Machine Learning with Python

Vocabulary

Machine Learning: Field of study that gives computers the ability to learn without being explicitly programmed [1959, Arthur Samuel]

Major machine learning techniques

• Regression/Estimation:

Predicting continuous values

• Classification:

Predicting the item class/category of a case

• Clustering:

Finding the structure of data; summarization

• Associations:

Associating frequent co-occurring items/events.

• Anomaly detection:

Discovering abnormal and unusual cases

• Sequence mining:

Predicting next events; click-stream (Markov Model, HMM)

• Dimension Reduction:

Reducing the size of data (PCA)

• Recommendation systems:

Python for Machine Learning

NumPy - SciPy - Scikit-learn - Pandas - Matplotlib.

Comparaison

Supervised	Unsupervised
Classification: Classifies labeled data	Clustering: Finds patterns and groupings from unlabeled data
Regression: Predicts trends using previous labeled data	Has fewer evaluation methods than supervised learning
Has more evaluation methods than unsupervised learning	Less controlled environment
Controlled environment	

Regression algorithms

Ordinal regression

Poisson regression

Fast forest quantile regression

Linear, Polynomial, Lasso, Stepwise, Ridge regression

Bayesian linear regression

Neural network regression

Decision forest regression

Boosted decision tree regression

KNN (K-nearest neighbors)

Simple Linear Regression

$$\hat{y} = \theta_0 + \theta_1 \mathbf{x}_1$$

$$MSE = \frac{1}{n} \sum_{i=1}^{n} (y_i - \hat{y}_i)^2$$

Estimating the parameters

$$\boldsymbol{\theta}_{1} = \frac{\sum_{i=1}^{s} (x_{i} - \bar{x}) (y_{i} - \bar{y})}{\sum_{i=1}^{s} (x_{i} - \bar{x})^{2}}$$

$$\boldsymbol{\theta}_0 = \overline{\boldsymbol{y}} - \boldsymbol{\theta}_1 \overline{\boldsymbol{x}}$$

Model Evaluation

- Train and Test on the Same Dataset
- \bullet Train/Test Split

$$Error = \frac{1}{n} \sum_{j=1}^{n} |y_j - \hat{y}_j|$$

Training accuracy is the percentage of correct predictions that the model makes when using the test dataset

High training accuracy isn't necessarily a good thing Result of over-fitting

• Over-fit: the model is overly trained to the dataset, which may capture noise and produce a non-generalized model

Out-of-sample accuracy is the percentage of correct predictions that the model makes on data that the model has not been trained on.

K-fold cross-validation

Metrics in Regression Models

$$MAE = \frac{1}{n} \sum_{j=1}^{n} |y_j - \hat{y}_j|$$

$$MSE = \frac{1}{n} \sum_{i=1}^{n} (y_i - \hat{y}_i)^2$$

$$RMSE = \sqrt{\frac{1}{n} \sum_{j=1}^{n} (y_j - \hat{y}_j)^2}$$

$$RAE = \frac{\sum_{j=1}^{n} |y_j - \hat{y}_j|}{\sum_{j=1}^{n} |y_j - \hat{y}_j|^2}$$

$$RSE = \frac{\sum_{j=1}^{n} (y_j - \hat{y}_j)^2}{\sum_{j=1}^{n} (y_j - \hat{y}_j)^2}$$

$$R^2 = 1 - RSE$$

Multiple Linear Regression

$$y = \theta_0 + \theta_1 x_1 + \theta_2 x_2 + \ldots + \theta_n x_n$$
$$\hat{y} = \theta^T X$$

$$heta^T = [heta_0, heta_1, heta_2, \ldots] \quad X = \left[egin{array}{c} 1 \\ x_1 \\ x_2 \\ \ldots \end{array}
ight]$$

Classification algorithms

Decision Trees (ID3, C4.5, C5.0)

Naive Baves

Linear Discriminant Analysis

k-Nearest Neighbor

Logistic Regression

Neural Networks

Support Vector Machines (SVM)

K-Nearest Neighbours

Based on: similar cases with same class labels are near each other.

K-nearest neighbor algorithm

- 1. Pick a value for K.
- 2. Calculate the distance of unknown case from all cases.
- 3. Select the K-observations in the training data that are "nearest" to the unknown data point.
- 4. Predict the response of the unknown data point using the most popular response value from the Knearest neighbors.

