Applications of Data Analytics in Manufacturing

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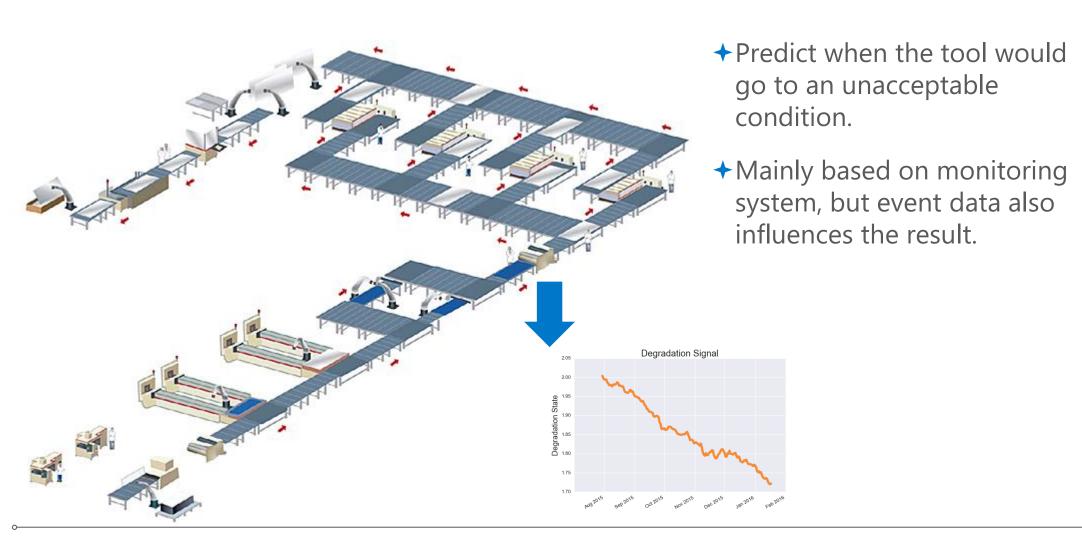
Rolls Royce Digital.

Overview of Production Line



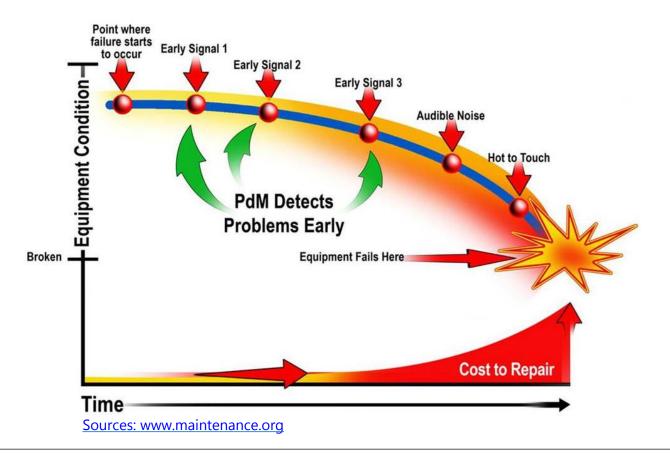
- →Event system: recipe changes, tool maintenance activities (PM), human actions.
- →Measurement system: deviation quality gates, yield output.
- ◆Investigation system: find quality deviation root causes

Event System Solution: Predictive Maintenance



Predictive Maintenance

→Predict when a tool would go bad.



Maintenance Model

Terminologies:

Degradation signal Y:

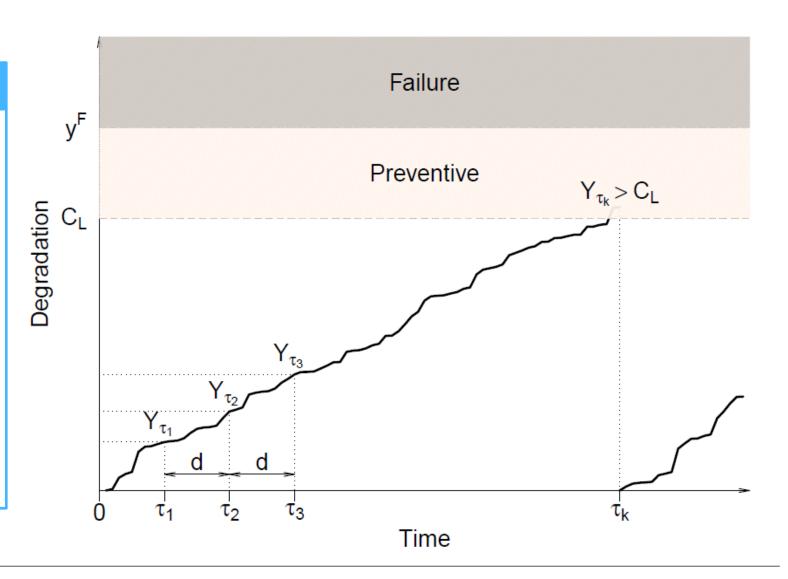
- Monotonically increasing or decreasing
- Recover after CM/PM

