Pipeline

MACHINE LEARNING WITH PYSPARK



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Leakage?

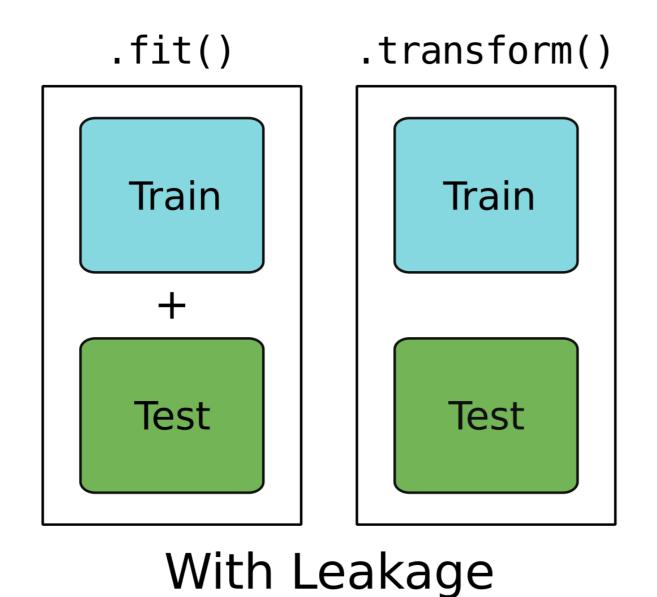
Only for training data.

.transform()

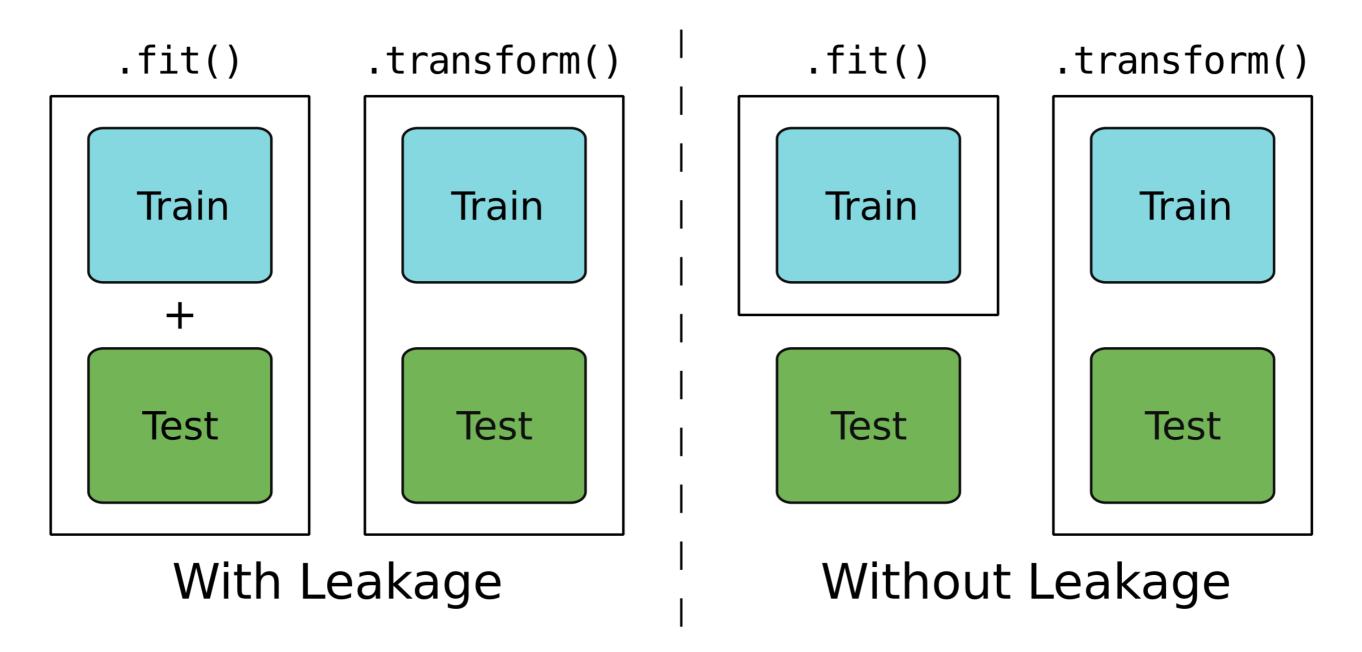
For testing and training data.



