• Boxplot:

$$-Q_1 = N \times 0.25$$

$$-Q_2 = N \times 0.5$$

$$-Q_3 = N \times 0.75$$

$$-IQR = Q_3 - Q_1$$

$$- Bounds = [Q_1 - 1.5 \times IQR, Q_3 + 1.5 \times IQR]$$

• Pearson =
$$\frac{\Sigma(y_{1i} - \bar{y}_1)(y_{2i} - \bar{y}_2)}{\sqrt{\Sigma(y_{1i} - \bar{y}_1)^2 \times (y_{2i} - \bar{y}_2)^2}}$$

• Spearman: Assign ranks and apply Pearson formula. Example: $[20, 10, 20, 30, 20] \rightarrow [3, 1, 3, 5, 3]$

• Normalization:

- MinMax:
$$\frac{y_i - min}{max - min}$$

- Standardization:
$$\frac{y_i - \mu}{\sigma}$$

• Binarization:

- Range (equal width): Depends on variable range Example: $y \in [-1, 1] : [0.2, -0.1, 0.6] \rightarrow [1, 0, 1]$
- Frequency (equal depth): Depends on variable mean Example: $\bar{y} = 25 : [10, 40, 30, 20] \rightarrow [0, 1, 1, 0]$

	True		
	A	В	С
_, A	TA	FA	FA
P F	FB	TB	FΒ
^L C	FC	FC	TC

TN

FPFN

• Metrics:

$$- \ {\rm Accuracy} = \frac{TP + TN}{total}$$

- Error rate =
$$1 - Accuracy = \frac{FP + FN}{total}$$

$$- \text{ Recall} = \frac{TP}{TP + FN} \text{ (Sensitivity)}$$

- Fallout =
$$\frac{TN}{TN + FP}$$
 (Specificity)

$$- \text{ Precision} = \frac{TP}{TP + FP}$$

$$- F_1 = \frac{TP}{TP + \frac{1}{2}(FP + FN)}$$

• Error:

- Sum of Squares Error: $SSE = \sum (Z \hat{Z})^2$
- Maen Squared Error: $MSE = \frac{1}{n}SSE$
- Root Maen Squared Error: $RMSE = \sqrt{MSE}$
- Mean Absolute Error: $MAE = \frac{1}{n} \sum |Z \hat{Z}|$
- Information Gain: $IG(y_{out}|y_i) = E(y_{out}) E(y_{out}|y_i)$
- Entropy: $E(y) = -\sum P(x_i) \log(P(x_i))$
- Decision trees:
 - 1. Choose feature with highest IG.
 - 2. Split dataset by that feature, create leaves if necessary.
 - 3. Repeat until unable to proceed.

Prune: (Given a twig)

1. Count it's leaves labels.

Example: $\#A = 5, \ \#B = 6$

- 2. Remove it's leaves.
- 3. Relabel twig as a leaf.

Example: B(6/11), #B > #A

• Vector Norm:

$$||x||_p = \sqrt[p]{\sum_{i=1}^n |x_i|^p} \qquad ||x||_\infty = \max |x_i|$$

• Matrix Multiplication:

$$\begin{bmatrix} \dots & \dots & n \\ \dots & \dots & \dots \\ m & \dots & \dots \end{bmatrix} \cdot \begin{bmatrix} \dots & \dots & l \\ \dots & \dots & \dots \\ n & \dots & \dots \end{bmatrix} = \begin{bmatrix} \dots & \dots & l \\ \dots & \dots & \dots \\ m & \dots & \dots \end{bmatrix}$$

• Gaussian Distribution

- Variance:
$$var = \frac{\sum (y_i - \mu)^2}{n(-1)}$$

- Standard Deviation: $\sigma = \sqrt{va}$

$$-P(y|\mu,\sigma) = \frac{1}{\sqrt{2\pi} \cdot \sigma} \cdot \exp\left(-\frac{1}{2} \left(\frac{y-\mu}{\sigma}\right)^2\right)$$

• Gaussian Mixture:

- Covariance:
$$cov(y_1, y_2) = \frac{\sum (y_{1i} - \mu_1)(y_{2i} - \mu_2)}{n(-1)}$$

– Covariance Matrix:
$$\Sigma(y_1, y_2) = \begin{bmatrix} var(y_1) & cov(y_1, y_2) \\ cov(y_1, y_2) & var(y_2) \end{bmatrix}$$

$$- |\Sigma| = \begin{vmatrix} a & b \\ c & d \end{vmatrix} = ad - bc$$

$$- P(y|\mu, \Sigma) = \frac{1}{(2\pi)^{d/2} |\Sigma|^{1/2}} \cdot exp\left(-\frac{1}{2}(y-\mu)^T \Sigma^{-1}(y-\mu)\right)$$

• Naive Bayes:

– MAP:
$$P(C|x) = \frac{P(C)P(x|C)}{P(x)}$$

- ML:
$$P(C|x) = P(x|C)$$

- 1. Calculate probability for each class.
- 2. Calculate P(x|C) for each attribute and each class.
- 3. TODO
- K-Nearest Neighbors:
 - 1. Distances: (for n variables)

- Manhattan:
$$\sum |y_{1i} - y_{2i}|$$

- Euclidean:
$$\sqrt{\sum (y_{1i} - y_{2i})^2}$$

- Cosine:
$$\frac{\sum \overline{y_{1i}} \ y_{2i}}{\sqrt{\sum y_{1i}^2} \sqrt{\sum y_{2i}^2}}$$

- Hamming: #Differences
- 2. Choose K nearest neighbors.
- 3. Classify using mean if variable is numeric, or mode if it is categoric.
- 4. If weighted, divide by weight.