Failure Limit y^F:

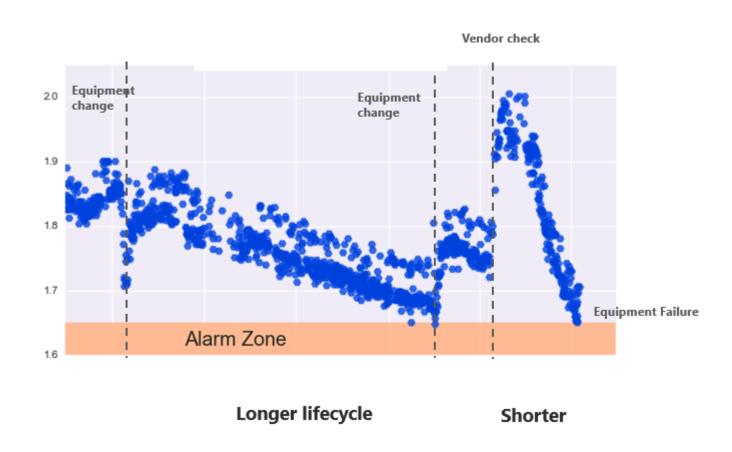
 Tool fails when degradation signal hits

Control Limit C_L:

 Perform PM when degradation signal hits



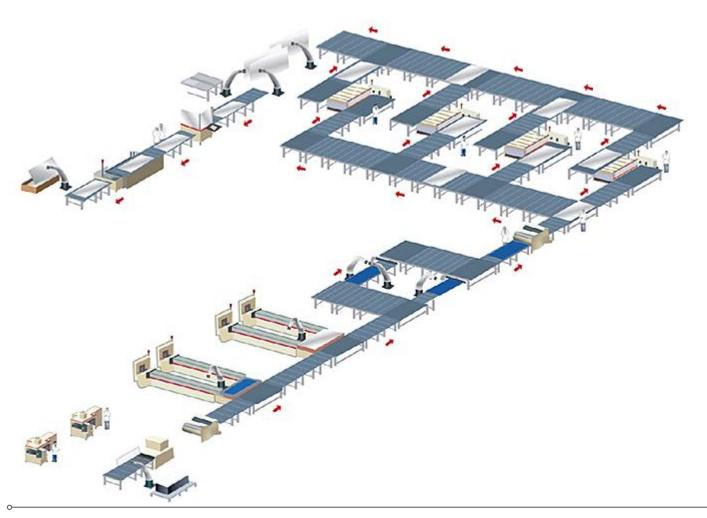
Real Scenario



Problem Formulation



Measurement Solution: Virtual Measurement



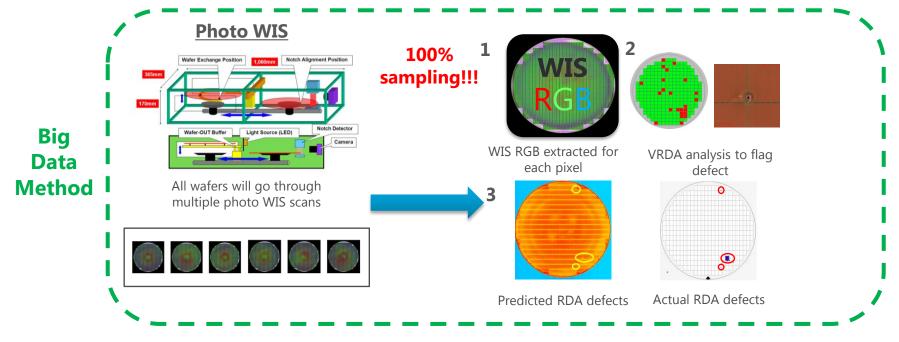
- → Measurement is an expensive operation, some destructive.
 Consequently, its sampling rate is low.
- ✦ However, some monitoring systems have 100% sampling rate, either through sensors or scanning images.

Virtual Measurement

→Motivation: Predict low-sampled measurement using 100% sampled images.

