can capture an arbitrarily complex classification criterion
- Actual structure of the tree depends on impurity calculation
- Danger of overfitting: model tied too closely to training set

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- In principle, a decision tree can capture an arbitrarily complex classification criterion
- Actual structure of the tree depends on impurity calculation
- Danger of overfitting: model tied too closely to training set
- Is there an alternative to pruning?

Ensemble models

- Sequence of independent training data sets D_1, D_2, \dots, D_k
- Generate models M_1, M_2, \dots, M_k
- Take this **ensemble** of models and “average” them
 - For regression, take the mean of the predictions
 - For classification, take a vote among the results and choose the most popular one

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- Take this **ensemble** of models and “average” them
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 - For classification, take a vote among the results and choose the most popular one
- **Challenge:** Infeasible to get large number of independent training samples
- Can we build independent models from a single training data set?
 - Strategy to build the model is fixed
 - Same data will produce same model

Bootstrap Aggregating = Bagging

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 - $TD = \{d_1, d_2, \dots, d_N\}$
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 - Repeat K times

Sample of size K

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 - Put it back into the set
 - Repeat K times
- Some items in the sample will be repeated
- If sample size is same as data size ($K = N$), expected number of distinct items is $(1 - \frac{1}{e}) \cdot N$
 - Approx 63.2%

Bootstrap Aggregating = Bagging

- Sample with replacement of size N : bootstrap sample
 - Approx 2/3 of full training data

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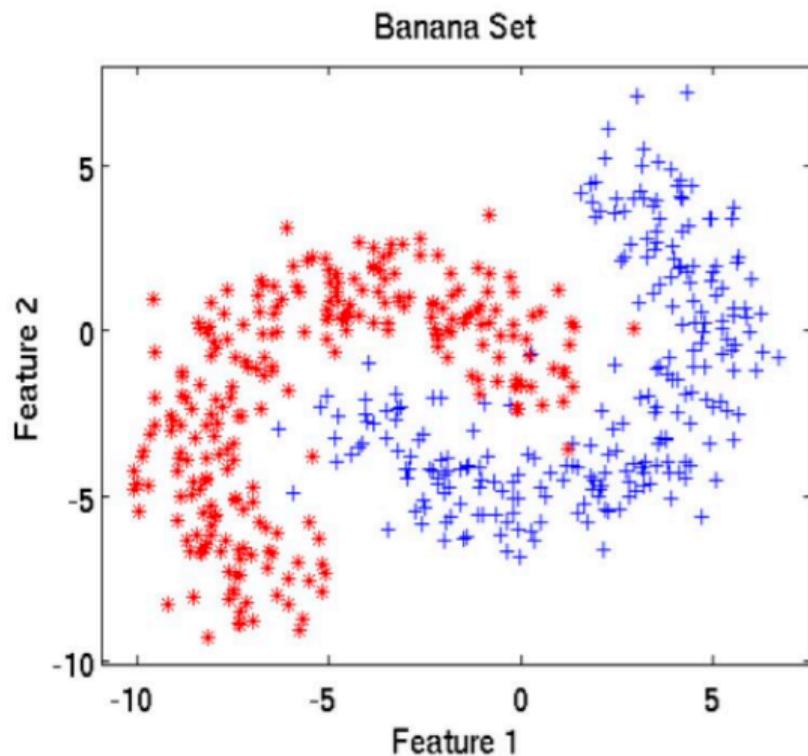
D_1, \dots, D_k

