Lecture 10 – Classification and Regression 2

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Objectives

Decision tree
Splitting rule: information gain
Splitting rule: Gini coefficient
Decision tree generation
Random forest
Image analysis example

Decision tree

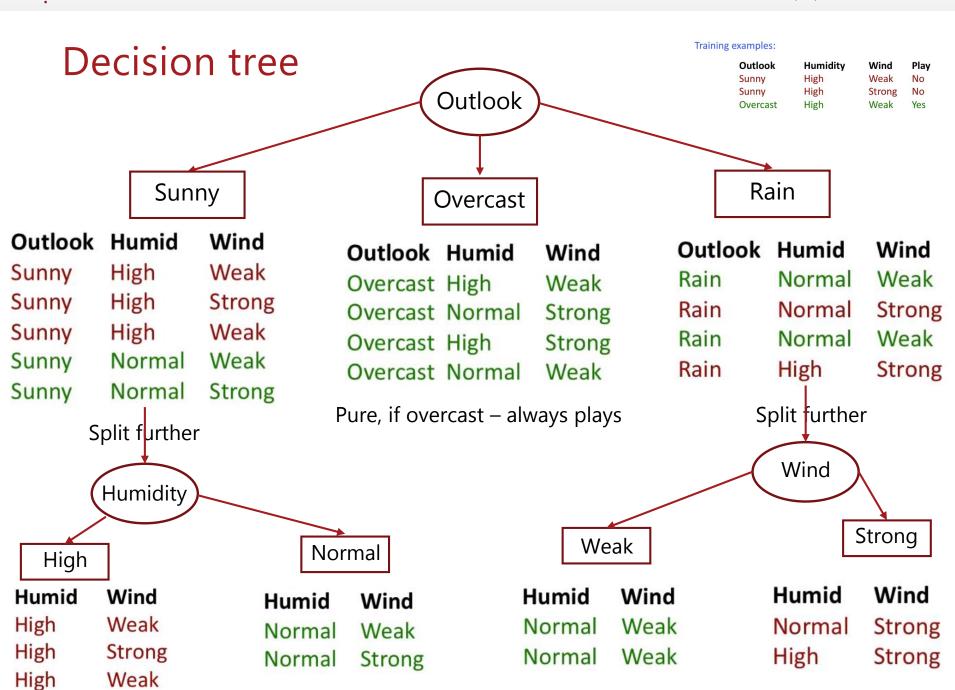
We want to predict if John plays tennis in a given day

 We collected history of 13 days with some factors we believe may be important

- We want make a prediction for new days:
 - D14 [Rain, High, Weak]
- Divide and conquer:
 - Split training samples according to some attribute
 - If the result is pure stop, overwise continue splitting

Training examples:

Outlook	Humidity	Wind	Play
Sunny	High	Weak	No
Sunny	High	Strong	No
Overcast	High	Weak	Yes
Rain	Normal	Weak	Yes
Rain	Normal	Strong	No
Overcast	Normal	Strong	Yes
Sunny	High	Weak	No
Sunny	Normal	Weak	Yes
Rain	Normal	Weak	Yes
Sunny	Normal	Strong	Yes
Overcast	High	Strong	Yes
Overcast	Normal	Weak	Yes
Rain	High	Strong	No

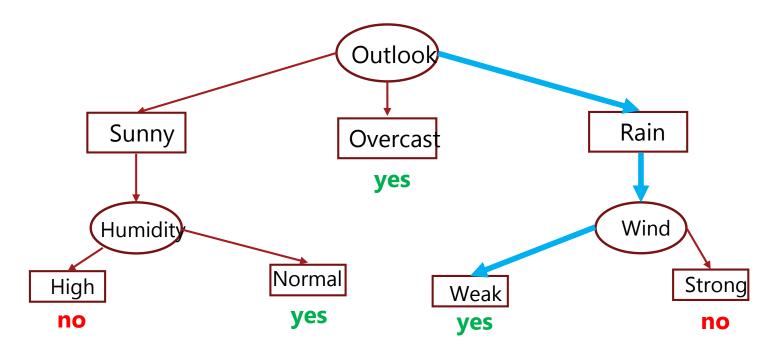


Decision tree

Training examples:

OutlookHumidityWindPlaySunnyHighWeakNoSunnyHighStrongNoOvercastHighWeakYes

We generated a decision tree



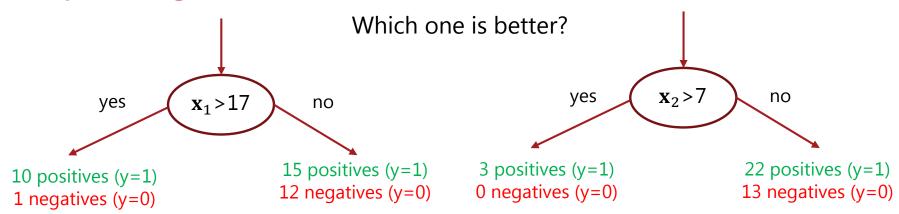
For our test example [Rain, High, Weak], we predict John to play

Binary decision tree: algorithm

- Training dataset $\mathbf{T} = \{(\mathbf{x}_1, \mathbf{y}_1), (\mathbf{x}_2, \mathbf{y}_2), ..., (\mathbf{x}_n, \mathbf{y}_n)\}; y \in \{0,1\}$:
- Select feature j using the current subset of examples T.
- Find the optimal split of T using feature j:
 - The new subsets T_0 and T_1 .
- Continue for each T_0 and T_1 :
 - If stopping criteria is not reached for T_k , go to step 1

How to select j? How to split T?

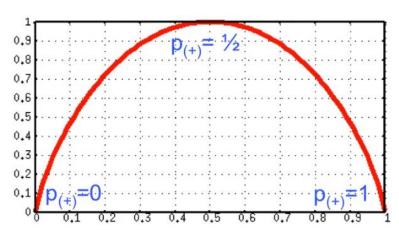
Splitting rule



- Intuitively:
 - Purity (more certain separation) :
 - 9 pos / 1 neg => very certain separation (90%)
 - 5 pos / 4 neg => not certain (56%)
 - Power of separation:
 - Let's say we have 30 samples
 - Separation with 10 pos / 2 neg is better than
 - Separation with 5 pos / 1 neg
 - Symmetric:
 - 10 pos / 2 neg is as good as 2 pos / 10 neg

Splitting rule: entropy

- Entropy $H(T) = -p_+ \log_2 p_+ p_- \log_2 p_-$:
 - *H*(9 pos, 1 neg):
 - $H(T) = -0.9 \log_2 0.9 0.1 \log_2 0.1 = 0.137 + 0.332 = \mathbf{0.469}$
 - *H*(5 pos, 4 neg):
 - $H(T) = -0.56 \log_2 0.56 0.44 \log_2 0.44 = 0.468 + 0.521 = 0.989$



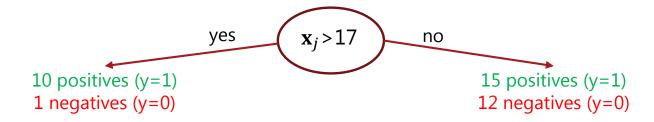
- *H*(8 class1, 7 class2, 5 class3):
 - $H(T) = -0.4 \log_2 0.4 0.35 \log_2 0.35 0.25 \log_2 0.25 = 0.137 + 0.332 = 1.56$
- H(11 class1, 9 class2, 0 class3):
 - $H(T) = -0.55 \log_2 0.55 0.45 \log_2 0.45 0 \log_2 0 = 0.468 + 0.521 = 0.993$



Splitting rule: information gain

Sets with good separation and large power:

$$Gain(T, j) = H(T) - \sum_{i=0,1} \frac{|T_i|}{|T|} H(T_i)$$



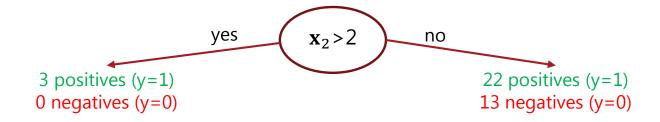
$$Gain(T, j)$$

= $H(25pos, 13pos) - \frac{11}{38}H(10pos, 1pos) - \frac{27}{38}H(15pos, 12pos)$
= $0.927 - \frac{11}{38}0.44 - \frac{27}{38}0.99 = 0.096$

Splitting rule: information gain

Sets with good separation and large power:

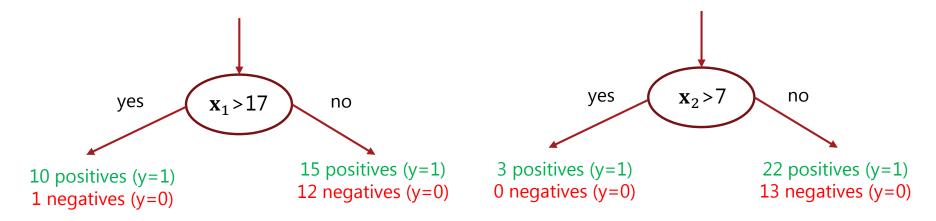
$$Gain(T,j) = H(T) - \sum_{i=0,1} \frac{|T_i|}{|T|} H(T_i)$$



$$Gain(T, j)$$

= $H(25pos, 13pos) - \frac{3}{38}H(3pos, 0pos) - \frac{35}{38}H(22pos, 13pos)$
= $0.927 - \frac{3}{38}0 - \frac{35}{38}0.95 = 0.05$

Splitting rule: information gain



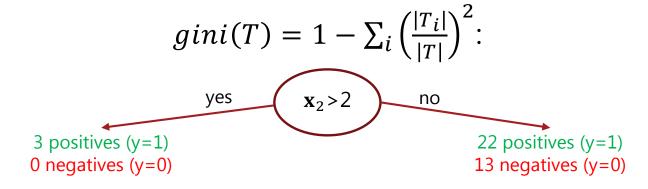
Information gain is 0.096

Information gain is 0.05

- We want to maximize the gain:
 - It is better to split using $x_1 > 17$
 - $Gain(T,j) = H(T) \sum_{i=0,1} \frac{|T_i|}{|T|} H(T_i)$

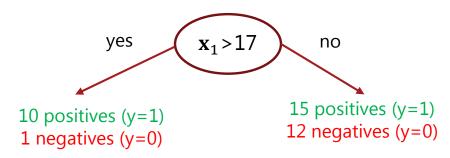
Splitting rule: Gini index

Alternative splitting rule:



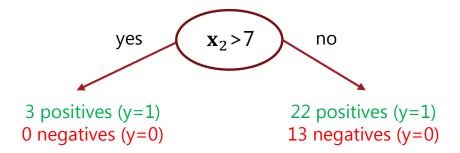
$$Gini(T,j) = \frac{3}{38}gini(T_0) + \frac{35}{38}gini(T_1)$$
$$= \frac{3}{38}(1-1-0) + \frac{35}{38}(1-0.395-0.138) = 0.509$$

Splitting rule: Gini

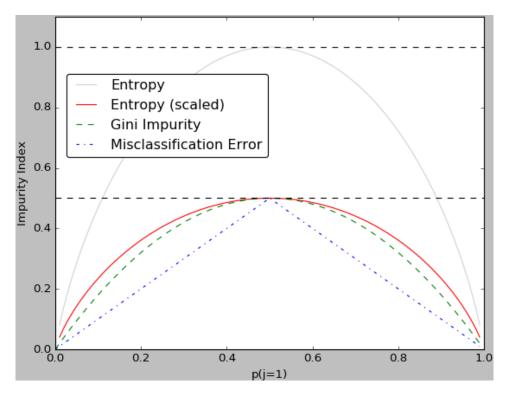


Gini index is 0.399

We want to minimize the Gini index



Gini index is 0.509

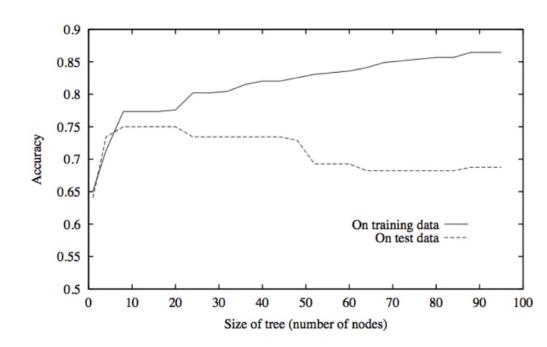


Stopping condition

Overfitting problem:

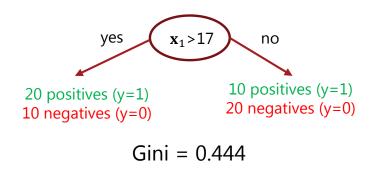
- If we allow the tree to grow until all terminal leaves are pure, the training classification will be perfect
- Leaves can contain only one element
- Testing accuracy may be unreliable