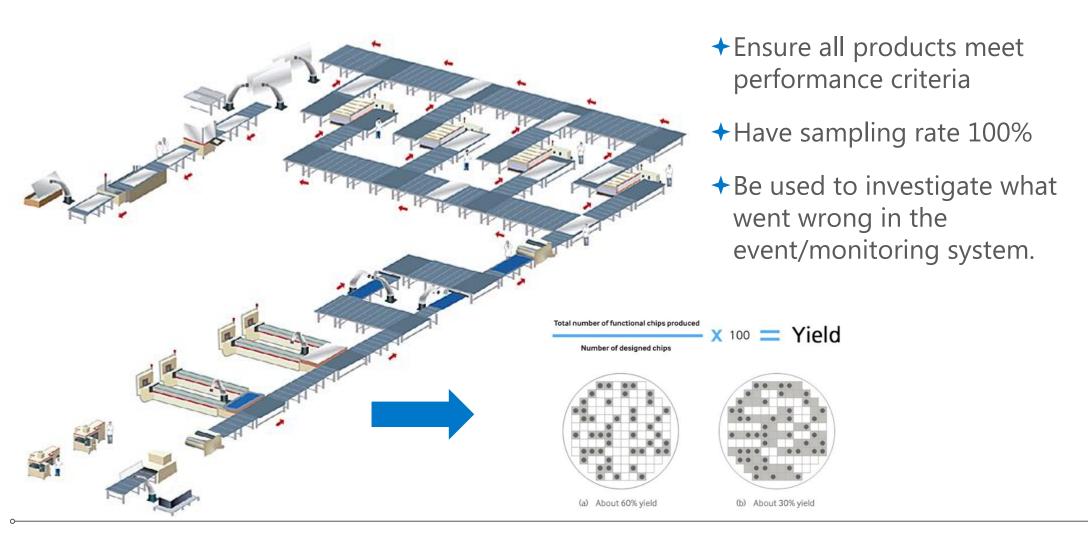
Virtual Measurement





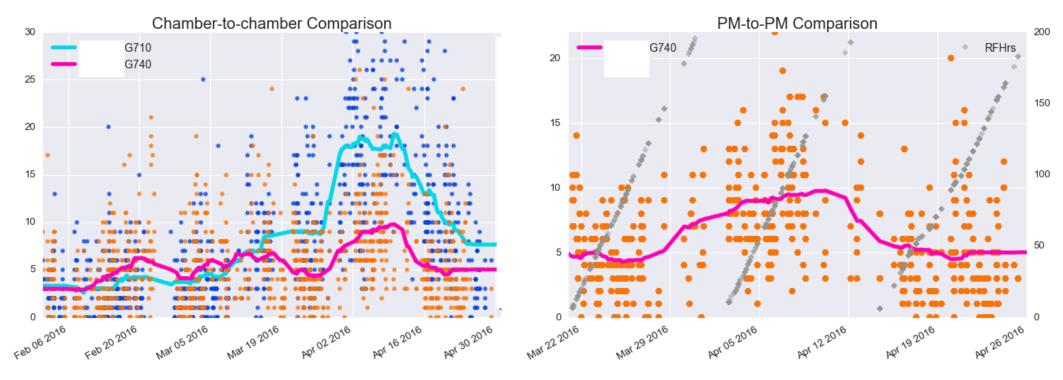
100% wafer sampling → Higher visibility to detect defective wafers

Investigation Solution: Yield Modelling



Yield Modelling

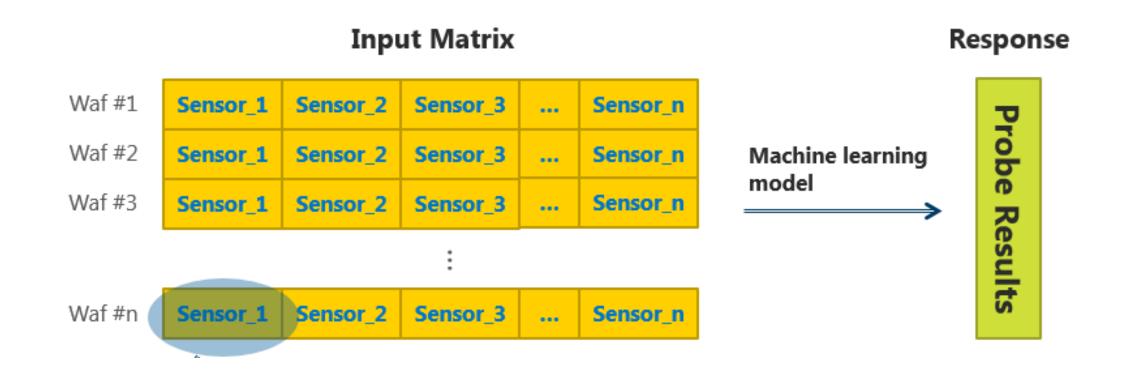
◆Investigate what went wrong with a bad tool



What are the root causes? How to improve yield?

Answer from tool sensor signals => Bad maintenance activity.

Yield Modelling



Modelling with Lasso and Random Forest

Purposes of Data Modelling

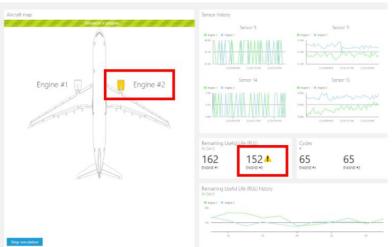
→Inference

- Explain the relationship among predictors and between predictors and responses.
- Tell data insights and trigger investigation

→Prediction

 Estimate responses given predictor values in a set of unobserved samples.





