

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

- 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: $\bar{y} = 25 : [10, 40, 30, 20] \rightarrow [0, 1, 1, 0]$

- Confusion Matrix:

		True			B
		A	B	C	
Pred	A	TA	FA	FA	TP
	B	FB	TB	FB	TN
	C	FC	FC	TC	FP
					FN

- Metrics:

- Accuracy = $\frac{TP + TN}{total}$
- Error rate = $1 - Accuracy = \frac{FP + FN}{total}$
- Recall = $\frac{TP}{TP + FN}$ (Sensitivity)
- Fallout = $\frac{TN}{TN + FP}$ (Specificity)
- Precision = $\frac{TP}{TP + FP}$
- $F_1 = \frac{TP}{TP + \frac{1}{2}(FP + FN)}$

- Error:

Sum of Squares Error: $\sum(z - \hat{z})^2$

Root Mean Squared Error: $\sqrt{\frac{1}{n} SSE}$

Mean Absolute (Percentage) Error: $\frac{1}{n} \sum \left| \frac{z - \hat{z}}{z} \right|$

- Information Gain: $IG(class|y_i) = E(class) - E(class|y_i)$

- Entropy: $E(y) = \sum_{i=1}^k \frac{|y_i|}{|y|} \sum_{j=1}^n -P(y_{ij}) \log(P(y_{ij}))$

- 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:

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

- Matrix Multiplication:

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

- Gaussian Distribution:

- Variance: $var = \frac{\sum(y_i - \mu)^2}{n(-1)}$
- Standard Deviation: $\sigma = \sqrt{var}$
- $P(y|\mu, \sigma) = \frac{1}{\sqrt{2\pi} \cdot \sigma} \cdot \exp\left(-\frac{1}{2} \left(\frac{y - \mu}{\sigma}\right)^2\right)$

- Gaussian Mixture:

- Covariance: $cov(y_1, y_2) = \frac{\sum(y_{1i} - \mu_1)(y_{2i} - \mu_2)}{n(-1)}$
- Covariance Matrix: $\Sigma(y_1, y_2) = \begin{bmatrix} var(y_1) & cov(y_1, y_2) \\ cov(y_1, y_2) & var(y_2) \end{bmatrix}$
- $det|\Sigma| = det \begin{vmatrix} a & b \\ c & d \end{vmatrix} = ad - bc$
- $P(y|\mu, \Sigma) = \frac{1}{(2\pi)^{d/2} |\Sigma|^{1/2}} \cdot exp\left(-\frac{1}{2}(y - \mu)^T \Sigma^{-1}(y - \mu)\right)$

- Naive Bayes:

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

1. Calculate $P(C)$ for each class.
2. Calculate $P(y|C)$ for each variable for each class.
3. Calculate likelihood: $P(x|C) = P(y_1|C) \times \dots \times P(y_d|C)$

- K-Nearest Neighbors:

1. Distances: (for n variables)
 - Manhattan: $\sum |y_{1i} - y_{2i}|$
 - Euclidean: $\sqrt{\sum (y_{1i} - y_{2i})^2}$
 - Cosine: $\frac{\sum y_{1i} y_{2i}}{\sqrt{\sum y_{1i}^2} \sqrt{\sum y_{2i}^2}}$
 - Hamming: #Differences
2. If weighted, for each variable multiply by weight.
3. Choose K nearest neighbors.
4. Classify using mean if variable is numeric, or mode if it is categorical.

- Regressions: $\hat{z} = Xw = (w^T X^T)^T$

- Linear: $y = w_0 + w_1 x$
- Polynomial: $w = (X^T X)^{-1} X^T z$
- Ridge: $w = (X^T X + \lambda I)^{-1} X^T z$

- Perceptron:

$\hat{z} = a(w^T x), a \leftarrow$ activation function

Gradient Descent:

$$w^{new} = w^{old} + \Delta w \quad \text{where} \quad \Delta w = -\eta \frac{\partial E}{\partial w}$$

- Neural Networks (MLP):

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

- Backward:

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

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

$$* w^{[i]'} = w^{[i]} - \eta \cdot \frac{\partial E}{\partial w^{[i]}}$$

$$* b^{[i]'} = b^{[i]} - \eta \cdot \frac{\partial E}{\partial b^{[i]}}$$

– Multiple observations:

1. Apply forward propagation for each observation.
2. Calculate $\delta_i^{[l]}$ for each observation.
3. $\frac{\partial E}{\partial w^{[l]}} = \sum_{i=1}^n \delta_i^{[l]} \cdot \left(x_i^{[l-1]}\right)^T$
4. $\frac{\partial E}{\partial b^{[l]}} = \sum_{i=1}^n \delta_i^{[l]}$