A leaky model

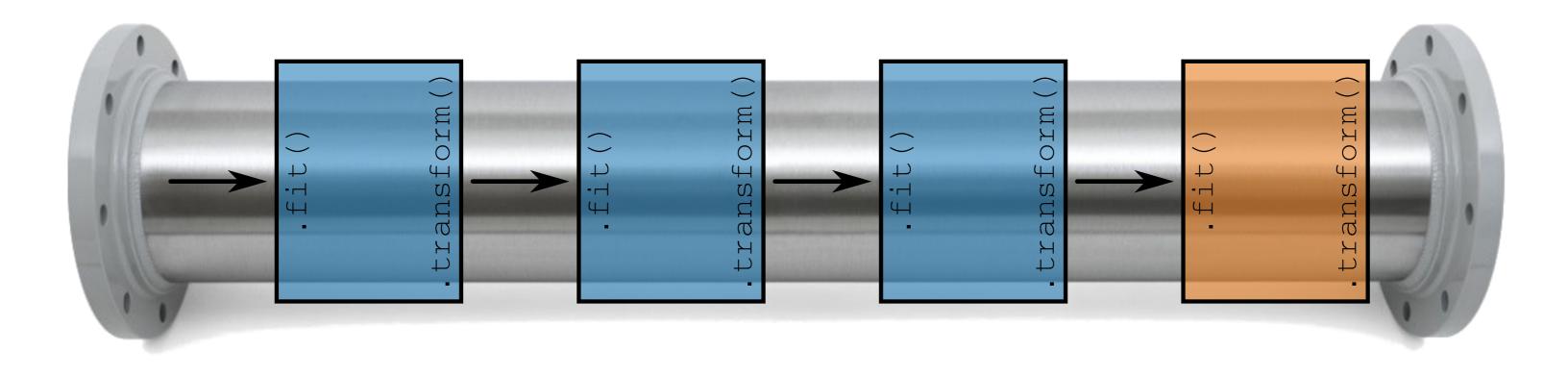


A watertight model



Pipeline

A pipeline consists of a series of operations.



You could apply each operation individually... or you could just apply the pipeline!

Cars model: Steps

```
indexer = StringIndexer(inputCol='type', outputCol='type_idx')
onehot = OneHotEncoder(inputCols=['type_idx'], outputCols=['type_dummy'])
assemble = VectorAssembler(
  inputCols=['mass', 'cyl', 'type_dummy'],
  outputCol='features'
)
regression = LinearRegression(labelCol='consumption')
```

Cars model: Applying steps

Training data

```
indexer = indexer.fit(cars_train)
cars_train = indexer.transform(cars_train)

onehot = onehot.fit(cars_train)
cars_train = onehot.transform(cars_train)
```

```
cars_train = assemble.transform(cars_train)
```

```
# Fit model to training data
regression = regression.fit(cars_train)
```

Testing data

```
cars_test = indexer.transform(cars_test)

cars_test = onehot.transform(cars_test)

cars_test = assemble.transform(cars_test)

# Make predictions on testing data
predictions = regression.transform(cars_test)
```

Cars model: Pipeline

Combine steps into a pipeline.

```
from pyspark.ml import Pipeline

pipeline = Pipeline(stages=[indexer, onehot, assemble, regression])
```

Training data

pipeline = pipeline.fit(cars_train)

Testing data

predictions = pipeline.transform(cars_test)

Cars model: Stages

Access individual stages using the .stages attribute.

```
# The LinearRegression object (fourth stage -> index 3)
pipeline.stages[3]
print(pipeline.stages[3].intercept)
```

```
4.19433571782916
```

```
print(pipeline.stages[3].coefficients)
```

```
DenseVector([0.0028, 0.2705, -1.1813, -1.3696, -1.1751, -1.1553, -1.8894])
```



Pipelines streamline workflow!

