

LECTURE 17

Classification

Building models of classification in sklearn

Data Science, Spring 2024 @ Knowledge Stream

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Outline

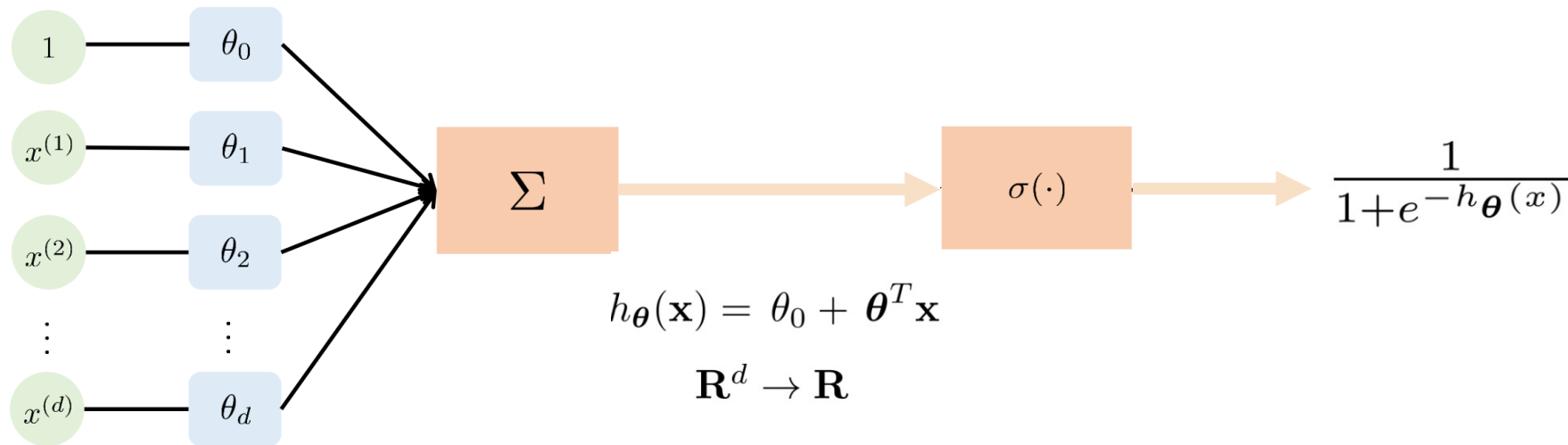
Lecture 17

- Introduction to Classification
- Types of Classification
- Classification Algorithms
- Performance Metrics
- Overfitting and Underfitting
- Conclusion

