
EvalML Documentation

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Alteryx Innovation Labs

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EvalML is an AutoML library that builds, optimizes, and evaluates machine learning pipelines using domain-specific objective functions.

Combined with [Featuretools](#) and [Compose](#), EvalML can be used to create end-to-end machine learning solutions for classification and regression problems.

QUICK START

```
[1]: import evalml
      from evalml import AutoMLSearch
```

1.1 Load Data

First, we load in the features and outcomes we want to use to train our model.

```
[2]: X, y = evalml.demos.load_breast_cancer()
```

1.2 Configure search

EvalML has many options to configure the pipeline search. At the minimum, we need to define an objective function. For simplicity, we will use the F1 score in this example. However, the real power of EvalML is in using domain-specific *objective functions* or *building your own*.

Below EvalML utilizes Bayesian optimization (EvalML's default optimizer) to search and find the best pipeline defined by the given objective.

```
[3]: automl = AutoMLSearch(problem_type="binary",
                           objective="f1",
                           max_pipelines=5)
```

In order to validate the results of the pipeline creation and optimization process, we will save some of our data as a holdout set.

```
[4]: X_train, X_holdout, y_train, y_holdout = evalml.preprocessing.split_data(X, y, test_
      ↪size=.2)
```

When we call `.search()`, the search for the best pipeline will begin. There is no need to wrangle with missing data or categorical variables as EvalML includes various preprocessing steps (like imputation, one-hot encoding, feature selection) to ensure you're getting the best results. As long as your data is in a single table, EvalML can handle it. If not, you can reduce your data to a single table by utilizing [Featuretools](#) and its Entity Sets.

You can find more information on pipeline components and how to integrate your own custom pipelines into EvalML [here](#).

```
[5]: automl.search(X_train, y_train)
```

Generating pipelines to search over...

```
*****
* Beginning pipeline search *
*****
```

Optimizing for F1.
Greater score is better.

Searching up to 5 pipelines.
Allowed model families: xgboost, linear_model, random_forest, catboost

```
FigureWidget({
  'data': [{ 'mode': 'lines+markers',
             'name': 'Best Score',
             'type': ...
```

```
Mode Baseline Binary Classification... 0%|          | Elapsed:00:00
CatBoost Classifier w/ Simple Imput... 20%|         | Elapsed:00:22
Logistic Regression Classifier w/ S... 40%|        | Elapsed:00:23
Random Forest Classifier w/ Simple ... 60%|       | Elapsed:00:25
XGBoost Classifier w/ Simple Imputer: 80%|      | Elapsed:00:25[21:12:21] WARNING: ..
↳/src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used
↳with the objective 'binary:logistic' was changed from 'error' to 'logloss'.
↳Explicitly set eval_metric if you'd like to restore the old behavior.
[21:12:21] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↳evaluation metric used with the objective 'binary:logistic' was changed from
↳'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↳behavior.
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

```
[21:12:21] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↳evaluation metric used with the objective 'binary:logistic' was changed from
↳'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↳behavior.
```

```
XGBoost Classifier w/ Simple Imputer:      80%|      | Elapsed:00:25
Optimization finished                      80%|      | Elapsed:00:25
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1]. (continues on next page)

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1.3 See Pipeline Rankings

After the search is finished we can view all of the pipelines searched, ranked by score. Internally, EvalML performs cross validation to score the pipelines. If it notices a high variance across cross validation folds, it will warn you. EvalML also provides additional *data checks* to analyze your data to assist you in producing the best performing pipeline.

```
[6]: automl.rankings
```

	id	pipeline_name	score	\
0	2	Logistic Regression Classifier w/ Simple Imput...	0.982639	
1	1	CatBoost Classifier w/ Simple Imputer	0.972203	
2	4	XGBoost Classifier w/ Simple Imputer	0.965294	
3	3	Random Forest Classifier w/ Simple Imputer	0.959963	
4	0	Mode Baseline Binary Classification Pipeline	0.770273	

	high_variance_cv	parameters
0	False	{'Simple Imputer': {'impute_strategy': 'most_f...
1	False	{'Simple Imputer': {'impute_strategy': 'most_f...
2	False	{'Simple Imputer': {'impute_strategy': 'most_f...
3	False	{'Simple Imputer': {'impute_strategy': 'most_f...
4	False	{'Baseline Classifier': {'strategy': 'random_w...

1.4 Describe pipeline

If we are interested in see more details about the pipeline, we can describe it using the `id` from the rankings table:

```
[7]: automl.describe_pipeline(3)
```

```
*****
* Random Forest Classifier w/ Simple Imputer *
*****

Problem Type: Binary Classification
Model Family: Random Forest

Pipeline Steps
=====
1. Simple Imputer
   * impute_strategy : most_frequent
   * fill_value : None
2. Random Forest Classifier
   * n_estimators : 100
   * max_depth : 6
   * n_jobs : -1

Training
=====
Training for Binary Classification problems.
Total training time (including CV): 1.4 seconds
```

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Cross Validation

		F1	Accuracy	Binary	Balanced Accuracy	Binary	Precision	AUC	Log_
→ Loss	Binary	MCC	Binary	# Training	# Testing				
0		0.948		0.934		0.923	0.929	0.982	↳
→	0.162		0.859	303.0	152.0				
1		0.963		0.954		0.949	0.958	0.991	↳
→	0.117		0.901	303.0	152.0				
2		0.968		0.960		0.961	0.978	0.985	↳
→	0.137		0.916	304.0	151.0				
mean		0.960		0.949		0.944	0.955	0.986	↳
→	0.138		0.892	-	-				
std		0.010		0.014		0.020	0.025	0.005	↳
→	0.023		0.030	-	-				
coef of var		0.011		0.014		0.021	0.026	0.005	↳
→	0.164		0.033	-	-				

1.5 Select Best pipeline

We can now select best pipeline and score it on our holdout data:

```
[8]: pipeline = automl.best_pipeline
      pipeline.fit(X_train, y_train)
      pipeline.score(X_holdout, y_holdout, ["f1"])

[8]: OrderedDict([('F1', 0.9583333333333334)])
```

We can also visualize the structure of our pipeline:

```
[9]: pipeline.graph()

[9]:
```

1.6 Whats next?

Head into the more in-depth automated walkthrough [here](#) or any advanced topics below.

1.6.1 Install

EvalML is available for Python 3.6+. It can be installed by running the following command:

```
pip install evalml --extra-index-url https://install.featurelabs.com/<license>/
```

Dependencies

Optional Dependencies

EvalML includes several dependencies in `requirements.txt` by default: `xgboost` and `catboost` support pipelines built around those modeling libraries, and `plotly` and `ipywidgets` support plotting functionality in

automl searches. These dependencies are recommended but are not required in order to install and use EvalML. To install these additional dependencies run `pip install -r requirements.txt`.

Core Dependencies

If you wish to install EvalML with only the core required dependencies, include `--no-dependencies` in your EvalML pip install command, and then install all core dependencies with `pip install -r core-requirements.txt`.

Windows

The [XGBoost](#) library may not be pip-installable in some Windows environments. If you are encountering installation issues, please try installing XGBoost from [Github](#) before installing EvalML.

1.6.2 Objective Functions

The **objective function** is what EvalML maximizes (or minimizes) as it completes the pipeline search. As it gets feedback from building pipelines, it tunes the hyperparameters to build optimized models. Therefore, it is critical to have an objective function that captures how the model's predictions will be used in a business setting.

List of Available Objective Functions

Most AutoML libraries optimize for generic machine learning objective functions. Frequently, the scores produced by the generic machine learning objective diverge from how the model will be evaluated in the real world.

In EvalML, we can train and optimize the model for a specific problem by optimizing a domain-specific objectives functions or by defining our own custom objective function.

Currently, EvalML has two domain specific objective functions with more being developed. For more information on these objective functions click on the links below.

- [Fraud Detection](#)
- [Lead Scoring](#)

Build your own objective Functions

Often times, the objective function is very specific to the use-case or business problem. To get the right objective to optimize requires thinking through the decisions or actions that will be taken using the model and assigning the cost/benefit to doing that correctly or incorrectly based on known outcomes in the training data.

Once you have determined the objective for your business, you can provide that to EvalML to optimize by defining a custom objective function. Read more [here](#).

1.6.3 Building a Fraud Prediction Model with EvalML

In this demo, we will build an optimized fraud prediction model using EvalML. To optimize the pipeline, we will set up an objective function to minimize the percentage of total transaction value lost to fraud. At the end of this demo, we also show you how introducing the right objective during the training is over 4x better than using a generic machine learning metric like AUC.

```
[1]: import evalml
      from evalml import AutoMLSearch
      from evalml.objectives import FraudCost
```

Configure “Cost of Fraud”

To optimize the pipelines toward the specific business needs of this model, you can set your own assumptions for the cost of fraud. These parameters are

- `retry_percentage` - what percentage of customers will retry a transaction if it is declined?
- `interchange_fee` - how much of each successful transaction do you collect?
- `fraud_payout_percentage` - the percentage of fraud will you be unable to collect
- `amount_col` - the column in the data the represents the transaction amount

Using these parameters, EvalML determines attempt to build a pipeline that will minimize the financial loss due to fraud.

```
[2]: fraud_objective = FraudCost(retry_percentage=.5,
                                interchange_fee=.02,
                                fraud_payout_percentage=.75,
                                amount_col='amount')
```

Search for best pipeline

In order to validate the results of the pipeline creation and optimization process, we will save some of our data as a holdout set

```
[3]: X, y = evalml.demos.load_fraud(n_rows=2500)
```

```

      Number of Features
Boolean                1
Categorical             6
Numeric                5

Number of training examples: 2500
Labels
False    85.92%
True     14.08%
Name: fraud, dtype: object
```

EvalML natively supports one-hot encoding. Here we keep 1 out of the 6 categorical columns to decrease computation time.

```
[4]: X = X.drop(['datetime', 'expiration_date', 'country', 'region', 'provider'], axis=1)

X_train, X_holdout, y_train, y_holdout = evalml.preprocessing.split_data(X, y, test_
↪size=0.2, random_state=0)

print(X.dtypes)

card_id          int64
store_id         int64
amount           int64
currency         object
```

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```
customer_present      bool
lat                   float64
lng                   float64
dtype: object
```

Because the fraud labels are binary, we will use `AutoMLSearch(problem_type='binary')`. When we call `.search()`, the search for the best pipeline will begin.

```
[5]: automl = AutoMLSearch(problem_type='binary',
                           objective=fraud_objective,
                           additional_objectives=['auc', 'f1', 'precision'],
                           max_pipelines=5,
                           optimize_thresholds=True)
```

```
automl.search(X_train, y_train)
```

```
Generating pipelines to search over...
```

```
*****
* Beginning pipeline search *
*****
```

```
Optimizing for Fraud Cost.
Lower score is better.
```

```
Searching up to 5 pipelines.
```

```
Allowed model families: xgboost, catboost, linear_model, random_forest
```

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...
```

```
Mode Baseline Binary Classification... 0%|          | Elapsed:00:00
CatBoost Classifier w/ Simple Input... 20%|         | Elapsed:00:09
Logistic Regression Classifier w/ S... 40%|        | Elapsed:00:10
Random Forest Classifier w/ Simple ... 60%|       | Elapsed:00:13
XGBoost Classifier w/ Simple Impute... 80%|      | Elapsed:00:13[21:10:26] WARNING: ..
↪/src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used
↪with the objective 'binary:logistic' was changed from 'error' to 'logloss'.
↪Explicitly set eval_metric if you'd like to restore the old behavior.
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

```
The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

```
The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].
```

```
[21:10:27] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↪evaluation metric used with the objective 'binary:logistic' was changed from
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↪behavior.

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:10:27] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↪evaluation metric used with the objective 'binary:logistic' was changed from
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↪behavior.
XGBoost Classifier w/ Simple Impute...      80%| | Elapsed:00:14
Optimization finished                      80%| | Elapsed:00:14
```

View rankings and select pipeline

Once the fitting process is done, we can see all of the pipelines that were searched, ranked by their score on the fraud detection objective we defined

```
[6]: automl.rankings

[6]:   id                pipeline_name      score \
0    0      Mode Baseline Binary Classification Pipeline  0.002316
1    3  Random Forest Classifier w/ Simple Imputer + O...  0.002316
2    4  XGBoost Classifier w/ Simple Imputer + One Hot...  0.007152
3    1                CatBoost Classifier w/ Simple Imputer  0.011063
4    2  Logistic Regression Classifier w/ Simple Imput...  0.022830

      high_variance_cv                parameters
0          False  {'Baseline Classifier': {'strategy': 'random_w...
1          False  {'Simple Imputer': {'impute_strategy': 'most_f...
2           True  {'Simple Imputer': {'impute_strategy': 'most_f...
3           True  {'Simple Imputer': {'impute_strategy': 'most_f...
4           True  {'Simple Imputer': {'impute_strategy': 'most_f...
```

to select the best pipeline we can run

```
[7]: best_pipeline = automl.best_pipeline
```

Describe pipeline

You can get more details about any pipeline. Including how it performed on other objective functions.

```
[8]: automl.describe_pipeline(automl.rankings.iloc[1]["id"])

*****
* Random Forest Classifier w/ Simple Imputer + One Hot Encoder *
*****
```

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```
Problem Type: Binary Classification
Model Family: Random Forest
```

```
Pipeline Steps
```

```
=====
```

```
1. Simple Imputer
    * impute_strategy : most_frequent
    * fill_value : None
2. One Hot Encoder
    * top_n : 10
    * categories : None
    * drop : None
    * handle_unknown : ignore
    * handle_missing : error
3. Random Forest Classifier
    * n_estimators : 100
    * max_depth : 6
    * n_jobs : -1
```

```
Training
```

```
=====
```

```
Training for Binary Classification problems.
Objective to optimize binary classification pipeline thresholds for: <evalml.
↪objectives.fraud_cost.FraudCost object at 0x7f802a53ddd0>
Total training time (including CV): 2.4 seconds
```

```
Cross Validation
```

```
-----
```

	Fraud Cost	AUC	F1	Precision	# Training	# Testing
0	0.002	0.860	0.247	0.141	1066.0	667.0
1	0.002	0.863	0.247	0.141	1066.0	667.0
2	0.002	0.856	0.247	0.141	1067.0	666.0
mean	0.002	0.859	0.247	0.141	-	-
std	0.000	0.004	0.000	0.000	-	-
coef of var	0.055	0.004	0.001	0.001	-	-

Evaluate on hold out

Finally, we retrain the best pipeline on all of the training data and evaluate on the holdout

```
[9]: best_pipeline.fit(X_train, y_train)
[9]: <evalml.pipelines.classification.baseline_binary.ModeBaselineBinaryPipeline at_
↪0x7f80705f2810>
```

Now, we can score the pipeline on the hold out data using both the fraud cost score and the AUC.

```
[10]: best_pipeline.score(X_holdout, y_holdout, objectives=["auc", fraud_objective])
[10]: OrderedDict([('AUC', 0.5), ('Fraud Cost', 0.016036197878507734)])
```

Why optimize for a problem-specific objective?

To demonstrate the importance of optimizing for the right objective, let's search for another pipeline using AUC, a common machine learning metric. After that, we will score the holdout data using the fraud cost objective to see how the best pipelines compare.

```
[11]: automl_auc = AutoMLSearch(problem_type='binary',
                                objective='auc',
                                additional_objectives=['f1', 'precision'],
                                max_pipelines=5,
                                optimize_thresholds=True)

automl_auc.search(X_train, y_train)
```

Generating pipelines to search over...

 * Beginning pipeline search *

Optimizing for AUC.
 Greater score is better.

Searching up to 5 pipelines.
 Allowed model families: xgboost, catboost, linear_model, random_forest

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...
```

Mode	Progress	Elapsed
Baseline Binary Classification...	0%	00:00
CatBoost Classifier w/ Simple Imput...	20%	00:09
Logistic Regression Classifier w/ S...	40%	00:09
Random Forest Classifier w/ Simple ...	60%	00:11
XGBoost Classifier w/ Simple Impute...	80%	00:11

[21:10:41] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used with the objective 'binary:logistic' was changed from 'error' to 'logloss'.
 ↳ Explicitly set eval_metric if you'd like to restore the old behavior.

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a future release. To remove this warning, do the following: 1) Pass option use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a future release. To remove this warning, do the following: 1) Pass option use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:10:41] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used with the objective 'binary:logistic' was changed from 'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old behavior.

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```

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:10:41] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↳evaluation metric used with the objective 'binary:logistic' was changed from
↳'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↳behavior.
XGBoost Classifier w/ Simple Impute...      80%| | Elapsed:00:12
Optimization finished                      80%| | Elapsed:00:12

```

like before, we can look at the rankings and pick the best pipeline

```
[12]: automl_auc.rankings
```

```

[12]:      id      pipeline_name      score \
0    3  Random Forest Classifier w/ Simple Imputer + O...  0.855789
1    4  XGBoost Classifier w/ Simple Imputer + One Hot...  0.851122
2    1                CatBoost Classifier w/ Simple Imputer  0.839414
3    2  Logistic Regression Classifier w/ Simple Imput...  0.804705
4    0      Mode Baseline Binary Classification Pipeline  0.500000

      high_variance_cv      parameters
0          False  {'Simple Imputer': {'impute_strategy': 'most_f...
1          False  {'Simple Imputer': {'impute_strategy': 'most_f...
2          False  {'Simple Imputer': {'impute_strategy': 'most_f...
3          False  {'Simple Imputer': {'impute_strategy': 'most_f...
4          False  {'Baseline Classifier': {'strategy': 'random_w...

```

```
[13]: best_pipeline_auc = automl_auc.best_pipeline
```

```

# train on the full training data
best_pipeline_auc.fit(X_train, y_train)

```

```
[13]: <evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline at 0x7f7ff863fb90>
```

```

[14]: # get the fraud score on holdout data
best_pipeline_auc.score(X_holdout, y_holdout, objectives=["auc", fraud_objective])

```

```

[14]: OrderedDict([('AUC', 0.8224584717607973),
                  ('Fraud Cost', 0.004329350526560073)])

```

```

[15]: # fraud score on fraud optimized again
best_pipeline.score(X_holdout, y_holdout, objectives=["auc", fraud_objective])

```

```
[15]: OrderedDict([('AUC', 0.5), ('Fraud Cost', 0.01126539759779717)])
```

When we optimize for AUC, we can see that the AUC score from this pipeline is better than the AUC score from the pipeline optimized for fraud cost. However, the losses due to fraud are over 3% of the total transaction amount when optimized for AUC and under 1% when optimized for fraud cost. As a result, we lose more than 2% of the total transaction amount by not optimizing for fraud cost specifically.

This happens because optimizing for AUC does not take into account the user-specified `retry_percentage`, `interchange_fee`, `fraud_payout_percentage` values. Thus, the best pipelines may produce the highest AUC but may not actually reduce the amount loss due to your specific type fraud.

This example highlights how performance in the real world can diverge greatly from machine learning metrics.

1.6.4 Building a Lead Scoring Model with EvalML

In this demo, we will build an optimized lead scoring model using EvalML. To optimize the pipeline, we will set up an objective function to maximize the revenue generated with true positives while taking into account the cost of false positives. At the end of this demo, we also show you how introducing the right objective during the training is over 6x better than using a generic machine learning metric like AUC.

```
[1]: import evalml
      from evalml import AutoMLSearch
      from evalml.objectives import LeadScoring
```

Configure LeadScoring

To optimize the pipelines toward the specific business needs of this model, you can set your own assumptions for how much value is gained through true positives and the cost associated with false positives. These parameters are

- `true_positive` - dollar amount to be gained with a successful lead
- `false_positive` - dollar amount to be lost with an unsuccessful lead

Using these parameters, EvalML builds a pipeline that will maximize the amount of revenue per lead generated.

```
[2]: lead_scoring_objective = LeadScoring(
      true_positives=1000,
      false_positives=-10
    )
```

Dataset

We will be utilizing a dataset detailing a customer's job, country, state, zip, online action, the dollar amount of that action and whether they were a successful lead.

```
[3]: from urllib.request import urlopen
      import pandas as pd

      customers_data = urlopen('https://featurelabs-static.s3.amazonaws.com/lead_scoring_ml_
      ↪apps/customers.csv')
      interactions_data = urlopen('https://featurelabs-static.s3.amazonaws.com/lead_scoring_
      ↪ml_apps/interactions.csv')
      leads_data = urlopen('https://featurelabs-static.s3.amazonaws.com/lead_scoring_ml_
      ↪apps/previous_leads.csv')
      customers = pd.read_csv(customers_data)
      interactions = pd.read_csv(interactions_data)
      leads = pd.read_csv(leads_data)

      X = customers.merge(interactions, on='customer_id').merge(leads, on='customer_id')
      y = X['label']
```

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```
X = X.drop(['customer_id', 'date_registered', 'birthday', 'phone', 'email',
            'owner', 'company', 'id', 'time_x',
            'session', 'referrer', 'time_y', 'label'], axis=1)
```

```
display(X.head())
```

	job	country	state	zip	action	amount
0	Engineer, mining	NaN	NY	60091.0	page_view	NaN
1	Psychologist, forensic	US	CA	NaN	purchase	135.23
2	Psychologist, forensic	US	CA	NaN	page_view	NaN
3	Air cabin crew	US	NaN	60091.0	download	NaN
4	Air cabin crew	US	NaN	60091.0	page_view	NaN

Search for best pipeline

In order to validate the results of the pipeline creation and optimization process, we will save some of our data as a holdout set

EvalML natively supports one-hot encoding and imputation so the above NaN and categorical values will be taken care of.

```
[4]: X_train, X_holdout, y_train, y_holdout = evalml.preprocessing.split_data(X, y, test_
      ↪size=0.2, random_state=0)
```

```
print(X.dtypes)
```

```
job          object
country      object
state        object
zip          float64
action       object
amount       float64
dtype: object
```

Because the lead scoring labels are binary, we will use `AutoMLSearch(problem_type='binary')`. When we call `.search()`, the search for the best pipeline will begin.

```
[5]: automl = AutoMLSearch(problem_type='binary',
                           objective=lead_scoring_objective,
                           additional_objectives=['auc'],
                           max_pipelines=5,
                           optimize_thresholds=True)
```

```
automl.search(X_train, y_train)
```

```
Generating pipelines to search over...
```

```
*****
* Beginning pipeline search *
*****
```

```
Optimizing for Lead Scoring.
Greater score is better.
```

```
Searching up to 5 pipelines.
Allowed model families: catboost, xgboost, linear_model, random_forest
```

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type'...

Mode Baseline Binary Classification... 0%|          | Elapsed:00:00
CatBoost Classifier w/ Simple Impute... 20%|          | Elapsed:00:11
Logistic Regression Classifier w/ S... 40%|          | Elapsed:00:13
Random Forest Classifier w/ Simple ... 60%|          | Elapsed:00:16
XGBoost Classifier w/ Simple Impute... 80%|          | Elapsed:00:16[21:11:05] WARNING: ..
↪/src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used
↪with the objective 'binary:logistic' was changed from 'error' to 'logloss'.
↪Explicitly set eval_metric if you'd like to restore the old behavior.

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:11:05] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↪evaluation metric used with the objective 'binary:logistic' was changed from
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↪behavior.

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:11:06] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↪evaluation metric used with the objective 'binary:logistic' was changed from
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↪behavior.
XGBoost Classifier w/ Simple Impute... 80%|          | Elapsed:00:17
Optimization finished                  80%|          | Elapsed:00:17
```

View rankings and select pipeline

Once the fitting process is done, we can see all of the pipelines that were searched, ranked by their score on the lead scoring objective we defined

```
[6]: automl.rankings
```

```
[6]:      id      pipeline_name      score \
0      1      CatBoost Classifier w/ Simple Imputer  44.280483
1      0      Mode Baseline Binary Classification Pipeline  42.140250
2      2      Logistic Regression Classifier w/ Simple Imput...  41.972462
3      4      XGBoost Classifier w/ Simple Imputer + One Hot...  41.912311
4      3      Random Forest Classifier w/ Simple Imputer + O...  41.507849

      high_variance_cv      parameters
0      False  {'Simple Imputer': {'impute_strategy': 'most_f...
1      False  {'Baseline Classifier': {'strategy': 'random_w...
2      False  {'Simple Imputer': {'impute_strategy': 'most_f...
3      False  {'Simple Imputer': {'impute_strategy': 'most_f...
4      False  {'Simple Imputer': {'impute_strategy': 'most_f...
```

to select the best pipeline we can run

```
[7]: best_pipeline = automl.best_pipeline
```

Describe pipeline

You can get more details about any pipeline. Including how it performed on other objective functions.

