scikit-learn

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Mines Linux Users Group

Introduction

Machine Learning - What is it really?

- Goal: Extract Knowledge from Data
- Sometimes called predictive analysis or statistical learning
- Given a large matrix of observations X, fit a function f(x)
 that maps observation x to a response variable y

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Important Terms

- **Classifiers** Algorithms that learn functions to map observations to a *discrete* response. E.g., is this tumor malignant or benign? Is this email spam or not?
- **Regressors** Algorithms that learn functions to map observations to a *continuous* response. E.g., how much should this house cost?
- **Underfitting** The learned function is too simple. "We barely studied for the exam."
- Overfitting The learned function is too complex. "We memorized all the practice problems, but don't understand the material."
- **Generalization** How well does the learned function extend to new observations?

- Provides many machine learning tools with a common Estimator interface¹
- Built in helpers for common ML tasks (e.g., metrics, preprocessing)
- Easily combine algorithms to make a complex pipeline²
- Relies heavily on numpy and scipy, often used with pandas

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Supervised Learning

Learning to Predict Breast Cancer

tree.fit(X_train, y_train) # Learn a Decision Function

from sklearn.tree import DecisionTreeClassifier
from sklearn.datasets import load_breast_cancer

Evaluating Accuracy of a Model

```
# How well did we do?
train_acc = tree.score(X_train, y_train)
test_acc = tree.score(X_test, y_test)
print("Training Accuracy: {:.3f}".format(train_acc))
print("Testing Accuracy: {:.3f}".format(test_acc))
# Training Accuracy: 1.000
# Testing Accuracy: 0.923
```

- Decision trees are a common first step, because they're easy to interpret and don't require much preprocessing
- Decision trees are prone to overfitting, so a good improvement is the RandomForest
- Support Vector Machines, Logistic/Linear Regression, and Artificial Neural Networks are commonly the first algorithms studied
- See the scikit-learn documentation for a comprehensive guide of available algorithms

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Becoming a "Data Scientist"

- 1. Get some (more) data
- 2. Pick an algorithm (or algorithm chain)
- 3. Train the model
- 4. Test generalization ability of trained model
- 5. Good enough? Done. Else, go back to step 1 or 2.

Then, tell people you're a genius . . . it's that easy!

Unsupervised Learning

Distinction from Supervised Learning

- **Supervised Learning** You tell the model what the correct answers are for training examples.
- **Unsupervised Learning** You ask the model to extract information from a dataset.
- **Unsupervised Clustering** Partition data into similar groups.

Example: K-Means Clustering

Unsupervised Transformations Create new representations of data. Example: Principal Component Analysis

Model Evaluation and Improvement

- Accuracy is not always the best metric for your system
- Plenty of others exist, pick the best for your business costs
- Look in the sklearn.metrics module for alternatives
- You can also use your own evaluation function!

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Cross Validation

Never Fit Models to Test Data! Ever!

Learning the parameters of a prediction function and testing it on the same data is a methodological mistake: a model that would just repeat the labels of the samples that it has just seen would have a perfect score but would fail to predict anything useful on yet-unseen data. This situation is called overfitting.

	◀ Total Number of Dataset →	
Experiment 1		
Experiment 2		Training Validation
Experiment 3		
Experiment 4		
Experiment 5		

Grid Search with Cross Validation

```
from sklearn.tree import DecisionTreeClassifier
from sklearn.datasets import load_breast_cancer
from sklearn.model_selection import train_test_split
from sklearn.model_selection import GridSearchCV
cancer = load breast cancer() # Get some data
X_train, X_test, y_train, y_test = train_test_split(
        cancer.data, cancer.target,
        stratify=cancer.target, random_state=1337)
tree = DecisionTreeClassifier(random_state=7331)
search_grid = {'criterion': ['gini', 'entropy'],
        'max_depth' : [5, 10, 15, 20]}
# search_grid could also be a list of dicts
search = GridSearchCV(tree, search_grid, cv=5)
search.fit(X_train, v_train)
print(search.best_params_)
```

Pipelines

Pipelines

Use Pipeline to combine multiple estimators into a single estimator. Two conveniences:

- 1. Convenience: You only have to call fit and predict once on your data to fit a whole sequence of estimators.
- 2. Joint parameter selection: You can grid search over parameters of all estimators in the pipeline at once.

A Simple Pipeline

Grid Search - Tuning a Complex Pipeline

```
from sklearn.pipeline import make_pipeline
from sklearn.svm import SVC
from sklearn.decomposition import PCA
from sklearn.preprocessing import StandardScaler
from sklearn.model_selection import GridSearchCV
pipe = make_pipeline(PCA(), StandardScaler(), SVC())
params = dict(pca_n_components=[2, 5, 10],
        svc__C=[0.1, 10, 100])
grid = GridSearchCV(pipe, param_grid=params)
# Next, call grid.fit on some training data
# This will use cross validation to estimation performance using each
# combination of parameters for pipeline in params dict
# With fitted model
print(grid.best_params_)
```

Questions?

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