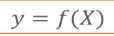
Data Structure

Sample training data ~20k rows, 100 unique engine id

Sample testing data ~13k rows, 100 unique engine id

id	cycle	setting1	setting2	setting3	s1	52	53		s19	s20	s21	RUL	label1	label2
	1 1	-0.0007	-0.0004	100	518.67	641.82	1589.7		100	39.06	23.419			
	1 2	0.0019	-0.0003	100	518.67	642.15	1591.82		100	39	23.4236			
	1 3	-0.0043	0.0003	100	518.67	642.35	1587.99		100	38.95	23.3442			
	1 191	. 0	-0.0004	100	518.67	643.34	1602.36		100	38.45	23.1295			
	1 192	0.0009	0	100	518.67	643.54	1601.41		100	38.48	22.9649			
	2 1	-0.0018	0.0006	100	518.67	641.89	1583.84		100	38.94	23.4585			
	2 2	0.0043	-0.0003	100	518.67	641.82	1587.05		100	39.06	23.4085			
	2 3	0.0018	0.0003	100	518.67	641.55	1588.32		100	39.11	23.425			
	2 286	-0.001	-0.0003	100	518.67	643.44	1603.63		100	38.33	23.0169			
	2 287	-0.0005	0.0006	100	518.67	643.85	1608.5		100	38.43	23.0848			
id	cycle	setting1	setting2	setting3	c1	s2	s 3		s19	s20	521			
		0.0023	_	-				_	100					
	_	2 -0.0027							100					
		0.0003						_	100					
													_	
	1 3	-0.0025	0.0004	100	518.6	7 642.7	9 1585.72	2	100	39.09	23.4069			
	1 3	-0.0006	0.0004	100	518.6	7 642.5	8 1581.2	2	100	38.81	23.3552			
	2	-0.0009	0.0004	100	518.6	7 642.6	6 1589.3	3	100	39	23.3923			
	2	-0.0011	0.0002	100	518.6	7 642.5	1 1588.43	3	100	38.84	23.2902			
	2	0.0002	0.0003	100	518.6	7 642.5	8 1595.0	5	100	39.02	23.4064			
	2 4	0.0011	-0.0001	100	518.6	7 642.6	4 1587.7	L	100	38.99	23.2918			
	2 4	0.0018	-0.0001	100	518.6	7 642.5	5 1586.59	9	100	38.81	23.2618			
	3	-0.0001	0.0001	100	518.6	7 642.0	3 1589.92	2	100	38.99	23.296			
	3	0.0039	-0.0003	100	518.6	7 642.2	3 1597.3	L	100	38.84	23.3191			
	3	0.0006	0.0003	100	518.6	7 642.9	8 1586.7	7	100	38.69	23.3774			
	3 12	0.0014	0.0002	100	518.6	7 643.2	4 1588.64	1	100	38.56	23.227			
	3 12	-0.0016	0.0004	100	518.6	7 642.8	8 1589.75	5	100	38.93	23.274			

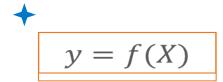
Data Structure



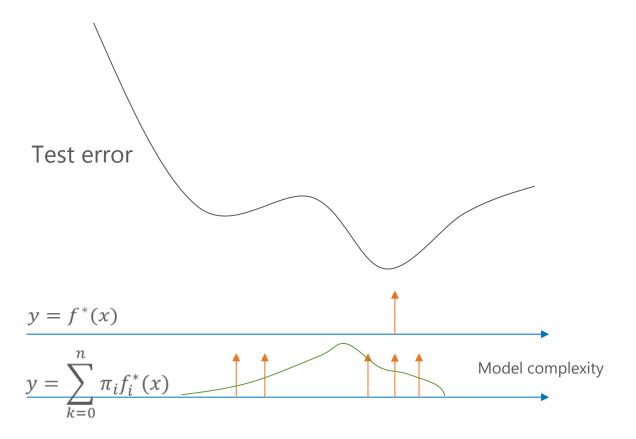


1. LASSO

Basic Model Assumptions



- Two approaches to estimate f
 - Underlying model style (*)
 - Bayesian style



Ref: Chapter 5. <u>Deep Learning</u>. Ian Goodfellow, Yoshua Bengio and Aaron Courville

Linear Regression

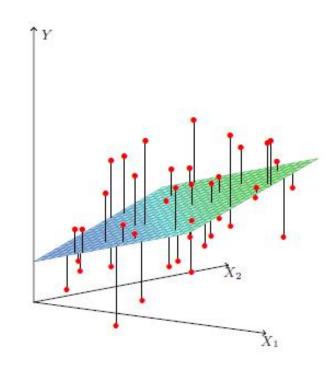
Assumption: $f(X) = B_0 + \sum_{j=1}^p X_j B_j$