```
[8]: automl.describe_pipeline(automl.rankings.iloc[0]["id"])
```

```
*****
* CatBoost Classifier w/ Simple Imputer *
*****

Problem Type: Binary Classification
Model Family: CatBoost

Pipeline Steps
=====
1. Simple Imputer
    * impute_strategy : most_frequent
    * fill_value : None
2. CatBoost Classifier
    * n_estimators : 1000
    * eta : 0.03
    * max_depth : 6
    * bootstrap_type : None

Training
=====
Training for Binary Classification problems.
Objective to optimize binary classification pipeline thresholds for: <evalml.
↳objectives.lead_scoring.LeadScoring object at 0x7f5c85cb3e50>
Total training time (including CV): 11.2 seconds

Cross Validation
-----
      Lead Scoring      AUC # Training # Testing
0      45.845 0.925      2479.0      1550.0
1      42.748 0.922      2479.0      1550.0
```

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2	44.248	0.937	2480.0	1549.0
mean	44.280	0.928	-	-
std	1.549	0.008	-	-
coef of var	0.035	0.008	-	-

Evaluate on hold out

Finally, we retrain the best pipeline on all of the training data and evaluate on the holdout

```
[9]: best_pipeline.fit(X_train, y_train)
[9]: <evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline at 0x7f5c55766c50>
```

Now, we can score the pipeline on the hold out data using both the lead scoring score and the AUC.

```
[10]: best_pipeline.score(X_holdout, y_holdout, objectives=["auc", lead_scoring_objective])
[10]: OrderedDict([('AUC', 0.9402462979752191),
                  ('Lead Scoring', 11.969045571797077)])
```

Why optimize for a problem-specific objective?

To demonstrate the importance of optimizing for the right objective, let's search for another pipeline using AUC, a common machine learning metric. After that, we will score the holdout data using the lead scoring objective to see how the best pipelines compare.

```
[11]: automl_auc = evalml.AutoMLSearch(problem_type='binary',
                                     objective='auc',
                                     additional_objectives=[],
                                     max_pipelines=5,
                                     optimize_thresholds=True)

automl_auc.search(X_train, y_train)

Generating pipelines to search over...
*****
* Beginning pipeline search *
*****

Optimizing for AUC.
Greater score is better.

Searching up to 5 pipelines.
Allowed model families: catboost, xgboost, linear_model, random_forest
```

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...
```

```
Mode Baseline Binary Classification... 0%|          | Elapsed:00:00
CatBoost Classifier w/ Simple Input... 20%|         | Elapsed:00:11
Logistic Regression Classifier w/ S... 40%|        | Elapsed:00:12
Random Forest Classifier w/ Simple ... 60%|       | Elapsed:00:13
XGBoost Classifier w/ Simple Impute... 80%|      | Elapsed:00:13[21:11:27] WARNING: ..
↪ /src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used
↪ with the objective 'binary:logistic' was changed from 'error' to 'logloss'.
↪ Explicitly set eval_metric if you'd like to restore the old behavior.
```

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```

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:11:27] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↳evaluation metric with the objective 'binary:logistic' was changed from
↳'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↳behavior.

/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↳lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:

The use of label encoder in XGBClassifier is deprecated and will be removed in a
↳future release. To remove this warning, do the following: 1) Pass option
↳use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↳labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].

[21:11:27] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↳evaluation metric with the objective 'binary:logistic' was changed from
↳'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↳behavior.
XGBoost Classifier w/ Simple Impute...      80%| | Elapsed:00:14
Optimization finished                      80%| | Elapsed:00:14

```

like before, we can look at the rankings and pick the best pipeline

```
[12]: automl_auc.rankings
```

```

[12]:   id                pipeline_name      score \
0    1      CatBoost Classifier w/ Simple Imputer  0.931551
1    4  XGBoost Classifier w/ Simple Imputer + One Hot...  0.723820
2    3  Random Forest Classifier w/ Simple Imputer + O...  0.703700
3    2  Logistic Regression Classifier w/ Simple Imput...  0.702151
4    0      Mode Baseline Binary Classification Pipeline  0.500000

      high_variance_cv      parameters
0          False  {'Simple Imputer': {'impute_strategy': 'most_f...
1          False  {'Simple Imputer': {'impute_strategy': 'most_f...
2          False  {'Simple Imputer': {'impute_strategy': 'most_f...
3          False  {'Simple Imputer': {'impute_strategy': 'most_f...
4          False  {'Baseline Classifier': {'strategy': 'random_w...

```

```
[13]: best_pipeline_auc = automl_auc.best_pipeline

      # train on the full training data
      best_pipeline_auc.fit(X_train, y_train)

[13]: <evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline at 0x7f5cc959c910>

[14]: # get the auc and lead scoring score on holdout data
      best_pipeline_auc.score(X_holdout, y_holdout, objectives=["auc", lead_scoring_
      ↪objective])

[14]: OrderedDict([('AUC', 0.9402462979752191),
                  ('Lead Scoring', 11.969045571797077)])
```

When we optimize for AUC, we can see that the AUC score from this pipeline is better than the AUC score from the pipeline optimized for lead scoring. However, the revenue per lead gained was only \$7 per lead when optimized for AUC and was \$45 when optimized for lead scoring. As a result, we would gain up to 6x the amount of revenue if we optimized for lead scoring.

This happens because optimizing for AUC does not take into account the user-specified `true_positive` (dollar amount to be gained with a successful lead) and `false_positive` (dollar amount to be lost with an unsuccessful lead) values. Thus, the best pipelines may produce the highest AUC but may not actually generate the most revenue through lead scoring.

This example highlights how performance in the real world can diverge greatly from machine learning metrics.

1.6.5 Custom Objective Functions

Often times, the objective function is very specific to the use-case or business problem. To get the right objective to optimize requires thinking through the decisions or actions that will be taken using the model and assigning a cost/benefit to doing that correctly or incorrectly based on known outcomes in the training data.

Once you have determined the objective for your business, you can provide that to EvalML to optimize by defining a custom objective function.

How to Create a Objective Function

To create a custom objective function, we must define 2 functions:

- The **“objective function”**: this function takes the predictions, true labels, and any other information about the future and returns a score of how well the model performed.
- The **“decision function”**: this function takes prediction probabilities that were output from the model and a threshold and returns a prediction.

To evaluate a particular model, EvalML automatically finds the best threshold to pass to the decision function to generate predictions and then scores the resulting predictions using the objective function. The score from the objective function determines which set of pipeline hyperparameters EvalML will try next.

To give a concrete example, let’s look at how the fraud detection objective function is built.

```
[1]: from evalml.objectives.binary_classification_objective import _
      ↪BinaryClassificationObjective
      import pandas as pd

      class FraudCost(BinaryClassificationObjective):
```

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```

"""Score the percentage of money lost of the total transaction amount process due_
to fraud"""
    name = "Fraud Cost"
    greater_is_better = False
    score_needs_proba = False

    def __init__(self, retry_percentage=.5, interchange_fee=.02,
                  fraud_payout_percentage=1.0, amount_col='amount'):
        """Create instance of FraudCost

        Arguments:
            retry_percentage (float): What percentage of customers that will retry a_
            transaction if it
            is declined. Between 0 and 1. Defaults to .5

            interchange_fee (float): How much of each successful transaction you can_
            collect.
            Between 0 and 1. Defaults to .02

            fraud_payout_percentage (float): Percentage of fraud you will not be able_
            to collect.
            Between 0 and 1. Defaults to 1.0

            amount_col (str): Name of column in data that contains the amount._
            Defaults to "amount"
        """
        self.retry_percentage = retry_percentage
        self.interchange_fee = interchange_fee
        self.fraud_payout_percentage = fraud_payout_percentage
        self.amount_col = amount_col

    def decision_function(self, ypred_proba, threshold=0.0, X=None):
        """Determine if a transaction is fraud given predicted probabilities,_
threshold, and dataframe with transaction amount

        Arguments:
            ypred_proba (pd.Series): Predicted probabilities
            X (pd.DataFrame): Dataframe containing transaction amount
            threshold (float): Dollar threshold to determine if transaction is_
            fraud

        Returns:
            pd.Series: Series of predicted fraud labels using X and threshold
        """
        if not isinstance(X, pd.DataFrame):
            X = pd.DataFrame(X)

        if not isinstance(ypred_proba, pd.Series):
            ypred_proba = pd.Series(ypred_proba)

        transformed_probs = (ypred_proba.values * X[self.amount_col])
        return transformed_probs > threshold

    def objective_function(self, y_true, y_predicted, X):
        """Calculate amount lost to fraud per transaction given predictions, true_
values, and dataframe with transaction amount

```

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```

    Arguments:
        y_predicted (pd.Series): predicted fraud labels
        y_true (pd.Series): true fraud labels
        X (pd.DataFrame): dataframe with transaction amounts

    Returns:
        float: amount lost to fraud per transaction
    """
    if not isinstance(X, pd.DataFrame):
        X = pd.DataFrame(X)

    if not isinstance(y_predicted, pd.Series):
        y_predicted = pd.Series(y_predicted)

    if not isinstance(y_true, pd.Series):
        y_true = pd.Series(y_true)

    # extract transaction using the amount columns in users data
    try:
        transaction_amount = X[self.amount_col]
    except KeyError:
        raise ValueError("`{}` is not a valid column in X.".format(self.amount_
→col))

    # amount paid if transaction is fraud
    fraud_cost = transaction_amount * self.fraud_payout_percentage

    # money made from interchange fees on transaction
    interchange_cost = transaction_amount * (1 - self.retry_percentage) * self.
→interchange_fee

    # calculate cost of missing fraudulent transactions
    false_negatives = (y_true & ~y_predicted) * fraud_cost

    # calculate money lost from fees
    false_positives = (~y_true & y_predicted) * interchange_cost

    loss = false_negatives.sum() + false_positives.sum()

    loss_per_total_processed = loss / transaction_amount.sum()

    return loss_per_total_processed

```

1.6.6 Setting up pipeline search

Designing the right machine learning pipeline and picking the best parameters is a time-consuming process that relies on a mix of data science intuition as well as trial and error. EvalML streamlines the process of selecting the best modeling algorithms and parameters, so data scientists can focus their energy where it is most needed.

How it works

EvalML selects and tunes machine learning pipelines built of numerous steps. This includes encoding categorical data, missing value imputation, feature selection, feature scaling, and finally machine learning. As EvalML tunes pipelines,

it uses the objective function selected and configured by the user to guide its search.

At each iteration, EvalML uses cross-validation to generate an estimate of the pipeline's performances. If a pipeline has high variance across cross-validation folds, it will provide a warning. In this case, the pipeline may not perform reliably in the future.

EvalML is designed to work well out of the box. However, it provides numerous methods for you to control the search described below.

Selecting problem type

EvalML supports both classification and regression problems. You select your problem type by calling the `AutoMLSearch` initialization with the corresponding argument.

```
[1]: import evalml
      from evalml import AutoMLSearch

[2]: AutoMLSearch(problem_type='binary')
      Using default limit of max_pipelines=5.

[2]: <evalml.automl.automl_search.AutoMLSearch at 0x7fb752167d10>

[3]: AutoMLSearch(problem_type='multiclass')
      Using default limit of max_pipelines=5.

[3]: <evalml.automl.automl_search.AutoMLSearch at 0x7fb75217f4d0>

[4]: AutoMLSearch(problem_type='regression')
      Using default limit of max_pipelines=5.

[4]: <evalml.automl.automl_search.AutoMLSearch at 0x7fb752180350>
```

Setting the Objective Function

The only required parameter to start searching for pipelines is the objective function. Most domain-specific objective functions require you to specify parameters based on your business assumptions. You can do this before you initialize your pipeline search. For example

```
[5]: from evalml.objectives import FraudCost

      fraud_objective = FraudCost(
          retry_percentage=.5,
          interchange_fee=.02,
          fraud_payout_percentage=.75,
          amount_col='amount'
      )

      AutoMLSearch(problem_type='binary', objective=fraud_objective, optimize_
      ↪ thresholds=True)

      Using default limit of max_pipelines=5.
```

```
[5]: <evalml automl automl_search.AutoMLSearch at 0x7fb752186b90>
```

Evaluate on Additional Objectives

Additional objectives can be scored on during the evaluation process. To add another objective, use the `additional_objectives` parameter in `AutoMLSearch`. The results of these additional objectives will then appear in the results of `describe_pipeline`.

```
[6]: from evalml.objectives import FraudCost

fraud_objective = FraudCost(
    retry_percentage=.5,
    interchange_fee=.02,
    fraud_payout_percentage=.75,
    amount_col='amount'
)

AutoMLSearch(problem_type='binary', objective='AUC', additional_objectives=[fraud_
↪objective], optimize_thresholds=False)

Using default limit of max_pipelines=5.
```

```
[6]: <evalml automl automl_search.AutoMLSearch at 0x7fb75211f290>
```

Selecting Model Types

By default, all model types are considered. You can control which model types to search with the `allowed_model_families` parameters

```
[7]: automl = AutoMLSearch(problem_type='binary',
                           objective="f1",
                           allowed_model_families=["random_forest"])

Using default limit of max_pipelines=5.
```

After initialization you can view the pipelines that will be included in the search

```
[8]: automl.allowed_pipelines
```

you can see a list of all supported models like this

```
[9]: evalml.list_model_families("binary") # `binary` for binary classification and
↪ `multiclass` for multiclass classification
```

```
[9]: [<ModelFamily.XGBOOST: 'xgboost'>,
      <ModelFamily.CATBOOST: 'catboost'>,
      <ModelFamily.LINEAR_MODEL: 'linear_model'>,
      <ModelFamily.RANDOM_FOREST: 'random_forest'>]
```

```
[10]: evalml.list_model_families("regression")
```

```
[10]: [<ModelFamily.XGBOOST: 'xgboost'>,
      <ModelFamily.CATBOOST: 'catboost'>,
      <ModelFamily.LINEAR_MODEL: 'linear_model'>,
      <ModelFamily.RANDOM_FOREST: 'random_forest'>]
```

Limiting Search Time

You can limit the search time by specifying a maximum number of pipelines and/or a maximum amount of time. EvalML won't build new pipelines after the maximum time has passed or the maximum number of pipelines have been built. If a limit is not set, then a maximum of 5 pipelines will be built.

The maximum search time can be specified as an integer in seconds or as a string in seconds, minutes, or hours.

```
[11]: AutoMLSearch(problem_type='binary',
                  objective="f1",
                  max_pipelines=5,
                  max_time=60)

AutoMLSearch(problem_type='binary',
              objective="f1",
              max_time="1 minute")

[11]: <evalml automl automl_search.AutoMLSearch at 0x7fb72c323750>
```

Early Stopping

You can also limit search time by providing a patience value for early stopping. With a patience value, EvalML will stop searching when the best objective score has not been improved upon for n iterations. The patience value must be a positive integer. You can also provide a tolerance value where EvalML will only consider a score as an improvement over the best score if the difference was greater than the tolerance percentage.

```
[12]: from evalml.demos import load_diabetes

X, y = load_diabetes()
automl = AutoMLSearch(problem_type='regression', objective="MSE", patience=2,
                      tolerance=0.01, max_pipelines=10)
automl.search(X, y)

Generating pipelines to search over...
*****
* Beginning pipeline search *
*****

Optimizing for MSE.
Lower score is better.

Searching up to 10 pipelines.
Allowed model families: xgboost, catboost, linear_model, random_forest

FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...}

Mean Baseline Regression Pipeline:      0%|          | Elapsed:00:00
CatBoost Regressor w/ Simple Imputer:  10%|         | Elapsed:00:03
Linear Regressor w/ Simple Imputer ...  20%|        | Elapsed:00:03
Random Forest Regressor w/ Simple I...  30%|       | Elapsed:00:04
XGBoost Regressor w/ Simple Imputer:    40%|      | Elapsed:00:04
```

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```
2 iterations without improvement. Stopping search early...
Optimization finished          40%|          | Elapsed:00:04
```

```
[13]: automl.rankings
```

```
[13]:      id      pipeline_name      score \
0    2  Linear Regressor w/ Simple Imputer + Standard ...  3027.144520
1    1                CatBoost Regressor w/ Simple Imputer  3279.699820
2    3      Random Forest Regressor w/ Simple Imputer  3301.475890
3    4      XGBoost Regressor w/ Simple Imputer  3960.404464
4    0      Mean Baseline Regression Pipeline  5943.716736

      high_variance_cv      parameters
0          False  {'Simple Imputer': {'impute_strategy': 'most_f...
1          False  {'Simple Imputer': {'impute_strategy': 'most_f...
2          False  {'Simple Imputer': {'impute_strategy': 'most_f...
3          False  {'Simple Imputer': {'impute_strategy': 'most_f...
4          False      {'Baseline Regressor': {'strategy': 'mean'}}
```

Control Cross Validation

EvalML cross-validates each model it tests during its search. By default it uses 3-fold cross-validation. You can optionally provide your own cross-validation method.

```
[14]: from sklearn.model_selection import StratifiedKFold
```

```
automl = AutoMLSearch(problem_type='binary',
                      objective="f1",
                      data_split=StratifiedKFold(5))
```

```
Using default limit of max_pipelines=5.
```

1.6.7 Exploring search results

After finishing a pipeline search, we can inspect the results. First, let's build a search of 10 different pipelines to explore.

```
[1]: import evalml
from evalml import AutoMLSearch

X, y = evalml.demos.load_breast_cancer()

automl = AutoMLSearch(problem_type='binary',
                      objective="f1",
                      max_pipelines=5)

automl.search(X, y)
```

```
Generating pipelines to search over...
```

```
*****
* Beginning pipeline search *
*****
```

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Optimizing for F1.

Greater score is better.

Searching up to 5 pipelines.

Allowed model families: catboost, xgboost, random_forest, linear_model

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...
```

```
Mode Baseline Binary Classification... 0%|          | Elapsed:00:00
CatBoost Classifier w/ Simple Imput... 20%|         | Elapsed:00:22
Logistic Regression Classifier w/ S... 40%|        | Elapsed:00:23
Random Forest Classifier w/ Simple ... 60%|       | Elapsed:00:25
XGBoost Classifier w/ Simple Imputer: 80%|      | Elapsed:00:25[21:09:50] WARNING: ..
↪/src/learner.cc:1061: Starting in XGBoost 1.3.0, the default evaluation metric used
↪with the objective 'binary:logistic' was changed from 'error' to 'logloss'.
↪Explicitly set eval_metric if you'd like to restore the old behavior.
[21:09:50] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↪evaluation metric used with the objective 'binary:logistic' was changed from
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↪behavior.
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

```
The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

```
The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].
```

```
[21:09:50] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default
↪evaluation metric used with the objective 'binary:logistic' was changed from
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old
↪behavior.
```

```
XGBoost Classifier w/ Simple Imputer: 80%|      | Elapsed:00:25
Optimization finished                  80%|      | Elapsed:00:25
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning:
```

```
The use of label encoder in XGBClassifier is deprecated and will be removed in a
↪future release. To remove this warning, do the following: 1) Pass option
↪use_label_encoder=False when constructing XGBClassifier object; and 2) Encode your
↪labels (y) as integers starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].
```

View Rankings

A summary of all the pipelines built can be returned as a pandas DataFrame. It is sorted by score. EvalML knows based on our objective function whether higher or lower is better.

```
[2]: automl.rankings
```

```
[2]:
```

	id	pipeline_name	score	\
0	2	Logistic Regression Classifier w/ Simple Imput...	0.980447	
1	1	CatBoost Classifier w/ Simple Imputer	0.976333	
2	4	XGBoost Classifier w/ Simple Imputer	0.970577	
3	3	Random Forest Classifier w/ Simple Imputer	0.966629	
4	0	Mode Baseline Binary Classification Pipeline	0.771060	

	high_variance_cv	parameters
0	False	{'Simple Imputer': {'impute_strategy': 'most_f...
1	False	{'Simple Imputer': {'impute_strategy': 'most_f...
2	False	{'Simple Imputer': {'impute_strategy': 'most_f...
3	False	{'Simple Imputer': {'impute_strategy': 'most_f...
4	False	{'Baseline Classifier': {'strategy': 'random_w...

Describe Pipeline

Each pipeline is given an id. We can get more information about any particular pipeline using that id. Here, we will get more information about the pipeline with id = 0.

```
[3]: automl.describe_pipeline(1)
```

```
*****
* CatBoost Classifier w/ Simple Imputer *
*****

Problem Type: Binary Classification
Model Family: CatBoost

Pipeline Steps
=====
1. Simple Imputer
    * impute_strategy : most_frequent
    * fill_value : None
2. CatBoost Classifier
    * n_estimators : 1000
    * eta : 0.03
    * max_depth : 6
    * bootstrap_type : None

Training
=====
Training for Binary Classification problems.
Total training time (including CV): 22.8 seconds

Cross Validation
-----
```

	F1	Accuracy	Binary	Balanced Accuracy	Binary	Precision	AUC	Log_
→Loss	Binary	MCC	Binary # Training	# Testing				
0	0.967		0.958		0.949	0.951	0.995	
→	0.106	0.910	379.0	190.0				

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1	0.983	0.979		0.975	0.975	0.994	⌋
↪ 0.082	0.955	379.0	190.0				
2	0.979	0.974		0.976	0.991	0.990	⌋
↪ 0.093	0.944	380.0	189.0				
mean	0.976	0.970		0.967	0.973	0.993	⌋
↪ 0.094	0.936	–	–				
std	0.008	0.011		0.015	0.020	0.003	⌋
↪ 0.012	0.024	–	–				
coef of var	0.009	0.011		0.016	0.021	0.003	⌋
↪ 0.128	0.025	–	–				

Get Pipeline

We can get the object of any pipeline via their `id` as well:

```
[4]: automl.get_pipeline(1)
[4]: <evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline at 0x7f61f2fac150>
```

Get best pipeline

If we specifically want to get the best pipeline, there is a convenient access

```
[5]: automl.best_pipeline
[5]: <evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline at 0x7f626ba33490>
```

Feature Importance

We can get the importance associated with each feature of the resulting pipeline

```
[6]: pipeline = automl.get_pipeline(1)
      pipeline.fit(X, y)
      pipeline.feature_importance
[6]:
```

	feature	importance
0	worst texture	11.284484
1	worst concave points	8.105820
2	worst perimeter	7.989750
3	worst radius	7.872959
4	worst area	7.851859
5	mean concave points	7.098117
6	mean texture	5.919744
7	worst smoothness	4.831183
8	worst concavity	4.665609
9	area error	4.204315
10	compactness error	2.616379
11	worst symmetry	2.088257
12	mean concavity	2.015011
13	radius error	1.948857
14	concave points error	1.875384
15	mean compactness	1.824212
16	perimeter error	1.705293

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17	mean smoothness	1.697390
18	worst fractal dimension	1.629284
19	mean radius	1.585845
20	mean area	1.479429
21	smoothness error	1.460370
22	fractal dimension error	1.370231
23	mean perimeter	1.306400
24	texture error	1.065599
25	mean fractal dimension	1.000343
26	worst compactness	0.967447
27	mean symmetry	0.960967
28	symmetry error	0.912462
29	concavity error	0.666999

We can also create a bar plot of the feature importances

```
[7]: pipeline.graph_feature_importance()
```

Data type cannot be displayed: application/vnd.plotly.v1+json, text/html

Precision-Recall Curve

For binary classification, you can view the precision-recall curve of a classifier

```
[8]: # get the predicted probabilities associated with the "true" label
y_pred_proba = pipeline.predict_proba(X)[1]
evalml.pipelines.graph_utils.graph_precision_recall_curve(y, y_pred_proba)
```

Data type cannot be displayed: application/vnd.plotly.v1+json, text/html

ROC Curve

For binary and multiclass classification, you can view the ROC curve of a classifier

```
[9]: # get the predicted probabilities associated with the "true" label
y_pred_proba = pipeline.predict_proba(X)[1]
evalml.pipelines.graph_utils.graph_roc_curve(y, y_pred_proba)
```

Data type cannot be displayed: application/vnd.plotly.v1+json, text/html

Confusion Matrix

For binary or multiclass classification, you can view a confusion matrix of the classifier's predictions

