# Contents

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This project can be useful to anyone who wishes to do supervised classification under interpretability constraints: explicit logical rules have to be used for classifying data. This project is particularly suitable for supervised anomaly detection, i.e. imbalanced classification. Application domains include fraud detection, predictive maintenance, intrusion detection, churn detection…

This project comes with a skrules module which contains a single estimator with unit tests, along with examples and benchmarks.
CHAPTER 1

API Documentation

- skope_rules
CHAPTER 2

General examples

Introductory examples.

2.1 SkopeRules example

An example using SkopeRules for imbalanced classification.

SkopeRules find logical rules with high precision and fuse them. Finding good rules is done by fitting classification and regression trees to sub-samples. A fitted tree defines a set of rules (each tree node defines a rule); rules are then tested out of the bag, and the ones with higher precision are selected and merged. This produces a real-valued decision function, reflecting for each new sample how many rules (each weighted by respective precision) have found it abnormal.

```python
import numpy as np
import matplotlib.pyplot as plt
from skrules import SkopeRules

print(__doc__)

rng = np.random.RandomState(42)

n_inliers = 1000
n_outliers = 50

# Generate train data
I = 0.5 * rng.randn(int(n_inliers / 2), 2)
X_inliers = np.r_[I + 2, I - 2]
O = 0.5 * rng.randn(n_outliers, 2)
X_outliers = O  # np.r_[O, O + [2, -2]]
X_train = np.r_[X_inliers, X_outliers]
y_train = [0] * n_inliers + [1] * n_outliers
```
### 2.1.1 Training the SkopeRules classifier

```python
# fit the model
clf = SkopeRules(random_state=rng, n_estimators=10)
clf.fit(X_train, y_train)

# plot the line, the samples, and the nearest vectors to the plane
xx, yy = np.meshgrid(np.linspace(-5, 5, 50), np.linspace(-5, 5, 50))
Z = clf.decision_function(np.c_[xx.ravel(), yy.ravel()])
Z = Z.reshape(xx.shape)
plt.title("Skope Rules, value of the decision_function method")
plt.contourf(xx, yy, Z, cmap=plt.cm.Blues)

a = plt.scatter(X_inliers[:, 0], X_inliers[:, 1], c='white',
                 s=20, edgecolor='k')
b = plt.scatter(X_outliers[:, 0], X_outliers[:, 1], c='red',
                 s=20, edgecolor='k')
plt.axis('tight')
plt.xlim((-5, 5))
plt.ylim((-5, 5))
plt.legend([a, b],
           ['inliers', 'outliers'],
           loc="upper left")
plt.show()
```

---

**Skope Rules, value of the decision_function method**

- **inliers**
- **outliers**
2.1.2 Extracting top rules

On the 4 following figures, the predict_top_rules method is used with several values of n_rules. n_rules = 2 means that the prediction is done using only the 2 best rules.

```python
print('The 4 most precise rules are the following:')
for rule in clf.rules_[:4]:
    print(rule[0])

fig, axes = plt.subplots(2, 2, figsize=(12, 5),
    sharex=True, sharey=True)
for i_ax, ax in enumerate(np.ravel(axes)):
    Z = clf.predict_top_rules(np.c_[xx.ravel(), yy.ravel()], i_ax+1)
    Z = Z.reshape(xx.shape)
    ax.set_title("Prediction with predict_top_rules, n_rules="+str(i_ax+1))
    ax.contourf(xx, yy, Z, cmap=plt.cm.Blues)
    a = ax.scatter(X_inliers[:, 0], X_inliers[:, 1], c='white',
        s=20, edgecolor='k')
    b = ax.scatter(X_outliers[:, 0], X_outliers[:, 1], c='red',
        s=20, edgecolor='k')
    ax.axis('tight')
plt.xlim((-5, 5))
plt.ylim((-5, 5))
plt.legend([a, b],
    ["inliers", "outliers"],
    loc="upper left")
plt.show()
```

Out:

The 4 most precise rules are the following:
c0 <= 1.15681171417 and c0 > -0.680330395699 and c1 <= 1.08434700966
c0 > -0.841694712639 and c1 <= 0.687212407589 and c1 > -1.21965646744
c0 <= 1.03912234306 and c0 > -0.663101434708 and c1 <= 1.27947068214
c0 > -0.841694712639 and c1 <= 0.687212407589 and c1 > -1.43224191666

Total running time of the script: ( 0 minutes 1.548 seconds)
2.2 Example: detecting defaults on retail credits

SkopeRules finds logical rules with high precision and fuse them. Finding good rules is done by fitting classification and regression trees to sub-samples. A fitted tree defines a set of rules (each tree node defines a rule); rules are then tested out of the bag, and the ones with higher precision are kept.