Expect each
original item
to be in
 $\approx 2/3$ of D_i 's

Bootstrap Aggregating = Bagging

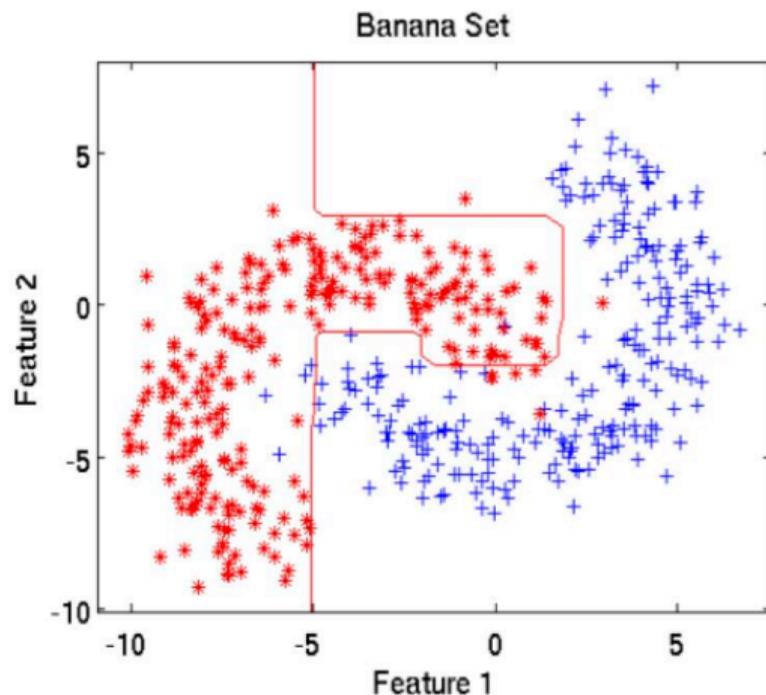
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 - Models will vary because each uses different training data
- Final classifier: report the majority answer
 - Assumptions: binary classifier, k odd
- Provably reduces variance