To find B

$$J_B = (y - XB)^T (y - XB)$$

$$\frac{\partial J_B}{\partial B} = -2X^T(y - XB)$$

$$\hat{B} = (X^T X)^{-1} X^T y$$



Ridge Regression

$$\mathbf{H}_B = (y - XB)^T (y - XB) + \lambda B^T B$$

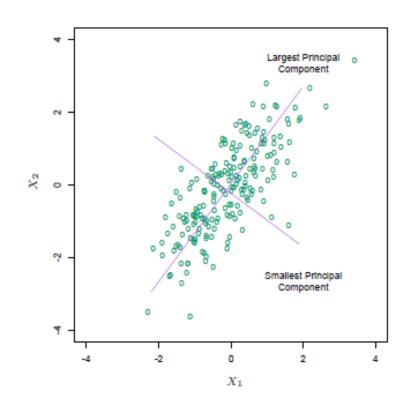
$$\hat{B} = (X^T X + \lambda I)^{-1} X^T y$$

With $X = UDV^T$

$$\widehat{B} = V(D^2 + \lambda I)^{-1}DU^T y$$

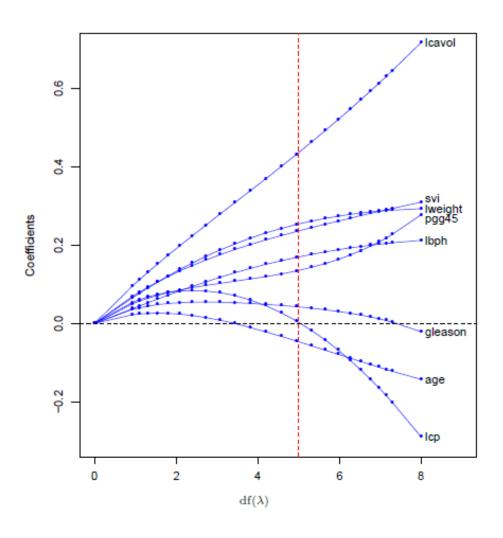
$$\hat{y} = X\hat{B} = \sum_{j=1}^{p} u_j \left(\frac{d_j^2}{d_j^2 + \lambda} \right) u_j^T y$$

 u_j columns of U, d_j diagonal elements of D



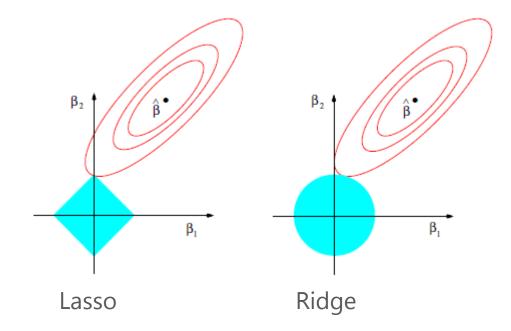
Ref: SVD Tutorial

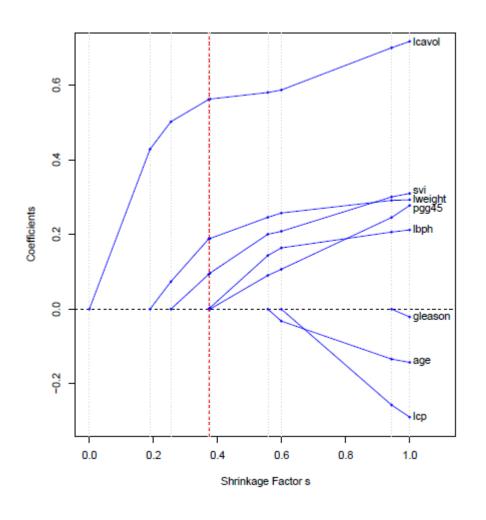
Ridge Regression



Lasso Regression

$${}^{\bullet}J_B = (y - XB)^T (y - XB) + \lambda \sum_{j=1}^p |B_j|$$

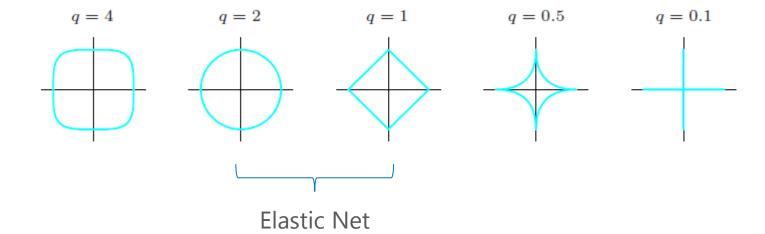




G-Formula

+

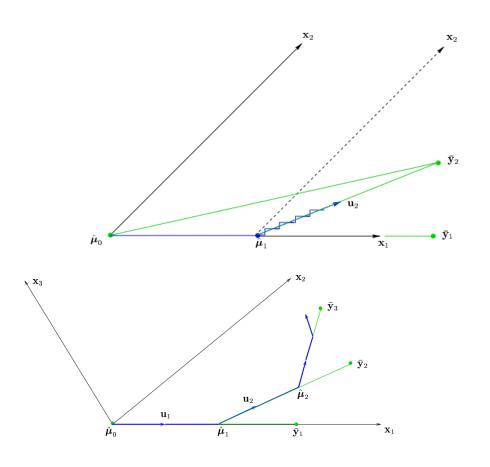
$$J_B = (y - XB)^T (y - XB) + \lambda \sum_{j=1}^p |B_j|^q$$



Lasso Solution – Least Angle Regression (LAR) *

Algorithm 3.2 Least Angle Regression.