– 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 $\sigma(x)$	$\frac{1}{1+e^{-x}}$	$x^{[i]} (1-x^{[i]})$
ArcTan $\arctan(x)$	$\arctan(x)$ or $\tan^{-1}(x)$	$\frac{1}{\left(x^{[i]}\right)^2 + 1}$
Hyperbolic tangent	$\frac{e^x - e^{-x}}{e^x + e^{-x}}$	$1 - \left(x^{[i]}\right)^2$
ReLU	0 if $x < 0$ x if $x \geq 0$	0 if $x^{[i]} < 0$ 1 if $x^{[i]} \geq 0$
Softmax	$\frac{e^{x_i}}{\sum_{j=1}^n e^{x_j}}$	$x^{[i]} (1-x^{[i]})$
Sign	1 if $x \geq 0$ 0 if $x < 0$	

NOTE:

- * When cross-entropy and softmax are combined:

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

• More Derivatives:

Function	Derivative	Function	Derivative
$x \pm y$	$x' \pm y'$	$f(x)^a$	$a f(x)^{a-1} f'(x)$
xy	$x'y + y'x$	$a^{f(x)}$	$a^{f(x)} f'(x) \ln a$
$\frac{x}{y}$	$\frac{xy' + y'x}{y^2}$	$\log_a f(x)$	$\frac{f'(x)}{f(x) \ln a}$

• K-Means:

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

• Silhouette: $\in [-1,1]$ (the closer to 1 the better)

– For an observation x_i :

- * a = average distance of x_i to the points in it's cluster
- * b = average distance of x_i 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

• EM:

– Initializaion: Initial mixture parameters

– Expectation (E-step):

Calculate weights for each datapoint x_i for each cluster 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)

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

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

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

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

• Quadratic Formula: $ax^2 + bx + c = 0 \Rightarrow x = \frac{-b \pm \sqrt{b^2 - 4ac}}{2a}$

• PCA: $\Sigma u = (\lambda I)u$

– Σ : Covariance Matrix

– I : Identity Matrix

– Eigenvalue (λ): $\det |\Sigma - \lambda I| = 0$ or $\lambda = \Sigma u \circ \frac{1}{u}$

Example:

$$\begin{aligned} \lambda &= \left(\begin{bmatrix} 2.917 & 2.667 \\ 2.667 & 2.667 \end{bmatrix} \begin{bmatrix} -0.690 \\ 0.723 \end{bmatrix} \right) \circ \begin{bmatrix} -0.690^{-1} \\ 0.723^{-1} \end{bmatrix} \\ &= \begin{bmatrix} -0.084 \\ 0.088 \end{bmatrix} \circ \begin{bmatrix} -0.690^{-1} \\ 0.723^{-1} \end{bmatrix} \\ &= \begin{bmatrix} 1.122 \\ 1.122 \end{bmatrix} \end{aligned}$$

– Eigenvector (u): $(\Sigma - \lambda I)u = \vec{0}$

Example:

$$\begin{aligned} u &\Leftrightarrow \left(\begin{bmatrix} 1.333 & 0.667 \\ 0.667 & 1.667 \end{bmatrix} - \begin{bmatrix} 2.187 & 0 \\ 0 & 2.187 \end{bmatrix} \right) \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \\ &\Leftrightarrow \left[\begin{array}{cc|c} -0.854 & 0.667 & 0 \\ 0.667 & -0.52 & 0 \end{array} \right] \Leftrightarrow \begin{matrix} L_1 \times -0.854^{-1} \\ L_2 \times 0.667^{-1} \end{matrix} \\ &\Leftrightarrow \left[\begin{array}{cc|c} 1 & -0.781 & 0 \\ 1 & -0.781 & 0 \end{array} \right] \Leftrightarrow L_2 - L_1 \\ &\Leftrightarrow \left[\begin{array}{cc|c} 1 & -0.781 & 0 \\ 0 & 0 & 0 \end{array} \right] \Leftrightarrow x_1 = 0.781x_2 \\ &= \begin{bmatrix} 0.781x \\ x \end{bmatrix} \end{aligned}$$

– Projecting (bivariate to univariate):

1. Choose highest λ (higher λ means more variation)
2. Calculate eigenvector (u)
3. Apply formula: $\phi = u^T X^T$

• Model complexity:

– Perceptron with d inputs: $d + 1$

– Tree with d features que tomam n valores: n^d

– MLP: estimated by the number of weights

– Bayesian Classifier: estimated by the number of parameters

- * With c classes: $c - 1$ priors

- * With d dimension: $(d \text{ averages} + \frac{d(d+1)}{2} \Sigma) \times c$