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Cross-Validation

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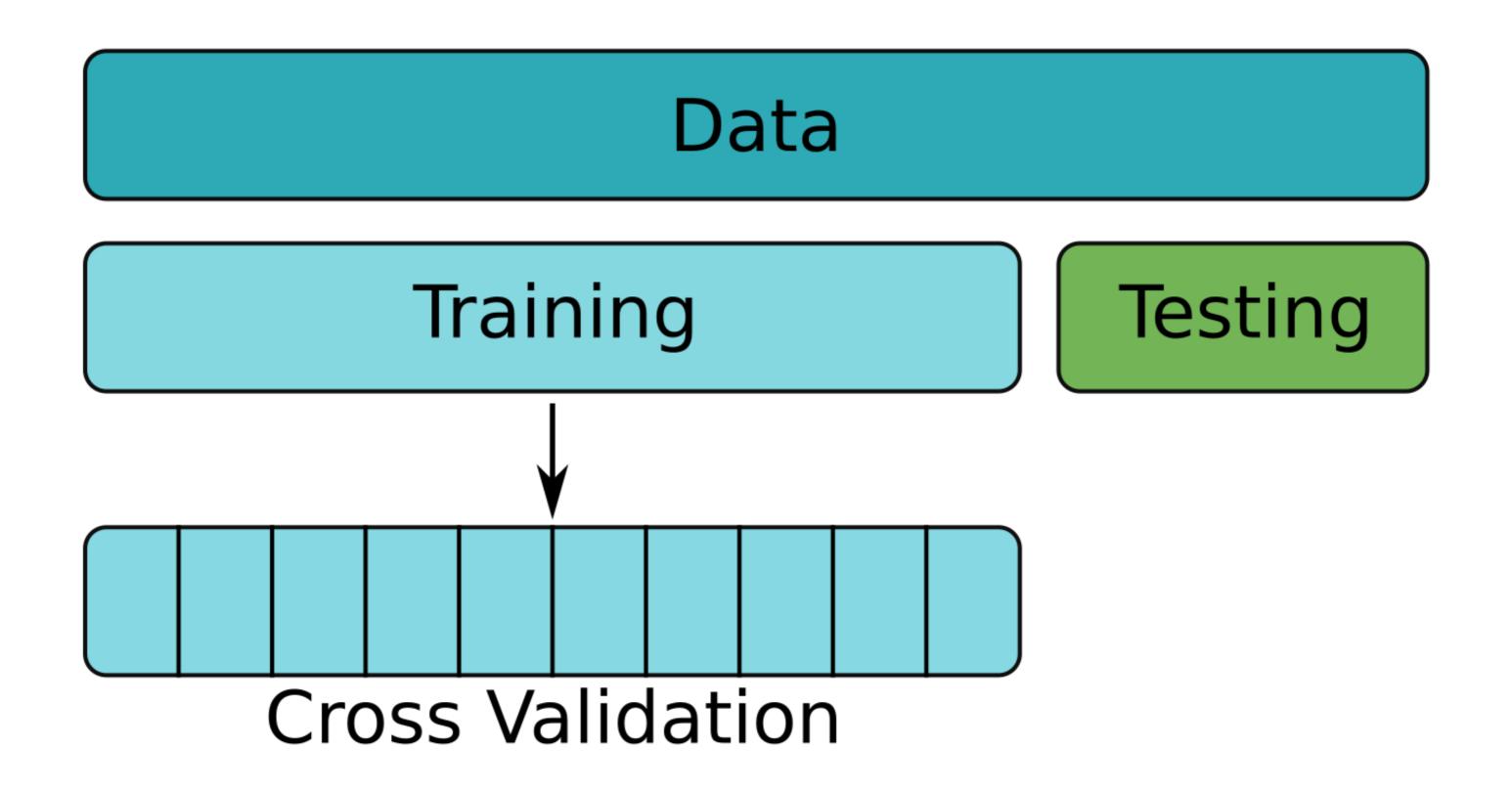
Data