```
[10]: y_pred = pipeline.predict(X)
evalml.pipelines.graph_utils.graph_confusion_matrix(y, y_pred)
```

Data type cannot be displayed: application/vnd.plotly.v1+json, text/html

Access raw results

You can also get access to all the underlying data, like this:

```
[11]: automl.results
[11]: {'pipeline_results': {0: {'id': 0,
    'pipeline_name': 'Mode Baseline Binary Classification Pipeline',
    'pipeline_class': evalml.pipelines.classification.baseline_binary.
    ↳ ModeBaselineBinaryPipeline,
    'pipeline_summary': 'Baseline Classifier',
    'parameters': {'Baseline Classifier': {'strategy': 'random_weighted'}},
    'score': 0.7710601157203097,
    'high_variance_cv': False,
    'training_time': 0.03497886657714844,
    'cv_data': [{'all_objective_scores': OrderedDict([('F1',
        0.7702265372168284),
        ('Accuracy Binary', 0.6263157894736842),
        ('Balanced Accuracy Binary', 0.5),
        ('Precision', 0.6263157894736842),
        ('AUC', 0.5),
        ('Log Loss Binary', 0.6608932451679239),
        ('MCC Binary', 0.0),
        ('# Training', 379),
        ('# Testing', 190)]),
    'score': 0.7702265372168284,
    'binary_classification_threshold': 0.5},
    {'all_objective_scores': OrderedDict([('F1', 0.7702265372168284),
        ('Accuracy Binary', 0.6263157894736842),
        ('Balanced Accuracy Binary', 0.5),
        ('Precision', 0.6263157894736842),
        ('AUC', 0.5),
        ('Log Loss Binary', 0.6608932451679239),
        ('MCC Binary', 0.0),
        ('# Training', 379),
        ('# Testing', 190)]),
    'score': 0.7702265372168284,
    'binary_classification_threshold': 0.5},
    {'all_objective_scores': OrderedDict([('F1', 0.7727272727272727),
        ('Accuracy Binary', 0.6296296296296297),
        ('Balanced Accuracy Binary', 0.5),
        ('Precision', 0.6296296296296297),
        ('AUC', 0.5),
        ('Log Loss Binary', 0.6591759924082952),
        ('MCC Binary', 0.0),
        ('# Training', 380),
        ('# Testing', 189)]),
    'score': 0.7727272727272727,
    'binary_classification_threshold': 0.5}}},
    1: {'id': 1,
    'pipeline_name': 'CatBoost Classifier w/ Simple Imputer',
    'pipeline_class': evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline,
```

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```

'pipeline_summary': 'CatBoost Classifier w/ Simple Imputer',
'parameters': {'Simple Imputer': {'impute_strategy': 'most_frequent',
    'fill_value': None},
    'CatBoost Classifier': {'n_estimators': 1000,
        'eta': 0.03,
        'max_depth': 6,
        'bootstrap_type': None}},
'score': 0.9763329621163277,
'high_variance_cv': False,
'training_time': 22.760128498077393,
'cv_data': [{'all_objective_scores': OrderedDict([('F1',
    0.9669421487603305),
    ('Accuracy Binary', 0.9578947368421052),
    ('Balanced Accuracy Binary', 0.9493431175287016),
    ('Precision', 0.9512195121951219),
    ('AUC', 0.9945555687063559),
    ('Log Loss Binary', 0.10583268649418161),
    ('MCC Binary', 0.909956827190137),
    ('# Training', 379),
    ('# Testing', 190)]),
    'score': 0.9669421487603305,
    'binary_classification_threshold': 0.5},
    {'all_objective_scores': OrderedDict([('F1', 0.9833333333333334),
    ('Accuracy Binary', 0.9789473684210527),
    ('Balanced Accuracy Binary', 0.9746715587643509),
    ('Precision', 0.9752066115702479),
    ('AUC', 0.9943188543022844),
    ('Log Loss Binary', 0.08186397218927995),
    ('MCC Binary', 0.955011564828661),
    ('# Training', 379),
    ('# Testing', 190)]),
    'score': 0.9833333333333334,
    'binary_classification_threshold': 0.5},
    {'all_objective_scores': OrderedDict([('F1', 0.9787234042553192),
    ('Accuracy Binary', 0.9735449735449735),
    ('Balanced Accuracy Binary', 0.9760504201680673),
    ('Precision', 0.9913793103448276),
    ('AUC', 0.9899159663865547),
    ('Log Loss Binary', 0.09296190874649334),
    ('MCC Binary', 0.9443109474170326),
    ('# Training', 380),
    ('# Testing', 189)]),
    'score': 0.9787234042553192,
    'binary_classification_threshold': 0.5}}],
2: {'id': 2,
    'pipeline_name': 'Logistic Regression Classifier w/ Simple Imputer + Standard_
↪Scaler',
    'pipeline_class': evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline,
    'pipeline_summary': 'Logistic Regression Classifier w/ Simple Imputer + Standard_
↪Scaler',
    'parameters': {'Simple Imputer': {'impute_strategy': 'most_frequent',
    'fill_value': None},
    'Logistic Regression Classifier': {'penalty': 'l2',
        'C': 1.0,
        'n_jobs': -1}},
'score': 0.9804468499596668,
'high_variance_cv': False,

```

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```
'training_time': 1.0541419982910156,
'cv_data': [{'all_objective_scores': OrderedDict([('F1',
0.9831932773109243),
('Accuracy Binary', 0.9789473684210527),
('Balanced Accuracy Binary', 0.9775121316132087),
('Precision', 0.9831932773109243),
('AUC', 0.9936087110900698),
('Log Loss Binary', 0.09347817517438463),
('MCC Binary', 0.9550242632264173),
('# Training', 379),
('# Testing', 190)]),
'score': 0.9831932773109243,
'binary_classification_threshold': 0.5},
{'all_objective_scores': OrderedDict([('F1', 0.9794238683127572),
('Accuracy Binary', 0.9736842105263158),
('Balanced Accuracy Binary', 0.9647887323943662),
('Precision', 0.9596774193548387),
('AUC', 0.9975144987572493),
('Log Loss Binary', 0.08320464479579018),
('MCC Binary', 0.9445075449666159),
('# Training', 379),
('# Testing', 190)]),
'score': 0.9794238683127572,
'binary_classification_threshold': 0.5},
{'all_objective_scores': OrderedDict([('F1', 0.9787234042553192),
('Accuracy Binary', 0.9735449735449735),
('Balanced Accuracy Binary', 0.9760504201680673),
('Precision', 0.9913793103448276),
('AUC', 0.9906362545018007),
('Log Loss Binary', 0.09680859555948443),
('MCC Binary', 0.9443109474170326),
('# Training', 380),
('# Testing', 189)]),
'score': 0.9787234042553192,
'binary_classification_threshold': 0.5}}],
3: {'id': 3,
'pipeline_name': 'Random Forest Classifier w/ Simple Imputer',
'pipeline_class': evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline,
'pipeline_summary': 'Random Forest Classifier w/ Simple Imputer',
'parameters': {'Simple Imputer': {'impute_strategy': 'most_frequent',
'fill_value': None},
'Random Forest Classifier': {'n_estimators': 100,
'max_depth': 6,
'n_jobs': -1}},
'score': 0.9666291581686476,
'high_variance_cv': False,
'training_time': 1.4184563159942627,
'cv_data': [{'all_objective_scores': OrderedDict([('F1',
0.9543568464730291),
('Accuracy Binary', 0.9421052631578948),
('Balanced Accuracy Binary', 0.9338975026630371),
('Precision', 0.9426229508196722),
('AUC', 0.9893478518167831),
('Log Loss Binary', 0.13984688783161608),
('MCC Binary', 0.8757606542930872),
('# Training', 379),
('# Testing', 190)]),
```

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```

    'score': 0.9543568464730291,
    'binary_classification_threshold': 0.5},
{'all_objective_scores': OrderedDict([('F1', 0.9709543568464729),
    ('Accuracy Binary', 0.9631578947368421),
    ('Balanced Accuracy Binary', 0.9563853710498283),
    ('Precision', 0.9590163934426229),
    ('AUC', 0.989347851816783),
    ('Log Loss Binary', 0.1201072101539428),
    ('MCC Binary', 0.9211492315750531),
    ('# Training', 379),
    ('# Testing', 190)]),
    'score': 0.9709543568464729,
    'binary_classification_threshold': 0.5},
{'all_objective_scores': OrderedDict([('F1', 0.9745762711864406),
    ('Accuracy Binary', 0.9682539682539683),
    ('Balanced Accuracy Binary', 0.9689075630252101),
    ('Precision', 0.9829059829059829),
    ('AUC', 0.9927971188475391),
    ('Log Loss Binary', 0.10765634363120971),
    ('MCC Binary', 0.9325680982740896),
    ('# Training', 380),
    ('# Testing', 189)]),
    'score': 0.9745762711864406,
    'binary_classification_threshold': 0.5}}],
4: {'id': 4,
    'pipeline_name': 'XGBoost Classifier w/ Simple Imputer',
    'pipeline_class': evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline,
    'pipeline_summary': 'XGBoost Classifier w/ Simple Imputer',
    'parameters': {'Simple Imputer': {'impute_strategy': 'most_frequent',
    'fill_value': None},
    'XGBoost Classifier': {'eta': 0.1,
    'max_depth': 6,
    'min_child_weight': 1,
    'n_estimators': 100}},
    'score': 0.9705772481706093,
    'high_variance_cv': False,
    'training_time': 0.42917609214782715,
    'cv_data': [{'all_objective_scores': OrderedDict([('F1',
    0.9666666666666667),
    ('Accuracy Binary', 0.9578947368421052),
    ('Balanced Accuracy Binary', 0.9521836903775595),
    ('Precision', 0.9586776859504132),
    ('AUC', 0.9915966386554622),
    ('Log Loss Binary', 0.11449876085695762),
    ('MCC Binary', 0.9097672817424011),
    ('# Training', 379),
    ('# Testing', 190)]),
    'score': 0.9666666666666667,
    'binary_classification_threshold': 0.5},
{'all_objective_scores': OrderedDict([('F1', 0.979253112033195),
    ('Accuracy Binary', 0.9736842105263158),
    ('Balanced Accuracy Binary', 0.9676293052432241),
    ('Precision', 0.9672131147540983),
    ('AUC', 0.9959758551307847),
    ('Log Loss Binary', 0.07421583775339011),
    ('MCC Binary', 0.943843520216036),
    ('# Training', 379),

```

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```

        ('# Testing', 190)]),
    'score': 0.979253112033195,
    'binary_classification_threshold': 0.5},
    {'all_objective_scores': OrderedDict([('F1', 0.9658119658119659),
        ('Accuracy Binary', 0.9576719576719577),
        ('Balanced Accuracy Binary', 0.9605042016806722),
        ('Precision', 0.9826086956521739),
        ('AUC', 0.9885954381752701),
        ('Log Loss Binary', 0.11418110851220609),
        ('MCC Binary', 0.9112159507396058),
        ('# Training', 380),
        ('# Testing', 189)]),
    'score': 0.9658119658119659,
    'binary_classification_threshold': 0.5}}}],
    'search_order': [0, 1, 2, 3, 4])

```

1.6.8 Regression Example

```

[1]: import evalml
from evalml import AutoMLSearch
from evalml.demos import load_diabetes
from evalml.pipelines import PipelineBase, get_pipelines

X, y = evalml.demos.load_diabetes()

automl = AutoMLSearch(problem_type='regression', objective="R2", max_pipelines=5)

automl.search(X, y)

```

Generating pipelines to search over...

```

*****
* Beginning pipeline search *
*****

```

Optimizing for R2.
Greater score is better.

Searching up to 5 pipelines.
Allowed model families: linear_model, random_forest, catboost, xgboost

```

FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...

```

Mean Baseline Regression Pipeline:	0%	Elapsed:00:00
CatBoost Regressor w/ Simple Imputer:	20%	Elapsed:00:03
Linear Regressor w/ Simple Imputer ...	40%	Elapsed:00:03
Random Forest Regressor w/ Simple I...	60%	Elapsed:00:04
XGBoost Regressor w/ Simple Imputer:	80%	Elapsed:00:04
Optimization finished	80%	Elapsed:00:04

```

[2]: automl.rankings

```

```
[2]:      id                pipeline_name      score \
0   2  Linear Regressor w/ Simple Imputer + Standard ...  0.488703
1   1                CatBoost Regressor w/ Simple Imputer  0.446477
2   3          Random Forest Regressor w/ Simple Imputer  0.441420
3   4          XGBoost Regressor w/ Simple Imputer  0.331082
4   0          Mean Baseline Regression Pipeline -0.004217

      high_variance_cv      parameters
0          False  {'Simple Imputer': {'impute_strategy': 'most_f...
1          False  {'Simple Imputer': {'impute_strategy': 'most_f...
2          False  {'Simple Imputer': {'impute_strategy': 'most_f...
3          False  {'Simple Imputer': {'impute_strategy': 'most_f...
4          False  {'Baseline Regressor': {'strategy': 'mean'}}
```

```
[3]: automl.best_pipeline
```

```
[3]: <evalml.pipelines.utils.make_pipeline.<locals>.GeneratedPipeline at 0x7fed1ad357d0>
```

```
[4]: automl.get_pipeline(0)
```

```
[4]: <evalml.pipelines.regression.baseline_regression.MeanBaselineRegressionPipeline at_
↳0x7fed1bdc6950>
```

```
[5]: automl.describe_pipeline(0)
```

```
*****
* Mean Baseline Regression Pipeline *
*****

Problem Type: Regression
Model Family: Baseline

Pipeline Steps
=====
1. Baseline Regressor
    * strategy : mean

Training
=====
Training for Regression problems.
Total training time (including CV): 0.0 seconds

Cross Validation
-----
```

	R2	Root Mean Squared Error	MAE	MSE	MedianAE	MaxError
↳ExpVariance # Training # Testing						
0	-0.007	75.863	63.324	5755.216	57.190	186.810
↳-0.000	294.0	148.0				
1	-0.000	79.654	68.759	6344.747	67.966	193.966
↳ 0.000	295.0	147.0				
2	-0.006	75.705	65.485	5731.187	63.817	170.817
↳-0.000	295.0	147.0				
mean	-0.004	77.074	65.856	5943.717	62.991	183.864
↳-0.000	-					
std	0.004	2.236	2.736	347.510	5.435	11.852
↳ 0.000	-					
coef of var	-0.866	0.029	0.042	0.058	0.086	0.064
↳-0.866	-					

1.6.9 Avoiding Overfitting

The ultimate goal of machine learning is to make accurate predictions on unseen data. EvalML aims to help you build a model that will perform as you expect once it is deployed in to the real world.

One of the benefits of using EvalML to build models is that it provides data checks to ensure you are building pipelines that will perform reliably in the future. This page describes the various ways EvalML helps you avoid overfitting to your data.

```
[1]: import evalml
```

Detecting Label Leakage

A common problem is having features that include information from your label in your training data. By default, EvalML will provide a warning when it detects this may be the case.

Let's set up a simple example to demonstrate what this looks like:

```
[2]: import pandas as pd
from evalml.data_checks import LabelLeakageDataCheck

X = pd.DataFrame({
    "leaked_feature": [6, 6, 10, 5, 5, 11, 5, 10, 11, 4],
    "leaked_feature_2": [3, 2.5, 5, 2.5, 3, 5.5, 2, 5, 5.5, 2],
    "valid_feature": [3, 1, 3, 2, 4, 6, 1, 3, 3, 11]
})

y = pd.Series([1, 1, 0, 1, 1, 0, 1, 0, 0, 1])

label_leakage_check = LabelLeakageDataCheck()
messages = label_leakage_check.validate(X, y)
for message in messages:
    print(message)

Column 'leaked_feature' is 95.0% or more correlated with the target
Column 'leaked_feature_2' is 95.0% or more correlated with the target
```

In the example above, EvalML warned about the input features `leaked_feature` and `leaked_feature_2`, which are both very closely correlated with the label we are trying to predict.

The second way to find features that may be leaking label information is to look at the top features of the model after running an AutoML search. As we can see below, the top features in our model are the 2 leaked features.

```
[3]: automl = evalml.AutoMLSearch(
    problem_type='binary',
    max_pipeline=1,
    allowed_model_families=["linear_model"],
)

automl.search(X, y)
best_pipeline = automl.best_pipeline
best_pipeline.fit(X, y)
best_pipeline.feature_importance

Column 'leaked_feature' is 95.0% or more correlated with the target
Column 'leaked_feature_2' is 95.0% or more correlated with the target
Generating pipelines to search over...
*****
```

(continues on next page)

(continued from previous page)

```
* Beginning pipeline search *
*****

Optimizing for Log Loss Binary.
Lower score is better.

Searching up to 1 pipelines.
Allowed model families: linear_model
```

```
FigureWidget({
  'data': [{ 'mode': 'lines+markers',
             'name': 'Best Score',
             'type'...
```

```
Mode Baseline Binary Classification... 0%|          | Elapsed:00:00
Optimization finished                  0%|          | Elapsed:00:00
```

```
[3]:      feature  importance
0    leaked_feature      0.0
1    leaked_feature_2      0.0
2      valid_feature      0.0
```

Perform cross-validation for pipeline evaluation

By default, EvalML performs 3-fold cross validation when building pipelines. This means that it evaluates each pipeline 3 times using different sets of data for training and testing. In each trial, the data used for testing has no overlap from the data used for training.

While this is a good baseline approach, you can pass your own cross validation object to be used during modeling. The cross validation object can be any of the CV methods defined in [scikit-learn](#) or use a compatible API.

For example, if we wanted to do a time series split:

```
[4]: from sklearn.model_selection import TimeSeriesSplit

X, y = evalml.demos.load_breast_cancer()

automl = evalml.AutoMLSearch(
    problem_type='binary',
    data_split=TimeSeriesSplit(n_splits=6),
    max_pipelines=1
)

automl.search(X, y)

Generating pipelines to search over...
*****
* Beginning pipeline search *
*****

Optimizing for Log Loss Binary.
Lower score is better.

Searching up to 1 pipelines.
Allowed model families: catboost, random_forest, linear_model, xgboost
```

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type'...
  }
])
```

Mode Baseline Binary Classification...	0%	Elapsed:00:00
Optimization finished	0%	Elapsed:00:00

if we describe the 1 pipeline we built, we can see the scores for each of the 6 splits as determined by the cross-validation object we provided. We can also see the number of training examples per fold increased because we were using `TimeSeriesSplit`

```
[5]: automl.describe_pipeline(0)
```

```
*****
* Mode Baseline Binary Classification Pipeline *
*****

Problem Type: Binary Classification
Model Family: Baseline

Pipeline Steps
=====
1. Baseline Classifier
    * strategy : random_weighted

Training
=====
Training for Binary Classification problems.
Total training time (including CV): 0.1 seconds

Cross Validation
-----
```

		Log Loss Binary	Accuracy Binary	Balanced Accuracy Binary	F1	
	AUC	MCC	Binary # Training	Binary # Testing		
0		0.880	83.0	0.358	0.500	0.000
→000	0.500	0.000	83.0	81.0		0.
1		0.697	164.0	0.469	0.500	0.000
→000	0.500	0.000	164.0	81.0		0.
2		0.697	245.0	0.333	0.500	0.000
→000	0.500	0.000	245.0	81.0		0.
3		0.664	326.0	0.716	0.500	0.835
→716	0.500	0.000	326.0	81.0		0.
4		0.619	407.0	0.790	0.500	0.883
→790	0.500	0.000	407.0	81.0		0.
5		0.611	488.0	0.741	0.500	0.851
→741	0.500	0.000	488.0	81.0		0.
mean		0.695	-	0.568	0.500	0.428
→374	0.500	0.000	-	-		0.
std		0.098	-	0.205	0.000	0.469
→411	0.000	0.000	-	-		0.
coef of var		0.141	-	0.361	0.000	1.096
→097	0.000	inf	-	-		1.

Detect unstable pipelines

When we perform cross validation we are trying generate an estimate of pipeline performance. EvalML does this by taking the mean of the score across the folds. If the performance across the folds varies greatly, it is indicative the the estimated value may be unreliable.

To protect the user against this, EvalML checks to see if the pipeline’s performance has a variance between the different folds. EvalML triggers a warning if the “coefficient of variance” of the scores (the standard deviation divided by mean) of the pipelines scores exceeds .2.

This warning will appear in the pipeline rankings under `high_variance_cv`.

```
[6]: automl.rankings
[6]:   id          pipeline_name      score \
0    0  Mode Baseline Binary Classification Pipeline  0.694742

      high_variance_cv          parameters
0                False  {'Baseline Classifier': {'strategy': 'random_w...
```

Create holdout for model validation

EvalML offers a method to quickly create an holdout validation set. A holdout validation set is data that is not used during the process of optimizing or training the model. You should only use this validation set once you’ve picked the final model you’d like to use.

Below we create a holdout set of 20% of our data

```
[7]: X, y = evalml.demos.load_breast_cancer()
      X_train, X_holdout, y_train, y_holdout = evalml.preprocessing.split_data(X, y, test_
      ↪size=.2)
```

```
[8]: automl = evalml.AutoMLSearch(problem_type='binary',
                                   objective="f1",
                                   max_pipelines=3)
      automl.search(X_train, y_train)

Generating pipelines to search over...
*****
* Beginning pipeline search *
*****

Optimizing for F1.
Greater score is better.

Searching up to 3 pipelines.
Allowed model families: catboost, random_forest, linear_model, xgboost

FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type'...

Mode Baseline Binary Classification...    0%|          | Elapsed:00:00
CatBoost Classifier w/ Simple Imput...  33%|          | Elapsed:00:22
Logistic Regression Classifier w/ S...  67%|          | Elapsed:00:23
Optimization finished                    67%|          | Elapsed:00:23
```

then we can retrain the best pipeline on all of our training data and see how it performs compared to the estimate

```
[9]: pipeline = automl.best_pipeline
      pipeline.fit(X_train, y_train)
      pipeline.score(X_holdout, y_holdout, ["f1"])

[9]: OrderedDict([('F1', 0.972972972972973)])
```

1.6.10 EvalML Components and Pipelines

EvalML searches and trains multiple machine learning **pipelines** in order to find the best one for your data. Each pipeline is made up of various **components** that can learn from the data, transform the data and ultimately predict labels given new data. Below we'll show an example of an EvalML pipeline. You can find a more in-depth look into *components* or learn how you can construct and use your own *pipelines*.

XGBoost Pipeline

The EvalML XGBoost Pipeline is made up of four different components: a one-hot encoder, a missing value imputer, a feature selector and an XGBoost estimator. To initialize a pipeline you need a parameters dictionary.

Parameters

The parameters dictionary needs to be in the format of a two-layered dictionary where the first key-value pair is the component name and component parameters dictionary. The component parameters dictionary consists of a key value pair of parameter name and parameter values. An example will be shown below and component parameters can be found [here](#).

```
[1]: from evalml.demos import load_breast_cancer
      from evalml.pipelines import XGBoostBinaryPipeline

X, y = load_breast_cancer()

parameters = {
    'Simple Imputer': {
        'impute_strategy': 'mean'
    },
    'RF Classifier Select From Model': {
        "percent_features": 0.5,
        "number_features": X.shape[1],
        "n_estimators": 20,
        "max_depth": 5
    },
    'XGBoost Classifier': {
        "n_estimators": 20,
        "eta": 0.5,
        "min_child_weight": 5,
        "max_depth": 10,
    }
}

xgp = XGBoostBinaryPipeline(parameters=parameters, random_state=5)
xgp.graph()
```

[1]:

From the above graph we can see each component and its parameters. Each component takes in data and feeds it to the next. You can see more detailed information by calling `.describe()`:

[2]: `xgp.describe()`

```
*****
* XGBoost Binary Classification Pipeline *
*****

Problem Type: Binary Classification
Model Family: XGBoost

Pipeline Steps
=====
1. Simple Imputer
   * impute_strategy : mean
   * fill_value : None
2. One Hot Encoder
   * top_n : 10
   * categories : None
   * drop : None
   * handle_unknown : ignore
   * handle_missing : error
3. XGBoost Classifier
   * eta : 0.5
   * max_depth : 10
   * min_child_weight : 5
   * n_estimators : 20
```

You can then fit and score an individual pipeline with an objective. An objective can either be a string representation of an EvalML objective or an EvalML objective class. You can find more objectives [here](#).