This example aims at finding logical rules to predict credit defaults. The analysis shows that setting.

2.2.1 Data import and preparation

There are 3 categorical variables (SEX, EDUCATION and MARRIAGE) and 20 numerical variables. The target (credit defaults) is transformed in a binary variable with integers 0 (no default) and 1 (default). From the 30000 credits, 50% are used for training and 50% are used for testing. The target is unbalanced with a 22%/78% ratio.

```python
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
from sklearn.metrics import roc_curve, precision_recall_curve
from sklearn.ensemble import RandomForestClassifier
from sklearn.model_selection import GridSearchCV
from sklearn.utils import shuffle
from skrules import SkopeRules
from skrules.datasets import load_credit_data

print(__doc__)
rng = np.random.RandomState(1)

# Importing data
dataset = load_credit_data()
X = dataset.data
y = dataset.target
# Shuffling data, preparing target and variables
data, y = shuffle(np.array(X), y, random_state=rng)
data = pd.DataFrame(data, columns=X.columns)
for col in ['ID']:
    del data[col]

# Quick feature engineering
data = data.rename(columns={'PAY_0': 'PAY_1'})
old_PAY = ['PAY_3', 'PAY_4', 'PAY_5', 'PAY_6']
data['PAY_old_mean'] = data[old_PAY].apply(lambda x: np.mean(x), axis=1)
old_BILL_AMT = ['BILL_AMT3', 'BILL_AMT4', 'BILL_AMT5', 'BILL_AMT6']
data['BILL_AMT_old_mean'] = data[old_BILL_AMT].apply(  
    lambda x: np.mean(x), axis=1)
data['BILL_AMT_old_std'] = data[old_BILL_AMT].apply(  
    lambda x: np.std(x), axis=1)
old_PAY_AMT = ['PAY_AMT3', 'PAY_AMT4', 'PAY_AMT5', 'PAY_AMT6']
data['PAY_AMT_old_mean'] = data[old_PAY_AMT].apply(  
    lambda x: np.mean(x), axis=1)
data['PAY_AMT_old_std'] = data[old_PAY_AMT].apply(  
    lambda x: np.std(x), axis=1)
```
```python
lambda x: np.std(x), axis=1)

data.drop(old_PAY_AMT + old_BILL_AMT + old_PAY, axis=1, inplace=True)

# Creating the train/test split
feature_names = list(data.columns)
print("List of variables used to train models : ", str(feature_names))
data = data.values
n_samples = data.shape[0]
n_samples_train = int(n_samples / 2)
y_train = y[:n_samples_train]
y_test = y[n_samples_train:]
X_train = data[:n_samples_train]
X_test = data[n_samples_train:]

Out:
List of variables used to train models : [u'LIMIT_BAL', u'SEX', u'EDUCATION', u'MARRIAGE', u'AGE', 'PAY_1', 'PAY_2', 'BILL_AMT1', 'BILL_AMT2', 'PAY_AMT1', 'PAY_AMT2', 'PAY_old_mean', 'BILL_AMT_old_mean', 'BILL_AMT_old_std', 'PAY_AMT_old_mean', 'PAY_AMT_old_std']

2.2.2 Benchmark with a Random Forest classifier

This part shows the training and performance evaluation of a random forest model. The objective remains to extract rules which targets credit defaults.

rf = GridSearchCV(
    RandomForestClassifier(
        random_state=rng,
        n_estimators=50,
        class_weight='balanced'),
    param_grid={'max_depth': range(3, 8, 1),
                 'max_features': np.linspace(0.1, 1., 5)},
    scoring={'AUC': 'roc_auc'}, cv=5,
    refit='AUC', n_jobs=-1)
rf.fit(X_train, y_train)
scoring_rf = rf.predict_proba(X_test)[:, 1]
print("Random Forest selected parameters : \$s\) % rf.best_params_

# Plot ROC and PR curves

fig, axes = plt.subplots(1, 2, figsize=(12, 5),
    sharex=True, sharey=True)
ax = axes[0]
fpr_RF, tpr_RF, _ = roc_curve(y_test, scoring_rf)
ax.step(fpr_RF, tpr_RF, linestyle='-.', c='g', lw=1, where='post')
ax.set_title("ROC", fontsize=20)
ax.legend(loc='upper center', fontsize=8)
ax.set_xlabel('False Positive Rate', fontsize=18)
ax.set_ylabel('True Positive Rate (Recall)', fontsize=18)
ax = axes[1]
```

