## 2. Greedy is good for $p \gg n$ : Boosting

data:  $(X_1, Y_1), \ldots, (X_n, Y_n)$  (i.i.d. or stationary), predictor variables  $X_i \in \mathbb{R}^p$ response variables  $Y_i \in \mathbb{R}$  or  $Y_i \in \{0, 1, \dots, J-1\}$ aim: estimation of function  $f(\cdot): \mathbb{R}^p \to \mathbb{R}$  (including feature selection) e.g.  $f(x) = \mathbf{E}[Y|X = x] \text{ or } f(x) = \mathbb{P}[Y = 1|X = x] \text{ with } Y \in \{0, 1\}$ or distribution of survival time Y given X depends on some function f(X) only our setting: typically p is very large historically: Boosting is an ensemble scheme (multiple predictions and averaging) 6 base procedure: algorithm A  $\hat{ heta}(\cdot)$  (a function estimate) data e.g.: simple linear regression, tree, MARS, "classical" smoothing, neural nets, ... generating multiple predictions: algorithm A  $\hat{\theta}_1(\cdot)$ weighted data 1 algorithm A  $\hat{\theta}_2(\cdot)$ weighted data 2 . . .  $\stackrel{\text{algorithm A}}{\longrightarrow} \quad \hat{\theta}_M(\cdot)$ weighted data M Aggregation:  $\hat{f}_A(\cdot) = \sum_{m=1}^M a_m \hat{\theta}_m(\cdot)$ data weights? averaging weights  $a_m$ ?

classification of 2 lymph nodal status in breast cancer using gene expressions from microarray data:

n = 33, p = 7129 (for CART: gene-preselection, reducing to p = 50)

method	test set error	gain over CART
CART	22.5%	_
LogitBoost with trees	16.3%	28%
LogitBoost with bagged trees	12.2%	46%

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**2.1. Boosting algorithms** 

 AdaBoost proposed for classification by Freund & Schapire (1996)

 data weights (rough original idea): large weights to previously heavily misclassified instances (sequential algorithm)

 averaging weights  $a_m$ : large if in-sample performance in mth round was good

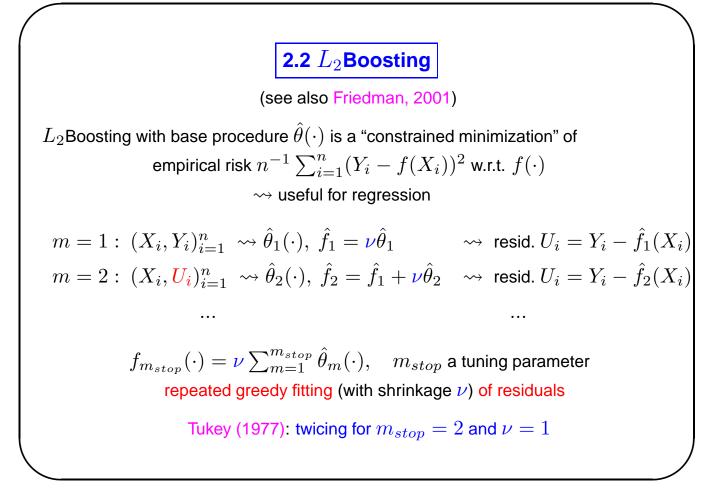
 Why should this be good?

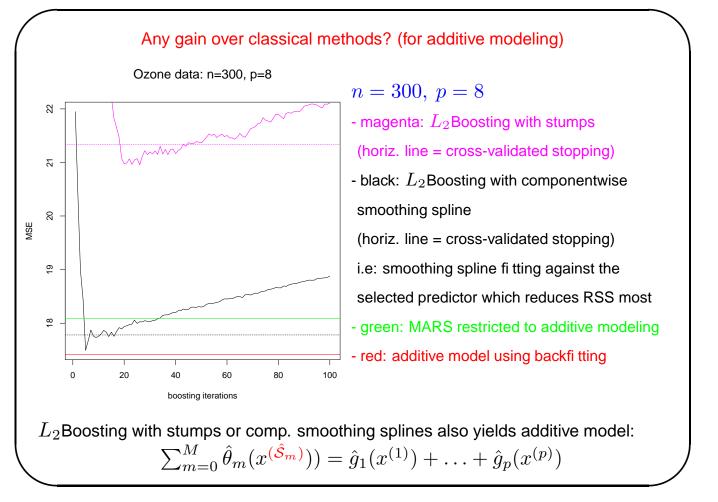
 (actually: other weighting schemes are equally good or better...)