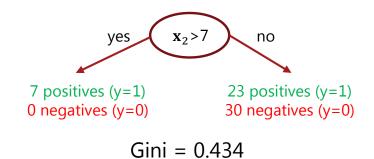
- How can we prevent it?
 - Validation data
 - Restrict tree depth
 - Restrict minimal leaf power
 - Pruning



Stopping condition

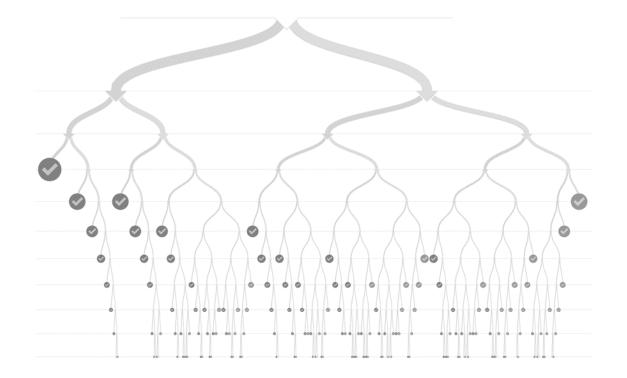
- Restrict the tree complexity:
 - No more than N splits are allowed before a leaf
 - Split cannot have less that x% of the original data
 - Allow only statistically significant restrictions
- What can be the problem with such restrictions?
 - Imbalanced splits





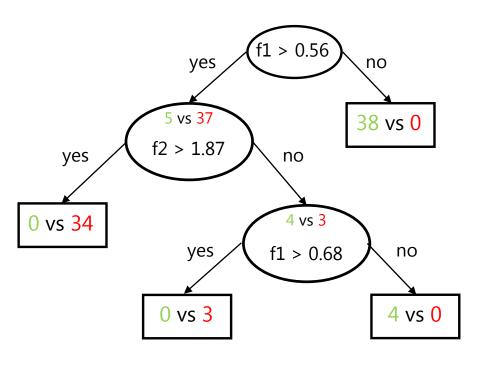
Pruning

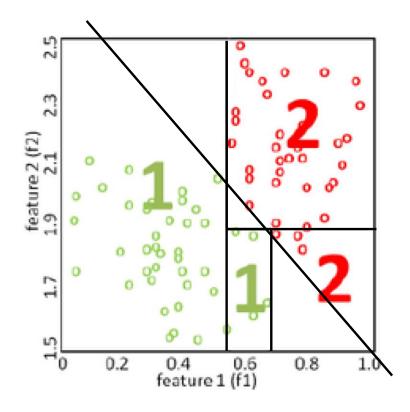
- Database is split into training and validation parts
- Go through every last level split:
 - Compare validation accuracy with the split and without (converted into a leaf)
 - If accuracy increases, replace the split with the leaf
- Go to next levels



Decision boundary

- Two classes in 2D space need to be separated
- Can decision tree generate such a visually-optimal border?





Decision boundary

How the pruned tree will perform on the following test examples:

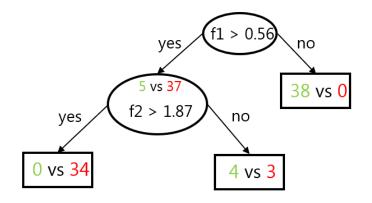
predictions

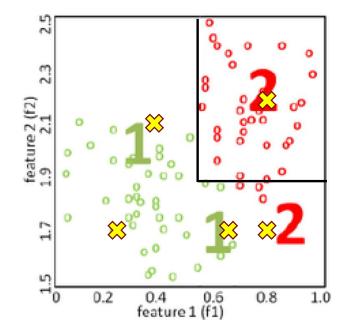
•
$$[0.25, 1.7]$$
 38 vs $0 = [1, 0]$

•
$$[0.65, 1.7]$$
 4 vs 3 => $[0.57, 0.43]$

•
$$[0.4, 2.1]$$
 38 vs $0 = [1, 0]$

•
$$[0.8, 2.2]$$
 0 vs 34 => $[0, 1]$





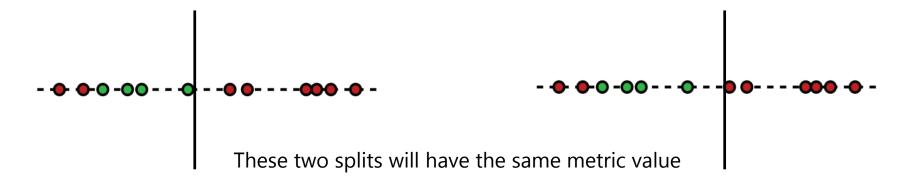
Selection of the threshold value

 What kind of intuitive rules we can establish during selection of the optimal threshold?



