# Putting it all together

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# **Objectives**

At the end of the day, you'll be able to:

- Use Pipeline for more compact and comprehensive code
- Build your own transformer and estimator
- Use FeatureUnion for running parallel processes for features extraction
- Use GridSearchCV for model tuning

### **Before**

We usually start modeling from simple steps like:

- Reading and splitting prepared data into train/test subsets
- Extracting target
- Extracting features
- In cycle: choose model, fit, predict, print the scores

Then we want more features, and more models, and more tuning...

Example: Classification of text documents using sparse features

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# Why Pipeline

Why use Pipeline instead of keeping the steps separate?

- It makes code more readable
- It keeps data during intermediate steps
- It makes it easy to change the order or add/remove steps
- You only have to call fit and predict once
- Joint parameter selection: grid search over parameters of all estimators at once

## scikit-learn Pipeline

Pipeline can be used to chain multiple estimators into one.

All estimators in a pipeline, except the last one, must be transformers. The last estimator may be any type (transformer, classifier, etc.).

# scikit-learn Pipeline

Figure 1:Simple pipeline

# make\_pipeline

```
make_pipeline(Binarizer(), MultinomialNB())

notation

make_pipeline(Binarizer(), MultinomialNB())

notation

make_pipeline(Binarizer(), MultinomialNB())

notation

n
```

Figure 2: Make pipeline

# **Custom Steps**

```
pipeline = Pipeline([('selector', ItemSelector(key='body')),
                          ('stats', TextStats()), # returns a list of dicts
                          ('vect', DictVectorizer())])
    class ItemSelector(BaseEstimator, TransformerMixin):
      ""For data grouped by feature, select subset of data at a provided key.'''
      def __init__(self, key):
        self.kev = kev
34
      def fit(self, x, y=None):
        return self
      def transform(self, data_dict):
        return data_dict[self.key]
```

Figure 3:Simple pipeline

#### **Custom Transformers**

A transformer is just an object that responds to fit, transform, and fit\_transform.

Inheriting from Transformer Mixin is not required, but helps to communicate intent, and gets you fit\_transform for free.

### **Custom Transformers**

```
def fit(self, x, y=None):
return self

def transform(self, posts):
return [{'length': len(text), 'num_sentences': text.count('.')}
for text in posts]
```

Figure 4:Simple pipeline

# Using FunctionTransformer

You can convert an existing Python function into a transformer with FunctionTransformer:

```
def all but first column(X):
  return X[:. 1:]
def drop first component(X, v):
  """ Create a pipeline with PCA and the column selector and use it to
  transform the dataset.
  111111
  pipeline = make pipeline(
    PCA(). FunctionTransformer(all but first column).
  X train, X test, v train, v test = train test split(X, v)
  pipeline.fit(X train, v train)
  return pipeline.transform(X test), v test
```

A estimator is an object that fits a model based on some training data and is capable of inferring some properties on new data. All estimators implement the fit method.

To create custom estimator, you need to implement the following interface:

- get\_params([deep]) Get parameters for this estimator
- $\bullet$  set\_params(\*\*params) Set the parameters of this estimator

You can inherit from BaseEstimator and optionally the mixin classes in sklearn.base.

```
class MajorityClassifier(BaseEstimator, ClassifierMixin):
      """Predicts the majority class of its training data."""
      def init (self):
45
        pass
      def fit(self, X, v):
        self.classes_, indices = np.unique(["foo", "bar", "foo"],
                                           return inverse=True)
        self.majority_ = np.argmax(np.bincount(indices))
        return self
      def predict(self, X):
        return np.repeat(self.classes_[self.majority_], len(X))
```

Figure 6: Custom Estimator

If you do not want to make your code dependent on scikit-learn, the easiest way to implement the interface is:

```
def get_params(self, deep=True):

# suppose this estimator has parameters "alpha" and "recursive"

return {"alpha": self.alpha, "recursive": self.recursive}

def set_params(self, **parameters):

for parameter, value in parameters.items():

self.setattr(parameter, value)

return self
```

Figure 7: Custom Estimator

You can check whether your estimator adheres to the scikit-learn interface and standards by running utils.estimator\_checks.check\_estimator on the class:

```
from sklearn.utils.estimator_checks import check_estimator
from sklearn.svm import LinearSVC
check_estimator(LinearSVC) # passes
69
```

Figure 8: Custom Estimator

## Estimator types

Some common functionality depends on the kind of estimator passed. This distinction is implemented using the \_estimator\_type attribute:

- "classifier" for classifiers
- "regressor" for regressors
- "clusterer" for clustering methods

Inheriting from ClassifierMixin, RegressorMixin or ClusterMixin will set the attribute automatically.

#### **FeatureUnion**

FeatureUnion combines several transformer objects into a new transformer that combines their output:

- each of transformer objects is fit to the data independently
- transformers are applied in parallel
- the sample vectors they output are concatenated end-to-end into larger vectors

FeatureUnion and Pipeline can be combined to create complex models.

### **FeatureUnion**

Figure 9:Feature Union

# Pipeline parameters

The purpose of the pipeline is to assemble several steps that can be cross-validated together while setting different parameters.

It enables setting parameters of the various steps using their names and the parameter name separated by a "\_\_\_" (<estimator>\_\_\_<parameter>).

```
pca = PCA(n_components=2)

selection = SelectKBest(k=1)

pipeline = Pipeline([("features", combined_features), ("svm", svm)])

param_grid = dict(features__pca__n_components=[1, 2, 3],

param_grid = dict(features__univ_select__k=[1, 2],

svm__C=[0.1, 1, 10])
```

### GridSearchCV

Parameters that are not directly learnt within estimators can be set by searching a parameter space for the best Cross-validation.

The grid search provided by GridSearchCV exhaustively generates candidates from a grid of parameter values specified with the param\_grid parameter.

### GridSearchCV

The following param\_grid specifies that two grids should be explored: one with a linear kernel and C values in [1, 10, 100, 1000], and the second one with an RBF kernel, and the cross-product of C values ranging in [1, 10, 100, 1000] and gamma values in [0.001, 0.0001].

# Pipeline with GridSearchCV

```
# define a pipeline combining a text feature extractor with a simple
    # classifier
    pipeline = Pipeline([
         ('vect', CountVectorizer()).
         ('tfidf', TfidfTransformer()),
         ('clf', SGDClassifier()),
139 1)
     # uncommenting more parameters will give better exploring power but will
     # increase processing time in a combinatorial way
     parameters = {
         'vect__max_df': (0.5, 0.75, 1.0),
         #'vect__max_features': (None, 5000, 10000, 50000),
         'vect ngram range': ((1, 1), (1, 2)), # unigrams or bigrams
         #'tfidf use idf': (True, False),
         #'tfidf norm': ('l1', 'l2'),
         'clf alpha': (0.00001, 0.000001).
         'clf__penalty': ('l2', 'elasticnet'),
         #'clf__n_iter': (10, 50, 80),
     # find the best parameters for both the feature extraction and the
     # classifier
     grid search = GridSearchCV(pipeline, parameters, n jobs=-1, verbose=1)
    grid_search.fit(data.data, data.target)
    print("Best score: %0.3f" % grid_search.best_score_)
     best parameters = grid search.best estimator .get params()
     for param name in sorted(parameters.kevs()):
         print("\t%s: %r" % (param_name, best_parameters[param_name]))
```