

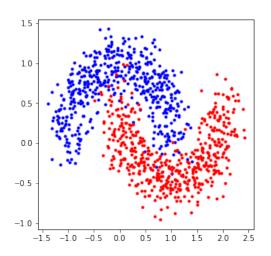
Performance Evaluation and Hyperparameter Estimation Machine Learning – Laboratory

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Running Example: Two-Moon Dataset

from sklearn.datasets import make moons

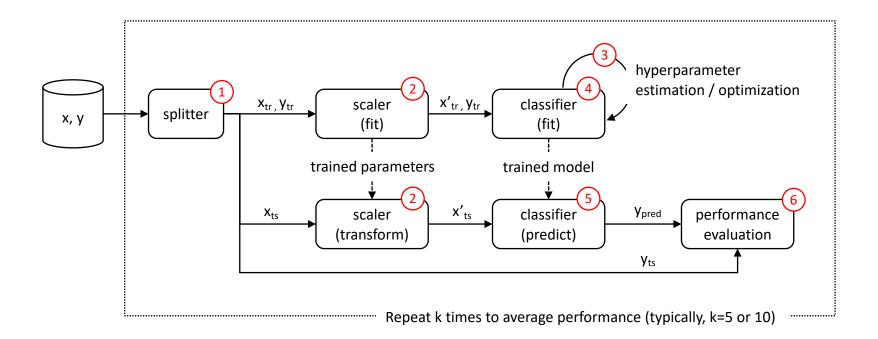


https://scikit-learn.org/stable/modules/generated/sklearn.datasets.make moons.html

Running Example: MNIST Digit Classification

- Dataset of handwritten digits
- 10 classes: 0, 1, ..., 9
- 28x28 images corresponding to 784 feature (pixel) values
- Pixel values in {0, 1, ..., 255}

ML System Design



ML System Design

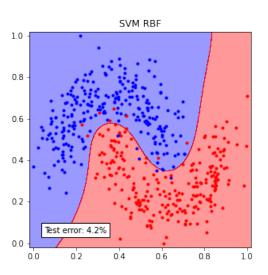
- 1. Sampling a training and a testing set (from the same underlying distribution)
 - e.g., splitting data x, y at random in x_tr, y_tr, x_ts, y_ts
- 2. Scaling training and test data (using parameters estimated on training data)
 - from sklearn.preprocessing import MinMaxScaler
- 3. Estimating classifier hyperparameters on training data
- 4. Fitting the classifier on training data
 - clf.fit(x_tr, y_tr)
- 5. Predicting the class labels of testing data
 - clf.predict(x ts)
- 6. Evaluating accuracy or classification error

Exercise 1 – Performance Evaluation

- Implement a machine-learning pipeline to process MNIST digit data as described before
- Use default hyperparameters for some sklearn classifiers (e.g., SVM, kNN)
- Estimate classification error on different runs (i.e., data splits) and average the results
- What happens if we change the classifier hyperparameters?

Solution

```
x, y = make moons(n samples=1000, noise=0.2)
splitter = ShuffleSplit(n splits=5, train size=0.5)
scaler = MinMaxScaler()
clf = svm.SVC(kernel='rbf', C=10, gamma=10.0)
clf name = 'SVM RBF'
for tr idx, ts idx in splitter.split(x, y):
    xtr = x[tr idx, :]
    ytr = y[tr idx]
    xts = x[ts idx, :]
    yts = y[ts idx]
    xtr = scaler.fit transform(xtr)
    xts = scaler.transform(xts)
    clf.fit(xtr, ytr)
    ypred = clf.predict(xts)
    error = (ypred != yts).mean()
    print("Test error: {:.1%}".format(error))
```



Test error: 4.0% Test error: 4.6%

...

Test error: 4.2%

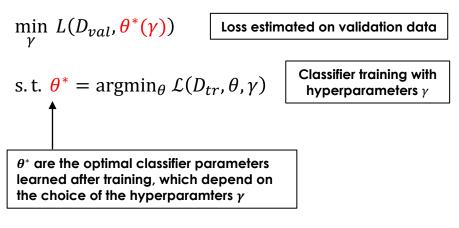
Mean test error: 4.2% +/- 2.7%

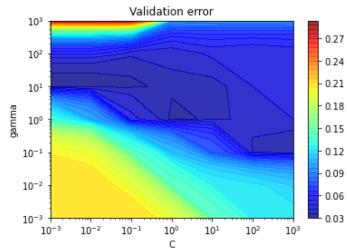
Hyperparameter Estimation / Optimization

