

# Assumptions of Machine Learning Classifiers

## 1. K-Nearest Neighbors (KNN)

KNN is a non-parametric, instance-based learning method. It does not assume a specific functional form, but it has several implicit assumptions:

### Assumptions

#### 1. Distance Meaningfulness

KNN assumes that data points that are close in feature space are similar.  
→ Therefore, features must be on similar scales (standardization needed).

#### 2. Local Smoothness

The target value changes smoothly in the neighborhood.  
Points near each other have similar labels.

#### 3. No Irrelevant Features

Too many irrelevant features reduce KNN performance ("curse of dimensionality").

#### 4. Balanced Classes

KNN assumes class distribution in neighborhoods is not heavily imbalanced.

## 2. Decision Tree Classifier

Decision Trees are flexible and non-linear, but still have some assumptions.

### Assumptions

#### 1. No Need for Feature Scaling

Tree splits are based on thresholding, so scaling does not matter.

#### 2. Features Are Sufficient to Separate Classes

The tree assumes the data can be separated using recursive binary splits.

### 3. Large Sample Size

Small datasets may cause highly unstable splits (high variance).

### 4. No Strong Multicollinearity

Highly correlated features may confuse split selection, though trees can still handle them better than linear models.

### 5. Data Purity Can Be Achieved

Tree assumes that splitting will eventually lead to homogeneous nodes.

## 3. Ensemble Learning

### A. Random Forest (Bagging)

Random Forests combine many decision trees trained on bootstrapped samples.

#### Assumptions

##### 1. Independence of Trees

Each tree should be different due to different samples and feature subsets.

##### 2. Low Bias Base Learners

Decision Trees must be strong enough to capture patterns.

##### 3. No Feature Scaling Required

Trees do not depend on distance.

##### 4. Sufficient Data

More data helps prevent random noise from dominating.

### B. Boosting (AdaBoost, Gradient Boosting, XGBoost)

Boosting builds models sequentially, where each new model corrects previous mistakes.

#### Assumptions

##### 1. Weak Learners Can Improve Sequentially

Boosting assumes each base learner performs slightly better than random.

##### 2. No Noise-Dominated Features

Boosting is sensitive to noise and outliers.

### 3. Features Don't Need Scaling

Trees are used; no distance metrics involved.

### 4. Additive Model Works

Boosting assumes errors can be reduced by combining many small models.

## 4. Support Vector Machines (SVM)

SVM finds the maximum-margin hyperplane to separate classes.

### Assumptions

#### 1. Data Is Separable (linearly or with kernel)

If linearly inseparable, kernel trick is used.

#### 2. Large Margin Exists

SVM assumes the best decision boundary is the one with maximum margin.

#### 3. Feature Scaling Is Required

SVM uses distance-based calculations; features must be standardized.

#### 4. Low Noise

SVM is sensitive to noise and overlapping classes.

#### 5. Balanced Classes

Imbalanced datasets can shift the margin unfavorably.

## 5. Linear Regression

### Assumptions

#### 1. Linearity

The relationship between features and target is linear.

#### 2. Independence of Errors

Residuals must not be correlated.

#### 3. Homoscedasticity (constant error variance)

Variance of residuals remains constant across predictions.

**4. Normality of Residuals**

Needed for statistical significance tests.

**5. No Perfect Multicollinearity**

Features must not be perfectly correlated.

**6. No Extreme Outliers**

Outliers affect slope drastically.