• Regressions:

- Linear:
$$W = (X^T X)^{-1} X^T Z$$

– Ridge:
$$W = (X^T X + \lambda I)^{-1} X^T Z$$

• Perceptron:

$$\hat{Z} = a(W^T X), \ a \leftarrow \text{activation function}$$

If $Z \neq \hat{Z} \longrightarrow W' = W + \eta(Z - \hat{Z})X$

• Neural Networks (MLP):

- Forward:
$$x^{[0]} \to z^{[1]} = w^{[1]}x^{[0]} + b^{[1]} \to x^{[1]} = a(z^{[1]}) \to \dots \to z^{[i]} = w^{[i]}x^{[i-1]} + b^{[i]} \to x^{[i]} = a(z^{[i]}) \to E$$

- Backward:

$$* \ \delta^{[last]} = \frac{\partial E}{\partial x^{[last]}} \circ \frac{\partial x^{[last]}}{\partial z^{[last]}}$$

$$* \ \delta^{[i]} = \left(w^{[i+1]}\right)^T \cdot \delta^{[i+1]} \circ \frac{\partial x^{[i]}}{\partial z^{[i]}}$$

$$*~b^{[i]'} = b^{[i]} - \eta rac{\partial E}{\partial b^{[i]}} \qquad *~rac{\partial E}{\partial b^{[i]}} = \delta^{[i]}$$

- Derivatives:

Derivatives.				
Name	Error function		$\frac{\partial E}{\partial x^{[i]}}$	
Squared Error	$\frac{1}{2} \left(x^{[i]} - t \right)^2$		$x^{[i]}-t$	
Cross- -entropy	$-\sum_{i=1}^{n} t_i \log \left(x_i^{[i]}\right)$		$-\frac{t}{x^{[i]}} + \frac{1-t}{1-x^{[i]}}$	
Name	Activation function	$\frac{\partial x^{[i]}}{\partial z^{[i]}}$		
Sigmoid $\sigma(x)$	$\frac{1}{1+e^{-x}} \qquad \sigma ($		$z^{[i]}$) $\left(1 - \sigma\left(z^{[i]}\right)\right)$	
Hyper. $tan.$ $tanh(x)$	$\frac{e^x - e^{-x}}{e^x + e^{-x}}$	$1 - \tanh^2\left(z^{[i]}\right)$		
$\begin{array}{c} \operatorname{ReLU} \\ R(x) \end{array}$	$ \begin{array}{c} 0 \text{ if } x < 0 \\ x \text{ if } x \ge 0 \end{array} $	$ \begin{array}{c c} 0 \text{ if } x < 0 \\ 1 \text{ if } x \ge 0 \end{array} $		
Softmax $S(x)$	$\frac{e^{x_i}}{\sum_{j=1}^n e^{x_j}}$	$S\left(z^{[i]}\right)\left(1-S\left(z^{[i]}\right)\right)$		
Name	Function		Derivative	
	$e^{f(x)}$		$e^x \times f'(x)$	

NOTE:

* When cross-entropy and softmax are combined:

$$\delta^{[last]} = \frac{\partial E}{\partial z^{[last]}} = x^{[last]} - t$$

• K-Means:

- 1. Assign each point to a cluster.
- 2. Update centroids: centroid $_{new}$ = mean of cluster points
- 3. Repeat until centroids don't change.

• EM:

- Initializaion: Initial mixture parameters
- Expectation (E-step):

Calculate weights for each datapoint
$$\mathbf{x}_i$$
 for each cluster \mathbf{c}_k :
$$\gamma_{ki} = \frac{\mathcal{N}(x_i | \mu_k, \Sigma_k) \cdot \pi_k}{\sum_{j=1}^k \mathcal{N}(x_i | \mu_j, \Sigma_j) \cdot \pi_j}$$

- Maximization (M-step):

Update parameters for each cluster: (for n observations)

$$* N_k = \sum_{i=1}^n \gamma_{ki}$$

$$* \mu_k = \frac{1}{N_k} \sum_{i=1}^n \gamma_{ki} \cdot x_i$$

*
$$\Sigma_k = \frac{1}{N_k} \sum_{i=1}^n \gamma_{ki} \cdot (x_i - \mu_k) \cdot (x_i - \mu_k)^T$$

$$* \ \pi_k = \frac{N_k}{N}$$

- Sillhouette: \in [-1,1] (the closer to 1 the better)
 - For an observation x_i :
 - * $a = \text{average distance of } x_i \text{ to the points in it's cluster}$
 - * $b = \text{average distance of } \mathbf{x}_i \text{ to points in closest cluster}$ * $s_{observation} = \frac{b-a}{max(a,b)}$
 - For a cluster:

Average of the cluster's observations silhouettes

- For the solution:

Average of the clusters silhouettes