

# Composite Models

Many methods can be seen as:

decision tree  
logistic function  
linear function  
... } of {  
decision trees  
logistic functions  
linear functions  
kernel functions  
lower dimensional subspace  
...

E.g., neural networks, regression trees, random forest, ...

Some combinations don't help.

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- Piecewise linear functions of input features  $\rightarrow$  neural networks (with ReLU)



# Boosting

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These weak learners are then boosted to be components in the ensemble that performs better than any of them.

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- At each stage, the base learner learns  $\hat{d}_i$  to minimize

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This is equivalent to learning from a modified dataset, where the previous prediction is subtracted from the actual value of the training set.
- Each learner is made to correct the errors of the previous prediction.



# Boosting\_learner

- 1: **procedure** *Boosting\_learner*( $Xs, Y, Es, L, K$ )
- 2:     **Inputs**
- 3:          $Xs$ : set of input features;  $Y$ : target feature;  $Es$ :  
training examples;  $L$ : base learner;  $K$ : number of components  
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- 12:     **return**  $p_k$

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- For regression, the loss is regularized squared error:

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The regularization is  $\Omega(f) = \gamma * |w| + \frac{1}{2} \lambda * \sum_j w_j^2$ , where  $w$  is vector of weights.  $\gamma$  and  $\lambda$  are nonnegative numbers.



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Optimize sum of log loss with the same regularization:

$$\left( \sum_e \text{logloss}(\hat{y}_e, y_e) \right) + \sum_{k=1}^K \Omega(f_k).$$

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- Gradient-boosted trees, the trees are built sequentially: each tree is learned assuming the previous trees are fixed.
- Two issues:
  - ▶ Selecting leaf values
  - ▶ Selecting splits
- For regression with squared error (or any loss based on the difference between the actual and predicted value), learn a tree for the difference between data and previous prediction.

## Selecting Leaf Values: Boolean Classification

- For the  $t$ th tree, optimize log loss with  $L2$  regularization:

$$\hat{y}_e^{(t)} = \text{sigmoid}\left(\sum_{k=1}^t f_k(x_e)\right)$$

$$\mathcal{L}^{(t)} = \sum_e \text{logloss}(\hat{y}_e^{(t)}, y_e) + \frac{1}{2}\lambda * \sum_j w_j^2 + \text{constant}$$

- Consider  $j$ th leaf, where  $I_j = \{e \mid q(x_e)=j\}$  is the set of training examples that map to it.
- Taking the derivative with respect to  $w_j$ :

$$\frac{\partial}{\partial w_j} \mathcal{L}^{(t)} = \lambda * w_j + \sum_{e \in I_j} (\hat{y}_e - y_e)$$

- A gradient descent step gives (Newton–Raphson method):

$$w_j = \frac{\sum_{e \in I_j} (y_e - \hat{y}_e^{(t-1)})}{\sum_{e \in I_j} \hat{y}_e^{(t-1)} * (1 - \hat{y}_e^{(t-1)}) + \lambda}$$