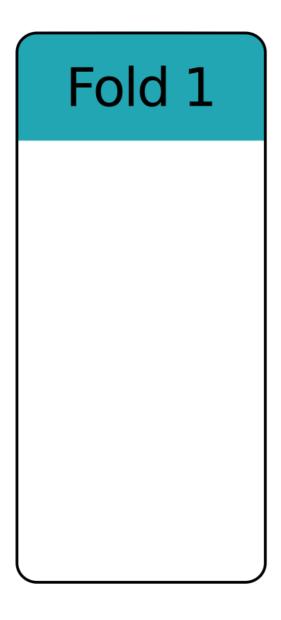
Data

Training

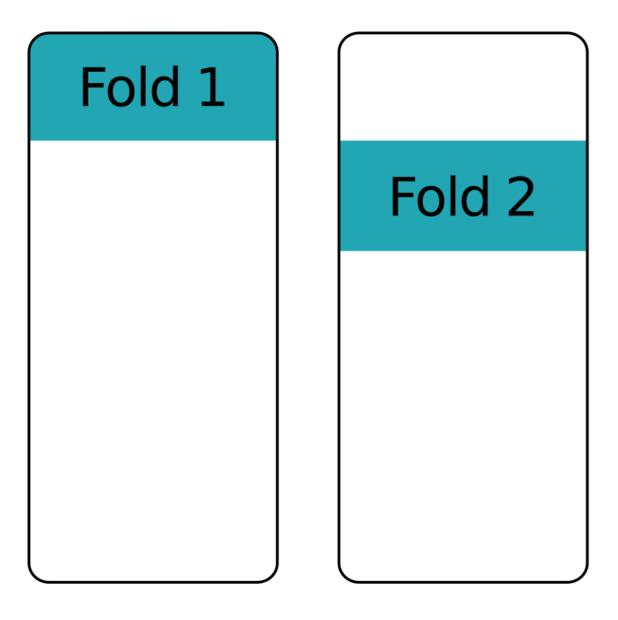
Testing



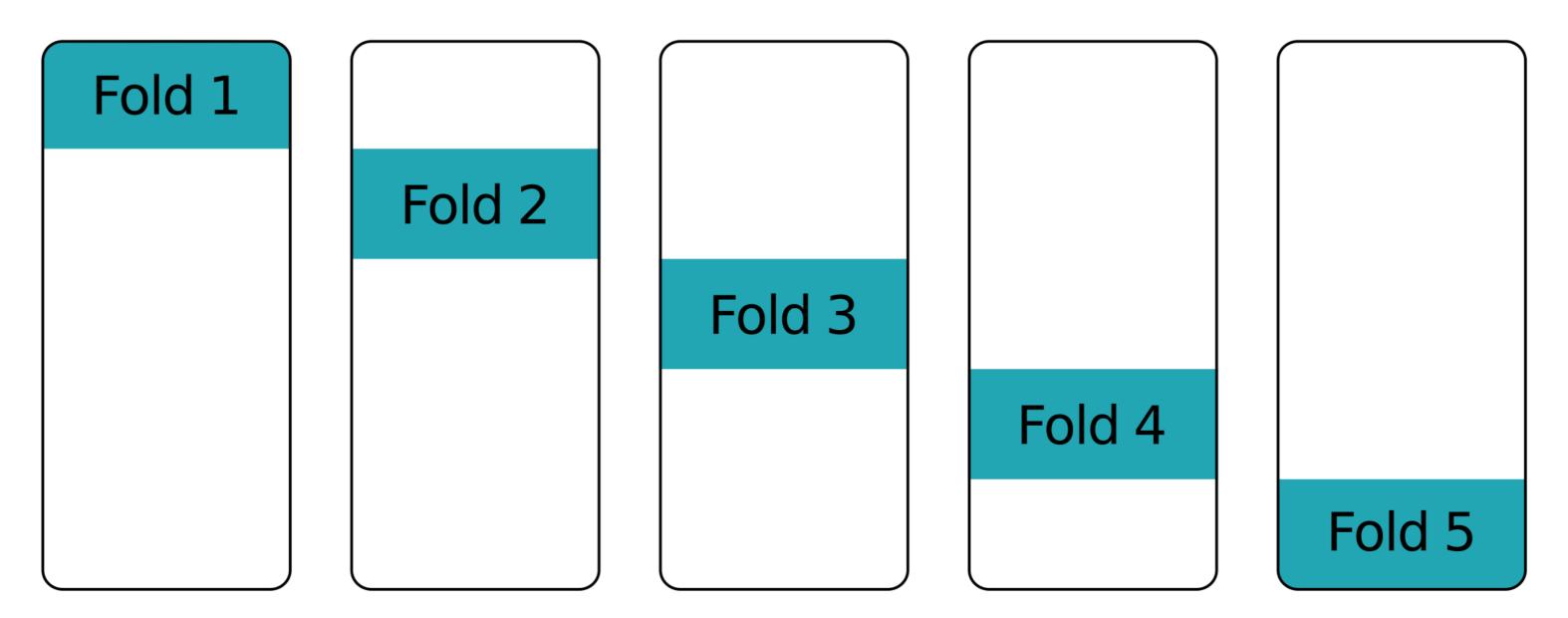
Fold upon fold - first fold



Fold upon fold - second fold



Fold upon fold - other folds



Cars revisited

```
cars.select('mass', 'cyl', 'consumption').show(5)
```

```
+----+
| mass|cyl|consumption|
+----+
|1451.0| 6| 9.05|
|1129.0| 4| 6.53|
|1399.0| 4| 7.84|
|1147.0| 4| 7.84|
|1111.0| 4| 9.05|
+----+
```

Estimator and evaluator

An object to build the model. This can be a pipeline.

```
regression = LinearRegression(labelCol='consumption')
```

An object to evaluate model performance.

```
evaluator = RegressionEvaluator(labelCol='consumption')
```



Grid and cross-validator

```
from pyspark.ml.tuning import CrossValidator, ParamGridBuilder
```

A grid of parameter values (empty for the moment).

```
params = ParamGridBuilder().build()
```

The cross-validation object.



Cross-validators need training too

Apply cross-validation to the training data.

```
cv = cv.fit(cars_train)
```

What's the average RMSE across the folds?

cv.avgMetrics

[0.800663722151572]



Cross-validators act like models

Make predictions on the original testing data.

```
evaluator.evaluate(cv.transform(cars_test))

# RMSE on testing data
0.745974203928479
```

Much smaller than the cross-validated RMSE.

```
# RMSE from cross-validation
0.800663722151572
```

A simple train-test split would have given an overly optimistic view on model performance.

Cross-validate all the models!