2.2. Example: detecting defaults on retail credits
precision_RF, recall_RF, _ = precision_recall_curve(y_test, scoring_rf)
ax.step(recall_RF, precision_RF, linestyle='-.', c='g', lw=1, where='post')
ax.set_title("Precision-Recall", fontsize=20)
ax.set_xlabel('Recall (True Positive Rate)', fontsize=18)
ax.set_ylabel('Precision', fontsize=18)
plt.show()

Out:

Random Forest selected parameters : {'max_features': 0.55, 'max_depth': 7}

The ROC and Precision-Recall curves illustrate the performance of Random Forests in this classification task. Suppose now that we add an interpretability contraint to this setting: Typically, we want to express our model in terms of logical rules detecting defaults. A random forest could be expressed in term of weighted sum of rules, but 1) such a large weighted sum, is hardly interpretable and 2) simplifying it by removing rules/weights is not easy, as optimality is targeted by the ensemble of weighted rules, not by each rule. In the following section, we show how SkopeRules can be used to produce a number of rules, each seeking for high precision on a potentially small area of detection (low recall).

### 2.2.3 Getting rules with skrules

This part shows how SkopeRules can be fitted to detect credit defaults. Performances are compared with the random forest model previously trained.

```python
# fit the model
clf = SkopeRules(
    max_depth_duplication=3, max_depth=3, max_features=0.5,
    max_samples_features=0.5, random_state=rng, n_estimators=20,
    feature_names=feature_names, recall_min=0.04, precision_min=0.6)
clf.fit(X_train, y_train)

# in the score_top_rules method, a score of k means that rule number k
# vote positively, but not rules 1, ..., k-1. It will allow us to plot
# performance of each rule separately on the ROC and PR plots.
scoring = clf.score_top_rules(X_test)
print(str(len(clf.rules_)) + ' rules have been built.')
print('The 5 most precise rules are the following: ')
```

Out:
for rule in clf.rules_[:5]:
    print(rule[0])

curves = [roc_curve, precision_recall_curve]
xlabels = ['False Positive Rate', 'Recall (True Positive Rate)']
ylabels = ['True Positive Rate (Recall)', 'Precision']

fig, axes = plt.subplots(1, 2, figsize=(12, 5),
                        sharex=True, sharey=True)

ax = axes[0]
fpr, tpr, _ = roc_curve(y_test, scoring)
fpr_rf, tpr_rf, _ = roc_curve(y_test, scoring_rf)
ax.scatter(fpr[:-1], tpr[:-1], c='b', s=10, label="rules of SkopeRules")
ax.step(fpr_RF, tpr_RF, linestyle='-.', c='g', lw=1, where='post',
        label="Random Forest")
ax.set_title("ROC", fontsize=20)
ax.legend(loc='upper center', fontsize=8)
ax.set_xlabel('False Positive Rate', fontsize=18)
ax.set_ylabel('True Positive Rate (Recall)', fontsize=18)

ax = axes[1]
precision, recall, _ = precision_recall_curve(y_test, scoring)
precision_rf, recall_rf, _ = precision_recall_curve(y_test, scoring_rf)
ax.scatter(recall[1:-1], precision[1:-1], c='b', s=10,
           label="rules of SkopeRules")
ax.step(recall_RF, precision_RF, linestyle='-.', c='g', lw=1, where='post',
        label="Random Forest")
ax.set_title("Precision-Recall", fontsize=20)
ax.set_xlabel('Recall (True Positive Rate)', fontsize=18)
ax.set_ylabel('Precision', fontsize=18)
plt.show()
The ROC and Precision-Recall curves show the performance of the rules generated by SkopeRules the (the blue points) and the performance of the Random Forest classifier fitted above. Each blue point represents the performance of a set of rules: Starting from the left on the precision-recall curve, the kth point represents the score associated to the concatenation (union) of the k first rules, etc. Thus, each blue point is associated with an interpretable classifier, which is a combination of a few rules. In terms of performance, each of these interpretable classifiers compare well with Random Forest, while offering complete interpretation. The range of recall and precision can be controlled by the precision_min and recall_min parameters. Here, setting precision_min to 0.6 force the rules to have a limited recall.

Total running time of the script: (1 minutes 53.372 seconds)

See the README for more information.