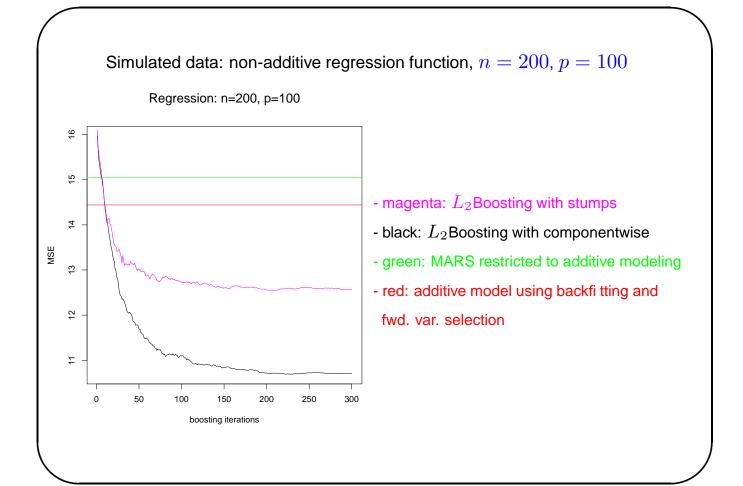
Breiman (1998/99):

AdaBoost is functional gradient descent (FGD) procedure

a mix of statistical estimation and numerical optimization...







similar for classification

very often: boosting performs comparatively well in high-dimensions (there is a lot of empirical evidence for this)

also SVM is often surprisingly accurate...

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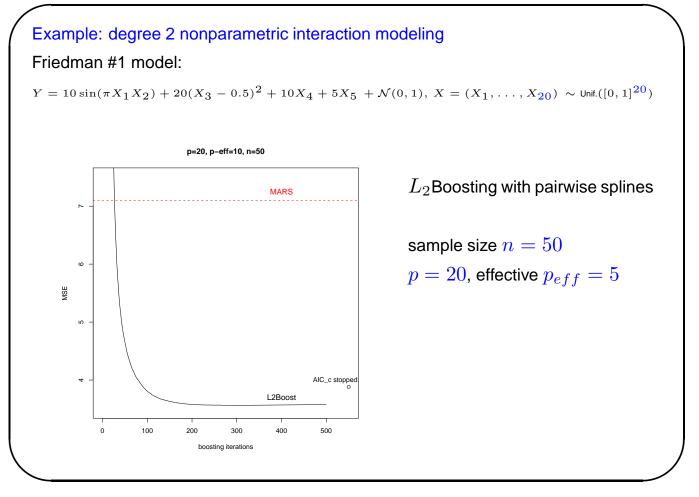
### 2.3. Choice of the base procedure

most popular in machine learning: tree algorithms (CART, C4.5) they do variable/feature selection

have seen: for componentwise smoothing splines or stumps

 $\longrightarrow$  boosting yields an additive model fit

→ we can use boosting for fitting in "quite many" structural models



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2.4.  $L_2 \mbox{Boosting}$  for high-dimensional linear models

linear model

$$Y = f(X) + \varepsilon,$$
  
$$f(x) = \sum_{j=1}^{p} \beta_j x^{(j)}, \quad p \gg n$$

or: a highly over-complete dictionary  $\{g_j(\cdot); j = 1, \ldots, p \gg n\}$ our approach:  $L_2$ Boosting with componentwise linear LS regression This base procedure fits a univariate linear regression model against the one predictor variable which reduces residual sum of squares most

first round of estimation: selected predictor variable 
$$X^{(\hat{S}_1)}$$
 (e.g. =  $X^{(3)}$ )  
corresponding ordinary least squares  $\hat{\beta}_{\hat{S}_1}$   
use shrunken fit  $\hat{f}_1 = \nu \hat{\beta}_{\hat{S}_1} X^{(\hat{S}_1)}$  (e.g.  $\nu = 0.1$ )  
second round of estimation: selected predictor variable  $X^{(\hat{S}_2)}$  (e.g. =  $X^{(21)}$ )  
corresponding OLS  $\hat{\beta}_{\hat{S}_2}$   
use shrunken fit  $\hat{f}_2 = \hat{f}_1 + \nu \hat{\beta}_{\hat{S}_2} X^{(\hat{S}_2)}$   
etc.  
very different from forward variable selection  
this method does variable selection and  
assigns variable amount of degrees of freedom for selected variables (shrinkage)  
not full OLS on selected variables (even with  $\nu = 1$ )  
For  $\nu = 1$ , this  $L_2$ Boosting is known as Matching Pursuit (Mallat and Zhang, 1993)