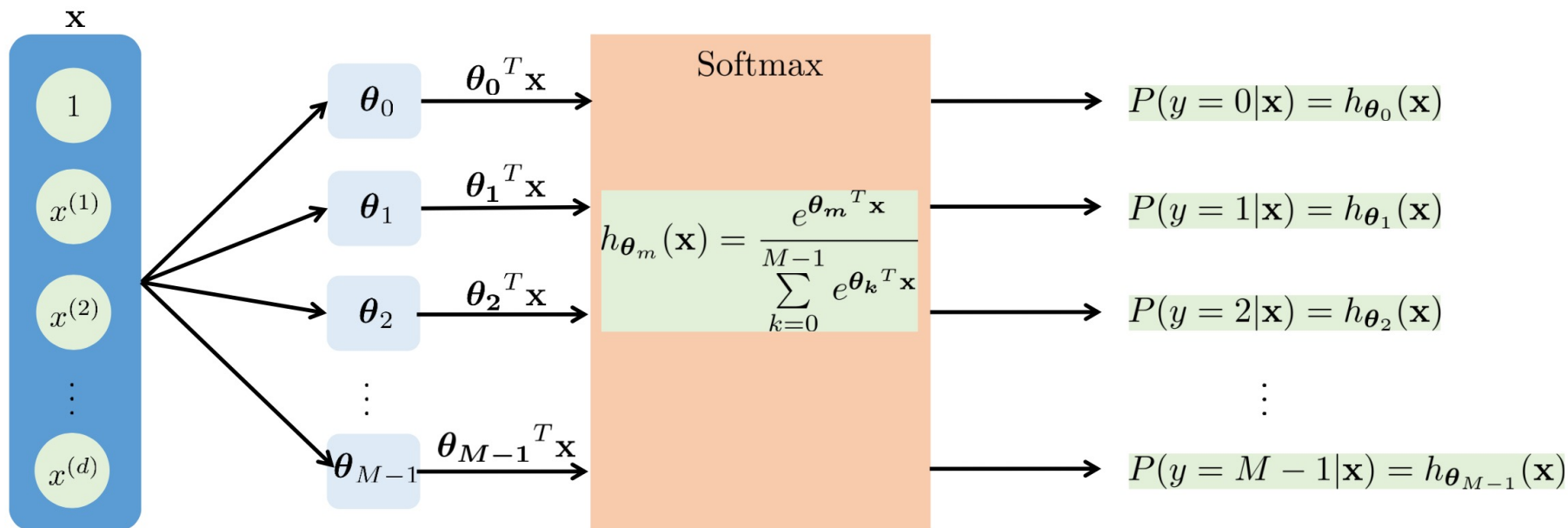
Classification

- Classification is defined as the process of recognition and grouping of objects
- Classification refers to a predictive modeling problem where a class label is predicted for a given example of input data
- For Classification, the training dataset must be sufficiently representative of the problem and have many examples of each class label.
 1. Logistic Regression
 2. Decision Trees
 3. Naive Bayes
 4. K-Nearest Neighbours
 5. Support Vector Machine
 6. Random Forest

Logistic Regression



Logistic Regression



Logistic Regression

```
X = df['data']
```

```
Y = df['target']
```

```
from sklearn.model_selection import train_test_split
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, random_state=16)
```

```
from sklearn.linear_model import LogisticRegression
```

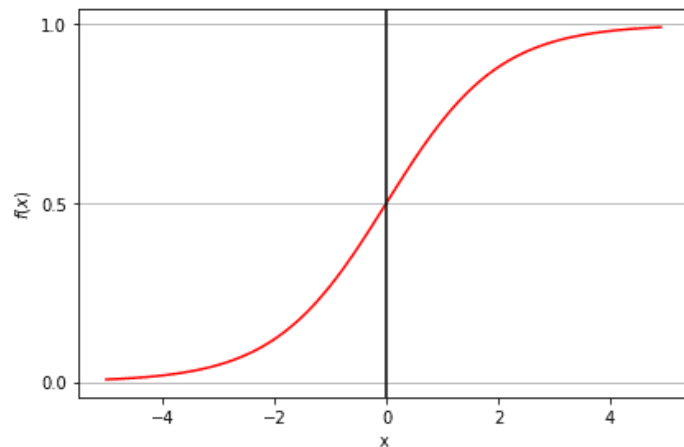
```
logreg = LogisticRegression(random_state=16)
```

```
logreg.fit(X_train, y_train)
```

```
y_pred = logreg.predict(X_test)
```

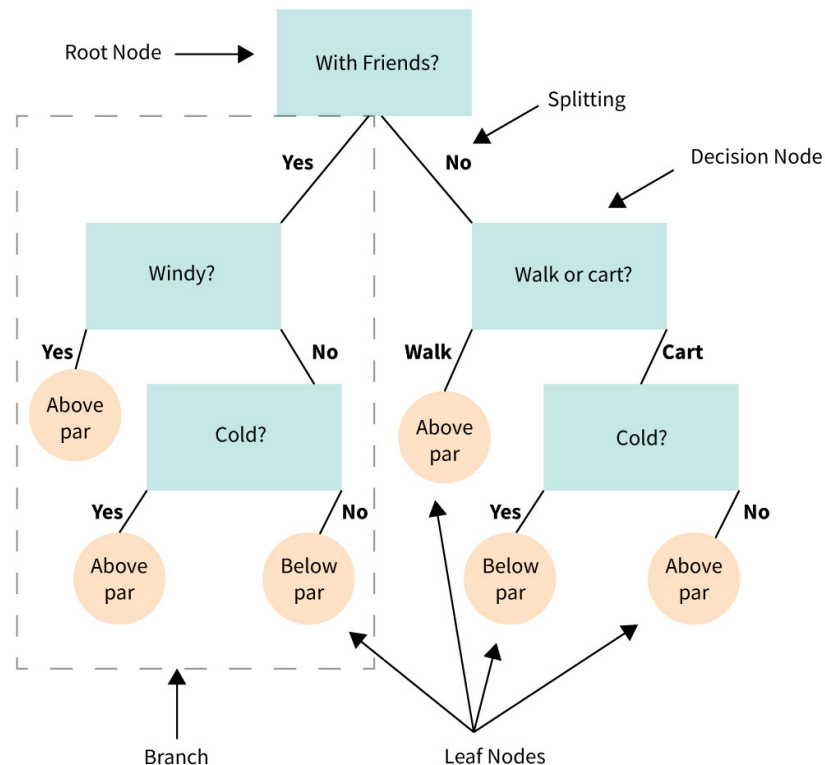
```
from sklearn import metrics
```

```
cnf_matrix = metrics.confusion_matrix(y_test, y_pred)
```