Evaluation Metrics in Classification

Jaccard index:

y: Actual labels, \hat{y} : Predicted labels

$$J(y,\hat{y}) = \frac{|y \cap \hat{y}|}{|y \cup \hat{y}|} = \frac{|y \cap \hat{y}|}{|y| + |\hat{y}| - |y \cap \hat{y}|}$$

confusion matrix:

TP: True Positive

FP: False Positive

$$\begin{array}{ll} {\rm Precision} &= {\rm TP}/({\rm TP} + {\rm FP}) \\ {\rm Recall} &= {\rm TP}/({\rm TP} + {\rm FN}) \end{array}$$

$$F1$$
-score = 2(prc * rec)/(prc+rec)

F1-score high accuracy $\rightarrow 1$

Log Loss:

Logarithmic loss measures the performance of a classifier where the predicted output is a probability value between 0 and 1.

LogLoss =
$$-\frac{1}{n}\sum (y \times \log(\hat{y}) + (1 - y) \times \log(1 - \hat{y}))$$

Log Loss high accuracy $\rightarrow 0$

Decision Trees

Each internal node corresponds to a test.

Each branch corresponds to a result of the test.

Each leaf node assians a classification

Decision Trees algorithm

- 1. Choose an attribute from your dataset.
- 2. Calculate the significance of attribute in splitting of data.
- 3. Split data based on the value of the best attribute.
- 4. Go to step 1.

Entropy: Measure of randomness or uncertainty.

The lower the Entropy, the less uniform the distribution, the purer the node.

Entropy =
$$-p(A)\log(p(A)) - p(B)\log(p(B))$$

Which attribute is the best?

The tree with the higher Information Gain after splitting.

Information Gain: is the information that can increase the level of certainty after splitting.

Information Gain = (Entropy before split) - (weighted entropy after split)

Logistic Regression

When is logistic regression suitable?

If your data is binary.

If you need probabilistic results.

When you need a linear decision boundary.

If you need to understand the impact of a feature.

$$\sigma\left(\theta^T X\right) = \frac{1}{1 + e^{-\theta^T X}}$$

General cost function

$$\sigma\left(\theta^T X\right) \longrightarrow P(y=1 \mid x)$$

Change the weight -> Reduce the cost

Cost function

$$Cost(\hat{y}, y) = \begin{cases} -\log(\hat{y}) & \text{if } y = 1\\ -\log(1 - \hat{y}) & \text{if } y = 0 \end{cases}$$

$$J(\theta) = -\frac{1}{m} \sum_{i=1}^{m} y^{i} \log(\hat{y}^{i}) + (1 - y^{i}) \log(1 - \hat{y}^{i})$$

$$\nabla J = \left[\frac{\partial J}{\partial \theta_1}, \frac{\partial J}{\partial \theta_2}, \frac{\partial J}{\partial \theta_3}, \dots, \frac{\partial J}{\partial \theta_k} \right]$$
$$\theta_{new} = \theta_{prv} - \eta \nabla J$$

Logistic Regression algorithm

- 1. initialize the parameters randomly.
- 2. Feed the cost function with training set, and calculate the error.
- 3. Calculate the gradient of cost function.
- 4. Update weights with new values.
- 5. Go to step 2 until cost is small enough.

Support Vector Machine

SVM is a supervised algorithm that classifies cases by finding a separator.

- 1. Mapping data to a high-dimensional feature space
- 2. Finding a separator

Kernelling in SVM is Mapping data into a higher dimensional space, in such a way that can change a linearly inseparable dataset into a linearly separable dataset. (Linear, Polynomial, RBF, Sigmoid) find hyperplane:

Find \mathbf{w} and b such that

$$\Phi(\mathbf{w}) = 1/2\mathbf{w}^{\mathrm{T}}\mathbf{w}$$
 is minimized;

and for all
$$\{(\mathbf{x_i}, y_i)\}: y_i(\mathbf{w}^T\mathbf{x_i} + b) \ge 1$$

Support Vector Machine

Advantages:

- Accurate in high-dimensional spaces
- Memory efficient

Disadvantages:

- Prone to over-fitting
- No probability estimation
- Small datasets

SVM applications

- Image recognition
- Text category assignment
- Detecting spam
- Sentiment analysis
- Gene Expression Classification
- Regression, outlier detection and clustering

Clustering

Cluster: A group of objects that are similar to other objects in the cluster, and dissimilar to data points in other clusters.

Reference

References