because of

variable selection and

assigning variable amount of degrees of freedom (shrinkage) for selected variables reminds to Lasso ( $\ell^1$ -penalized regression) (Tibshirani, 1996)

$$\hat{\beta}_{Lasso} = \operatorname{argmin}_{\beta} \sum_{i=1}^{n} (Y_i - \sum_{j=1}^{p} \beta_j X_i^{(j)})^2 + \underbrace{\lambda}_{\geq 0} \sum_{j=1}^{p} |\beta_j|$$

and indeed: there is a relation (Efron, Hastie, Johnstone, Tibshirani, 2004) but: the algorithms and estimates are not the same

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### Theorem for high dimensions (PB, 2004)

 $L_2$ Boosting with comp. linear LS regression is consistent (for suitable number of boosting iterations) if:

•  $p_n = O(\exp(Cn^{1-\xi})) \ (0 < \xi < 1)$ 

essentially exponentially many variables relative to n

•  $\sup_n \sum_{j=1}^{p_n} |\beta_{j,n}| < \infty \ \ell_1$ -sparseness of true function

i.e. for suitable, slowly growing  $m = m_n$ :

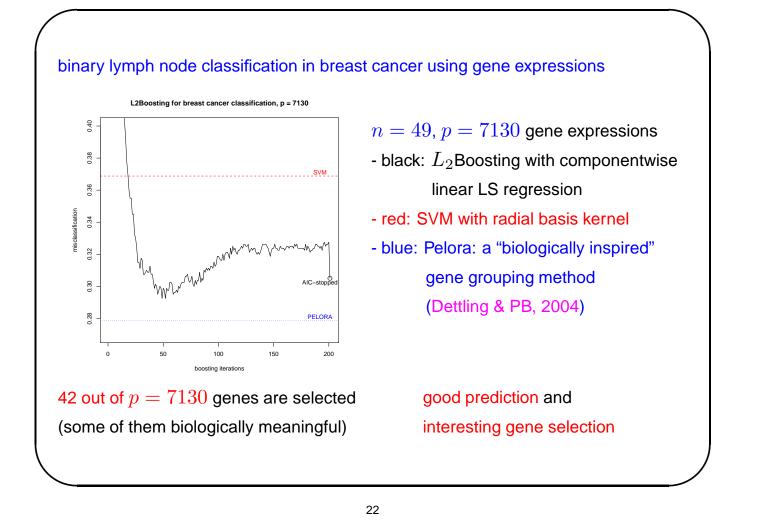
$$\mathbf{E}_X |\hat{f}_{m_n,n}(X) - f_n(X)|^2 = o_P(1) \ (n \to \infty)$$

"no" assumptions about the predictor variables/design matrix

in other words:

consistency for de-noising sparse signal with highly over-complete dictionaries

similar result has been given for the Lasso by Greenshtein and Ritov (2004)



**3.** 
$$L_2$$
**Boosting, Lasso and LARS**  
Efron et al. (2004): intriguing relation between  $L_2$ Boosting and  
Lasso:  $\hat{\beta}_{Lasso} = \operatorname{argmin}_{\beta} \sum_{i=1}^{n} (Y_i - \sum_{j=1}^{p} \beta_j X_i^{(j)})^2 + \lambda \sum_{j=1}^{p} |\beta_j|$   
for some special cases, roughly:  
iterations of " $L_2$ Boosting with "infinitesimally" small  $\nu$   
yield all Lasso solutions when varying  $\lambda$ "  
 $\rightsquigarrow$  computationally interesting to produce all Lasso solutions in  
one sweep of boosting  
Least Angle Regression LARS (Efron et al., 2004) is computationally even more  
clever and efficient than  $L_2$ Boosting

### for $p \gg n$

# both: Lasso/LARS and $L_2 \mbox{Boosting}$ are very useful

and LARS is really fast