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Grid Search

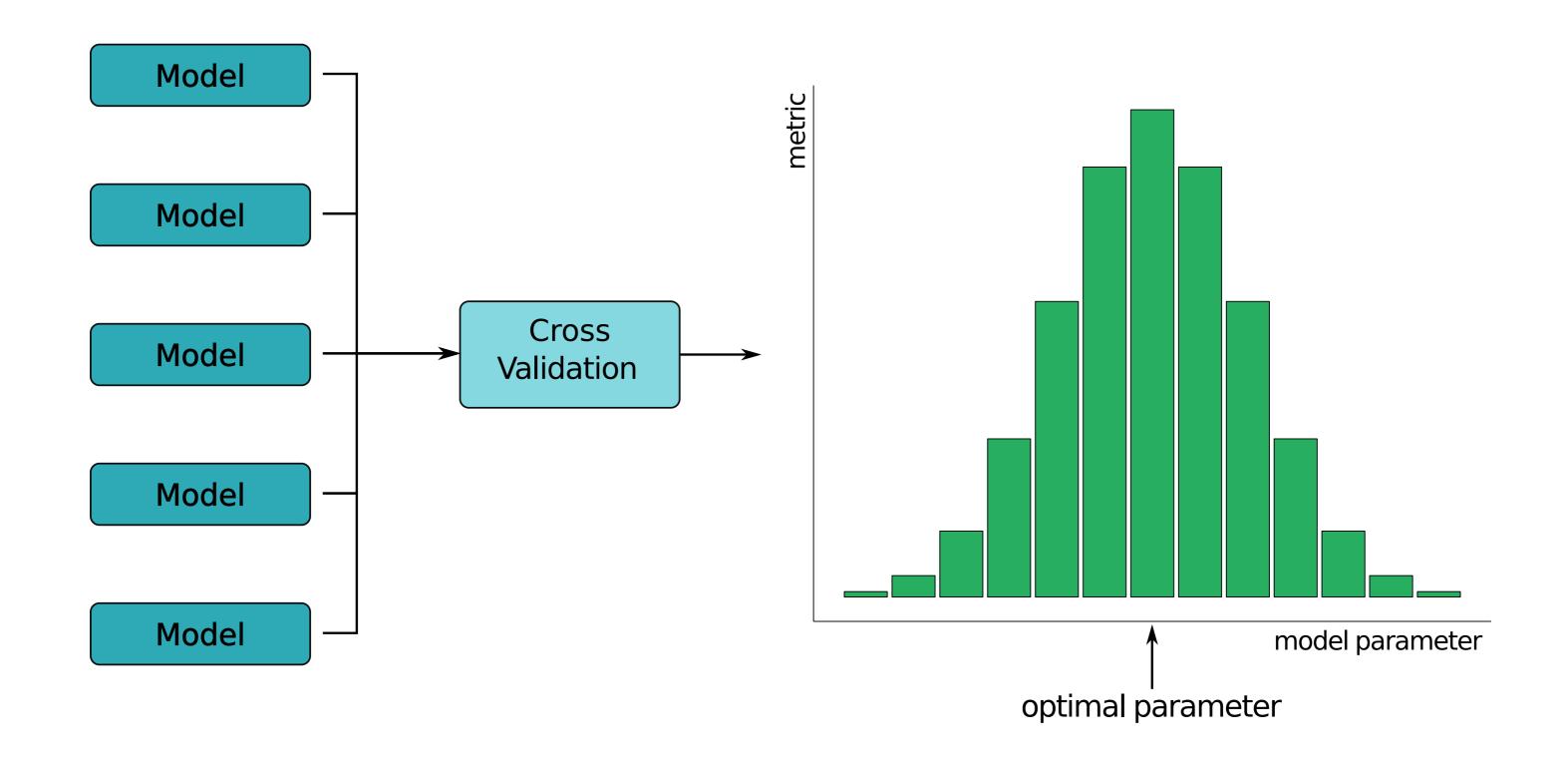
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Cars revisited (again)

```
cars.select('mass', 'cyl', 'consumption').show(5)
```

```
| mass|cyl|consumption|

+-----+

|1451.0| 6| 9.05|

|1129.0| 4| 6.53|

|1399.0| 4| 7.84|

|1147.0| 4| 7.84|

|1111.0| 4| 9.05|

+-----+
```

Fuel consumption with intercept

Linear regression with an intercept. Fit to training data.

```
regression = LinearRegression(labelCol='consumption', fitIntercept=True)
regression = regression.fit(cars_train)
```

Calculate the RMSE on the testing data.

```
evaluator.evaluate(regression.transform(cars_test))
```

```
# RMSE for model with an intercept 0.745974203928479
```



Fuel consumption without intercept

Linear regression without an intercept. Fit to training data.

```
regression = LinearRegression(labelCol='consumption', fitIntercept=False)
regression = regression.fit(cars_train)
```

Calculate the RMSE on the testing data.

```
# RMSE for model without an intercept (second model)
0.852819012439

# RMSE for model with an intercept (first model)
```

0.745974203928

Parameter grid

```
from pyspark.ml.tuning import ParamGridBuilder
# Create a parameter grid builder
params = ParamGridBuilder()
# Add grid points
params = params.addGrid(regression.fitIntercept, [True, False])
# Construct the grid
params = params.build()
# How many models?
print('Number of models to be tested: ', len(params))
```

```
Number of models to be tested: 2
```



Grid search with cross-validation

Create a cross-validator and fit to the training data.