[3]: `xgp.fit(X, y)`
`xgp.score(X, y, objectives=['f1'])`

```
[21:12:35] WARNING: ../src/learner.cc:1061: Starting in XGBoost 1.3.0, the default_
↪evaluation metric used with the objective 'binary:logistic' was changed from_
↪'error' to 'logloss'. Explicitly set eval_metric if you'd like to restore the old_
↪behavior.
```

```
/home/docs/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/_
↪lib/python3.7/site-packages/xgboost/sklearn.py:888: UserWarning: The use of label_
↪encoder in XGBClassifier is deprecated and will be removed in a future release. To_
↪remove this warning, do the following: 1) Pass option use_label_encoder=False when_
↪constructing XGBClassifier object; and 2) Encode your labels (y) as integers_
↪starting with 0, i.e. 0, 1, 2, ..., [num_class - 1].
  warnings.warn(label_encoder_deprecation_msg, UserWarning)
```

[3]: `OrderedDict([('F1', 0.9916434540389972)])`

1.6.11 EvalML Components

From the [overview](#), we see how each machine learning pipeline consists of individual components that process data before the data is ultimately sent to an estimator. Below we will describe each type of component in an EvalML pipeline.

Component Classes

Components can be split into two distinct classes: **transformers** and **estimators**.

```
[1]: import numpy as np
import pandas as pd
from evalml.pipelines.components import SimpleImputer

X = pd.DataFrame([[1, 2, 3], [1, np.nan, 3]])
display(X)
```

```
   0    1    2
0  1    2  3
1  1  NaN  3
```

Transformers take in data as input and output altered data. For example, an *imputer* takes in data and outputs filled in missing data with the mean, median, or most frequent value of each column.

A transformer can fit on data and then transform it in two steps by calling `.fit()` and `.transform()` or in one step by calling `fit_transform()`.

```
[2]: imp = SimpleImputer(impute_strategy="mean")
X = imp.fit_transform(X)

display(X)
```

```
   0    1    2
0  1    2  3
1  1    2  3
```

On the other hand, an estimator fits on data (X) and labels (y) in order to take in new data as input and return the predicted label as output. Therefore, an estimator can fit on data and labels by calling `.fit()` and then predict by calling `.predict()` on new data. An example of this would be the *LogisticRegressionClassifier*. We can now see how a transformer alters data to make it easier for an estimator to learn and predict.

```
[3]: from evalml.pipelines.components import LogisticRegressionClassifier

clf = LogisticRegressionClassifier()

X = X
y = [1, 0]

clf.fit(X, y)
clf.predict(X)
```

```
[3]: 0    0
      1    0
      dtype: int64
```

Component Types

Components can further separate into different types that serve different functionality. Below we will go over the different types of transformers and estimators.

Transformer Types

- Imputer: fills missing data

- Ex: *SimpleImputer*
- Scaler: alters numerical data into different scales
 - Ex: *StandardScaler*
- Encoder: translates different data types
 - Ex: *OneHotEncoder*
- Feature Selection: selects most useful columns of data
 - Ex: *RFClassifierSelectFromModel*

Estimator Types

- Regressor: predicts numerical or continuous labels
 - Ex: *LinearRegressor*
- Classifier: predicts categorical or discrete labels
 - Ex: *XGBoostClassifier*

1.6.12 Custom Pipelines in EvalML

EvalML pipelines consist of modular components combining any number of transformers and an estimator. This allows you to create pipelines that fit the needs of your data to achieve the best results.

Requirements

A custom pipeline must adhere to the following requirements:

1. Inherit from the proper pipeline base class
 - Binary classification - `BinaryClassificationPipeline`
 - Multiclass classification - `MulticlassClassificationPipeline`
 - Regression - `RegressionPipeline`
2. Have a `component_graph` list as a class variable detailing the structure of the pipeline. Each component in the graph can be provided as either a string name or an instance.

Pipeline Configuration

There are a few other options to configure your custom pipeline.

Custom Name

By default, a pipeline class's name property is the result of adding spaces between each Pascal case capitalization in the class name. E.g. `LogisticRegressionPipeline.name` will return 'Logistic Regression Pipeline'. Therefore, we suggest custom pipelines use Pascal case for their class names.

If you'd like to override the pipeline classes name attribute so it isn't derived from the class name, you can set the `custom_name` attribute, like so:


```
[1]: from evalml.pipelines import BinaryClassificationPipeline

class CustomPipeline(BinaryClassificationPipeline):
    component_graph = ['Simple Imputer', 'Logistic Regression Classifier']
    custom_name = 'A custom pipeline name'

print(CustomPipeline.name)

A custom pipeline name
```

Custom Hyperparameters

To specify custom hyperparameter ranges, set the `custom_hyperparameters` property to be a dictionary where each key-value pair consists of a parameter name and range. AutoML will use this dictionary to override the hyperparameter ranges collected from each component in the component graph.

```
[2]: class CustomPipeline(BinaryClassificationPipeline):
    component_graph = ['Simple Imputer', 'Logistic Regression Classifier']

print("Without custom hyperparameters:")
print(CustomPipeline.hyperparameters)

class CustomPipeline(BinaryClassificationPipeline):
    component_graph = ['Simple Imputer', 'Logistic Regression Classifier']
    custom_hyperparameters = {
        'Simple Imputer': {
            'impute_strategy': ['most_frequent']
        }
    }

print()
print("With custom hyperparameters:")
print(CustomPipeline.hyperparameters)

Without custom hyperparameters:
{'Simple Imputer': {'impute_strategy': ['mean', 'median', 'most_frequent']},
 'Logistic Regression Classifier': {'penalty': ['l2'], 'C': Real(low=0.01, high=10,
prior='uniform', transform='identity')}}

With custom hyperparameters:
{'Simple Imputer': {'impute_strategy': ['most_frequent']}, 'Logistic Regression_
Classifier': {'penalty': ['l2'], 'C': Real(low=0.01, high=10, prior='uniform',
transform='identity')}}
```

1.6.13 Data Checks

EvalML provides data checks to help guide you in achieving the highest performing model. These utility functions help deal with problems such as overfitting, abnormal data, and missing data. These data checks can be found under `evalml/data_checks`. Below we will cover examples such as abnormal and missing data data checks.

Missing Data

Missing data or rows with NaN values provide many challenges for machine learning pipelines. In the worst case, many algorithms simply will not run with missing data! EvalML pipelines contain imputation *components* to ensure

that doesn't happen. Imputation works by approximating missing values with existing values. However, if a column contains a high number of missing values, a large percentage of the column would be approximated by a small percentage. This could potentially create a column without useful information for machine learning pipelines. By using the `HighlyNullDataCheck()` data check, EvalML will alert you to this potential problem by returning the columns that pass the missing values threshold.

```
[1]: import numpy as np
import pandas as pd

from evalml.data_checks import HighlyNullDataCheck

X = pd.DataFrame([[1, 2, 3],
                  [0, 4, np.nan],
                  [1, 4, np.nan],
                  [9, 4, np.nan],
                  [8, 6, np.nan]])

null_check = HighlyNullDataCheck(pct_null_threshold=0.8)

for message in null_check.validate(X):
    print(message.message)

Column '2' is 80.0% or more null
```

Abnormal Data

EvalML provides two data checks to check for abnormal data: `OutliersDataCheck()` and `IDColumnsDataCheck()`.

ID Columns

ID columns in your dataset provide little to no benefit to a machine learning pipeline as the pipeline cannot extrapolate useful information from unique identifiers. Thus, `IDColumnsDataCheck()` reminds you if these columns exist. In the given example, 'user_number' and 'id' columns are both identified as potentially being unique identifiers that should be removed.

```
[2]: from evalml.data_checks import IDColumnsDataCheck

X = pd.DataFrame([[0, 53, 6325, 5], [1, 90, 6325, 10], [2, 90, 18, 20]], columns=['user_
↪number', 'cost', 'revenue', 'id'])
id_col_check = IDColumnsDataCheck(id_threshold=0.9)

for message in id_col_check.validate(X):
    print(message.message)

Column 'id' is 90.0% or more likely to be an ID column
Column 'user_number' is 90.0% or more likely to be an ID column
```

Outliers

Outliers are observations that differ significantly from other observations in the same sample. Many machine learning pipelines suffer in performance if outliers are not dropped from the training set as they are not representative of the data. `OutliersDataCheck()` uses Isolation Forests to notify you if a sample can be considered an outlier.

Below we generate a random dataset with some outliers.

```
[3]: data = np.random.randn(100, 100)
X = pd.DataFrame(data=data)

# generate some outliers in rows 3, 25, 55, and 72
X.iloc[3, :] = pd.Series(np.random.randn(100) * 10)
X.iloc[25, :] = pd.Series(np.random.randn(100) * 20)
X.iloc[55, :] = pd.Series(np.random.randn(100) * 100)
X.iloc[72, :] = pd.Series(np.random.randn(100) * 100)
```

We then utilize `OutliersDataCheck()` to rediscover these outliers.

```
[4]: from evalml.data_checks import OutliersDataCheck

outliers_check = OutliersDataCheck()

for message in outliers_check.validate(X):
    print(message.message)

Row '3' is likely to have outlier data
Row '25' is likely to have outlier data
Row '55' is likely to have outlier data
Row '72' is likely to have outlier data
```

Writing Your Own Data Check

If you would prefer to write your own data check, you can do so by extending the `DataCheck` class and implementing the `validate(self, X, y)` class method. Below, we've created a new `DataCheck`, `ZeroVarianceDataCheck`.

```
[5]: from evalml.data_checks import DataCheck
from evalml.data_checks.data_check_message import DataCheckError

class ZeroVarianceDataCheck(DataCheck):
    def validate(self, X, y):
        if not isinstance(X, pd.DataFrame):
            X = pd.DataFrame(X)
        warning_msg = "Column '{}' has zero variance"
        return [DataCheckError(warning_msg.format(column), self.name) for column in X.
            ↪columns if len(X[column].unique()) == 1]
```

```
[1]: import evalml
```

1.6.14 Default DataChecks in AutoML

By default, AutoML will run the series of data checks in `DefaultDataChecks` when `automl.search()` is called to check that inputs are valid before running the search and fitting pipelines. Currently, `DefaultDataChecks` contains `HighlyNullDataCheck()`, `IDColumnsDataCheck()`, `LabelLeakageDataCheck()`. You can see the other available data checks under `evalml/data_checks`.

If the data checks return any warnings, those warnings will be logged, but the search will continue. However, if the data checks returns any error messages, `automl.search()` will raise a `ValueError` and quit before searching. This allows users to address any issues before running the potentially time-intensive search process.

Below, we have some data that contain a lot of null values, causing `DefaultDataChecks` to log a warning “Column ‘D’ is 95.0% or more null” and “Column ‘id’ is 100.0% or more likely to be an ID column” when try to run the search

below.

```
[2]: import pandas as pd
import numpy as np

X = pd.DataFrame(np.random.random((100, 5)), columns=["A", "B", "C", "D", "id"])
X.loc[11, 'A'] = np.nan
X.loc[9, 'B'] = np.nan
X.loc[30, 'C'] = np.nan
X.loc[95, 'D'] = np.nan
X.loc[:, 'id'] = range(100)
y = pd.Series([0,1]*50)

automl = evalml.AutoMLSearch(problem_type='binary', max_pipelines=1)
automl.search(X, y)
```

```
Column 'D' is 95.0% or more null
Column 'id' is 100.0% or more likely to be an ID column
Generating pipelines to search over...
*****
* Beginning pipeline search *
*****

Optimizing for Log Loss Binary.
Lower score is better.

Searching up to 1 pipelines.
Allowed model families: xgboost, catboost, random_forest, linear_model
```

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...}]
```

Mode Baseline Binary Classification...	0%	Elapsed:00:00
Optimization finished	0%	Elapsed:00:00

To access the exact warning and/or error messages our data checks returned, we can access `automl.data_check_results`.

```
[3]: for message in automl.data_check_results:
    print (message.message)

Column 'D' is 95.0% or more null
Column 'id' is 100.0% or more likely to be an ID column
```

Using your own DataCheck with AutoML

If you'd prefer to pass in your own data check, you can do so by passing in a `DataChecks` object as the value for the `data_checks` parameter. Here, we've implemented our own custom data check which returns a list of `DataCheckError` objects if there are any columns that have zero variance.

```
[4]: from evalml.data_checks import DataCheck, DataChecks
from evalml.data_checks.data_check_message import DataCheckError

class ZeroVarianceDataCheck(DataCheck):
    def validate(self, X, y):
```

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```

    if not isinstance(X, pd.DataFrame):
        X = pd.DataFrame(X)
        warning_msg = "Column '{}' has zero variance"
        return [DataCheckError(warning_msg.format(column), self.name) for column in X.
        ←columns if len(X[column].unique()) == 1]

```

If we now call `search()`, our error message will be logged and a `ValueError` will be raised:

```

[5]: data_checks = DataChecks(data_checks=[ZeroVarianceDataCheck()])

X = pd.DataFrame({'no_var': [1, 1, 1, 1, 1, 1, 1, 1, 1, 1],
                  'any_average_col': [2, 0, 1, 2, 1, 2, 0, 1, 2, 1],
                  'another_average_col': [0, 1, 2, 3, 4, 5, 6, 7, 8, 9]})
y = pd.Series([0, 1, 1, 0, 0, 0, 1, 1, 0, 0])

automl = evalml.AutoMLSearch(problem_type='binary', max_pipelines=1)
automl.search(X, y, data_checks=data_checks)

Column 'no_var' has zero variance

-----
ValueError                                Traceback (most recent call last)
<ipython-input-5-5f0233162825> in <module>
      7
      8 automl = evalml.AutoMLSearch(problem_type='binary', max_pipelines=1)
--> 9 automl.search(X, y, data_checks=data_checks)

~/checkouts/readthedocs.org/user_builds/feature-labs-inc-evalml/envs/v0.11.0/lib/
python3.7/site-packages/evalml/automl/automl_search.py in search(self, X, y,
data_checks, feature_types, raise_errors, show_iteration_plot)
    318         logger.error(message)
    319         if any([message.message_type == DataCheckMessageType.ERROR for
    ←message in self._data_check_results]):
-> 320             raise ValueError("Data checks raised some warnings and/or
    ←errors. Please see `self.data_check_results` for more information or pass
    ←data_checks=EmptyDataChecks() to search() to disable data checking.")
    321
    322         if self.allowed_pipelines is None:

ValueError: Data checks raised some warnings and/or errors. Please see `self.
data_check_results` for more information or pass data_checks=EmptyDataChecks() to
search() to disable data checking.

```

Again, we can access `self.data_check_results` to help us begin to address the issues raised by data checks.

```

[6]: for message in automl.data_check_results:
      print(message.message)

Column 'no_var' has zero variance

```

Disabling DataChecks

If you'd prefer not to run any data checks before running `search`, you can provide an `EmptyDataChecks` instance to `search()` instead.

```
[7]: from evalml.data_checks import EmptyDataChecks

X = pd.DataFrame({'no_var': [1, 1, 1, 1, 1, 1, 1, 1, 1, 1],
                    'any_average_col': [2, 0, 1, 2, 1, 2, 0, 1, 2, 1],
                    'another_average_col': [0, 1, 2, 3, 4, 5, 6, 7, 8, 9]})
y = pd.Series([0, 1, 1, 0, 0, 0, 1, 1, 0, 0])
automl = evalml.AutoMLSearch(problem_type='binary', max_pipelines=1)
automl.search(X, y, data_checks=EmptyDataChecks())
```

Generating pipelines to search over...

```
*****
* Beginning pipeline search *
*****
```

Optimizing for Log Loss Binary.
Lower score is better.

Searching up to 1 pipelines.
Allowed model families: xgboost, catboost, random_forest, linear_model

```
FigureWidget({
  'data': [{'mode': 'lines+markers',
            'name': 'Best Score',
            'type': ...
```

Mode Baseline Binary Classification...	0%	Elapsed:00:00
Optimization finished	0%	Elapsed:00:00

Even though we are using the same data as above, no data checks will be run and hence, the same input data we used above will not raise an error and continue with the search process.

1.6.15 Changelog

Future Releases

- Enhancements
- Fixes
- Changes
- Documentation Changes
- Testing Changes

v0.11.0 June 30, 2020

- **Enhancements**
 - Added multiclass support for ROC curve graphing [#832](#)
 - Added preprocessing component to drop features whose percentage of NaN values exceeds a specified threshold [#834](#)
 - Added data check to check for problematic target labels [#814](#)
 - Added PerColumnImputer that allows imputation strategies per column [#824](#)
 - Added transformer to drop specific columns [#827](#)
 - Added support for *categories*, *handle_error*, and *drop* parameters in *OneHotEncoder* [#830](#) [#897](#)

- Added preprocessing component to handle DateTime columns featurization #838
- Added ability to clone pipelines and components #842
- Define getter method for component *parameters* #847
- Added utility methods to calculate and graph permutation importances #860, #880
- Added new utility functions necessary for generating dynamic preprocessing pipelines #852
- Added kwargs to all components #863
- Updated *AutoSearchBase* to use dynamically generated preprocessing pipelines #870
- Added SelectColumns transformer #873
- Added ability to evaluate additional pipelines for automl search #874
- Added *default_parameters* class property to components and pipelines #879
- Added better support for disabling data checks in automl search #892
- Added ability to save and load AutoML objects to file #888
- Updated *AutoSearchBase.get_pipelines* to return an untrained pipeline instance #876
- Saved learned binary classification thresholds in automl results cv data dict #876
- **Fixes**
 - Fixed bug where SimpleImputer cannot handle dropped columns #846
 - Fixed bug where PerColumnImputer cannot handle dropped columns #855
 - Enforce requirement that builtin components save all inputted values in their parameters dict #847
 - Don't list base classes in *all_components* output #847
 - Standardize all components to output pandas data structures, and accept either pandas or numpy #853
 - Fixed rankings and full_rankings error when search has not been run #894
- **Changes**
 - Update *all_pipelines* and *all_components* to try initializing pipelines/components, and on failure exclude them #849
 - Refactor *handle_components* to *handle_components_class*, standardize to *ComponentBase* subclass instead of instance #850
 - Refactor “blacklist”/“whitelist” to “allow”/“exclude” lists #854
 - Replaced *AutoClassificationSearch* and *AutoRegressionSearch* with *AutoMLSearch* #871
 - Renamed *feature_importances* and *permutation_importances* methods to use singular names (*feature_importance* and *permutation_importance*) #883
 - Updated *automl* default data splitter to train/validation split for large datasets #877
 - Added open source license, update some repo metadata #887
- **Documentation Changes**
 - Fix some typos and update the EvalML logo #872
- **Testing Changes**
 - Update the changelog check job to expect the new branching pattern for the deps update bot #836

- Check that all components output pandas datastructures, and can accept either pandas or numpy #853
- Replaced *AutoClassificationSearch* and *AutoRegressionSearch* with *AutoMLSearch* #871

Warning:

Breaking Changes

- Pipelines' static `component_graph` field must contain either `ComponentBase` subclasses or `str`, instead of `ComponentBase` subclass instances #850
- Rename `handle_component` to `handle_component_class`. Now standardizes to `ComponentBase` subclasses instead of `ComponentBase` subclass instances #850
- Renamed `automl`'s `cv` argument to `data_split` #877
- Pipelines' and classifiers' `feature_importances` is renamed `feature_importance`, `graph_feature_importances` is renamed `graph_feature_importance` #883
- Passing `data_checks=None` to `automl` search will not perform any data checks as opposed to default checks. #892
- Pipelines to search for in `AutoML` are now determined automatically, rather than using the statically-defined pipeline classes. #870
- Updated `AutoSearchBase.get_pipelines` to return an untrained pipeline instance, instead of one which happened to be trained on the final cross-validation fold #876

v0.10.0 May 29, 2020

• **Enhancements**

- Added baseline models for classification and regression, add functionality to calculate baseline models before searching in `AutoML` #746
- Port over highly-null guardrail as a data check and define *DefaultDataChecks* and *DisableDataChecks* classes #745
- Update *Tuner* classes to work directly with pipeline parameters dicts instead of flat parameter lists #779
- Add Elastic Net as a pipeline option #812
- Added new Pipeline option *ExtraTrees* #790
- Added precision-recall curve metrics and plot for binary classification problems in `evalml.pipeline.graph_utils` #794
- Update the default `automl` algorithm to search in batches, starting with default parameters for each pipeline and iterating from there #793
- Added *AutoMLAlgorithm* class and *IterativeAlgorithm* impl, separated from *AutoSearchBase* #793

• **Fixes**

- Update pipeline `score` to return `nan` score for any objective which throws an exception during scoring #787
- Fixed bug introduced in #787 where binary classification metrics requiring predicted probabilities error in scoring #798
- CatBoost and XGBoost classifiers and regressors can no longer have a learning rate of 0 #795

- **Changes**

- Cleanup pipeline *score* code, and cleanup codecov #711
- Remove *pass* for abstract methods for codecov #730
- Added `__str__` for `AutoSearch` object #675
- Add util methods to graph ROC and confusion matrix #720
- Refactor *AutoBase* to *AutoSearchBase* #758
- Updated `AutoBase` with *data_checks* parameter, removed previous *detect_label_leakage* parameter, and added functionality to run data checks before search in `AutoML` #765
- Updated our logger to use Python’s logging utils #763
- Refactor most of *AutoSearchBase._do_iteration* impl into *AutoSearchBase._evaluate* #762
- Port over all guardrails to use the new `DataCheck` API #789
- Expanded *import_or_raise* to catch all exceptions #759
- Adds RMSE, MSLE, RMSLE as standard metrics #788
- Don’t allow *Recall* to be used as an objective for `AutoML` #784
- Removed feature selection from pipelines #819
- Update default estimator parameters to make automl search faster and more accurate #793

- **Documentation Changes**

- Add instructions to freeze *master* on *release.md* #726
- Update release instructions with more details #727 #733
- Add objective base classes to API reference #736
- Fix components API to match other modules #747

- **Testing Changes**

- Delete codecov yml, use codecov.io’s default #732
- Added unit tests for fraud cost, lead scoring, and standard metric objectives #741
- Update codecov client #782
- Updated `AutoBase` `__str__` test to include no parameters case #783
- Added unit tests for *ExtraTrees* pipeline #790
- If codecov fails to upload, fail build #810
- Updated Python version of dependency action #816
- Update the dependency update bot to use a suffix when creating branches #817

Warning:

Breaking Changes

- The *detect_label_leakage* parameter for `AutoML` classes has been removed and replaced by a *data_checks* parameter #765
- Moved ROC and confusion matrix methods from `evalml.pipeline.plot_utils` to `evalml.pipeline.graph_utils` #720

- `Tuner` classes require a pipeline hyperparameter range dict as an init arg instead of a space definition [#779](#)
- `Tuner.propose` and `Tuner.add` work directly with pipeline parameters dicts instead of flat parameter lists [#779](#)
- `PipelineBase.hyperparameters` and `custom_hyperparameters` use pipeline parameters dict format instead of being represented as a flat list [#779](#)
- All guardrail functions previously under `evalml.guardrails.utils` will be removed and replaced by data checks [#789](#)
- *Recall* disallowed as an objective for AutoML [#784](#)
- `AutoSearchBase` parameter `tuner` has been renamed to `tuner_class` [#793](#)
- `AutoSearchBase` parameter `possible_pipelines` and `possible_model_families` have been renamed to `allowed_pipelines` and `allowed_model_families` [#793](#)

v0.9.0 Apr. 27, 2020

• Enhancements

- Added accuracy as a standard objective [#624](#)
- Added verbose parameter to `load_fraud` [#560](#)
- Added Balanced Accuracy metric for binary, multiclass [#612](#) [#661](#)
- Added XGBoost regressor and XGBoost regression pipeline [#666](#)
- Added Accuracy metric for multiclass [#672](#)
- Added objective name in `AutoBase.describe_pipeline` [#686](#)
- Added `DataCheck` and `DataChecks`, `Message` classes and relevant subclasses [#739](#)