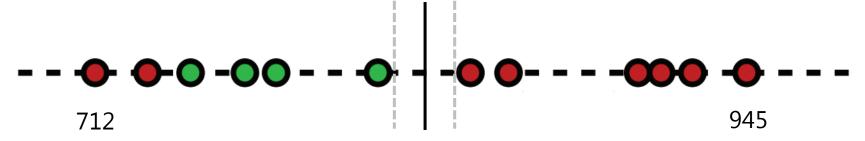
- Do not look outside interval of feature values [712, 945]
- The split quality metrics assign binary values for each sample.

Entropy =
$$-p_+ \log_2 p_+ - p_- \log_2 p_-$$
 Gini = $1 - \sum_{i=0,1} \left(\frac{|T_i|}{|T|}\right)^2$



Selection of the threshold value

 Only average points between neighboring samples need to be considered



Only check different-class neighbors



There is at least one green point.

Moving the threshold ever right or left will definitely improve separation

21

Selection of the threshold value

Algorithm:

Sort input samples according to a feature of interest X:

$$\{(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)\}$$

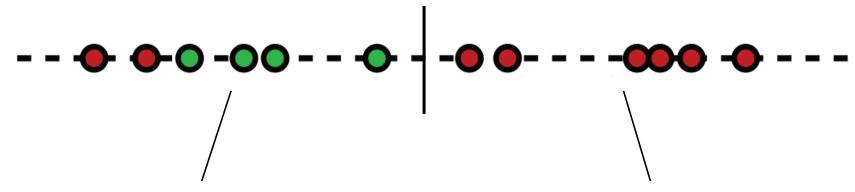
Consider the following split points:

$$x_i + \frac{(x_{i+1} - x_i)}{2}$$
: $y_{i+1} \neq x_i$

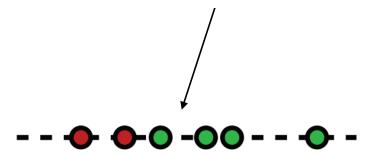
Find the optimal split point by computing splitting metrics

Selection of the threshold value

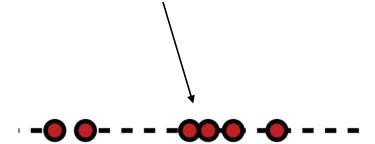
 Sorting is very expensive step, but it needs to be done only once at the beginning for every feature:



After splitting, we got to already sorted pieces



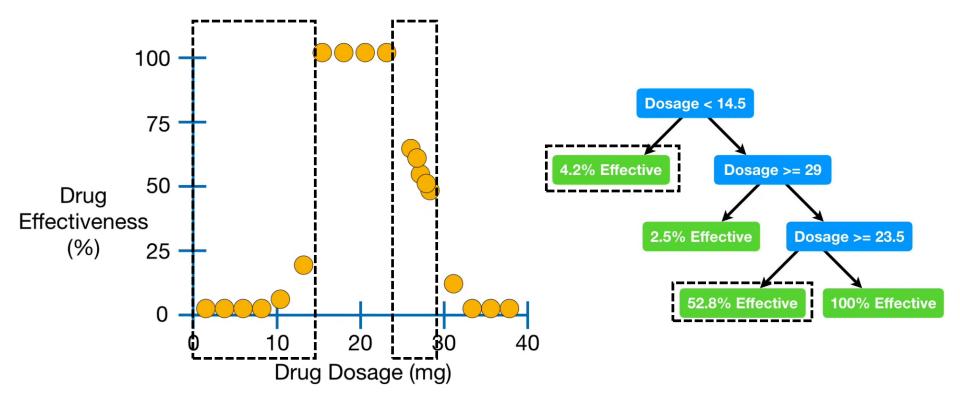
After splitting, we got to already sorted pieces