Decision Trees (DTs)

- A hierarchical, tree structure, consists of a root node, branches, internal nodes and leaf nodes
- Algorithms for binary classification
 1. Logistic Regression
 2. **Decision Trees**
 3. Naive Bayes
 4. k-Nearest Neighbours
 5. Support Vector Machine
 6. Random Forest



Decision Tree

```
X = df['data']
```

```
Y = df['target']
```

```
from sklearn.model_selection import train_test_split
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, random_state=16)
```

```
from sklearn.tree import DecisionTreeClassifier
```

```
dt = DecisionTreeClassifier ()
```

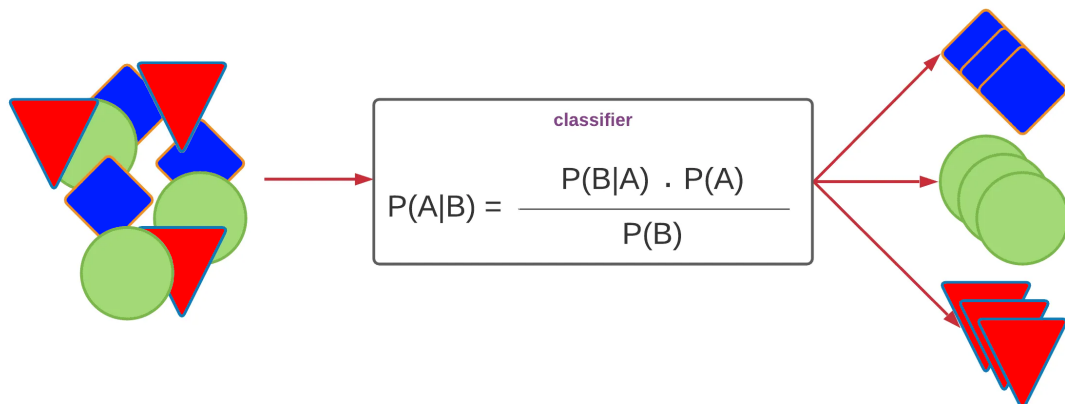
```
dt.fit(X_train, y_train)
```

```
y_pred = dt.predict(X_test)
```

```
from sklearn import metrics
```

```
cnf_matrix = metrics.confusion_matrix(y_test, y_pred)
```


- The Naïve Bayes classifier is used for classification tasks, like text classification.
- Algorithms for binary classification
 1. Logistic Regression
 2. Decision Trees
 3. Naive Bayes
 4. k-Nearest Neighbours
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- The Naive Bayes classifier is a family of probabilistic machine learning models based on [Bayes' theorem](#).
- The "naive" assumption is of [feature independence](#).
- It's commonly used for [text classification](#) tasks, particularly for document categorization and spam email filtering.
- Estimate $P(y|X)$ from the data using the Bayes Theorem

$$P(y \mid \mathbf{x}) = \underset{y \in \mathcal{Y}}{\text{maximize}} \quad \frac{P(\mathbf{x} \mid y) P(y)}{P(\mathbf{x})}$$

Naive Bayes

```
X = df['data']
```

```
Y = df['target']
```

```
from sklearn.model_selection import train_test_split
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, random_state=16)
```

```
from sklearn.naive_bayes import GaussianNB
```

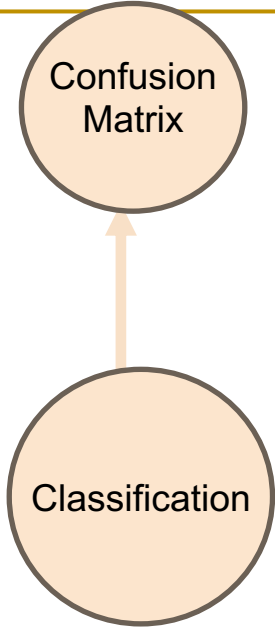
```
nb = GaussianNB ()
```

```
nb.fit(X_train, y_train)
```

```
y_pred =nb.predict(X_test)
```