- 1. Standardize the predictors to have mean zero and unit norm. Start with the residual $\mathbf{r} = \mathbf{y} \bar{\mathbf{y}}, \, \beta_1, \beta_2, \dots, \beta_p = 0$.
- 2. Find the predictor \mathbf{x}_i most correlated with \mathbf{r} .
- 3. Move β_j from 0 towards its least-squares coefficient $\langle \mathbf{x}_j, \mathbf{r} \rangle$, until some other competitor \mathbf{x}_k has as much correlation with the current residual as does \mathbf{x}_j .
- 4. Move β_j and β_k in the direction defined by their joint least squares coefficient of the current residual on $(\mathbf{x}_j, \mathbf{x}_k)$, until some other competitor \mathbf{x}_l has as much correlation with the current residual.
- 5. Continue in this way until all p predictors have been entered. After $\min(N-1,p)$ steps, we arrive at the full least-squares solution.

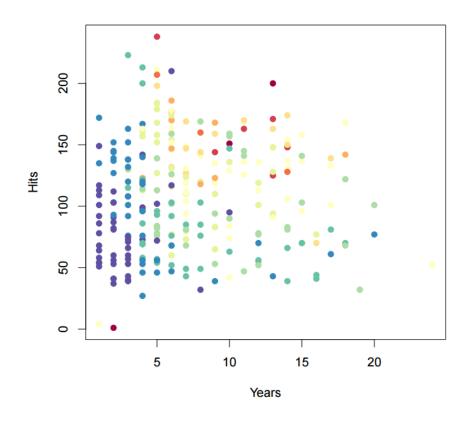


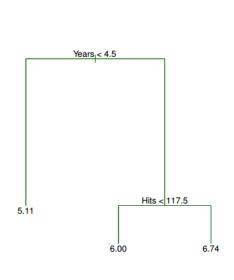
→Questions?

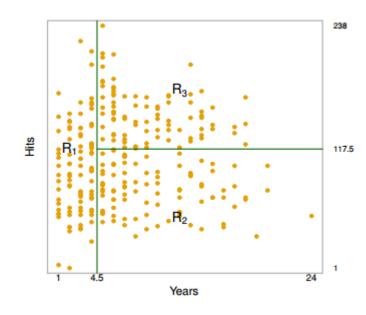
2. RANDOM FOREST

Decision Tree - Baseball Player Salary

Salary is color-coded from low (blue, green) to high (yellow,red)







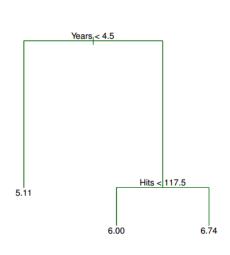
Decision Tree

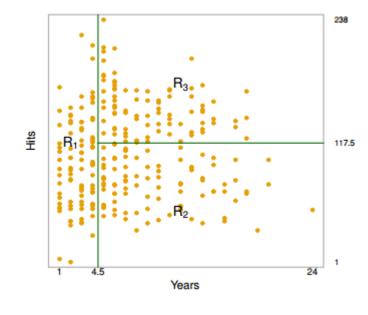
An approach that is known as recursive binary splitting

→Top-down,

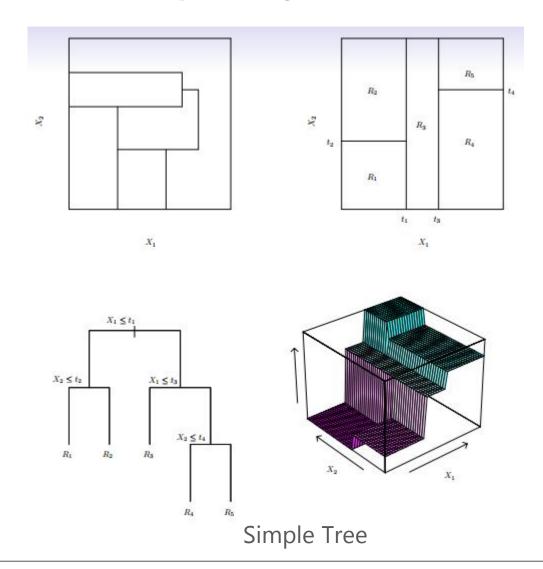
+Greedy

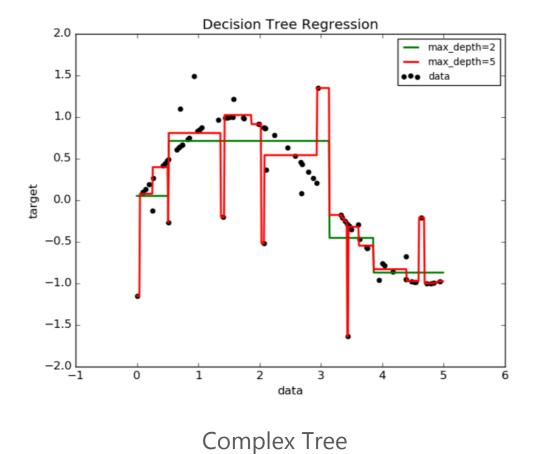
Predict the response for a given test observation <u>using the mean</u> of the training observations in the region