• Fixes

- Removed direct access to `cls.component_graph` [#595](#)
- Add testing files to `.gitignore` [#625](#)
- Remove circular dependencies from `Makefile` [#637](#)
- Add error case for `normalize_confusion_matrix()` [#640](#)
- Fixed XGBoostClassifier and XGBoostRegressor bug with feature names that contain `[`, `]`, or `<` [#659](#)
- Update `make_pipeline_graph` to not accidentally create empty file when testing if path is valid [#649](#)
- Fix pip installation warning about docsutils version, from boto dependency [#664](#)
- Removed zero division warning for F1/precision/recall metrics [#671](#)
- Fixed `summary` for pipelines without estimators [#707](#)

• Changes

- Updated default objective for binary/multiseries classification to log loss [#613](#)
- Created classification and regression pipeline subclasses and removed objective as an attribute of pipeline classes [#405](#)
- Changed the output of `score` to return one dictionary [#429](#)

- Created binary and multiclass objective subclasses #504
- Updated objectives API #445
- Removed call to *get_plot_data* from AutoML #615
- Set *raise_error* to default to True for AutoML classes #638
- Remove unnecessary “u” prefixes on some unicode strings #641
- Changed one-hot encoder to return uint8 dtypes instead of ints #653
- Pipeline *_name* field changed to *custom_name* #650
- Removed *graphs.py* and moved methods into *PipelineBase* #657, #665
- Remove s3fs as a dev dependency #664
- Changed requirements-parser to be a core dependency #673
- Replace *supported_problem_types* field on pipelines with *problem_type* attribute on base classes #678
- Changed AutoML to only show best results for a given pipeline template in *rankings*, added *full_rankings* property to show all #682
- Update *ModelFamily* values: don’t list xgboost/catboost as classifiers now that we have regression pipelines for them #677
- Changed AutoML’s *describe_pipeline* to get problem type from pipeline instead #685
- Standardize *import_or_raise* error messages #683
- Updated argument order of objectives to align with sklearn’s #698
- Renamed *pipeline.feature_importance_graph* to *pipeline.graph_feature_importances* #700
- Moved ROC and confusion matrix methods to *evalml.pipelines.plot_utils* #704
- Renamed *MultiClassificationObjective* to *MulticlassClassificationObjective*, to align with pipeline naming scheme #715

- **Documentation Changes**

- Fixed some sphinx warnings #593
- Fixed docstring for *AutoClassificationSearch* with correct command #599
- Limit readthedocs formats to pdf, not htmlzip and epub #594 #600
- Clean up objectives API documentation #605
- Fixed function on Exploring search results page #604
- Update release process doc #567
- *AutoClassificationSearch* and *AutoRegressionSearch* show inherited methods in API reference #651
- Fixed improperly formatted code in breaking changes for changelog #655
- Added configuration to treat Sphinx warnings as errors #660
- Removed separate plotting section for pipelines in API reference #657, #665
- Have leads example notebook load S3 files using https, so we can delete s3fs dev dependency #664
- Categorized components in API reference and added descriptions for each category #663

- Fixed Sphinx warnings about BalancedAccuracy objective #669
- Updated API reference to include missing components and clean up pipeline docstrings #689
- Reorganize API ref, and clarify pipeline sub-titles #688
- Add and update preprocessing utils in API reference #687
- Added inheritance diagrams to API reference #695
- Documented which default objective AutoML optimizes for #699
- Create separate install page #701
- Include more utils in API ref, like *import_or_raise* #704
- Add more color to pipeline documentation #705
- **Testing Changes**
 - Matched install commands of *check_latest_dependencies* test and it's GitHub action #578
 - Added Github app to auto assign PR author as assignee #477
 - Removed unneeded conda installation of xgboost in windows checkin tests #618
 - Update graph tests to always use tmpfile dir #649
 - Changelog checkin test workaround for release PRs: If 'future release' section is empty of PR refs, pass check #658
 - Add changelog checkin test exception for *dep-update* branch #723

Warning: Breaking Changes

- Pipelines will now no longer take an objective parameter during instantiation, and will no longer have an objective attribute.
- *fit()* and *predict()* now use an optional objective parameter, which is only used in binary classification pipelines to fit for a specific objective.
- *score()* will now use a required *objectives* parameter that is used to determine all the objectives to score on. This differs from the previous behavior, where the pipeline's objective was scored on regardless.
- *score()* will now return one dictionary of all objective scores.
- ROC and ConfusionMatrix plot methods via *Auto(*).plot* have been removed by #615 and are replaced by *roc_curve* and *confusion_matrix* in *evalml.pipelines.plot_utils* in #704
- *normalize_confusion_matrix* has been moved to *evalml.pipelines.plot_utils* #704
- Pipelines *_name* field changed to *custom_name*
- Pipelines *supported_problem_types* field is removed because it is no longer necessary #678
- Updated argument order of objectives' *objective_function* to align with sklearn #698
- *pipeline.feature_importance_graph* has been renamed to *pipeline.graph_feature_importances* in #700
- Removed unsupported MSLE objective #704

v0.8.0 Apr. 1, 2020

- **Enhancements**
 - Add normalization option and information to confusion matrix #484

- Add util function to drop rows with NaN values [#487](#)
- Renamed *PipelineBase.name* as *PipelineBase.summary* and redefined *PipelineBase.name* as class property [#491](#)
- Added access to parameters in Pipelines with *PipelineBase.parameters* (used to be return of *PipelineBase.describe*) [#501](#)
- Added *fill_value* parameter for SimpleImputer [#509](#)
- Added functionality to override component hyperparameters and made pipelines take hyperparameters from components [#516](#)
- Allow `numpy.random.RandomState` for *random_state* parameters [#556](#)
- **Fixes**
 - Removed unused dependency *matplotlib*, and move *category_encoders* to test reqs [#572](#)
- **Changes**
 - Undo version cap in XGBoost placed in [#402](#) and allowed all released of XGBoost [#407](#)
 - Support pandas 1.0.0 [#486](#)
 - Made all references to the logger static [#503](#)
 - Refactored *model_type* parameter for components and pipelines to *model_family* [#507](#)
 - Refactored *problem_types* for pipelines and components into *supported_problem_types* [#515](#)
 - Moved *pipelines/utils.save_pipeline* and *pipelines/utils.load_pipeline* to *PipelineBase.save* and *PipelineBase.load* [#526](#)
 - Limit number of categories encoded by OneHotEncoder [#517](#)
- **Documentation Changes**
 - Updated API reference to remove PipelinePlot and added moved PipelineBase plotting methods [#483](#)
 - Add code style and github issue guides [#463](#) [#512](#)
 - Updated API reference for to surface class variables for pipelines and components [#537](#)
 - Fixed README documentation link [#535](#)
 - Unhid PR references in changelog [#656](#)
- **Testing Changes**
 - Added automated dependency check PR [#482](#), [#505](#)
 - Updated automated dependency check comment [#497](#)
 - Have build_docs job use python executor, so that env vars are set properly [#547](#)
 - Added simple test to make sure OneHotEncoder's top_n works with large number of categories [#552](#)
 - Run windows unit tests on PRs [#557](#)

Warning: Breaking Changes

- `AutoClassificationSearch` and `AutoRegressionSearch`'s `model_types` parameter has been refactored into `allowed_model_families`

- `ModelTypes` enum has been changed to `ModelFamily`
- Components and Pipelines now have a `model_family` field instead of `model_type`
- `get_pipelines` utility function now accepts `model_families` as an argument instead of `model_types`
- `PipelineBase.name` no longer returns structure of pipeline and has been replaced by `PipelineBase.summary`
- `PipelineBase.problem_types` and `Estimator.problem_types` has been renamed to `supported_problem_types`
- `pipelines/utils.save_pipeline` and `pipelines/utils.load_pipeline` moved to `PipelineBase.save` and `PipelineBase.load`

v0.7.0 Mar. 9, 2020

- **Enhancements**

- Added emacs buffers to `.gitignore` #350
- Add CatBoost (gradient-boosted trees) classification and regression components and pipelines #247
- Added Tuner abstract base class #351
- Added `n_jobs` as parameter for `AutoClassificationSearch` and `AutoRegressionSearch` #403
- Changed colors of confusion matrix to shades of blue and updated axis order to match scikit-learn's #426
- Added `PipelineBase.graph` and `feature_importance_graph` methods, moved from previous location #423
- Added support for python 3.8 #462

- **Fixes**

- Fixed ROC and confusion matrix plots not being calculated if user passed own additional_objectives #276
- Fixed `ReadtheDocs` `FileNotFoundError` exception for fraud dataset #439

- **Changes**

- Added `n_estimators` as a tunable parameter for `XGBoost` #307
- Remove unused parameter `ObjectiveBase.fit_needs_proba` #320
- Remove extraneous parameter `component_type` from all components #361
- Remove unused `rankings.csv` file #397
- Downloaded demo and test datasets so unit tests can run offline #408
- Remove `_needs_fitting` attribute from Components #398
- Changed `plot.feature_importance` to show only non-zero feature importances by default, added optional parameter to show all #413
- Refactored `PipelineBase` to take in parameter dictionary and moved pipeline metadata to class attribute #421
- Dropped support for Python 3.5 #438

- Removed unused *apply.py* file [#449](#)
- Clean up requirements.txt to remove unused deps [#451](#)
- Support installation without all required dependencies [#459](#)
- **Documentation Changes**
 - Update release.md with instructions to release to internal license key [#354](#)
- **Testing Changes**
 - Added tests for utils (and moved current utils to gen_utils) [#297](#)
 - Moved XGBoost install into it's own separate step on Windows using Conda [#313](#)
 - Rewind pandas version to before 1.0.0, to diagnose test failures for that version [#325](#)
 - Added dependency update checkin test [#324](#)
 - Rewind XGBoost version to before 1.0.0 to diagnose test failures for that version [#402](#)
 - Update dependency check to use a whitelist [#417](#)
 - Update unit test jobs to not install dev deps [#455](#)

Warning: Breaking Changes

- Python 3.5 will not be actively supported.

v0.6.0 Dec. 16, 2019

- **Enhancements**
 - Added ability to create a plot of feature importances [#133](#)
 - Add early stopping to AutoML using patience and tolerance parameters [#241](#)
 - Added ROC and confusion matrix metrics and plot for classification problems and introduce PipelineSearchPlots class [#242](#)
 - Enhanced AutoML results with search order [#260](#)
 - Added utility function to show system and environment information [#300](#)
- **Fixes**
 - Lower botocore requirement [#235](#)
 - Fixed decision_function calculation for FraudCost objective [#254](#)
 - Fixed return value of Recall metrics [#264](#)
 - Components return *self* on fit [#289](#)
- **Changes**
 - Renamed automl classes to AutoRegressionSearch and AutoClassificationSearch [#287](#)
 - Updating demo datasets to retain column names [#223](#)
 - Moving pipeline visualization to PipelinePlots class [#228](#)
 - Standarizing inputs as pd.DataFrame / pd.Series [#130](#)
 - Enforcing that pipelines must have an estimator as last component [#277](#)
 - Added ipywidgets as a dependency in requirements.txt [#278](#)

- Added Random and Grid Search Tuners #240
- **Documentation Changes**
 - Adding class properties to API reference #244
 - Fix and filter FutureWarnings from scikit-learn #249, #257
 - Adding Linear Regression to API reference and cleaning up some Sphinx warnings #227
- **Testing Changes**
 - Added support for testing on Windows with CircleCI #226
 - Added support for doctests #233

Warning: Breaking Changes

- The `fit()` method for `AutoClassifier` and `AutoRegressor` has been renamed to `search()`.
- `AutoClassifier` has been renamed to `AutoClassificationSearch`
- `AutoRegressor` has been renamed to `AutoRegressionSearch`
- `AutoClassificationSearch.results` and `AutoRegressionSearch.results` now is a dictionary with `pipeline_results` and `search_order` keys. `pipeline_results` can be used to access a dictionary that is identical to the old `.results` dictionary. Whereas, `search_order` returns a list of the search order in terms of `pipeline_id`.
- Pipelines now require an estimator as the last component in `component_list`. Slicing pipelines now throws an `NotImplementedError` to avoid returning pipelines without an estimator.

v0.5.2 Nov. 18, 2019

- **Enhancements**
 - Adding basic pipeline structure visualization #211
- **Documentation Changes**
 - Added notebooks to build process #212

v0.5.1 Nov. 15, 2019

- **Enhancements**
 - Added basic outlier detection guardrail #151
 - Added basic ID column guardrail #135
 - Added support for unlimited pipelines with a `max_time` limit #70
 - Updated `.readthedocs.yaml` to successfully build #188
- **Fixes**
 - Removed MSLE from default additional objectives #203
 - Fixed `random_state` passed in pipelines #204
 - Fixed slow down in `RFRegressor` #206
- **Changes**
 - Pulled information for `describe_pipeline` from pipeline's new `describe` method #190
 - Refactored pipelines #108

- Removed guardrails from Auto(*) #202, #208

- **Documentation Changes**

- Updated documentation to show max_time enhancements #189
- Updated release instructions for RTD #193
- Added notebooks to build process #212
- Added contributing instructions #213
- Added new content #222

v0.5.0 Oct. 29, 2019

- **Enhancements**

- Added basic one hot encoding #73
- Use enums for model_type #110
- Support for splitting regression datasets #112
- Auto-infer multiclass classification #99
- Added support for other units in max_time #125
- Detect highly null columns #121
- Added additional regression objectives #100
- Show an interactive iteration vs. score plot when using fit() #134

- **Fixes**

- Reordered *describe_pipeline* #94
- Added type check for model_type #109
- Fixed *s* units when setting string max_time #132
- Fix objectives not appearing in API documentation #150

- **Changes**

- Reorganized tests #93
- Moved logging to its own module #119
- Show progress bar history #111
- Using cloudpickle instead of pickle to allow unloading of custom objectives #113
- Removed render.py #154

- **Documentation Changes**

- Update release instructions #140
- Include additional_objectives parameter #124
- Added Changelog #136

- **Testing Changes**

- Code coverage #90
- Added CircleCI tests for other Python versions #104
- Added doc notebooks as tests #139

- Test metadata for CircleCI and 2 core parallelism [#137](#)

v0.4.1 Sep. 16, 2019

- **Enhancements**

- Added AutoML for classification and regressor using Autobase and Skopt [#7](#) [#9](#)
- Implemented standard classification and regression metrics [#7](#)
- Added logistic regression, random forest, and XGBoost pipelines [#7](#)
- Implemented support for custom objectives [#15](#)
- Feature importance for pipelines [#18](#)
- Serialization for pipelines [#19](#)
- Allow fitting on objectives for optimal threshold [#27](#)
- Added detect label leakage [#31](#)
- Implemented callbacks [#42](#)
- Allow for multiclass classification [#21](#)
- Added support for additional objectives [#79](#)

- **Fixes**

- Fixed feature selection in pipelines [#13](#)
- Made random_seed usage consistent [#45](#)

- **Documentation Changes**

- Documentation Changes
- Added docstrings [#6](#)
- Created notebooks for docs [#6](#)
- Initialized readthedocs EvalML [#6](#)
- Added favicon [#38](#)

- **Testing Changes**

- Added testing for loading data [#39](#)

v0.2.0 Aug. 13, 2019

- **Enhancements**

- Created fraud detection objective [#4](#)

v0.1.0 July. 31, 2019

- *First Release*

- **Enhancements**

- Added lead scoring objective [#1](#)
- Added basic classifier [#1](#)

- **Documentation Changes**

- Initialized Sphinx for docs [#1](#)

1.6.16 API Reference

Demo Datasets

<code>load_fraud</code>	Load credit card fraud dataset.
<code>load_wine</code>	Load wine dataset.
<code>load_breast_cancer</code>	Load breast cancer dataset.
<code>load_diabetes</code>	Load diabetes dataset.

`evalml.demos.load_fraud`

`evalml.demos.load_fraud(n_rows=None, verbose=True)`

Load credit card fraud dataset. The fraud dataset can be used for binary classification problems.

Parameters

- **n_rows** (*int*) – number of rows from the dataset to return
- **verbose** (*bool*) – whether to print information about features and labels

Returns X, y

Return type pd.DataFrame, pd.Series

`evalml.demos.load_wine`

`evalml.demos.load_wine()`

Load wine dataset. Multiclass problem

Returns X, y

Return type pd.DataFrame, pd.Series

`evalml.demos.load_breast_cancer`

`evalml.demos.load_breast_cancer()`

Load breast cancer dataset. Multiclass problem

Returns X, y

Return type pd.DataFrame, pd.Series

`evalml.demos.load_diabetes`

`evalml.demos.load_diabetes()`

Load diabetes dataset. Regression problem

Returns X, y

Return type pd.DataFrame, pd.Series

Preprocessing

Utilities to preprocess data before using evalml.

<code>drop_nan_target_rows</code>	Drops rows in X and y when row in the target y has a value of NaN.
<code>label_distribution</code>	Get the label distributions.
<code>load_data</code>	Load features and labels from file.
<code>number_of_features</code>	Get the number of features for specific dtypes.
<code>split_data</code>	Splits data into train and test sets.

evalml.preprocessing.drop_nan_target_rows

evalml.preprocessing.drop_nan_target_rows(X, y)

Drops rows in X and y when row in the target y has a value of NaN.

Parameters

- **X** (*pd.DataFrame*) – Data to transform
- **y** (*pd.Series*) – Target values

Returns Transformed X (and y, if passed in) with rows that had a NaN value removed.

Return type *pd.DataFrame*

evalml.preprocessing.label_distribution

evalml.preprocessing.label_distribution(labels)

Get the label distributions.

Parameters **labels** (*pd.Series*) – Label values

Returns Label values and their frequency distribution as percentages.

Return type *pd.Series*

evalml.preprocessing.load_data

evalml.preprocessing.load_data(path, index, label, n_rows=None, drop=None, verbose=True, ***kwargs*)

Load features and labels from file.

Parameters

- **path** (*str*) – Path to file or a http/ftp/s3 URL
- **index** (*str*) – Column for index
- **label** (*str*) – Column for labels
- **n_rows** (*int*) – Number of rows to return
- **drop** (*list*) – List of columns to drop
- **verbose** (*bool*) – If True, prints information about features and labels

Returns features and labels

Return type `pd.DataFrame, pd.Series`

`evalml.preprocessing.number_of_features`

`evalml.preprocessing.number_of_features(dtypes)`

Get the number of features for specific dtypes.

Parameters `dtypes` (`pd.Series`) – dtypes to get the number of features for

Returns dtypes and the number of features for each input type

Return type `pd.Series`

`evalml.preprocessing.split_data`

`evalml.preprocessing.split_data(X, y, regression=False, test_size=0.2, random_state=None)`

Splits data into train and test sets.

Parameters

- **X** (`pd.DataFrame` or `np.array`) – data of shape `[n_samples, n_features]`
- **y** (`pd.Series`) – labels of length `[n_samples]`
- **regression** (`bool`) – if true, do not use stratified split
- **test_size** (`float`) – percent of train set to holdout for testing
- **random_state** (`int`, `np.random.RandomState`) – seed for the random number generator

Returns features and labels each split into train and test sets

Return type `pd.DataFrame, pd.DataFrame, pd.Series, pd.Series`

AutoML

AutoML Search Classes

AutoMLSearch

Automated Pipeline search.

evalml automl AutoMLSearch

evalml.automl.automl_search.AutoMLSearch

```
class evalml.automl.AutoMLSearch(problem_type=None, objective='auto',
                                max_pipelines=None, max_time=None, pa-
                                tience=None, tolerance=None, data_split=None, al-
                                lowed_pipelines=None, allowed_model_families=None,
                                start_iteration_callback=None, add_result_callback=None,
                                additional_objectives=None, random_state=0,
                                n_jobs=-1, tuner_class=None, verbose=True, opti-
                                mize_thresholds=False)
```

Automated Pipeline search.

Methods

<code>__init__</code>	Automated pipeline search
<code>add_to_rankings</code>	Fits and evaluates a given pipeline then adds the results to the automl rankings with the requirement that automl search has been run.
<code>describe_pipeline</code>	Describe a pipeline
<code>get_pipeline</code>	Given the ID of a pipeline training result, returns an untrained instance of the specified pipeline initialized with the parameters used to train that pipeline during automl search.
<code>load</code>	Loads AutoML object at file path
<code>save</code>	Saves AutoML object at file path
<code>search</code>	Find best classifier

evalml.automl.AutoMLSearch.__init__

```
AutoMLSearch.__init__(problem_type=None, objective='auto', max_pipelines=None,
                       max_time=None, patience=None, tolerance=None, data_split=None,
                       allowed_pipelines=None, allowed_model_families=None,
                       start_iteration_callback=None, add_result_callback=None,
                       additional_objectives=None, random_state=0, n_jobs=-1,
                       tuner_class=None, verbose=True, optimize_thresholds=False)
```

Automated pipeline search

Parameters

- **problem_type** (*str* or `ProblemTypes`) – Choice of ‘regression’, ‘binary’, or ‘multiclass’, depending on the desired problem type.
- **objective** (*str*, `ObjectiveBase`) – The objective to optimize for. When set to auto, chooses: `LogLossBinary` for binary classification problems, `LogLossMulticlass` for multiclass classification problems, and `R2` for regression problems.
- **max_pipelines** (*int*) – Maximum number of pipelines to search. If `max_pipelines` and `max_time` is not set, then `max_pipelines` will default to `max_pipelines` of 5.
- **max_time** (*int*, *str*) – Maximum time to search for pipelines. This will not start a new pipeline search after the duration has elapsed. If it is an integer, then the time will be in seconds. For strings, time can be specified as seconds, minutes, or hours.
- **patience** (*int*) – Number of iterations without improvement to stop search early. Must be positive. If `None`, early stopping is disabled. Defaults to `None`.
- **tolerance** (*float*) – Minimum percentage difference to qualify as score improvement for early stopping. Only applicable if `patience` is not `None`. Defaults to `None`.
- **allowed_pipelines** (*list(class)*) – A list of `PipelineBase` subclasses indicating the pipelines allowed in the search. The default of `None` indicates all pipelines for this problem type are allowed. Setting this field will cause `allowed_model_families` to be ignored.
- **allowed_model_families** (*list(str, ModelFamily)*) – The model families to search. The default of `None` searches over all model families. Run `evalml.list_model_families("binary")` to see options. Change *binary* to *multiclass* or *regression* depending on the problem type. Note that if `allowed_pipelines` is provided, this parameter will be ignored.
- **data_split** (`sklearn.model_selection.BaseCrossValidator`) – data splitting method to use. Defaults to `StratifiedKFold`.
- **tuner_class** – the tuner class to use. Defaults to `scikit-optimize` tuner
- **start_iteration_callback** (*callable*) – function called before each pipeline training iteration. Passed two parameters: `pipeline_class`, `parameters`.
- **add_result_callback** (*callable*) – function called after each pipeline training iteration. Passed two parameters: `results`, `trained_pipeline`.
- **additional_objectives** (*list*) – Custom set of objectives to score on. Will override default objectives for problem type if not empty.
- **random_state** (*int*, `np.random.RandomState`) – The random seed/state. Defaults to 0.
- **n_jobs** (*int* or `None`) – Non-negative integer describing level of parallelism used for pipelines. `None` and 1 are equivalent. If set to -1, all CPUs are used. For `n_jobs` below -1, `(n_cpus + 1 + n_jobs)` are used.
- **verbose** (*boolean*) – If `True`, turn verbosity on. Defaults to `True`

evalml.automl.AutoMLSearch.add_to_rankings

`AutoMLSearch.add_to_rankings(pipeline, X, y)`

Fits and evaluates a given pipeline then adds the results to the automl rankings with the requirement that automl search has been run. Please use the same data as previous runs of automl search. If pipeline already exists in rankings this method will return `None`.

Parameters

- **pipeline** (`PipelineBase`) – pipeline to train and evaluate.
- **x** (`pd.DataFrame`) – the input training data of shape `[n_samples, n_features]`.
- **y** (`pd.Series`) – the target training labels of length `[n_samples]`.

evalml automl.AutoMLSearch.describe_pipeline

`AutoMLSearch.describe_pipeline(pipeline_id, return_dict=False)`

Describe a pipeline

Parameters

- **pipeline_id** (`int`) – pipeline to describe
- **return_dict** (`bool`) – If True, return dictionary of information about pipeline. Defaults to False.

Returns Description of specified pipeline. Includes information such as type of pipeline components, problem, training time, cross validation, etc.

evalml automl.AutoMLSearch.get_pipeline

`AutoMLSearch.get_pipeline(pipeline_id, random_state=0)`

Given the ID of a pipeline training result, returns an untrained instance of the specified pipeline initialized with the parameters used to train that pipeline during automl search.