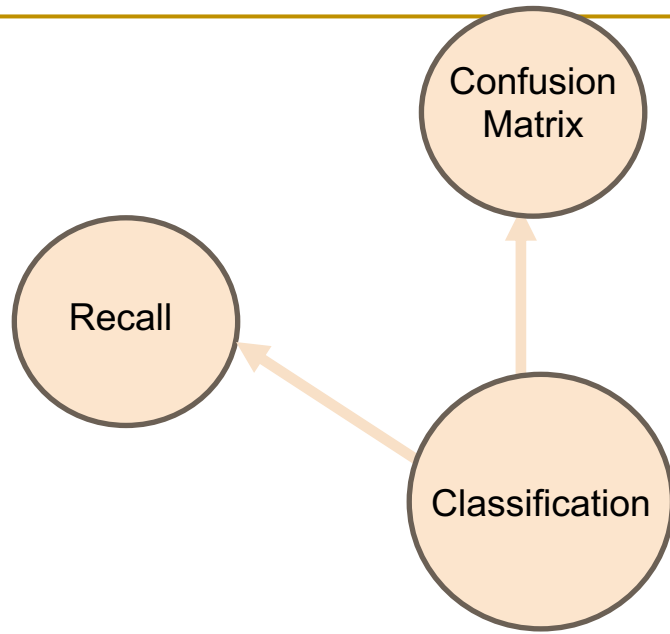
```
from sklearn import metrics
```

```
cnf_matrix = metrics.confusion_matrix(y_test, y_pred)
```



| | Predicted 0 | Predicted 1 |
|-------------|----------------|----------------|
| Actual 0 | TN | FP |
| Actual 1 | FN | TP |

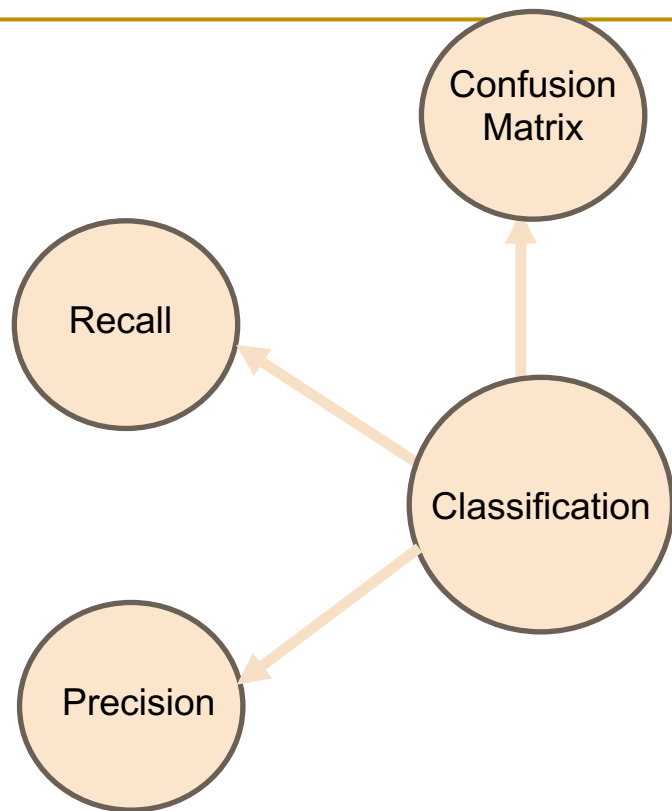
Performance Matrices



| | Predicted 0 | Predicted 1 |
|--------------------|-----------------------|-----------------------|
| Actual 0 | TN | FP |
| Actual 1 | FN | TP |