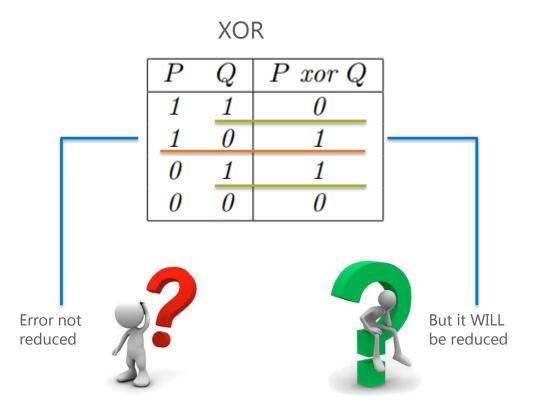
Tree Complexity and Overfitting





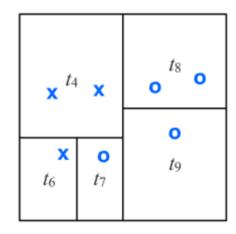
Tree Pruning

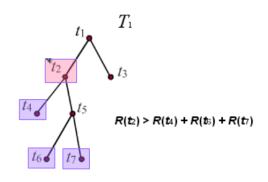
- →Motivation: Simple tree is too biased. Complex tree is overfitting.
- Naïve solution: grow the tree only so long as the decrease in the Residual-Sum-of-Square due to each split exceeds some (high) threshold.
 - Result in smaller trees, but is too short-sighted:
 a seemingly worthless split early on in the tree
 might be followed by a very good split



Tree Pruning

- →Motivation: Simple tree is too biased. Complex tree is overfitting.
- →Better solution: grow a large tree, then merge back nodes to obtain a smaller tree of the right size <u>link</u>





$$0.25 > 0 + 0 + 0$$

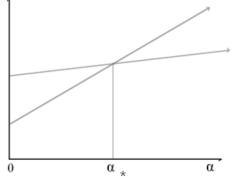
Cost complexity pruning (or Weakest link pruning) *

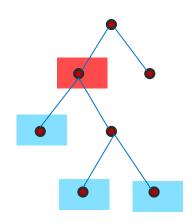
- Define T_t a branch rooted at a node t, and and \tilde{T}_t is its set of terminal nodes.
 - Let α be a regularization parameter.

$$R_{\alpha}(t) = R(t) + \alpha * 1$$

$$R_{\alpha}(T_t) = \sum_{t' \text{ in } \tilde{T}_t} R(t') + \alpha * |\tilde{T}_t|$$

$$\alpha_* = \frac{R(t) - R(T_t)}{\left|\tilde{T}_t\right| - 1}$$





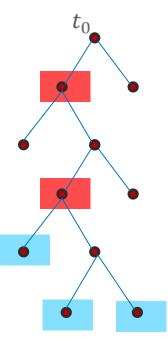
Cost complexity pruning (or Weakest link pruning) *

- \uparrow 1. Construct a large tree T_0 .
 - 2. Find a node t that minimizes the function $g(t) = \frac{R(t) R(T_t)}{|\tilde{T}_t| 1}$. Let $\alpha_1 = g(t)$ and remove all sub-nodes under t to produce T_1 .
 - 3. Repeat step 2 to find two sequences

$$\alpha_1 < \alpha_2 < \alpha_3 < \dots < \alpha_{|T|}$$

$$T_1 > T_2 > T_3 > \dots > t_0$$

Note: R(t) can be stored.