Parameters

- **pipeline_id** (`int`) – pipeline to retrieve
- **random_state** (`int`, `np.random.RandomState`) – The random seed/state. Defaults to 0.

Returns untrained pipeline instance associated with the provided ID

Return type *PipelineBase*

evalml automl.AutoMLSearch.load

static `AutoMLSearch.load(file_path)`

Loads AutoML object at file path

Parameters **file_path** (`str`) – location to find file to load

Returns AutoSearchBase object

evalml automl.AutoMLSearch.save

`AutoMLSearch.save(file_path)`

Saves AutoML object at file path

Parameters **file_path** (`str`) – location to save file

Returns None

evalml.automl.AutoMLSearch.search

`AutoMLSearch.search(X, y, data_checks='auto', feature_types=None, raise_errors=True, show_iteration_plot=True)`

Find best classifier

Parameters

- **X** (*pd.DataFrame*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]
- **feature_types** (*list, optional*) – list of feature types, either numerical or categorical. Categorical features will automatically be encoded
- **raise_errors** (*boolean*) – If True, raise errors and exit search if a pipeline errors during fitting. If False, set scores for the errored pipeline to NaN and continue search. Defaults to True.
- **show_iteration_plot** (*boolean, True*) – Shows an iteration vs. score plot in Jupyter notebook. Disabled by default in non-Jupyter environments.
- **data_checks** (*DataChecks, list(Datacheck), str, None*) – A collection of data checks to run before automl search. If data checks produce any errors, an exception will be thrown before the search begins. If “disabled” or None, no data checks will be done. If set to “auto”, DefaultDataChecks will be done. Default value is set to “auto”.

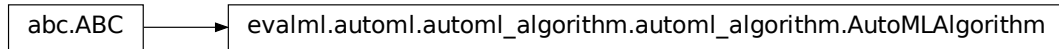
Returns self

Attributes

<code>best_pipeline</code>	Returns an untrained instance of the best pipeline and parameters found during automl search.
<code>data_check_results</code>	
<code>full_rankings</code>	Returns a pandas.DataFrame with scoring results from all pipelines searched
<code>has_searched</code>	Returns <i>True</i> if search has been ran and <i>False</i> if not
<code>rankings</code>	Returns a pandas.DataFrame with scoring results from the highest-scoring set of parameters used with each pipeline.

AutoML Algorithm Classes

<code>AutoMLAlgorithm</code>	Base class for the automl algorithms which power evalml.
<code>IterativeAlgorithm</code>	An automl algorithm which first fits a base round of pipelines with default parameters, then does a round of parameter tuning on each pipeline in order of performance.

evalml.automl.automl_algorithm.AutoMLAlgorithm

```

class evalml.automl.automl_algorithm.AutoMLAlgorithm(allowed_pipelines=None,
                                                    max_pipelines=None,
                                                    tuner_class=None,      ran-
                                                    dom_state=0)

```

Base class for the automl algorithms which power evalml.

Methods

<code>__init__</code>	This class represents an automated machine learning (AutoML) algorithm.
<code>add_result</code>	Register results from evaluating a pipeline
<code>next_batch</code>	Get the next batch of pipelines to evaluate

evalml.automl.automl_algorithm.AutoMLAlgorithm.__init__

```

AutoMLAlgorithm.__init__(allowed_pipelines=None, max_pipelines=None, tuner_class=None,
                        random_state=0)

```

This class represents an automated machine learning (AutoML) algorithm. It encapsulates the decision-making logic behind an automl search, by both deciding which pipelines to evaluate next and by deciding what set of parameters to configure the pipeline with.

To use this interface, you must define a `next_batch` method which returns the next group of pipelines to evaluate on the training data. That method may access state and results recorded from the previous batches, although that information is not tracked in a general way in this base class. Overriding `add_result` is a convenient way to record pipeline evaluation info if necessary.

Parameters

- **allowed_pipelines** (*list(class)*) – A list of PipelineBase subclasses indicating the pipelines allowed in the search. The default of None indicates all pipelines for this problem type are allowed.
- **max_pipelines** (*int*) – The maximum number of pipelines to be evaluated.
- **tuner_class** (*class*) – A subclass of Tuner, to be used to find parameters for each pipeline. The default of None indicates the SKOptTuner will be used.
- **random_state** (*int, np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.automl.automl_algorithm.AutoMLAlgorithm.add_result`AutoMLAlgorithm.add_result(score_to_minimize, pipeline)`

Register results from evaluating a pipeline

Parameters

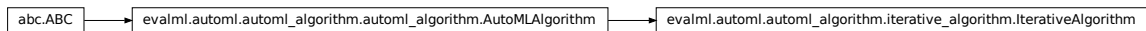
- **score_to_minimize** (*float*) – The score obtained by this pipeline on the primary objective, converted so that lower values indicate better pipelines.
- **pipeline** (*PipelineBase*) – The trained pipeline object which was used to compute the score.

evalml.automl.automl_algorithm.AutoMLAlgorithm.next_batch`AutoMLAlgorithm.next_batch()`

Get the next batch of pipelines to evaluate

Returns a list of instances of PipelineBase subclasses, ready to be trained and evaluated.**Return type** list(*PipelineBase*)**Attributes**

<code>batch_number</code>	Returns the number of batches which have been recommended so far.
<code>pipeline_number</code>	Returns the number of pipelines which have been recommended so far.

evalml.automl.automl_algorithm.IterativeAlgorithm

```

class evalml.automl.automl_algorithm.IterativeAlgorithm(allowed_pipelines=None,
                                                         max_pipelines=None,
                                                         tuner_class=None,
                                                         random_state=0,
                                                         pipelines_per_batch=5,
                                                         n_jobs=-1,          num-
                                                         ber_features=None)

```

An automl algorithm which first fits a base round of pipelines with default parameters, then does a round of parameter tuning on each pipeline in order of performance.

Methods

<code>__init__</code>	An automl algorithm which first fits a base round of pipelines with default parameters, then does a round of parameter tuning on each pipeline in order of performance.
<code>add_result</code>	Register results from evaluating a pipeline
<code>next_batch</code>	Get the next batch of pipelines to evaluate

`evalml.automl.automl_algorithm.IterativeAlgorithm.__init__`

`IterativeAlgorithm.__init__` (*allowed_pipelines=None*, *max_pipelines=None*,
tuner_class=None, *random_state=0*, *pipelines_per_batch=5*,
n_jobs=-1, *number_features=None*)

An automl algorithm which first fits a base round of pipelines with default parameters, then does a round of parameter tuning on each pipeline in order of performance.

Parameters

- **`allowed_pipelines`** (*list (class)*) – A list of PipelineBase subclasses indicating the pipelines allowed in the search. The default of None indicates all pipelines for this problem type are allowed.
- **`max_pipelines`** (*int*) – The maximum number of pipelines to be evaluated.
- **`tuner_class`** (*class*) – A subclass of Tuner, to be used to find parameters for each pipeline. The default of None indicates the SKOptTuner will be used.
- **`random_state`** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.
- **`pipelines_per_batch`** (*int*) – the number of pipelines to be evaluated in each batch, after the first batch.
- **`n_jobs`** (*int or None*) – Non-negative integer describing level of parallelism used for pipelines.
- **`number_features`** (*int*) – The number of columns in the input features.

`evalml.automl.automl_algorithm.IterativeAlgorithm.add_result`

`IterativeAlgorithm.add_result` (*score_to_minimize*, *pipeline*)

Register results from evaluating a pipeline

Parameters

- **`score_to_minimize`** (*float*) – The score obtained by this pipeline on the primary objective, converted so that lower values indicate better pipelines.
- **`pipeline`** (*PipelineBase*) – The trained pipeline object which was used to compute the score.

`evalml.automl.automl_algorithm.IterativeAlgorithm.next_batch`

`IterativeAlgorithm.next_batch` ()

Get the next batch of pipelines to evaluate

Returns a list of instances of PipelineBase subclasses, ready to be trained and evaluated.

Return type `list(PipelineBase)`

Attributes

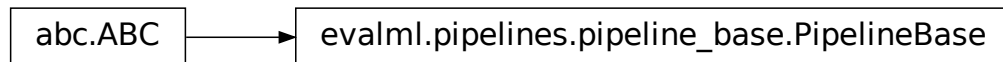
<code>batch_number</code>	Returns the number of batches which have been recommended so far.
<code>pipeline_number</code>	Returns the number of pipelines which have been recommended so far.

Pipelines

Pipeline Base Classes

<i>PipelineBase</i>	Base class for all pipelines.
<i>ClassificationPipeline</i>	Pipeline subclass for all classification pipelines.
<i>BinaryClassificationPipeline</i>	Pipeline subclass for all binary classification pipelines.
<i>MulticlassClassificationPipeline</i>	Pipeline subclass for all multiclass classification pipelines.
<i>RegressionPipeline</i>	Pipeline subclass for all regression pipelines.

evalml.pipelines.PipelineBase



class `evalml.pipelines.PipelineBase` (*parameters*, *random_state=0*)
Base class for all pipelines.

Methods

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph

Continued on next page

Table 12 – continued from previous page

<code>graph_feature_importance</code>	Generate a bar graph of the pipeline’s feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

`evalml.pipelines.PipelineBase.__init__`

`PipelineBase.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary `{ }` implies using all default values for component parameters.
- **random_state** (*int*, `np.random.RandomState`) – The random seed/state. Defaults to 0.

`evalml.pipelines.PipelineBase.clone`

`PipelineBase.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.PipelineBase.describe`

`PipelineBase.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.PipelineBase.fit`

`PipelineBase.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.PipelineBase.get_component

`PipelineBase.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.PipelineBase.graph

`PipelineBase.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters **filepath** (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.PipelineBase.graph_feature_importance

`PipelineBase.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.PipelineBase.load

static `PipelineBase.load(file_path)`

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.PipelineBase.predict

`PipelineBase.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.PipelineBase.save`

`PipelineBase.save` (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns `None`

`evalml.pipelines.PipelineBase.score`

`PipelineBase.score` (*X*, *y*, *objectives*)

Evaluate model performance on current and additional objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type `dict`

`evalml.pipelines.ClassificationPipeline`



class `evalml.pipelines.ClassificationPipeline` (*parameters*, *random_state=0*)

Pipeline subclass for all classification pipelines.

Methods

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters

Continued on next page

Table 13 – continued from previous page

<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline’s feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.ClassificationPipeline.__init__`

`ClassificationPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.ClassificationPipeline.clone`

`ClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.ClassificationPipeline.describe`

`ClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.ClassificationPipeline.fit

`ClassificationPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.ClassificationPipeline.get_component

`ClassificationPipeline.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ClassificationPipeline.graph

`ClassificationPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ClassificationPipeline.graph_feature_importance

`ClassificationPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ClassificationPipeline.load

static `ClassificationPipeline.load(file_path)`

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ClassificationPipeline.predict

`ClassificationPipeline.predict` (*X*, *objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.ClassificationPipeline.predict_proba

`ClassificationPipeline.predict_proba` (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.ClassificationPipeline.save

`ClassificationPipeline.save` (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns *None*

evalml.pipelines.ClassificationPipeline.score

`ClassificationPipeline.score` (*X*, *y*, *objectives*)

Evaluate model performance on objectives

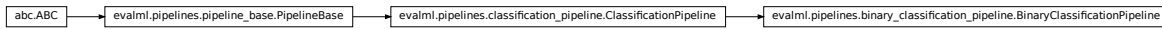
Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type *dict*

evalml.pipelines.BinaryClassificationPipeline



class evalml.pipelines.**BinaryClassificationPipeline** (*parameters*, *random_state=0*)
 Pipeline subclass for all binary classification pipelines.

Methods

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.BinaryClassificationPipeline.__init__

BinaryClassificationPipeline.**__init__** (*parameters*, *random_state=0*)
 Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or ComponentBase subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.BinaryClassificationPipeline.clone

BinaryClassificationPipeline.**clone** (*random_state=0*)
 Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.BinaryClassificationPipeline.describe`

`BinaryClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.BinaryClassificationPipeline.fit`

`BinaryClassificationPipeline.fit(X, y)`

Build a model

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

`evalml.pipelines.BinaryClassificationPipeline.get_component`

`BinaryClassificationPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.BinaryClassificationPipeline.graph`

`BinaryClassificationPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.BinaryClassificationPipeline.graph_feature_importance

BinaryClassificationPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.BinaryClassificationPipeline.load

static BinaryClassificationPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.BinaryClassificationPipeline.predict

BinaryClassificationPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.BinaryClassificationPipeline.predict_proba

BinaryClassificationPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.BinaryClassificationPipeline.save

BinaryClassificationPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.BinaryClassificationPipeline.score

BinaryClassificationPipeline.**score**(*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.MulticlassClassificationPipeline

class evalml.pipelines.**MulticlassClassificationPipeline**(*parameters*, *random_state=0*)

Pipeline subclass for all multiclass classification pipelines.

Methods

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.MulticlassClassificationPipeline.__init__

MulticlassClassificationPipeline.**__init__**(*parameters*, *random_state=0*)

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or ComponentBase subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.MulticlassClassificationPipeline.clone

`MulticlassClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.MulticlassClassificationPipeline.describe

`MulticlassClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.MulticlassClassificationPipeline.fit

`MulticlassClassificationPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.MulticlassClassificationPipeline.get_component

`MulticlassClassificationPipeline.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.MulticlassClassificationPipeline.graph

`MulticlassClassificationPipeline.graph` (*filepath=None*)

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type `graphviz.Digraph`

evalml.pipelines.MulticlassClassificationPipeline.graph_feature_importance

`MulticlassClassificationPipeline.graph_feature_importance` (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

evalml.pipelines.MulticlassClassificationPipeline.load

static `MulticlassClassificationPipeline.load` (*file_path*)

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

evalml.pipelines.MulticlassClassificationPipeline.predict

`MulticlassClassificationPipeline.predict` (*X, objective=None*)

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.MulticlassClassificationPipeline.predict_proba

`MulticlassClassificationPipeline.predict_proba` (*X*)

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type `pd.DataFrame`

evalml.pipelines.MulticlassClassificationPipeline.save

MulticlassClassificationPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.MulticlassClassificationPipeline.score

MulticlassClassificationPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on objectives

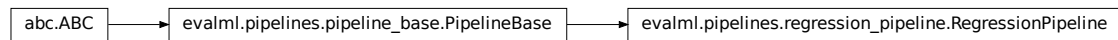
Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.RegressionPipeline



class evalml.pipelines.**RegressionPipeline** (*parameters*, *random_state=0*)

Pipeline subclass for all regression pipelines.

Methods

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.

Continued on next page

Table 16 – continued from previous page

<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

evalml.pipelines.RegressionPipeline.__init__

`RegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.RegressionPipeline.clone

`RegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.RegressionPipeline.describe

`RegressionPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.RegressionPipeline.fit

`RegressionPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.RegressionPipeline.get_component

`RegressionPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.RegressionPipeline.graph

`RegressionPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.RegressionPipeline.graph_feature_importance

`RegressionPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.RegressionPipeline.load

static `RegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.RegressionPipeline.predict

`RegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.RegressionPipeline.save

`RegressionPipeline.save` (*file_path*)

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns None

evalml.pipelines.RegressionPipeline.score

`RegressionPipeline.score` (*X*, *y*, *objectives*)

Evaluate model performance on current and additional objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type dict

Classification Pipelines

<code>CatBoostBinaryClassificationPipeline</code>	CatBoost Pipeline for binary classification.
<code>CatBoostMulticlassClassificationPipeline</code>	CatBoost Pipeline for multiclass classification.
<code>ENBinaryPipeline</code>	Elastic Net Pipeline for binary classification problems.
<code>ENMulticlassPipeline</code>	Elastic Net Pipeline for multiclass classification problems.
<code>ETBinaryClassificationPipeline</code>	Extra Trees Pipeline for binary classification.
<code>ETMulticlassClassificationPipeline</code>	Extra Trees Pipeline for multiclass classification.
<code>LogisticRegressionBinaryPipeline</code>	Logistic Regression Pipeline for binary classification.
<code>LogisticRegressionMulticlassPipeline</code>	Logistic Regression Pipeline for multiclass classification.
<code>RFBinaryClassificationPipeline</code>	Random Forest Pipeline for binary classification.
<code>RFMulticlassClassificationPipeline</code>	Random Forest Pipeline for multiclass classification.
<code>XGBoostBinaryPipeline</code>	XGBoost Pipeline for binary classification.
<code>XGBoostMulticlassPipeline</code>	XGBoost Pipeline for multiclass classification.
<code>BaselineBinaryPipeline</code>	Baseline Pipeline for binary classification.
<code>BaselineMulticlassPipeline</code>	Baseline Pipeline for multiclass classification.
<code>ModeBaselineBinaryPipeline</code>	Mode Baseline Pipeline for binary classification.
<code>ModeBaselineMulticlassPipeline</code>	Mode Baseline Pipeline for multiclass classification.

evalml.pipelines.CatBoostBinaryClassificationPipeline

```

class evalml.pipelines.CatBoostBinaryClassificationPipeline(parameters, random_state=0)
    CatBoost Pipeline for binary classification. CatBoost is an open-source library and natively supports categorical features.

    For more information, check out https://catboost.ai/ Note: impute_strategy must support both string and numeric data

    name = 'Cat Boost Binary Classification Pipeline'
    custom_name = None
    summary = 'CatBoost Classifier w/ Simple Imputer'
    component_graph = ['Simple Imputer', 'CatBoost Classifier']
    problem_type = 'binary'
    model_family = 'catboost'
    hyperparameters = {'CatBoost Classifier': {'eta': Real(low=1e-06, high=1, prior='uniform')}}
    custom_hyperparameters = {'Simple Imputer': {'impute_strategy': ['most_frequent']}}
    default_parameters = {'CatBoost Classifier': {'bootstrap_type': None, 'eta': 0.03,

```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.CatBoostBinaryClassificationPipeline.__init__

`CatBoostBinaryClassificationPipeline.__init__(parameters, random_state=0)`
Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.CatBoostBinaryClassificationPipeline.clone`

`CatBoostBinaryClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.CatBoostBinaryClassificationPipeline.describe`

`CatBoostBinaryClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.CatBoostBinaryClassificationPipeline.fit`

`CatBoostBinaryClassificationPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

`evalml.pipelines.CatBoostBinaryClassificationPipeline.get_component`

`CatBoostBinaryClassificationPipeline.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.CatBoostBinaryClassificationPipeline.graph

`CatBoostBinaryClassificationPipeline.graph (filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.CatBoostBinaryClassificationPipeline.graph_feature_importance

`CatBoostBinaryClassificationPipeline.graph_feature_importance (show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.CatBoostBinaryClassificationPipeline.load

static `CatBoostBinaryClassificationPipeline.load (file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.CatBoostBinaryClassificationPipeline.predict

`CatBoostBinaryClassificationPipeline.predict (X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.CatBoostBinaryClassificationPipeline.predict_proba

`CatBoostBinaryClassificationPipeline.predict_proba (X)`

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type `pd.DataFrame`

`evalml.pipelines.CatBoostBinaryClassificationPipeline.save`

`CatBoostBinaryClassificationPipeline.save(file_path)`

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

`evalml.pipelines.CatBoostBinaryClassificationPipeline.score`

`CatBoostBinaryClassificationPipeline.score(X, y, objectives)`

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- **y** (*pd.Series*) – true labels of length `[n_samples]`
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type `dict`

`evalml.pipelines.CatBoostMulticlassClassificationPipeline`



```
class evalml.pipelines.CatBoostMulticlassClassificationPipeline(parameters,
                                                                random_state=0)
```

CatBoost Pipeline for multiclass classification. CatBoost is an open-source library and natively supports categorical features.

For more information, check out <https://catboost.ai/> Note: `impute_strategy` must support both string and numeric data

```
name = 'Cat Boost Multiclass Classification Pipeline'
```

```
custom_name = None
```

```
summary = 'CatBoost Classifier w/ Simple Imputer'
```

```
component_graph = ['Simple Imputer', 'CatBoost Classifier']
```

```
problem_type = 'multiclass'
```

```
model_family = 'catboost'
```

```
hyperparameters = {'CatBoost Classifier': {'eta': Real(low=1e-06, high=1, prior='uniform', min_value=1e-06, max_value=1.0)}}
```

```
custom_hyperparameters = {'Simple Imputer': {'impute_strategy': ['most_frequent']}}
```

```
default_parameters = {'CatBoost Classifier': {'bootstrap_type': None, 'eta': 0.03,
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.__init__`

`CatBoostMulticlassClassificationPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.clone`

`CatBoostMulticlassClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.CatBoostMulticlassClassificationPipeline.describe`CatBoostMulticlassClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline.
Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.CatBoostMulticlassClassificationPipeline.fit`CatBoostMulticlassClassificationPipeline.fit(X, y)`

Build a model

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- `y` (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.CatBoostMulticlassClassificationPipeline.get_component`CatBoostMulticlassClassificationPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.CatBoostMulticlassClassificationPipeline.graph`CatBoostMulticlassClassificationPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.CatBoostMulticlassClassificationPipeline.graph_feature_importance`CatBoostMulticlassClassificationPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.load`

static `CatBoostMulticlassClassificationPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.predict`

`CatBoostMulticlassClassificationPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `objective` (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.predict_proba`

`CatBoostMulticlassClassificationPipeline.predict_proba(X)`

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`

Returns probability estimates

Return type `pd.DataFrame`

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.save`

`CatBoostMulticlassClassificationPipeline.save(file_path)`

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

`evalml.pipelines.CatBoostMulticlassClassificationPipeline.score`

`CatBoostMulticlassClassificationPipeline.score(X, y, objectives)`

Evaluate model performance on objectives

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – true labels of length `[n_samples]`

- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.ENBinaryPipeline



class evalml.pipelines.ENBinaryPipeline(*parameters*, *random_state=0*)

Elastic Net Pipeline for binary classification problems.

name = 'ENBinary Pipeline'

custom_name = None

summary = 'Elastic Net Classifier w/ One Hot Encoder + Simple Imputer'

component_graph = ['One Hot Encoder', 'Simple Imputer', 'Elastic Net Classifier']

problem_type = 'binary'

model_family = 'linear_model'

hyperparameters = {'Elastic Net Classifier': {'alpha': Real(low=0, high=1, prior='un

custom_hyperparameters = None

default_parameters = {'Elastic Net Classifier': {'alpha': 0.5, 'l1_ratio': 0.5, 'ma

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.

Continued on next page

Table 23 – continued from previous page

<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.ENBinaryPipeline.__init__`

`ENBinaryPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.ENBinaryPipeline.clone`

`ENBinaryPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.ENBinaryPipeline.describe`

`ENBinaryPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.ENBinaryPipeline.fit`

`ENBinaryPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.ENBinaryPipeline.get_component

`ENBinaryPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ENBinaryPipeline.graph

`ENBinaryPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ENBinaryPipeline.graph_feature_importance

`ENBinaryPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ENBinaryPipeline.load

static `ENBinaryPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ENBinaryPipeline.predict

`ENBinaryPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.ENBinaryPipeline.predict_proba

ENBinaryPipeline.**predict_proba**(*X*)

Make probability estimates for labels.

Parameters *X* (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.ENBinaryPipeline.save

ENBinaryPipeline.**save**(*file_path*)

Saves pipeline at file path

Parameters *file_path* (*str*) – location to save file

Returns None

evalml.pipelines.ENBinaryPipeline.score

ENBinaryPipeline.**score**(*X*, *y*, *objectives*)

Evaluate model performance on objectives

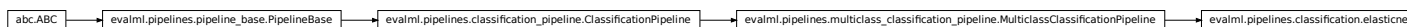
Parameters

- *X* (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- *y* (*pd.Series*) – true labels of length [n_samples]
- *objectives* (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.ENMulticlassPipeline



class evalml.pipelines.**ENMulticlassPipeline**(*parameters*, *random_state=0*)

Elastic Net Pipeline for multiclass classification problems.

name = 'ENMulticlass Pipeline'

custom_name = None

summary = 'Elastic Net Classifier w/ One Hot Encoder + Simple Imputer'

component_graph = ['One Hot Encoder', 'Simple Imputer', 'Elastic Net Classifier']

problem_type = 'multiclass'

model_family = 'linear_model'

hyperparameters = {'Elastic Net Classifier': {'alpha': Real(low=0, high=1, prior='un