Regression tree

Instead of classification, decision trees can do regression. Two challenges:

- Computing prediction:
 - Average values of the leaf
 - Some simple model of the values on the final leaf



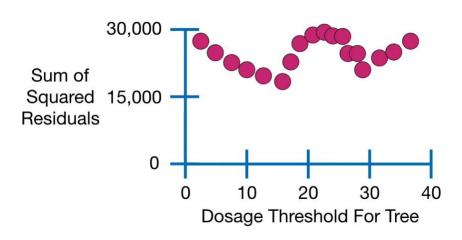
Splitting rule: reduction of variance

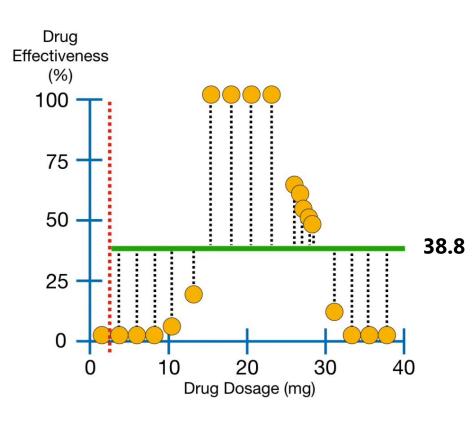
Splitting rule:

$$F(T) = \sum_{x \in T} (x - \bar{x})^2$$

$$(0 - 0)^2 + (0 - 38.8)^2 + (0 - 38.8)^2 + (0 - 38.8)^2$$

+ $(5 - 38.8)^2 + (20 - 38.8)^2 + (100 - 38.8)^2$
+ $(100 - 38.8)^2 + ... + (0 - 38.8)^2$
= 27465.5





How would you quantify the quality of this split?

Splitting rule: categorical variables

If a feature is not numerical but categorical:

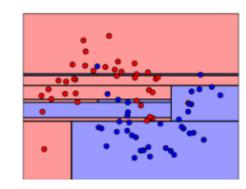
- Binary {male/female}, {had history of disease/ no history}
- Non-binary utilized {drug A/drug B/drug C}

For non-binary categorical features:

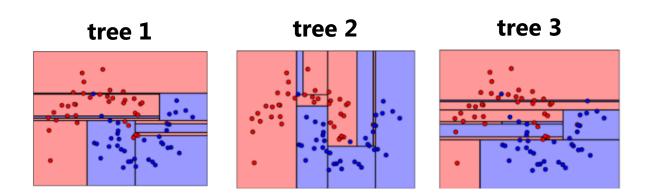
Check binary splits in the form drug A vs. drug not A

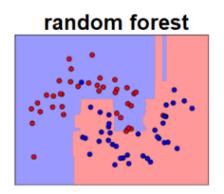
Random Forests

- Single tree cannot generalize well on complex data
- Idea is to generate many weaker trees instead of one full tree



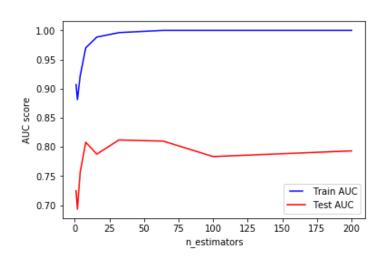
- Introduce randomness during training of each tree:
 - Train each tree on a random subset of training samples S_r
 - During split generation, only consider a random subsample of features $d \ll D$

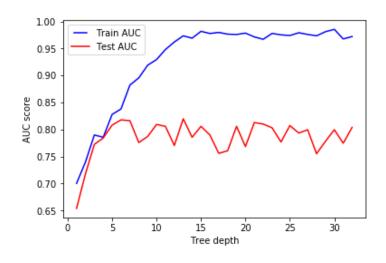




Random forests

Tree depth vs number of trees



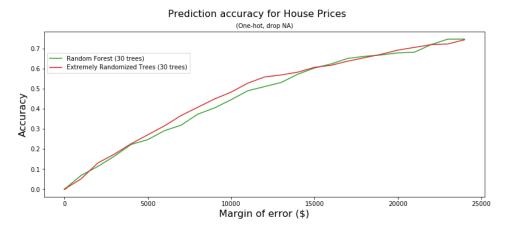


• How important are ratios $|S_r|/|S|$ and |d|/|D|?