$$\text{Recall} = \frac{TP}{TP + FN}$$

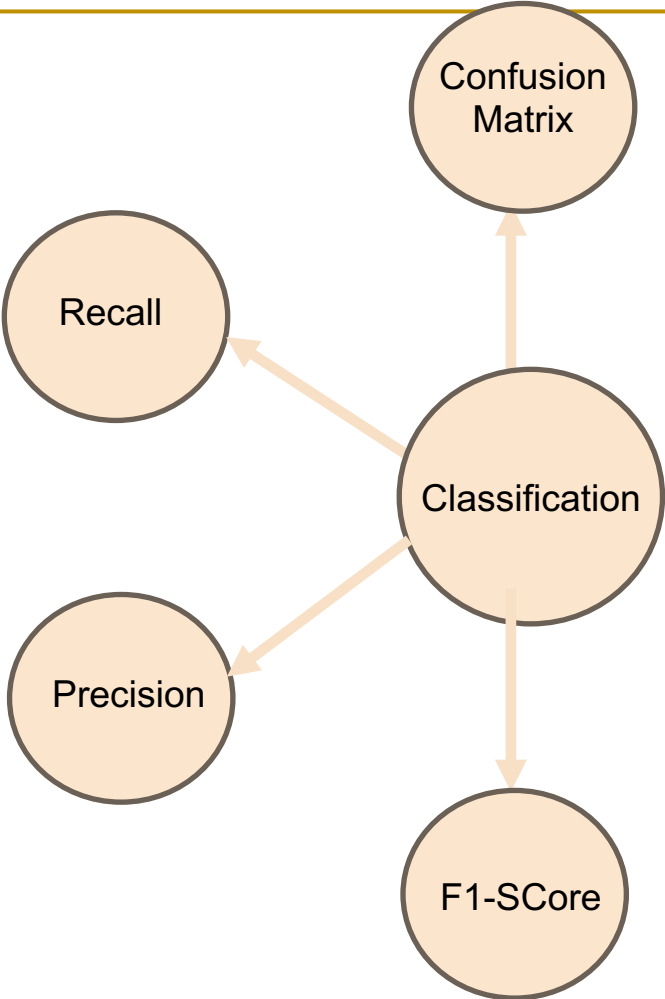
Performance Matrices



| | Predicted 0 | Predicted 1 |
|--------------------|-----------------------|-----------------------|
| Actual 0 | TN | FP |
| Actual 1 | FN | TP |

$$\text{Recall} = \frac{TP}{TP + FN}$$

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



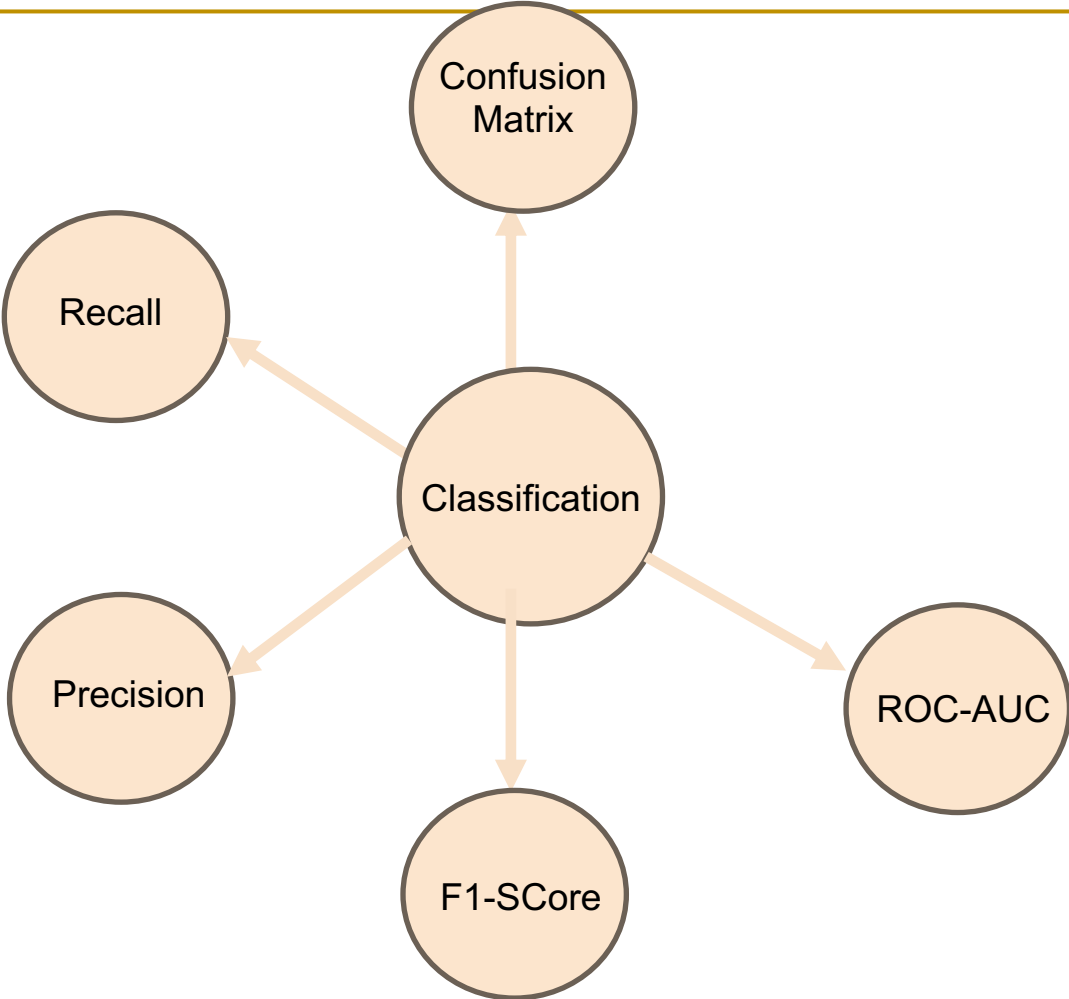
| | Predicted 0 | Predicted 1 |
|--------------------|-----------------------|-----------------------|
| Actual 0 | TN | FP |
| Actual 1 | FN | TP |