```
custom_hyperparameters = None
```

```
default_parameters = {'Elastic Net Classifier': {'alpha': 0.5, 'l1_ratio': 0.5, 'ma
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.ENMulticlassPipeline.__init__

```
ENMulticlassPipeline.__init__(parameters, random_state=0)
```

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.ENMulticlassPipeline.clone

```
ENMulticlassPipeline.clone(random_state=0)
```

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.ENMulticlassPipeline.describe`

`ENMulticlassPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline.
Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.ENMulticlassPipeline.fit`

`ENMulticlassPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

`evalml.pipelines.ENMulticlassPipeline.get_component`

`ENMulticlassPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.ENMulticlassPipeline.graph`

`ENMulticlassPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ENMulticlassPipeline.graph_feature_importance

`ENMulticlassPipeline.graph_feature_importance` (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

evalml.pipelines.ENMulticlassPipeline.load

static `ENMulticlassPipeline.load` (*file_path*)

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

evalml.pipelines.ENMulticlassPipeline.predict

`ENMulticlassPipeline.predict` (*X, objective=None*)

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.ENMulticlassPipeline.predict_proba

`ENMulticlassPipeline.predict_proba` (*X*)

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type `pd.DataFrame`

evalml.pipelines.ENMulticlassPipeline.save

`ENMulticlassPipeline.save` (*file_path*)

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

evalml.pipelines.ENMulticlassPipeline.score

ENMulticlassPipeline.**score** (*X*, *y*, *objectives*)
Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.ETBinaryClassificationPipeline



```
class evalml.pipelines.ETBinaryClassificationPipeline(parameters, random_state=0)
    Extra Trees Pipeline for binary classification.

    name = 'Extra Trees Binary Classification Pipeline'
    custom_name = 'Extra Trees Binary Classification Pipeline'
    summary = 'Extra Trees Classifier w/ One Hot Encoder + Simple Imputer'
    component_graph = ['One Hot Encoder', 'Simple Imputer', 'Extra Trees Classifier']
    problem_type = 'binary'
    model_family = 'extra_trees'
    hyperparameters = {'Extra Trees Classifier': {'max_depth': Integer(low=4, high=10, p
    custom_hyperparameters = None
    default_parameters = {'Extra Trees Classifier': {'max_depth': 6, 'max_features': 'a
```

Instance attributes

feature_importance	Return importance associated with each feature.
parameters	Returns parameter dictionary for this pipeline
threshold	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
-----------------------	---

Continued on next page

Table 27 – continued from previous page

<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.ETBinaryClassificationPipeline.__init__`

`ETBinaryClassificationPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.ETBinaryClassificationPipeline.clone`

`ETBinaryClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.ETBinaryClassificationPipeline.describe`

`ETBinaryClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.ETBinaryClassificationPipeline.fit

`ETBinaryClassificationPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.ETBinaryClassificationPipeline.get_component

`ETBinaryClassificationPipeline.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ETBinaryClassificationPipeline.graph

`ETBinaryClassificationPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ETBinaryClassificationPipeline.graph_feature_importance

`ETBinaryClassificationPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ETBinaryClassificationPipeline.load

static `ETBinaryClassificationPipeline.load(file_path)`

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ETBinaryClassificationPipeline.predict

ETBinaryClassificationPipeline.**predict** (*X*, *objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.ETBinaryClassificationPipeline.predict_proba

ETBinaryClassificationPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.ETBinaryClassificationPipeline.save

ETBinaryClassificationPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns *None*

evalml.pipelines.ETBinaryClassificationPipeline.score

ETBinaryClassificationPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type *dict*

evalml.pipelines.ETMulticlassClassificationPipeline



```

class evalml.pipelines.ETMulticlassClassificationPipeline(parameters, random_state=0)
    Extra Trees Pipeline for multiclass classification.

    name = 'Extra Trees Multiclass Classification Pipeline'
    custom_name = 'Extra Trees Multiclass Classification Pipeline'
    summary = 'Extra Trees Classifier w/ One Hot Encoder + Simple Imputer'
    component_graph = ['One Hot Encoder', 'Simple Imputer', 'Extra Trees Classifier']
    problem_type = 'multiclass'
    model_family = 'extra_trees'
    hyperparameters = {'Extra Trees Classifier': {'max_depth': Integer(low=4, high=10, prior=6)}}
    custom_hyperparameters = None
    default_parameters = {'Extra Trees Classifier': {'max_depth': 6, 'max_features': 'auto'}}
  
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.ETMulticlassClassificationPipeline.__init__

ETMulticlassClassificationPipeline.__init__(parameters, random_state=0)

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: component_graph (list): List of components in order. Accepts strings or ComponentBase subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.ETMulticlassClassificationPipeline.clone

ETMulticlassClassificationPipeline.clone(random_state=0)

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a RandomState instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.ETMulticlassClassificationPipeline.describe

ETMulticlassClassificationPipeline.describe()

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if return_dict is True, else None

Return type dict

evalml.pipelines.ETMulticlassClassificationPipeline.fit

ETMulticlassClassificationPipeline.fit(X, y)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.ETMulticlassClassificationPipeline.get_component

ETMulticlassClassificationPipeline.**get_component** (*name*)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ETMulticlassClassificationPipeline.graph

ETMulticlassClassificationPipeline.**graph** (*filepath=None*)

Generate an image representing the pipeline graph

Parameters **filepath** (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ETMulticlassClassificationPipeline.graph_feature_importance

ETMulticlassClassificationPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ETMulticlassClassificationPipeline.load

static ETMulticlassClassificationPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ETMulticlassClassificationPipeline.predict

ETMulticlassClassificationPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.ETMulticlassClassificationPipeline.predict_probaETMulticlassClassificationPipeline.**predict_proba**(*X*)

Make probability estimates for labels.

Parameters *X* (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]**Returns** probability estimates**Return type** *pd.DataFrame***evalml.pipelines.ETMulticlassClassificationPipeline.save**ETMulticlassClassificationPipeline.**save**(*file_path*)

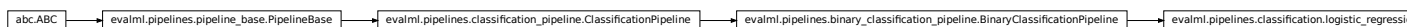
Saves pipeline at file path

Parameters *file_path* (*str*) – location to save file**Returns** None**evalml.pipelines.ETMulticlassClassificationPipeline.score**ETMulticlassClassificationPipeline.**score**(*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- *X* (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- *y* (*pd.Series*) – true labels of length [n_samples]
- *objectives* (*list*) – list of objectives to score

Returns ordered dictionary of objective scores**Return type** dict**evalml.pipelines.LogisticRegressionBinaryPipeline**

```
class evalml.pipelines.LogisticRegressionBinaryPipeline(parameters, random_state=0)
```

Logistic Regression Pipeline for binary classification.

name = 'Logistic Regression Binary Pipeline'**custom_name** = None**summary** = 'Logistic Regression Classifier w/ Simple Imputer + One Hot Encoder + Standard Scaler'**component_graph** = ['Simple Imputer', 'One Hot Encoder', 'Standard Scaler', 'Logistic Regression Classifier']**problem_type** = 'binary'**model_family** = 'linear_model'

```
hyperparameters = {'Logistic Regression Classifier': {'C': Real(low=0.01, high=10, pr
custom_hyperparameters = None
default_parameters = {'Logistic Regression Classifier': {'C': 1.0, 'n_jobs': -1, 'pe
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.LogisticRegressionBinaryPipeline.__init__`

`LogisticRegressionBinaryPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.LogisticRegressionBinaryPipeline.clone`

`LogisticRegressionBinaryPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.LogisticRegressionBinaryPipeline.describe`

`LogisticRegressionBinaryPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.LogisticRegressionBinaryPipeline.fit`

`LogisticRegressionBinaryPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

`evalml.pipelines.LogisticRegressionBinaryPipeline.get_component`

`LogisticRegressionBinaryPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.LogisticRegressionBinaryPipeline.graph`

`LogisticRegressionBinaryPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.LogisticRegressionBinaryPipeline.graph_feature_importance

LogisticRegressionBinaryPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.LogisticRegressionBinaryPipeline.load

static LogisticRegressionBinaryPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.LogisticRegressionBinaryPipeline.predict

LogisticRegressionBinaryPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.LogisticRegressionBinaryPipeline.predict_proba

LogisticRegressionBinaryPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.LogisticRegressionBinaryPipeline.save

LogisticRegressionBinaryPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.LogisticRegressionBinaryPipeline.score

`LogisticRegressionBinaryPipeline.score` (*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.LogisticRegressionMulticlassPipeline

```
class evalml.pipelines.LogisticRegressionMulticlassPipeline(parameters, random_state=0)
```

Logistic Regression Pipeline for multiclass classification.

name = 'Logistic Regression Multiclass Pipeline'

custom_name = None

summary = 'Logistic Regression Classifier w/ One Hot Encoder + Simple Imputer + Standard Scaler'

component_graph = ['One Hot Encoder', 'Simple Imputer', 'Standard Scaler', 'Logistic Regression Classifier']

problem_type = 'multiclass'

model_family = 'linear_model'

hyperparameters = {'Logistic Regression Classifier': {'C': Real(low=0.01, high=10, prior=1.0, steps=1000)}}

custom_hyperparameters = None

default_parameters = {'Logistic Regression Classifier': {'C': 1.0, 'n_jobs': -1, 'penalty': 'l2'}}

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.

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Table 33 – continued from previous page

<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline’s feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.LogisticRegressionMulticlassPipeline.__init__`

`LogisticRegressionMulticlassPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.LogisticRegressionMulticlassPipeline.clone`

`LogisticRegressionMulticlassPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.LogisticRegressionMulticlassPipeline.describe`

`LogisticRegressionMulticlassPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.LogisticRegressionMulticlassPipeline.fit

LogisticRegressionMulticlassPipeline.**fit** (*X*, *y*)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.LogisticRegressionMulticlassPipeline.get_component

LogisticRegressionMulticlassPipeline.**get_component** (*name*)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.LogisticRegressionMulticlassPipeline.graph

LogisticRegressionMulticlassPipeline.**graph** (*filepath=None*)

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.LogisticRegressionMulticlassPipeline.graph_feature_importance

LogisticRegressionMulticlassPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.LogisticRegressionMulticlassPipeline.load

static LogisticRegressionMulticlassPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.LogisticRegressionMulticlassPipeline.predict

LogisticRegressionMulticlassPipeline.**predict** (*X*, *objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.LogisticRegressionMulticlassPipeline.predict_proba

LogisticRegressionMulticlassPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.LogisticRegressionMulticlassPipeline.save

LogisticRegressionMulticlassPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns *None*

evalml.pipelines.LogisticRegressionMulticlassPipeline.score

LogisticRegressionMulticlassPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on objectives

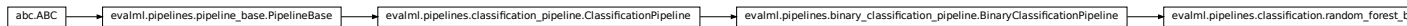
Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type *dict*

evalml.pipelines.RFBinaryClassificationPipeline



```

class evalml.pipelines.RFBinaryClassificationPipeline(parameters, random_state=0)
    Random Forest Pipeline for binary classification.

    name = 'Random Forest Binary Classification Pipeline'
    custom_name = 'Random Forest Binary Classification Pipeline'
    summary = 'Random Forest Classifier w/ Simple Imputer + One Hot Encoder'
    component_graph = ['Simple Imputer', 'One Hot Encoder', 'Random Forest Classifier']
    problem_type = 'binary'
    model_family = 'random_forest'
    hyperparameters = {'One Hot Encoder': {}, 'Random Forest Classifier': {'max_depth':
    custom_hyperparameters = None
    default_parameters = {'One Hot Encoder': {'categories': None, 'drop': None, 'handle

```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.RFBinaryClassificationPipeline.__init__`

`RFBinaryClassificationPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.RFBinaryClassificationPipeline.clone`

`RFBinaryClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.RFBinaryClassificationPipeline.describe`

`RFBinaryClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.RFBinaryClassificationPipeline.fit`

`RFBinaryClassificationPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- **y** (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

evalml.pipelines.RFBinaryClassificationPipeline.get_component

`RFBinaryClassificationPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.RFBinaryClassificationPipeline.graph

`RFBinaryClassificationPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.RFBinaryClassificationPipeline.graph_feature_importance

`RFBinaryClassificationPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.RFBinaryClassificationPipeline.load

static `RFBinaryClassificationPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.RFBinaryClassificationPipeline.predict

`RFBinaryClassificationPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.RFBinaryClassificationPipeline.predict_proba`RFBinaryClassificationPipeline.predict_proba(X)`

Make probability estimates for labels.

Parameters *X* (`pd.DataFrame` or `np.array`) – data of shape [n_samples, n_features]**Returns** probability estimates**Return type** `pd.DataFrame`**evalml.pipelines.RFBinaryClassificationPipeline.save**`RFBinaryClassificationPipeline.save(file_path)`

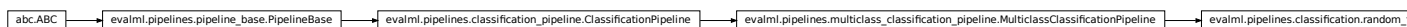
Saves pipeline at file path

Parameters *file_path* (`str`) – location to save file**Returns** `None`**evalml.pipelines.RFBinaryClassificationPipeline.score**`RFBinaryClassificationPipeline.score(X, y, objectives)`

Evaluate model performance on objectives

Parameters

- *X* (`pd.DataFrame` or `np.array`) – data of shape [n_samples, n_features]
- *y* (`pd.Series`) – true labels of length [n_samples]
- **objectives** (`list`) – list of objectives to score

Returns ordered dictionary of objective scores**Return type** `dict`**evalml.pipelines.RFMulticlassClassificationPipeline**

```
class evalml.pipelines.RFMulticlassClassificationPipeline(parameters, random_state=0)
    Random Forest Pipeline for multiclass classification.

    name = 'Random Forest Multiclass Classification Pipeline'
    custom_name = 'Random Forest Multiclass Classification Pipeline'
    summary = 'Random Forest Classifier w/ One Hot Encoder + Simple Imputer'
    component_graph = ['One Hot Encoder', 'Simple Imputer', 'Random Forest Classifier']
    problem_type = 'multiclass'
    model_family = 'random_forest'
```

```
hyperparameters = {'One Hot Encoder': {}, 'Random Forest Classifier': {'max_depth':
custom_hyperparameters = None
default_parameters = {'One Hot Encoder': {'categories': None, 'drop': None, 'handle
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.RFMulticlassClassificationPipeline.__init__`

`RFMulticlassClassificationPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.RFMulticlassClassificationPipeline.clone`

`RFMulticlassClassificationPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.RFMulticlassClassificationPipeline.describe`

`RFMulticlassClassificationPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.RFMulticlassClassificationPipeline.fit`

`RFMulticlassClassificationPipeline.fit(X, y)`

Build a model

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

`evalml.pipelines.RFMulticlassClassificationPipeline.get_component`

`RFMulticlassClassificationPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.RFMulticlassClassificationPipeline.graph`

`RFMulticlassClassificationPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.RFMulticlassClassificationPipeline.graph_feature_importance

RFMulticlassClassificationPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.RFMulticlassClassificationPipeline.load

static RFMulticlassClassificationPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.RFMulticlassClassificationPipeline.predict

RFMulticlassClassificationPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.RFMulticlassClassificationPipeline.predict_proba

RFMulticlassClassificationPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.RFMulticlassClassificationPipeline.save

RFMulticlassClassificationPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.RFMulticlassClassificationPipeline.score

`RFMulticlassClassificationPipeline.score` (*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.XGBoostBinaryPipeline



class evalml.pipelines.XGBoostBinaryPipeline (*parameters*, *random_state=0*)

XGBoost Pipeline for binary classification.

name = 'XGBoost Binary Classification Pipeline'

custom_name = 'XGBoost Binary Classification Pipeline'

summary = 'XGBoost Classifier w/ Simple Imputer + One Hot Encoder'

component_graph = ['Simple Imputer', 'One Hot Encoder', 'XGBoost Classifier']

problem_type = 'binary'

model_family = 'xgboost'

hyperparameters = {'One Hot Encoder': {}, 'Simple Imputer': {'impute_strategy': ['m', 'f', 'a']}}

custom_hyperparameters = None

default_parameters = {'One Hot Encoder': {'categories': None, 'drop': None, 'handle_unknown': 'ignore'}}

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
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Table 39 – continued from previous page

<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.XGBoostBinaryPipeline.__init__`

`XGBoostBinaryPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.XGBoostBinaryPipeline.clone`

`XGBoostBinaryPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.XGBoostBinaryPipeline.describe`

`XGBoostBinaryPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.XGBoostBinaryPipeline.fit`

`XGBoostBinaryPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

`evalml.pipelines.XGBoostBinaryPipeline.get_component`

`XGBoostBinaryPipeline.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.XGBoostBinaryPipeline.graph`

`XGBoostBinaryPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

`evalml.pipelines.XGBoostBinaryPipeline.graph_feature_importance`

`XGBoostBinaryPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

`evalml.pipelines.XGBoostBinaryPipeline.load`

static `XGBoostBinaryPipeline.load(file_path)`

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.XGBoostBinaryPipeline.predict

XGBoostBinaryPipeline.**predict** (*X*, *objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.XGBoostBinaryPipeline.predict_proba

XGBoostBinaryPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.XGBoostBinaryPipeline.save

XGBoostBinaryPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.XGBoostBinaryPipeline.score

XGBoostBinaryPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.XGBoostMulticlassPipeline



```

class evalml.pipelines.XGBoostMulticlassPipeline(parameters, random_state=0)
    XGBoost Pipeline for multiclass classification.

    name = 'XGBoost Multiclass Classification Pipeline'
    custom_name = 'XGBoost Multiclass Classification Pipeline'
    summary = 'XGBoost Classifier w/ One Hot Encoder + Simple Imputer'
    component_graph = ['One Hot Encoder', 'Simple Imputer', 'XGBoost Classifier']
    problem_type = 'multiclass'
    model_family = 'xgboost'
    hyperparameters = {'One Hot Encoder': {}, 'Simple Imputer': {'impute_strategy': ['m
    custom_hyperparameters = None
    default_parameters = {'One Hot Encoder': {'categories': None, 'drop': None, 'handle
  
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.XGBoostMulticlassPipeline.__init__

XGBoostMulticlassPipeline.**__init__**(*parameters*, *random_state*=0)

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: *component_graph* (list): List of components in order. Accepts strings or ComponentBase subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.XGBoostMulticlassPipeline.clone

XGBoostMulticlassPipeline.**clone**(*random_state*=0)

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a RandomState instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.XGBoostMulticlassPipeline.describe

XGBoostMulticlassPipeline.**describe**()

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if return_dict is True, else None

Return type dict

evalml.pipelines.XGBoostMulticlassPipeline.fit

XGBoostMulticlassPipeline.**fit**(*X*, *y*)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.XGBoostMulticlassPipeline.get_component

`XGBoostMulticlassPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.XGBoostMulticlassPipeline.graph

`XGBoostMulticlassPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.XGBoostMulticlassPipeline.graph_feature_importance

`XGBoostMulticlassPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.XGBoostMulticlassPipeline.load

static `XGBoostMulticlassPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.XGBoostMulticlassPipeline.predict

`XGBoostMulticlassPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.XGBoostMulticlassPipeline.predict_proba

XGBoostMulticlassPipeline.**predict_proba**(X)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.XGBoostMulticlassPipeline.save

XGBoostMulticlassPipeline.**save**(file_path)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.XGBoostMulticlassPipeline.score

XGBoostMulticlassPipeline.**score**(X, y, objectives)

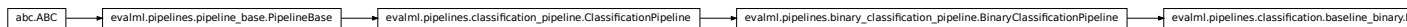
Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.BaselineBinaryPipeline

```
class evalml.pipelines.BaselineBinaryPipeline(parameters, random_state=0)
```

Baseline Pipeline for binary classification.

```
name = 'Baseline Classification Pipeline'
```

```
custom_name = 'Baseline Classification Pipeline'
```

```
summary = 'Baseline Classifier'
```

```
component_graph = ['Baseline Classifier']
```

```
problem_type = 'binary'
```

```
model_family = 'baseline'
```

```
hyperparameters = {'Baseline Classifier': {}}
```

```
custom_hyperparameters = None
default_parameters = {'Baseline Classifier': {'strategy': 'mode'}}
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.BaselineBinaryPipeline.__init__`

`BaselineBinaryPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int, np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.BaselineBinaryPipeline.clone`

`BaselineBinaryPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.BaselineBinaryPipeline.describe`

`BaselineBinaryPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.BaselineBinaryPipeline.fit`

`BaselineBinaryPipeline.fit(X, y)`

Build a model

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

`evalml.pipelines.BaselineBinaryPipeline.get_component`

`BaselineBinaryPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.BaselineBinaryPipeline.graph`

`BaselineBinaryPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

`evalml.pipelines.BaselineBinaryPipeline.graph_feature_importance`

`BaselineBinaryPipeline.graph_feature_importance` (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

`evalml.pipelines.BaselineBinaryPipeline.load`

static `BaselineBinaryPipeline.load` (*file_path*)

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

`evalml.pipelines.BaselineBinaryPipeline.predict`

`BaselineBinaryPipeline.predict` (*X, objective=None*)

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.BaselineBinaryPipeline.predict_proba`

`BaselineBinaryPipeline.predict_proba` (*X*)

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type `pd.DataFrame`

`evalml.pipelines.BaselineBinaryPipeline.save`

`BaselineBinaryPipeline.save` (*file_path*)

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

evalml.pipelines.BaselineBinaryPipeline.score

`BaselineBinaryPipeline.score` (*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.BaselineMulticlassPipeline

class evalml.pipelines.**BaselineMulticlassPipeline** (*parameters*, *random_state=0*)

Baseline Pipeline for multiclass classification.

name = 'Baseline Multiclass Classification Pipeline'

custom_name = 'Baseline Multiclass Classification Pipeline'

summary = 'Baseline Classifier'

component_graph = ['Baseline Classifier']

problem_type = 'multiclass'

model_family = 'baseline'

hyperparameters = {'Baseline Classifier': {}}

custom_hyperparameters = None

default_parameters = {'Baseline Classifier': {'strategy': 'mode'}}

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.

Continued on next page

Table 45 – continued from previous page

<i>describe</i>	Outputs pipeline details including component parameters
<i>fit</i>	Build a model
<i>get_component</i>	Returns component by name
<i>graph</i>	Generate an image representing the pipeline graph
<i>graph_feature_importance</i>	Generate a bar graph of the pipeline’s feature importance
<i>load</i>	Loads pipeline at file path
<i>predict</i>	Make predictions using selected features.
<i>predict_proba</i>	Make probability estimates for labels.
<i>save</i>	Saves pipeline at file path
<i>score</i>	Evaluate model performance on objectives

`evalml.pipelines.BaselineMulticlassPipeline.__init__`

`BaselineMulticlassPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.BaselineMulticlassPipeline.clone`

`BaselineMulticlassPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.BaselineMulticlassPipeline.describe`

`BaselineMulticlassPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.BaselineMulticlassPipeline.fit`BaselineMulticlassPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self**evalml.pipelines.BaselineMulticlassPipeline.get_component**`BaselineMulticlassPipeline.get_component(name)`

Returns component by name

Parameters **name** (*str*) – name of component**Returns** component to return**Return type** Component**evalml.pipelines.BaselineMulticlassPipeline.graph**`BaselineMulticlassPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.**Returns** Graph object that can be directly displayed in Jupyter notebooks.**Return type** graphviz.Digraph**evalml.pipelines.BaselineMulticlassPipeline.graph_feature_importance**`BaselineMulticlassPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.**Returns** plotly.Figure, a bar graph showing features and their corresponding importance**evalml.pipelines.BaselineMulticlassPipeline.load****static** `BaselineMulticlassPipeline.load(file_path)`

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file**Returns** PipelineBase object

evalml.pipelines.BaselineMulticlassPipeline.predict

BaselineMulticlassPipeline.**predict** (*X*, *objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.BaselineMulticlassPipeline.predict_proba

BaselineMulticlassPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.BaselineMulticlassPipeline.save

BaselineMulticlassPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.BaselineMulticlassPipeline.score

BaselineMulticlassPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.ModeBaselineBinaryPipeline



```

class evalml.pipelines.ModeBaselineBinaryPipeline(parameters, random_state=0)
    Mode Baseline Pipeline for binary classification.

    name = 'Mode Baseline Binary Classification Pipeline'
    custom_name = 'Mode Baseline Binary Classification Pipeline'
    summary = 'Baseline Classifier'
    component_graph = ['Baseline Classifier']
    problem_type = 'binary'
    model_family = 'baseline'
    hyperparameters = {'Baseline Classifier': {}}
    custom_hyperparameters = {'strategy': ['mode']}
    default_parameters = {'Baseline Classifier': {'strategy': 'mode'}}

```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline
<code>threshold</code>	

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

evalml.pipelines.ModeBaselineBinaryPipeline.__init__

ModeBaselineBinaryPipeline.__init__(parameters, random_state=0)

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: component_graph (list): List of components in order. Accepts strings or ComponentBase subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.ModeBaselineBinaryPipeline.clone

ModeBaselineBinaryPipeline.clone(random_state=0)

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a RandomState instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.ModeBaselineBinaryPipeline.describe

ModeBaselineBinaryPipeline.describe()

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if return_dict is True, else None

Return type dict

evalml.pipelines.ModeBaselineBinaryPipeline.fit

ModeBaselineBinaryPipeline.fit(X, y)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.ModeBaselineBinaryPipeline.get_component

`ModeBaselineBinaryPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ModeBaselineBinaryPipeline.graph

`ModeBaselineBinaryPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ModeBaselineBinaryPipeline.graph_feature_importance

`ModeBaselineBinaryPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ModeBaselineBinaryPipeline.load

static `ModeBaselineBinaryPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ModeBaselineBinaryPipeline.predict

`ModeBaselineBinaryPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

`evalml.pipelines.ModeBaselineBinaryPipeline.predict_proba`

`ModeBaselineBinaryPipeline.predict_proba(X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type `pd.DataFrame`

`evalml.pipelines.ModeBaselineBinaryPipeline.save`

`ModeBaselineBinaryPipeline.save(file_path)`

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns `None`

`evalml.pipelines.ModeBaselineBinaryPipeline.score`

`ModeBaselineBinaryPipeline.score(X, y, objectives)`

Evaluate model performance on objectives

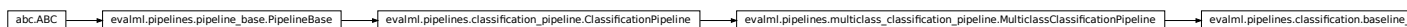
Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

Returns ordered dictionary of objective scores

Return type `dict`

`evalml.pipelines.ModeBaselineMulticlassPipeline`



```
class evalml.pipelines.ModeBaselineMulticlassPipeline(parameters, random_state=0)
```

Mode Baseline Pipeline for multiclass classification.