•
$$|d| = \sqrt{|D|}$$

•
$$|d| = |D|$$

•
$$|d| = 1$$



Random forests: kernels

• We previously observed axis-aligned kernels:

$$h(\mathbf{x}, i, \tau) = [\tau_1 \ge x_i \ge \tau_2]$$
 e. g. $[\infty \ge x_i \ge 15.5]$



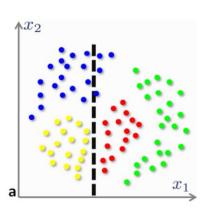
$$h(\mathbf{x}, \phi, \mathbf{\psi}, \tau) = [\tau_1 \ge \phi(\mathbf{x})^T \cdot \mathbf{\psi} \ge \tau_2]$$

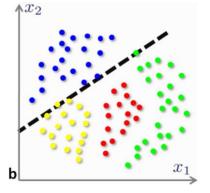
e. g. $\phi(\mathbf{x}) = [x_i, x_j, 1]; \quad \mathbf{\psi} = [\psi_1, \psi_2, \psi_3]$

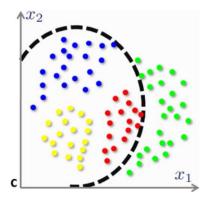
Non-linear kernel:

$$h(\mathbf{x}, \phi, \mathbf{\psi}, \tau) = [\tau_1 \ge \phi(\mathbf{x})^T \cdot \mathbf{\psi} \cdot \phi(\mathbf{x}) \ge \tau_2]$$

e. g. $\phi(\mathbf{x}) = [x_i, x_j, 1]; \quad \mathbf{\psi} = \mathbb{R}^{3\times3}$

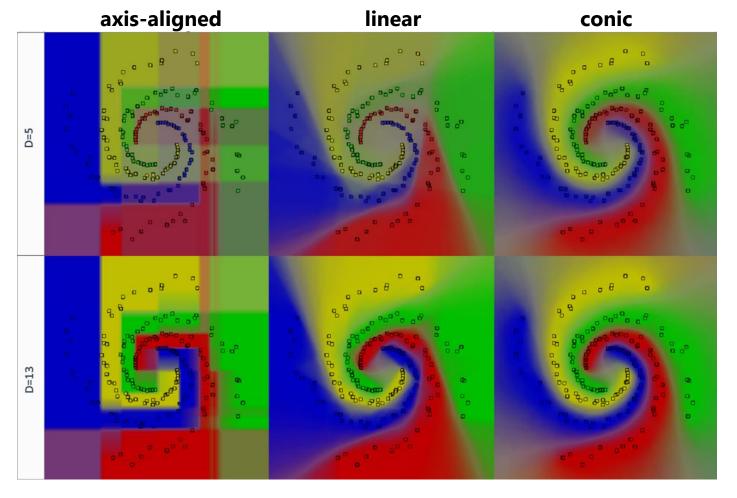






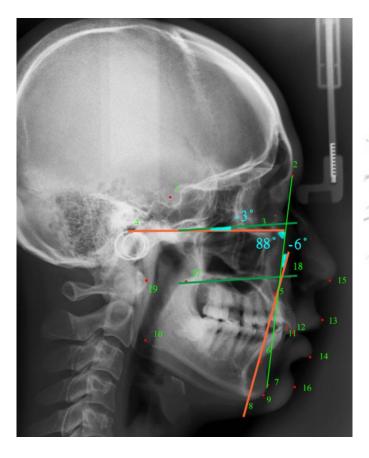
Random forests: kernels

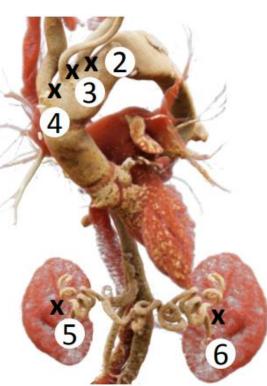
• Kernels are computationally expansive, the number of different combinations of i, j, ... in $\phi(\mathbf{x}) = [x_i, x_j, 1]$ grows exponentially