What's the cross-validated RMSE for each model?

```
cv.avgMetrics
```

[0.800663722151, 0.907977823182]



The best model & parameters

```
# Access the best model cv.bestModel
```

Or just use the cross-validator object.

```
predictions = cv.transform(cars_test)
```

Retrieve the best parameter.

```
cv.bestModel.explainParam('fitIntercept')
```

```
'fitIntercept: whether to fit an intercept term (default: True, current: True)'
```



A more complicated grid

How many models now?

```
print ('Number of models to be tested: ', len(params))
```

```
Number of models to be tested: 50
```

Find the best parameters!

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Ensemble

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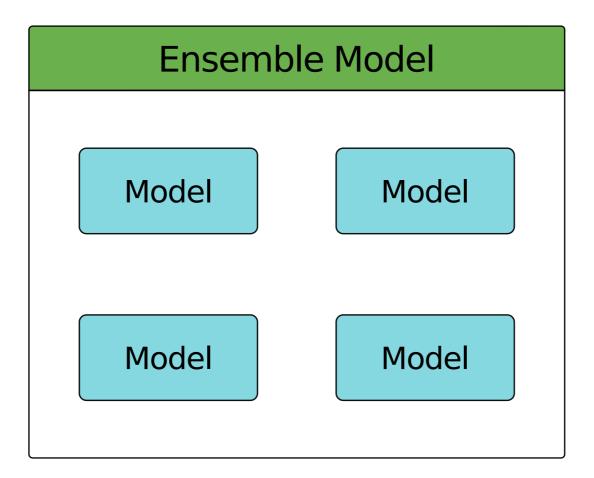
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What's an ensemble?

It's a collection of models.



Wisdom of the Crowd — collective opinion of a group better than that of a single expert.

Ensemble diversity

Diversity and **independence** are important because the best collective decisions are the product of disagreement and contest, not consensus or compromise.

? James Surowiecki, *The Wisdom of Crowds*

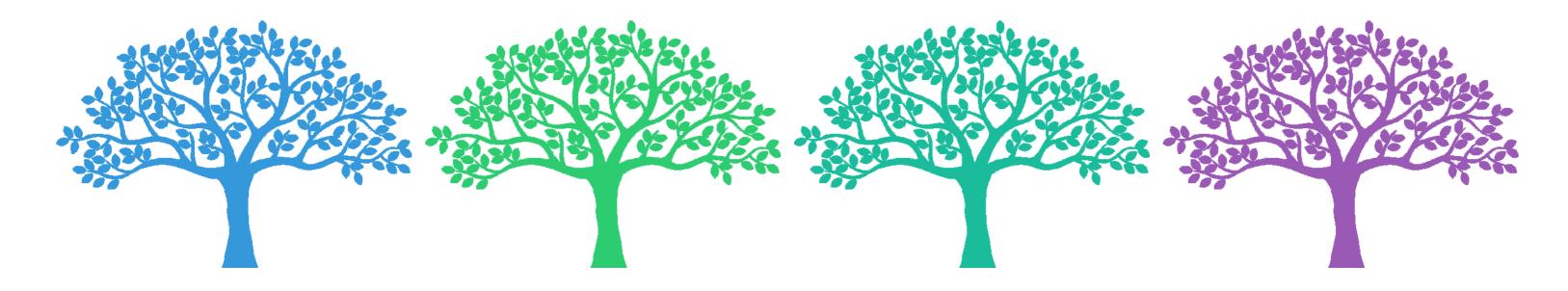
Random Forest

Random Forest — an ensemble of Decision Trees

Creating model diversity:

- each tree trained on random subset of data
- random subset of features used for splitting at each node

No two trees in the forest should be the same.