$$\text{Recall} = \frac{TP}{TP + FN}$$

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

$$\text{F1 Score} = \frac{TP}{TP + \frac{1}{2}(FP + FN)}$$

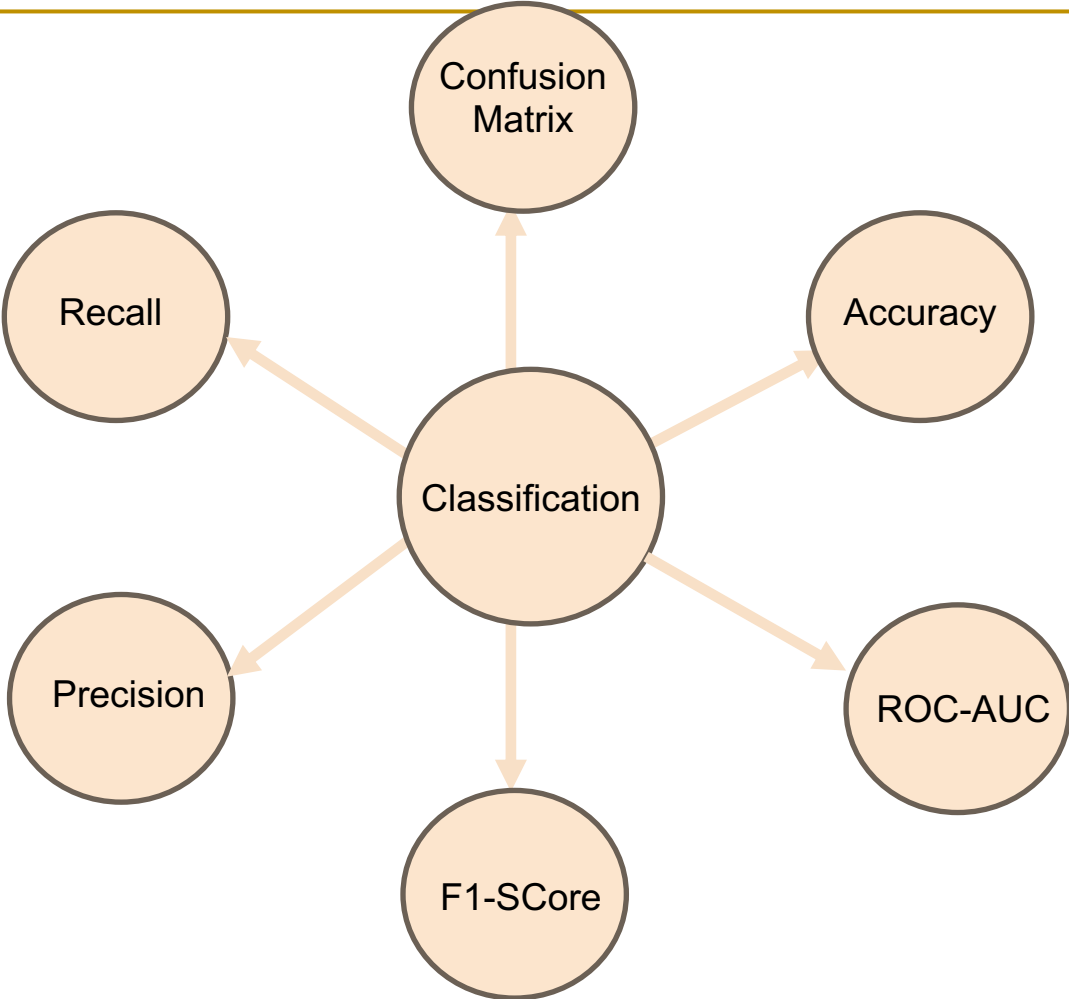
Performance Matrices



| | Predicted 0 | Predicted 1 |
|--------------------|-----------------------|-----------------------|
| Actual 0 | TN | FP |
| Actual 1 | FN | TP |

