

α by cross-validation

	$\alpha = \alpha_1$	$\alpha = \alpha_2$...	$\alpha = \alpha_B$	
$x^{(j)}_{K=1}$	$C_{\alpha_1}^{(j)}$				$C_{\alpha}(j)$
$K=2$	\vdots	\vdots		\vdots	
\vdots					
$K=K$	$C_{\alpha_K}^{(j)}$				
	$CV(\alpha_1)$	$CV(\alpha_2)$...	$CV(\alpha_B)$	
	\min		α which minimize the CV		

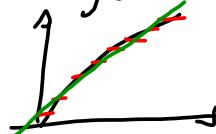
Advantages of trees

- simplicity and ease of communication;
- compact representation;
- speed of computation (easy to parallelize)
- easy to handle both continuous and categorical variables (mixture of them)
- easily implement different loss functions;
- easy to handle missing data;
- automatic variable selection



Disadvantages

- instability of the results \rightarrow bagging, boosting and random forest
- difficulties with online computation (difficult to integrate new information)
- very hard to approximate very steep functions
- there are no formal statistical procedures (tests of hypotheses, confidence intervals, ...)
- difficult to evaluate variable importance



Classification trees

Start $p=1$, $k=2$

Instead of estimating $P(x)$, we try to estimate $\hat{P}_k(x) = \Pr[Y=1|x=x]$

$$\hat{P}_k(x) = \sum_{h=1}^J P_h \mathbb{1}(x \in R_h)$$

where P_h is the probability that $Y=1$ given $x \in R_h$. We allocate each observation to 0 or 1 depending $\hat{P}_k(x)$, if $\hat{P}_k(x) > t$, t being a threshold (usually $\frac{1}{2}$)

To estimate P_h , we can simply use the frequency of 1 in R_h

$$\hat{P}_h = \frac{1}{n_j} \sum_{i:x_i \in R_h} \mathbb{1}(y_i = 1)$$

Given the binary nature of Y , we need to use a specific version of the deviance (binomial deviance, see notes of lecture 5) as the loss function

$$D = -2 \sum_{i=1}^n y_i \log \hat{P}_h + (1-y_i) \log (1-\hat{P}_h)$$

which can be rewritten as

$$D = -2 \sum_{h=1}^J n_h (\hat{P}_h \log \hat{P}_h + (1-\hat{P}_h) \log (1-\hat{P}_h)) = \sum_{h=1}^J D_h$$

because the probability of $Y=1$ is constantly equal to P_h in each region R_h

Consider D in the form

$$D = 2n \sum_{h=1}^J \frac{n_h}{n} Q(\hat{P}_h)$$

weight proportional
 to the size of leaves

measure of impurity
 how much the elements of a region
 are non-homogeneous
 $\hookrightarrow Q=0$ if all observations are 0 or 1
 $\hookrightarrow Q=\frac{1}{2}$ in the case of maximum
 entropy (Soil/Sox)

Where

$$Q(P_h) = - \sum_k P_{jk} \log P_{jk} = - (P_{j1} \log P_{j1} + P_{j0} \log P_{j0}) \quad k=2$$

$$= - (P_{j1} \log P_{j1} + (1-P_{j1}) \log (1-P_{j1}))$$

We can then use different measures of impurity, e.g.

Gini index $Q(P_h) = \sum_{k \in \{0,1\}} P_{jk} (1-P_{jk})$

missclassification error $\sum_{h=1}^J \frac{1}{n_h} \sum_{i:x_i \in R_h} \mathbb{1}(y_i \neq k_h)$