- How can we select the best hyperparameters for our classifier, i.e., the hyperparameters that we expect to perform best on the test set?
- Remark: we cannot look at the test set more than once. We cannot optimize the
 hyperparameters by iteratively testing how they perform on the test set.
 - This will lead us to overestimate the performance of the system in production!
- Validation data. The standard procedure is to optimize these hyperparameters using a (separate) validation set and compute their expected performance on such data

Hyperparameter Estimation / Optimization

 In principle, hyperparameter optimization aims to solving the following (bi-level) optimization problem:



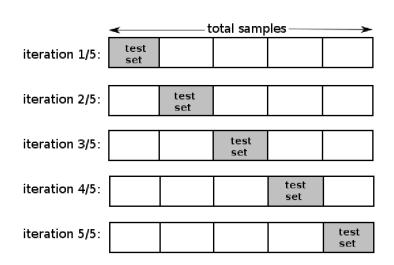


Hyperparameter Estimation / Optimization

- Hyperparameter optimization is normally performed using grid search with k-fold cross validation (even though more complex optimization techniques can be used)
- The procedure works as follows:
 - Set different hyperparameter values to test (the grid space) $C \in \{10, 100, 1000\}$ $\gamma \in \{0.1, 0.2, 0.5, 1.0\}$
 - For each set of hyperparameter values (e.g., C and gamma pairs):
 - Train the classifier on the training set with the current hyperparameters
 - Evaluate the performance on the validation set
 - Repeat K times on different training-validation splits, and average the performance
- At the end of this procedure, for each set of hyperparameters we will get an averaged performance estimation (e.g., the mean validation loss/error)
- We can thus select the best set of hyperparameters as those that minimize the mean loss/error (or that maximize the mean validation accuracy)

Hyperparameter Optimization with k-Fold Cross Validation

- Normally k-fold cross validation is used to create training-validation sets for hyperparameter estimation
 - Split data into K folds
 - K-1 folds are used for training, 1 for validation
 - the process is repeated K times to estimate performance with the given hyperparameters (changing the validation fold at every iteration)
- 1. The process is repeated for every set of hyperparameters
- 2. The best set of hyperparameters is eventually selected
- 3. Classifier is trained using such hyperparameters on the full training set



Exercise 2 – Hyperparameter Optimization

- Solve again Exercise 1 but including hyperparameter optimization
 - For the linear SVM, optimize C
 - For the SVM with the RBF kernel, optimize C and gamma
 - For kNN, optimize k
- Use grid search with 5-fold cross validation to select the best hyperparameters that maximize the mean validation accuracy
- GridSearchCV from sklearn already implements the whole procedure
 - https://scikit-learn.org/stable/modules/generated/sklearn.model_selection.GridSearchCV.html

Solution

```
from sklearn.model selection import GridSearchCV
# let's go back to the two-moon dataset
x, y = make moons (n samples=1000, noise=0.2)
clf = GridSearchCV(estimator=svm.SVC(kernel='linear'),
                   param grid={'C': [0.01, 0.1, 1, 10, 100]})
acc = run(x, y, splitter, scaler, clf)
print("Hyperparameter estimation (5-fold xval)")
print(" - Best parameters set found on development set:", clf.best params )
print(" - Grid scores on development set:")
means = clf.cv results ['mean test score']
stds = clf.cv results ['std test score']
for mean, std, params in zip(means, stds, clf.cv results ['params']):
   print("
                  \$0.3f (+/-\$0.03f) for \$r" \% (mean, std * 2, params))
print("Mean test accuracy: {:.1%} +/- {:.1%}\n".format(acc.mean(), 2*acc.std()))
```

Solution

- Optimization of the hyperparameter C of a linear SVM on the two-moon data
 - Note that we have mean validation accuracy estimated for each C value
 - The best value C=1 is selected and used to fit the classifier on the whole training set
 - Mean test accuracy is reasonably close to the best mean validation accuracy

```
Hyperparameter estimation (5-fold xval)

- Best parameters set found on development set: {'C': 1}

- Grid scores on development set:

0.510 (+/-0.000) for {'C': 0.01}

0.842 (+/-0.081) for {'C': 0.1}

0.858 (+/-0.085) for {'C': 1}

0.858 (+/-0.082) for {'C': 10}

0.852 (+/-0.091) for {'C': 100}

Mean test accuracy: 85.3% +/- 1.2%
```