ROC curve is a graphical representation of the performance of the binary classifier

$$TPR = \frac{TP}{FN + TP}$$
$$FPR = \frac{FP}{TN + FP}$$



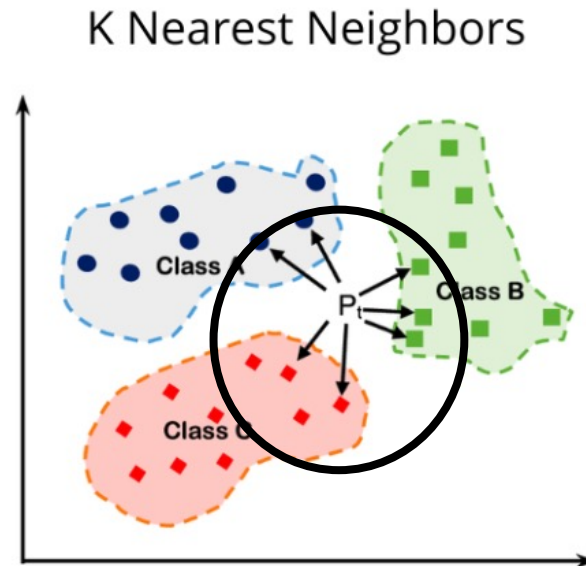
| | Predicted 0 | Predicted 1 |
|--------------------|-----------------------|-----------------------|
| Actual 0 | TN | FP |
| Actual 1 | FN | TP |

$$\text{Accuracy} = \frac{(TP + TN)}{(TP + FP + TN + FN)}$$

- `from sklearn.metrics import roc_curve`
- `from sklearn.metrics import roc_auc_score`
- `fpr, tpr, thresholds = roc_curve(y_true, y_scores)`
- `auc = roc_auc_score(y_true, y_scores)`

K-Nearest Neighbor

- K nearest neighbours is a simple algorithm that stores all available cases and classifies new cases based on a similarity measure.
- Algorithms for classification
 1. Logistic Regression
 2. Decision Trees
 3. Naive Bayes
 4. **k-Nearest Neighbours**
 5. Support Vector Machine
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- The k-nearest neighbours algorithm **stores** all the available data
- **Classifies** a new data point based on the **similarity measure** (e.g., distance functions).
- The data point is classified by a **majority vote** of its neighbours, with the data point being assigned to the class most common amongst its **K nearest neighbours** measured by a distance function.

- Loading the training and the test data.
- Choose the nearest data points (the value of K). K can be any integer.
- Do the following, for each test data point
 - Use Euclidean distance $\sqrt{\sum_{i=1}^k (x_i - y_i)^2}$ or Manhattan distance $\sum_{i=1}^k |x_i - y_i|$
to calculate the distance between test data and each row of training.
 - Sort the data set in ascending order based on the distance value.
 - From the sorted array, choose the top K rows.
 - Based on the most appearing class of these rows, it will assign a class to the test point.
 - End

Companies Using KNN

- Companies like [Amazon or Netflix](#) use [KNN](#) when recommending books to buy or movies to watch.
- How do these companies make recommendations?

Well, these companies gather data on the books you have read or movies you have watched on their website and apply KNN.

The companies will input your available customer data and compare that to other customers who have purchased similar books or have watched similar movies.

KNN Algorithm

```
X = df['data']
```

```
Y = df['target']
```

```
from sklearn.model_selection import train_test_split
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, random_state=16)
```

```
from sklearn.neighbors import KNeighborsClassifier
```

```
knn = KNeighborsClassifier(n_neighbors = k)
```

```
knn.fit(X_train, y_train)
```

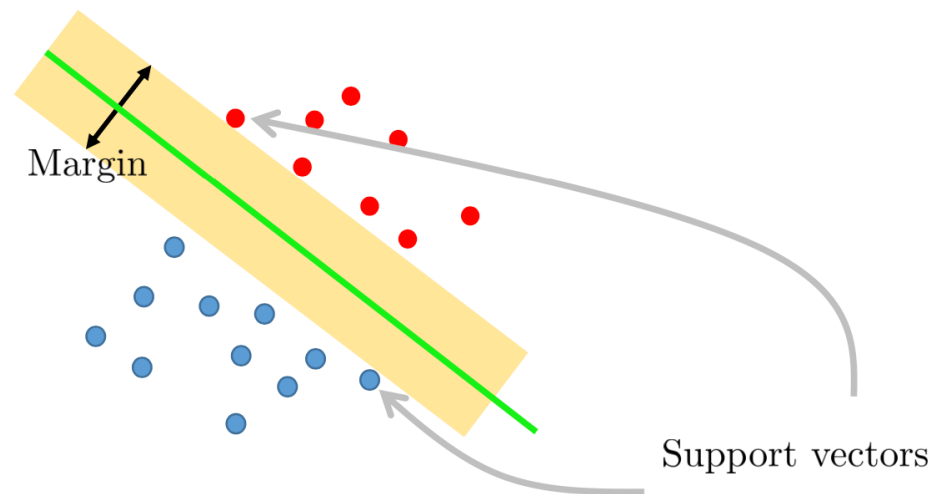
```
y_pred = knn.predict(X_test)
```

```
from sklearn import metrics
```

```
cnf_matrix = metrics.confusion_matrix(y_test, y_pred)
```