Bagging with decision trees



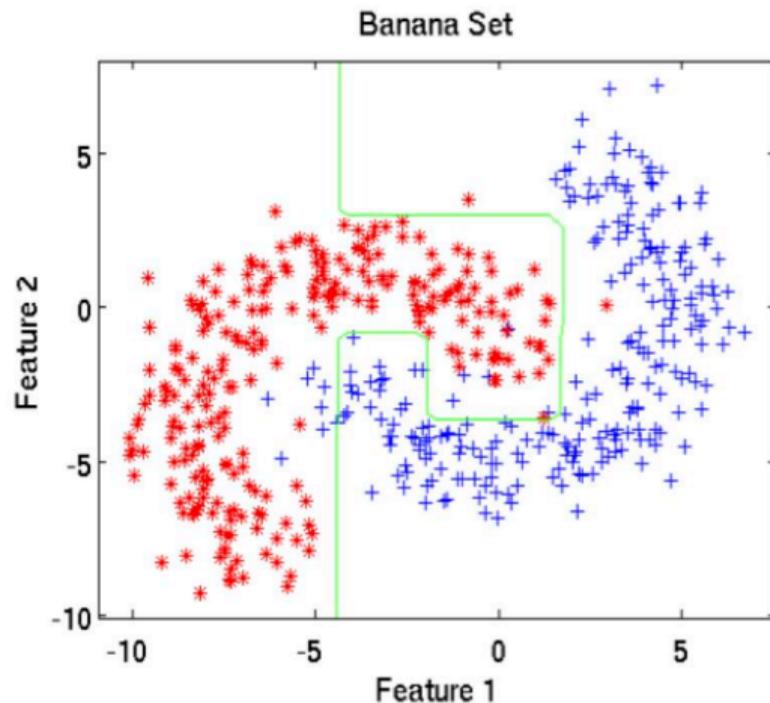
Training data

Bagging with decision trees



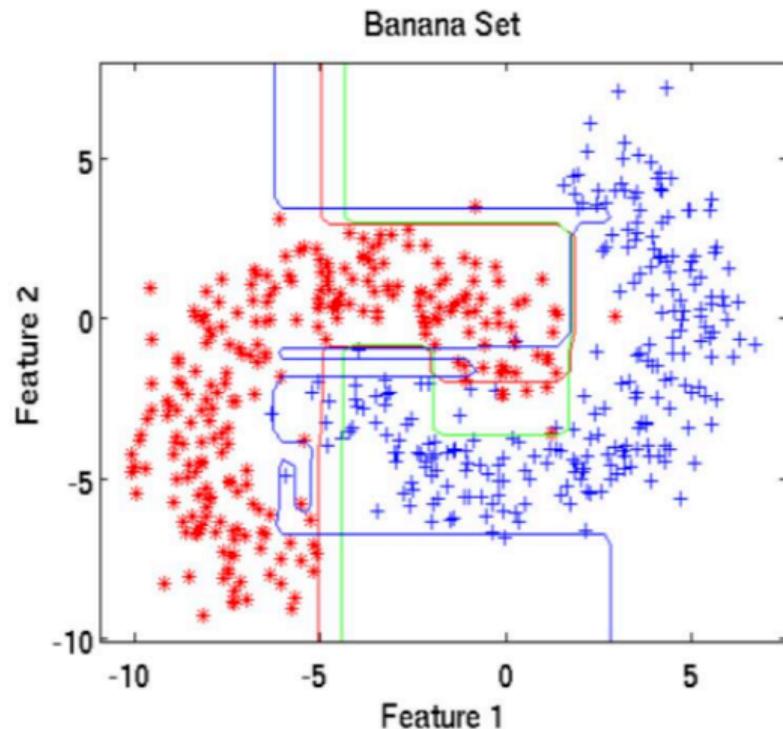
Decision boundary produced
by one tree

Bagging with decision trees



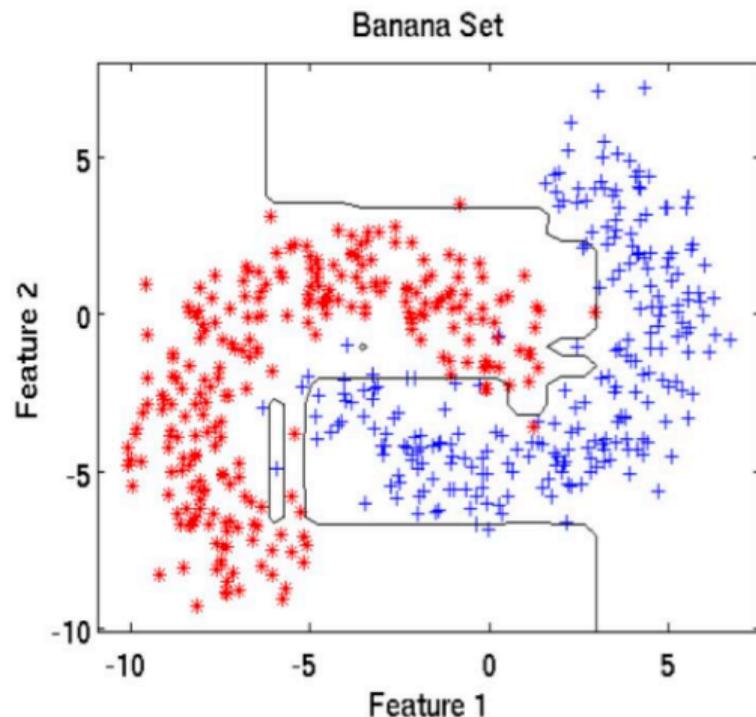
Decision boundary produced by a second tree

Bagging with decision trees



Three trees and final boundary overlaid

Bagging with decision trees



Final result from bagging all trees.

When to use bagging

- Bagging improves performance when there is high variance
 - Independent samples produce sufficiently different models

When to use bagging

- Bagging improves performance when there is high variance
 - Independent samples produce sufficiently different models
- A model with low variance will not show improvement
 - **k-nearest neighbour** classifier
 - Given an unknown input, find k nearest neighbours and choose majority
 - Across different subsets of training data, variation in k nearest neighbours is relatively small
 - Bootstrap samples will produce similar models

Random Forest

- Applying bagging to decision trees with a further twist

Graph theory

Forest =
Collection of
trees

Random Forest

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 - Instead, fix a small limit $m < M$ — say $m = \log_2 M + 1$
 - At each level, choose a random subset of available attributes of size m
 - Evaluate only these m attributes to choose next query

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 - At each level, choose a random subset of available attributes of size m
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 - No pruning — build each tree to the maximum
- Final classifier: vote on the results returned by T_1, T_2, \dots, T_k

- Theoretically, overall error rate depends on two factors
 - **Correlation** between pairs of trees — higher correlation results in higher overall error rate
 - **Strength (accuracy)** of each tree — higher strength of individual trees results in lower overall error rate

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- Reducing m , the number of attributes examined at each level, reduces correlation and strength
 - Both changes influence the error rate in opposite directions

m out of M
features

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Random Forest ...

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- Increasing m increases both correlation and strength
- Search for a value of m that optimizes overall error rate

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- **Oob classification** — for each d , vote only among those T_i where d is oob for D_i
- Use oob samples to validate the model
 - Estimate generalization error rate of overall model based on error rate of oob classification
 - Do not require a separate test data set

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- Compute weighted average of impurity gain
 - Weight is given by number of training samples at the node

Regression

Discover line $l = \theta_0 + \theta_1 x_1 + \theta_2 x_2 + \dots + \theta_m x_m$

Higher $\theta_i \Rightarrow$ higher impact of A_i

Requires normalization