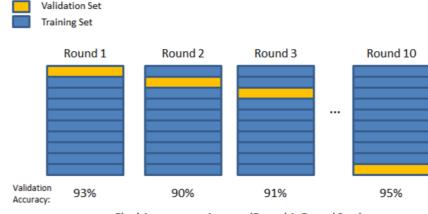
Final Decision Tree

Algorithm 8.1 Building a Regression Tree

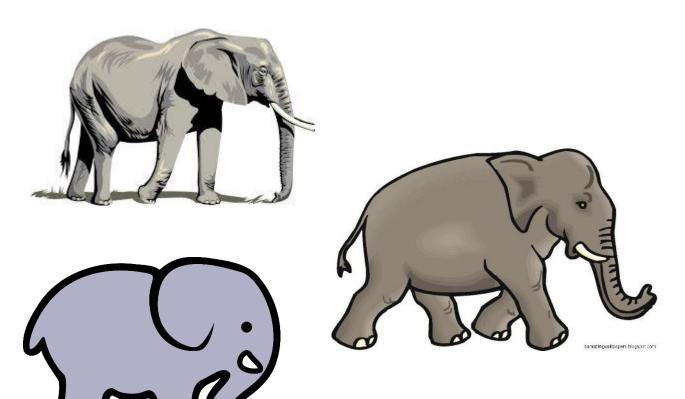
- Use recursive binary splitting to grow a large tree on the training data, stopping only when each terminal node has fewer than some minimum number of observations.
- Apply cost complexity pruning to the large tree in order to obtain a sequence of best subtrees, as a function of α.
- 3. Use K-fold cross-validation to choose α . That is, divide the training observations into K folds. For each $k = 1, \ldots, K$:
 - (a) Repeat Steps 1 and 2 on all but the kth fold of the training data.
 - (b) Evaluate the mean squared prediction error on the data in the left-out kth fold, as a function of α .

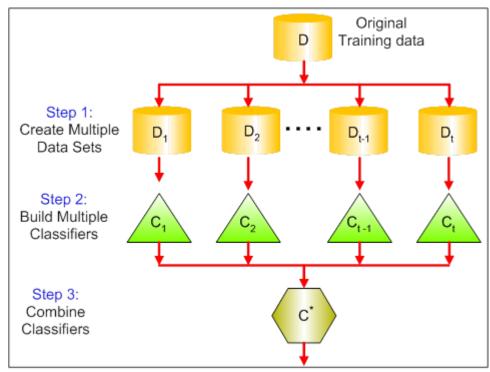
Average the results for each value of α , and pick α to minimize the average error.

4. Return the subtree from Step 2 that corresponds to the chosen value of α .



Bagging

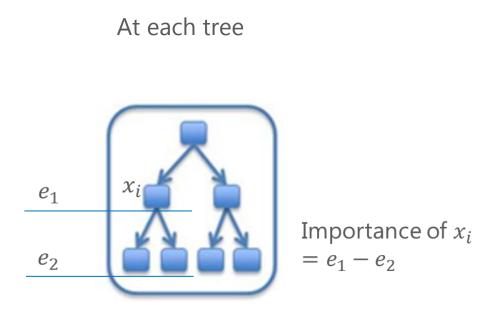


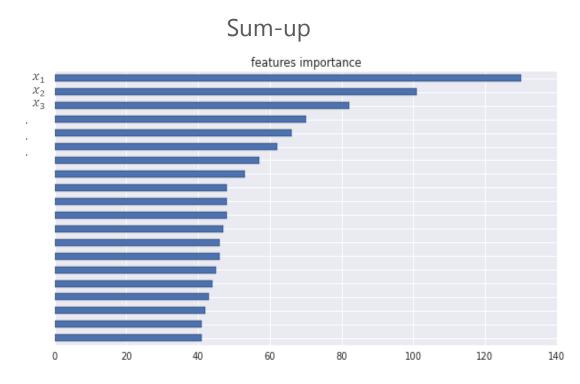


Bagging – Prediction

→Averaging predictions from multiple trees, each is constructed on one part of a training data set. Famous method: Bootstrap (sampling with replacement).

Bagging – Variable Importance

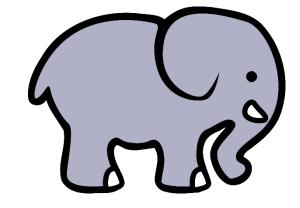




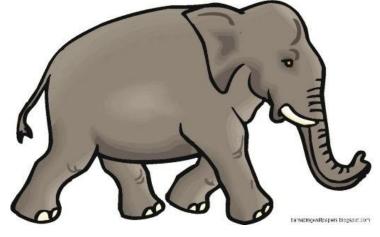
Random Forest







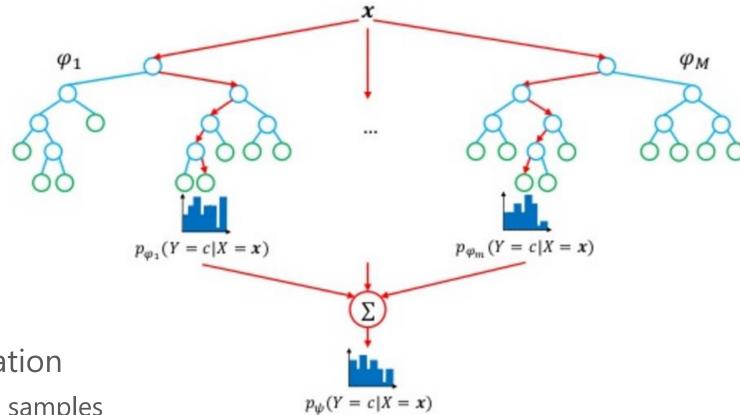








Random Forest



- **→**Randomization
 - Bootstrap samples
 - Feature subsets are different for these trees