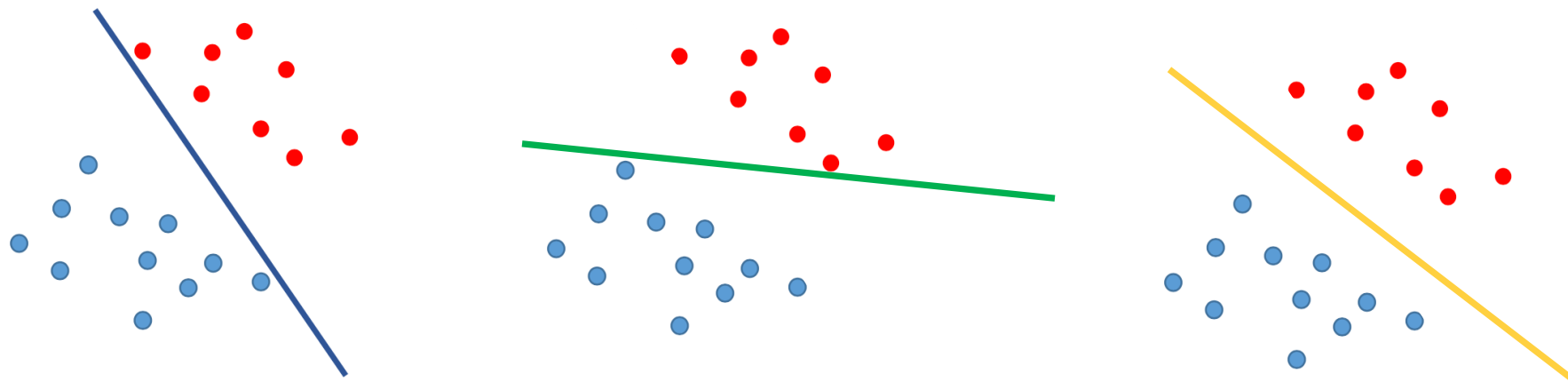
Support Vector Machine (SVM)

- **SVM** is used to classify data by finding the optimal decision boundary that maximally separates different classes
- Algorithms for binary classification
 1. Logistic Regression
 2. Decision Trees
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 4. k-Nearest Neighbours
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Support Vector Machine (SVM)

- We have linear separable classes.
- We can have multiple hyperplanes separating these classes.

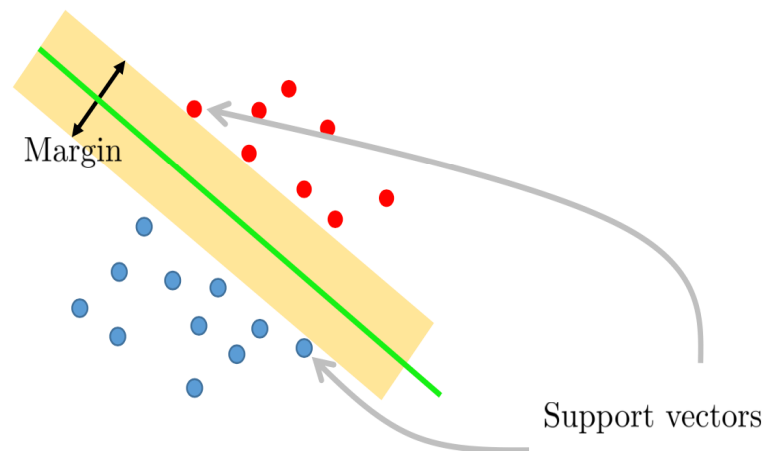


Q: Which one is the best decision boundary?

A: Maximum Margin Classifier (e.g., Support Vector Machine)

Support Vector Machine (SVM)

- Idea: Choose a fat separator
- The best boundary is the one that maximizes the **margin** or distance between the boundary and the “**difficult points**” close to the decision boundary.
- Margin against the data points are called **support vectors**.
- **Hard margin idea:** Find a maximum margin classifier with no errors on the training data.
- **Soft margin idea:** Find the maximum margin classifier while minimizing the number of training errors.



- ```
from sklearn.svm import SVC
svm_classifier = SVC(kernel = 'linear', C=1.0)
svm_classifier.fit(X_train, y_train)
y_pred = svm_classifier.predict(X_test)
```

You can choose different kernels (e.g., 'linear', 'poly', 'rbf') by specifying the kernel parameter. The C parameter controls the trade-off between maximizing the margin and minimizing the classification error.