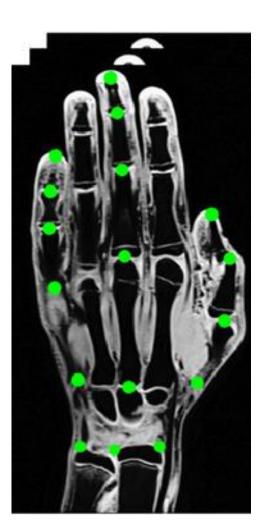
Random forests example: landmarking

Landmark detection in medical imaging

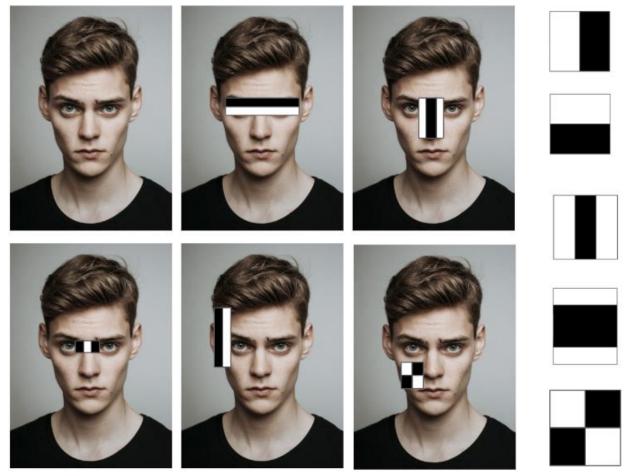




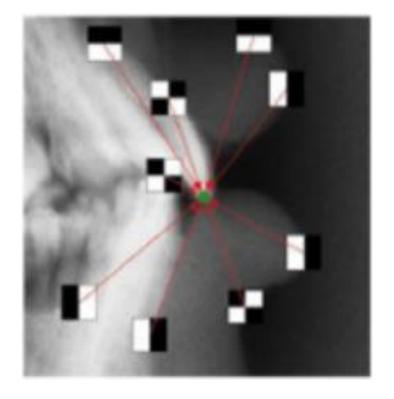




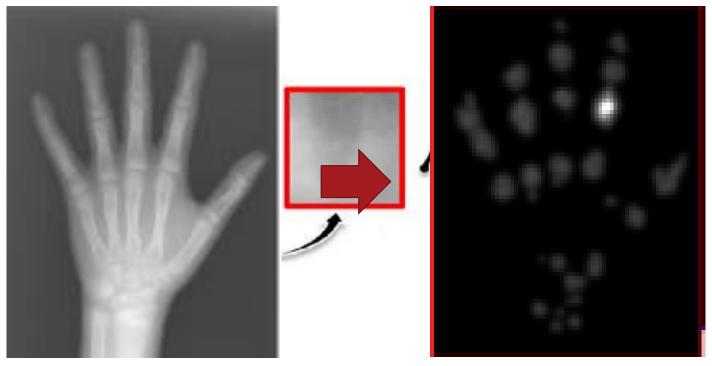
- Single pixel intensity
- Box feature (or Haar features)



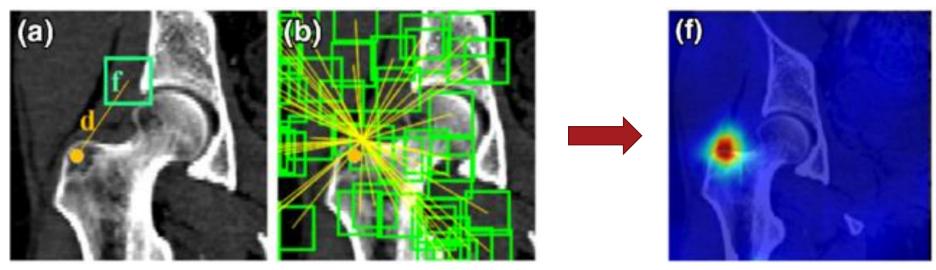
- How to find out which features are important to detect an object?
- Collect many images with the object
- Extract all features around the object examples in the images to form positive samples
- Extract all features around random pixels* in the image to form negative samples
- Generate random forest that will recognize important features and create rules for the object recognition



- For a new image:
 - Extract features for all pixels, and classify these pixels using the random forest
 - Find the pixels with the highest probability to be the landmark predicted by the random forest



- What type of the random forest was used for landmark detection example?
 - Classification / Regression
- Can we reformulate it as a regression?
 - Hint: if classification forest predicts pixels to be non-landmark, we just throw this pixel away. Can we still extract something useful from it?



Predict the displacement from pixel f to landmark d

Questions?