Create a forest of trees

Returning to cars data: manufactured in USA (0.0) or not (1.0).

Create Random Forest classifier.

```
from pyspark.ml.classification import RandomForestClassifier

forest = RandomForestClassifier(numTrees=5)
```

Fit to the training data.

```
forest = forest.fit(cars_train)
```

Seeing the trees

How to access trees within forest?

```
forest.trees
```

```
[DecisionTreeClassificationModel (uid=dtc_aa66702a4ce9) of depth 5 with 17 nodes,
DecisionTreeClassificationModel (uid=dtc_99f7efedafe9) of depth 5 with 31 nodes,
DecisionTreeClassificationModel (uid=dtc_9306e4a5fa1d) of depth 5 with 21 nodes,
DecisionTreeClassificationModel (uid=dtc_d643bd48a8dd) of depth 5 with 23 nodes,
DecisionTreeClassificationModel (uid=dtc_a2d5abd67969) of depth 5 with 27 nodes]
```

These can each be used to make individual predictions.

Predictions from individual trees

What predictions are generated by each tree?

```
|tree 0|tree 1|tree 2|tree 3|tree 4|label|
  0.0
        0.0
             0.0
                  0.0| 0.0| 0.0| <- perfect agreement
             0.0
                        0.0 0.0
  1.0
        1.0
                  1.0|
  0.0
        0.0
             0.0
                  1.0| 1.0| 1.0|
  0.0
        0.0
             0.0
                  1.0
                         0.0 0.0
  0.0
        1.0|
             1.0|
                  1.0|
                         0.0| 1.0|
       1.0|
             0.0 1.0 1.0 1.0
  1.0
        1.0|
             1.0 | 1.0 | 1.0 | <- perfect agreement
  1.0
```

Consensus predictions

Use the .transform() method to generate consensus predictions.

```
|label|probability
                                               |prediction|
     [[0.8,0.2]
                                               0.0
     [0.4, 0.6]
0.0
                                               11.0
     |[0.533333333333333, 0.4666666666666666] |0.0
     [0.71777777777778,0.2822222222222226] [0.0
     [0.39396825396825397, 0.606031746031746] [1.0
     [0.17660818713450294, 0.823391812865497] [1.0
     [0.053968253968253964, 0.946031746031746] [1.0
```

Feature importances

The model uses these features: cyl, size, mass, length, rpm and consumption.

Which of these is most or least important?

```
forest.featureImportances
```

```
SparseVector(6, {0: 0.0205, 1: 0.2701, 2: 0.108, 3: 0.1895, 4: 0.2939, 5: 0.1181})
```

Looks like:

- rpm is most important
- cyl is least important.

Gradient-Boosted Trees

Iterative boosting algorithm:

- 1. Build a Decision Tree and add to ensemble.
- 2. Predict label for each training instance using ensemble.
- 3. Compare predictions with known labels.
- 4. Emphasize training instances with incorrect predictions.
- 5. Return to 1.

Model improves on each iteration.

Boosting trees

Create a Gradient-Boosted Tree classifier.

```
from pyspark.ml.classification import GBTClassifier

gbt = GBTClassifier(maxIter=10)
```

Fit to the training data.

```
gbt = gbt.fit(cars_train)
```

Comparing trees

Let's compare the three types of tree models on testing data.

```
# AUC for Decision Tree
0.5875

# AUC for Random Forest
0.65

# AUC for Gradient-Boosted Tree
0.65
```

Both of the ensemble methods perform better than a plain Decision Tree.

Ensemble all of the models!

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Closing thoughts

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Things you've learned

- Load & prepare data
- Classifiers
 - Decision Tree
 - Logistic Regression
- Regression
 - Linear Regression
 - Penalized Regression
- Pipelines
- Cross-validation & grid search
- Ensembles







Learning more

Documentation at https://spark.apache.org/docs/latest/.



Congratulations!

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