```
name = 'Mode Baseline Multiclass Classification Pipeline'
```

```
custom_name = 'Mode Baseline Multiclass Classification Pipeline'
```

```
summary = 'Baseline Classifier'
```

```
component_graph = ['Baseline Classifier']
```

```
problem_type = 'multiclass'
```

```
model_family = 'baseline'
```

```
hyperparameters = {'Baseline Classifier': {}}
custom_hyperparameters = {'strategy': ['mode']}
default_parameters = {'Baseline Classifier': {'strategy': 'mode'}}
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on objectives

`evalml.pipelines.ModeBaselineMulticlassPipeline.__init__`

`ModeBaselineMulticlassPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.ModeBaselineMulticlassPipeline.clone`

`ModeBaselineMulticlassPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.ModeBaselineMulticlassPipeline.describe`

`ModeBaselineMulticlassPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.ModeBaselineMulticlassPipeline.fit`

`ModeBaselineMulticlassPipeline.fit(X, y)`

Build a model

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

`evalml.pipelines.ModeBaselineMulticlassPipeline.get_component`

`ModeBaselineMulticlassPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.ModeBaselineMulticlassPipeline.graph`

`ModeBaselineMulticlassPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ModeBaselineMulticlassPipeline.graph_feature_importance

ModeBaselineMulticlassPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ModeBaselineMulticlassPipeline.load

static ModeBaselineMulticlassPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ModeBaselineMulticlassPipeline.predict

ModeBaselineMulticlassPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.ModeBaselineMulticlassPipeline.predict_proba

ModeBaselineMulticlassPipeline.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.ModeBaselineMulticlassPipeline.save

ModeBaselineMulticlassPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.ModeBaselineMulticlassPipeline.score

ModeBaselineMulticlassPipeline.**score**(*X*, *y*, *objectives*)

Evaluate model performance on objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – list of objectives to score

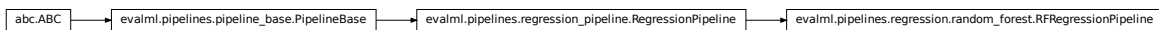
Returns ordered dictionary of objective scores

Return type dict

Regression Pipelines

<i>RFRegressionPipeline</i>	Random Forest Pipeline for regression problems.
<i>CatBoostRegressionPipeline</i>	CatBoost Pipeline for regression problems.
<i>ENRegressionPipeline</i>	Elastic Net Pipeline for regression problems.
<i>ETRegressionPipeline</i>	Extra Trees Pipeline for regression problems.
<i>LinearRegressionPipeline</i>	Linear Regression Pipeline for regression problems.
<i>XGBoostRegressionPipeline</i>	XGBoost Pipeline for regression problems.
<i>BaselineRegressionPipeline</i>	Baseline Pipeline for regression problems.
<i>MeanBaselineRegressionPipeline</i>	Baseline Pipeline for regression problems.

evalml.pipelines.RFRegressionPipeline



class evalml.pipelines.**RFRegressionPipeline**(*parameters*, *random_state=0*)

Random Forest Pipeline for regression problems.

name = 'Random Forest Regression Pipeline'

custom_name = 'Random Forest Regression Pipeline'

summary = 'Random Forest Regressor w/ One Hot Encoder + Simple Imputer'

component_graph = ['One Hot Encoder', 'Simple Imputer', 'Random Forest Regressor']

problem_type = 'regression'

model_family = 'random_forest'

hyperparameters = {'One Hot Encoder': {}, 'Random Forest Regressor': {'max_depth':

custom_hyperparameters = None

default_parameters = {'One Hot Encoder': {'categories': None, 'drop': None, 'handle

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

`evalml.pipelines.RFRegressionPipeline.__init__`

`RFRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.RFRegressionPipeline.clone`

`RFRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.RFRegressionPipeline.describe`

`RFRegressionPipeline.describe()`

Outputs pipeline details including component parameters

Parameters `return_dict` (*bool*) – If True, return dictionary of information about pipeline.
Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.RFRegressionPipeline.fit`

`RFRegressionPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

`evalml.pipelines.RFRegressionPipeline.get_component`

`RFRegressionPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

`evalml.pipelines.RFRegressionPipeline.graph`

`RFRegressionPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

`evalml.pipelines.RFRegressionPipeline.graph_feature_importance`

`RFRegressionPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

`evalml.pipelines.RFRegressionPipeline.load`

static `RFRegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

`evalml.pipelines.RFRegressionPipeline.predict`

`RFRegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `objective` (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.RFRegressionPipeline.save`

`RFRegressionPipeline.save(file_path)`

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

`evalml.pipelines.RFRegressionPipeline.score`

`RFRegressionPipeline.score(X, y, objectives)`

Evaluate model performance on current and additional objectives

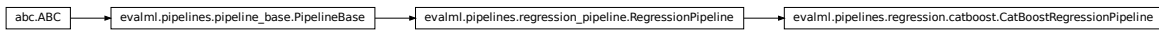
Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – true labels of length `[n_samples]`
- `objectives` (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type `dict`

evalml.pipelines.CatBoostRegressionPipeline



class evalml.pipelines.CatBoostRegressionPipeline (parameters, random_state=0)

CatBoost Pipeline for regression problems. CatBoost is an open-source library and natively supports categorical features.

For more information, check out <https://catboost.ai/>

Note: impute_strategy must support both string and numeric data

name = 'Cat Boost Regression Pipeline'

custom_name = None

summary = 'CatBoost Regressor w/ Simple Imputer'

component_graph = ['Simple Imputer', 'CatBoost Regressor']

problem_type = 'regression'

model_family = 'catboost'

hyperparameters = {'CatBoost Regressor': {'eta': Real(low=1e-06, high=1, prior='unif

custom_hyperparameters = {'Simple Imputer': {'impute_strategy': ['most_frequent']}}

default_parameters = {'CatBoost Regressor': {'bootstrap_type': None, 'eta': 0.03, 'l

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path

Continued on next page

Table 54 – continued from previous page

<i>score</i>	Evaluate model performance on current and additional objectives
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evalml.pipelines.CatBoostRegressionPipeline.__init__

`CatBoostRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.CatBoostRegressionPipeline.clone

`CatBoostRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.CatBoostRegressionPipeline.describe

`CatBoostRegressionPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.CatBoostRegressionPipeline.fit

`CatBoostRegressionPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.CatBoostRegressionPipeline.get_component

`CatBoostRegressionPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.CatBoostRegressionPipeline.graph

`CatBoostRegressionPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.CatBoostRegressionPipeline.graph_feature_importance

`CatBoostRegressionPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.CatBoostRegressionPipeline.load

static `CatBoostRegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.CatBoostRegressionPipeline.predict

`CatBoostRegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.CatBoostRegressionPipeline.save

`CatBoostRegressionPipeline.save` (*file_path*)

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns None

evalml.pipelines.CatBoostRegressionPipeline.score

`CatBoostRegressionPipeline.score` (*X*, *y*, *objectives*)

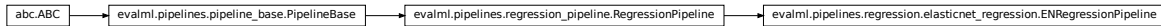
Evaluate model performance on current and additional objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.ENRegressionPipeline

class `evalml.pipelines.ENRegressionPipeline` (*parameters*, *random_state=0*)

Elastic Net Pipeline for regression problems.

name = 'ENRegression Pipeline'

custom_name = None

summary = 'Elastic Net Regressor w/ One Hot Encoder + Simple Imputer'

component_graph = ['One Hot Encoder', 'Simple Imputer', 'Elastic Net Regressor']

problem_type = 'regression'

model_family = 'linear_model'

hyperparameters = {'Elastic Net Regressor': {'alpha': Real(low=0, high=1, prior='uni

custom_hyperparameters = None

default_parameters = {'Elastic Net Regressor': {'alpha': 0.5, 'l1_ratio': 0.5, 'max

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
---------------------------------	---

Continued on next page

Table 55 – continued from previous page

<code>parameters</code>	Returns parameter dictionary for this pipeline
Methods:	
<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline’s feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

evalml.pipelines.ENRegressionPipeline.__init__

ENRegressionPipeline.**__init__**(*parameters*, *random_state=0*)

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or ComponentBase subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.ENRegressionPipeline.clone

ENRegressionPipeline.**clone**(*random_state=0*)

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a RandomState instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.ENRegressionPipeline.describe

ENRegressionPipeline.**describe**()

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline.
Defaults to false

Returns dictionary of all component parameters if return_dict is True, else None

Return type dict

evalml.pipelines.ENRegressionPipeline.fit

ENRegressionPipeline.**fit**(*X*, *y*)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.ENRegressionPipeline.get_component

ENRegressionPipeline.**get_component**(*name*)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ENRegressionPipeline.graph

ENRegressionPipeline.**graph**(*filepath=None*)

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ENRegressionPipeline.graph_feature_importance

ENRegressionPipeline.**graph_feature_importance**(*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

`evalml.pipelines.ENRegressionPipeline.load`

static `ENRegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

`evalml.pipelines.ENRegressionPipeline.predict`

`ENRegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `objective` (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.ENRegressionPipeline.save`

`ENRegressionPipeline.save(file_path)`

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

`evalml.pipelines.ENRegressionPipeline.score`

`ENRegressionPipeline.score(X, y, objectives)`

Evaluate model performance on current and additional objectives

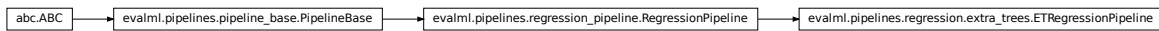
Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – true labels of length `[n_samples]`
- `objectives` (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type `dict`

evalml.pipelines.ETRegressionPipeline



```

class evalml.pipelines.ETRegressionPipeline(parameters, random_state=0)
    Extra Trees Pipeline for regression problems.

    name = 'Extra Trees Regression Pipeline'
    custom_name = 'Extra Trees Regression Pipeline'
    summary = 'Extra Trees Regressor w/ One Hot Encoder + Simple Imputer'
    component_graph = ['One Hot Encoder', 'Simple Imputer', 'Extra Trees Regressor']
    problem_type = 'regression'
    model_family = 'extra_trees'

    hyperparameters = {'Extra Trees Regressor': {'max_depth': Integer(low=4, high=10, pr
    custom_hyperparameters = None
    default_parameters = {'Extra Trees Regressor': {'max_depth': 6, 'max_features': 'au
  
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

`evalml.pipelines.ETRegressionPipeline.__init__`

`ETRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.ETRegressionPipeline.clone`

`ETRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.ETRegressionPipeline.describe`

`ETRegressionPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.ETRegressionPipeline.fit`

`ETRegressionPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- **y** (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

evalml.pipelines.ETRegressionPipeline.get_component

ETRegressionPipeline.**get_component** (*name*)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.ETRegressionPipeline.graph

ETRegressionPipeline.**graph** (*filepath=None*)

Generate an image representing the pipeline graph

Parameters **filepath** (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.ETRegressionPipeline.graph_feature_importance

ETRegressionPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.ETRegressionPipeline.load

static ETRegressionPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.ETRegressionPipeline.predict

ETRegressionPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.ETRegressionPipeline.save

ETRegressionPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.ETRegressionPipeline.score

ETRegressionPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on current and additional objectives

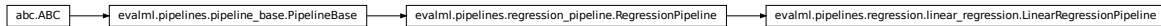
Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.LinearRegressionPipeline



class evalml.pipelines.**LinearRegressionPipeline** (*parameters*, *random_state=0*)

Linear Regression Pipeline for regression problems.

name = 'Linear Regression Pipeline'

custom_name = None

summary = 'Linear Regressor w/ One Hot Encoder + Simple Imputer + Standard Scaler'

component_graph = ['One Hot Encoder', 'Simple Imputer', 'Standard Scaler', 'Linear Reg'

problem_type = 'regression'

model_family = 'linear_model'

hyperparameters = {'Linear Regressor': {'fit_intercept': [True, False], 'normalize':

custom_hyperparameters = None

default_parameters = {'Linear Regressor': {'fit_intercept': True, 'n_jobs': -1, 'no

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
---------------------------------	---

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Table 59 – continued from previous page

<code>parameters</code>	Returns parameter dictionary for this pipeline
Methods:	
<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

evalml.pipelines.LinearRegressionPipeline.__init__

`LinearRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.LinearRegressionPipeline.clone

`LinearRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.LinearRegressionPipeline.describe

LinearRegressionPipeline.**describe**()

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline.
Defaults to false

Returns dictionary of all component parameters if return_dict is True, else None

Return type dict

evalml.pipelines.LinearRegressionPipeline.fit

LinearRegressionPipeline.**fit**(X, y)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.LinearRegressionPipeline.get_component

LinearRegressionPipeline.**get_component**(name)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.LinearRegressionPipeline.graph

LinearRegressionPipeline.**graph**(filepath=None)

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.LinearRegressionPipeline.graph_feature_importance

LinearRegressionPipeline.**graph_feature_importance**(show_all_features=False)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

`evalml.pipelines.LinearRegressionPipeline.load`

static `LinearRegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

`evalml.pipelines.LinearRegressionPipeline.predict`

`LinearRegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `objective` (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.LinearRegressionPipeline.save`

`LinearRegressionPipeline.save(file_path)`

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

`evalml.pipelines.LinearRegressionPipeline.score`

`LinearRegressionPipeline.score(X, y, objectives)`

Evaluate model performance on current and additional objectives

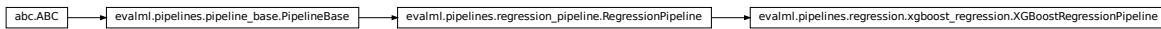
Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – true labels of length `[n_samples]`
- `objectives` (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type `dict`

evalml.pipelines.XGBoostRegressionPipeline



```

class evalml.pipelines.XGBoostRegressionPipeline(parameters, random_state=0)
    XGBoost Pipeline for regression problems.

    name = 'XGBoost Regression Pipeline'

    custom_name = None

    summary = 'XGBoost Regressor w/ One Hot Encoder + Simple Imputer'

    component_graph = ['One Hot Encoder', 'Simple Imputer', 'XGBoost Regressor']

    problem_type = 'regression'

    model_family = 'xgboost'

    hyperparameters = {'One Hot Encoder': {}, 'Simple Imputer': {'impute_strategy': ['m
    custom_hyperparameters = None

    default_parameters = {'One Hot Encoder': {'categories': None, 'drop': None, 'handle
  
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

evalml.pipelines.XGBoostRegressionPipeline.__init__

`XGBoostRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.XGBoostRegressionPipeline.clone

`XGBoostRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.XGBoostRegressionPipeline.describe

`XGBoostRegressionPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

evalml.pipelines.XGBoostRegressionPipeline.fit

`XGBoostRegressionPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.XGBoostRegressionPipeline.get_component

`XGBoostRegressionPipeline.get_component(name)`

Returns component by name

Parameters `name` (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.XGBoostRegressionPipeline.graph

`XGBoostRegressionPipeline.graph(filepath=None)`

Generate an image representing the pipeline graph

Parameters `filepath` (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.XGBoostRegressionPipeline.graph_feature_importance

`XGBoostRegressionPipeline.graph_feature_importance(show_all_features=False)`

Generate a bar graph of the pipeline's feature importance

Parameters `show_all_features` (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.XGBoostRegressionPipeline.load

static `XGBoostRegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.XGBoostRegressionPipeline.predict

`XGBoostRegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- `objective` (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.XGBoostRegressionPipeline.save

XGBoostRegressionPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.XGBoostRegressionPipeline.score

XGBoostRegressionPipeline.**score** (*X*, *y*, *objectives*)

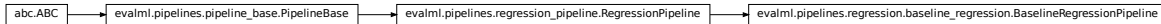
Evaluate model performance on current and additional objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type dict

evalml.pipelines.BaselineRegressionPipeline

class evalml.pipelines.**BaselineRegressionPipeline** (*parameters*, *random_state=0*)

Baseline Pipeline for regression problems.

name = 'Baseline Regression Pipeline'

custom_name = None

summary = 'Baseline Regressor'

component_graph = ['Baseline Regressor']

problem_type = 'regression'

model_family = 'baseline'

hyperparameters = {'Baseline Regressor': {}}

custom_hyperparameters = None

default_parameters = {'Baseline Regressor': {'strategy': 'mean'}}

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
---------------------------------	---

Continued on next page

Table 63 – continued from previous page

<code>parameters</code>	Returns parameter dictionary for this pipeline
Methods:	
<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline’s feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

evalml.pipelines.BaselineRegressionPipeline.__init__

`BaselineRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component’s parameters as values. An empty dictionary {} implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

evalml.pipelines.BaselineRegressionPipeline.clone

`BaselineRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

evalml.pipelines.BaselineRegressionPipeline.describe

BaselineRegressionPipeline.**describe**()

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline.
Defaults to false

Returns dictionary of all component parameters if return_dict is True, else None

Return type dict

evalml.pipelines.BaselineRegressionPipeline.fit

BaselineRegressionPipeline.**fit**(X, y)

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.BaselineRegressionPipeline.get_component

BaselineRegressionPipeline.**get_component**(name)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.BaselineRegressionPipeline.graph

BaselineRegressionPipeline.**graph**(filepath=None)

Generate an image representing the pipeline graph

Parameters **filepath** (*str*, *optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.BaselineRegressionPipeline.graph_feature_importance

BaselineRegressionPipeline.**graph_feature_importance**(show_all_features=False)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool*, *optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns `plotly.Figure`, a bar graph showing features and their corresponding importance

`evalml.pipelines.BaselineRegressionPipeline.load`

static `BaselineRegressionPipeline.load(file_path)`

Loads pipeline at file path

Parameters `file_path` (*str*) – location to load file

Returns `PipelineBase` object

`evalml.pipelines.BaselineRegressionPipeline.predict`

`BaselineRegressionPipeline.predict(X, objective=None)`

Make predictions using selected features.

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `objective` (*Object* or *string*) – the objective to use to make predictions

Returns estimated labels

Return type `pd.Series`

`evalml.pipelines.BaselineRegressionPipeline.save`

`BaselineRegressionPipeline.save(file_path)`

Saves pipeline at file path

Parameters `file_path` (*str*) – location to save file

Returns `None`

`evalml.pipelines.BaselineRegressionPipeline.score`

`BaselineRegressionPipeline.score(X, y, objectives)`

Evaluate model performance on current and additional objectives

Parameters

- `X` (*pd.DataFrame* or *np.array*) – data of shape `[n_samples, n_features]`
- `y` (*pd.Series*) – true labels of length `[n_samples]`
- `objectives` (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type `dict`

evalml.pipelines.MeanBaselineRegressionPipeline



```
class evalml.pipelines.MeanBaselineRegressionPipeline(parameters, random_state=0)
```

Baseline Pipeline for regression problems.

```
name = 'Mean Baseline Regression Pipeline'
```

```
custom_name = None
```

```
summary = 'Baseline Regressor'
```

```
component_graph = ['Baseline Regressor']
```

```
problem_type = 'regression'
```

```
model_family = 'baseline'
```

```
hyperparameters = {'Baseline Regressor': {}}
```

```
custom_hyperparameters = {'strategy': ['mean']}
```

```
default_parameters = {'Baseline Regressor': {'strategy': 'mean'}}
```

Instance attributes

<code>feature_importance</code>	Return importance associated with each feature.
<code>parameters</code>	Returns parameter dictionary for this pipeline

Methods:

<code>__init__</code>	Machine learning pipeline made out of transformers and a estimator.
<code>clone</code>	Constructs a new pipeline with the same parameters and components.
<code>describe</code>	Outputs pipeline details including component parameters
<code>fit</code>	Build a model
<code>get_component</code>	Returns component by name
<code>graph</code>	Generate an image representing the pipeline graph
<code>graph_feature_importance</code>	Generate a bar graph of the pipeline's feature importance
<code>load</code>	Loads pipeline at file path
<code>predict</code>	Make predictions using selected features.
<code>save</code>	Saves pipeline at file path
<code>score</code>	Evaluate model performance on current and additional objectives

`evalml.pipelines.MeanBaselineRegressionPipeline.__init__`

`MeanBaselineRegressionPipeline.__init__(parameters, random_state=0)`

Machine learning pipeline made out of transformers and a estimator.

Required Class Variables: `component_graph` (list): List of components in order. Accepts strings or `ComponentBase` subclasses in the list

Parameters

- **parameters** (*dict*) – dictionary with component names as keys and dictionary of that component's parameters as values. An empty dictionary `{}` implies using all default values for component parameters.
- **random_state** (*int*, *np.random.RandomState*) – The random seed/state. Defaults to 0.

`evalml.pipelines.MeanBaselineRegressionPipeline.clone`

`MeanBaselineRegressionPipeline.clone(random_state=0)`

Constructs a new pipeline with the same parameters and components.

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this pipeline with identical parameters and components

`evalml.pipelines.MeanBaselineRegressionPipeline.describe`

`MeanBaselineRegressionPipeline.describe()`

Outputs pipeline details including component parameters

Parameters **return_dict** (*bool*) – If True, return dictionary of information about pipeline. Defaults to false

Returns dictionary of all component parameters if `return_dict` is True, else None

Return type dict

`evalml.pipelines.MeanBaselineRegressionPipeline.fit`

`MeanBaselineRegressionPipeline.fit(X, y)`

Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape `[n_samples, n_features]`
- **y** (*pd.Series*) – the target training labels of length `[n_samples]`

Returns self

evalml.pipelines.MeanBaselineRegressionPipeline.get_component

MeanBaselineRegressionPipeline.**get_component** (*name*)

Returns component by name

Parameters **name** (*str*) – name of component

Returns component to return

Return type Component

evalml.pipelines.MeanBaselineRegressionPipeline.graph

MeanBaselineRegressionPipeline.**graph** (*filepath=None*)

Generate an image representing the pipeline graph

Parameters **filepath** (*str, optional*) – Path to where the graph should be saved. If set to None (as by default), the graph will not be saved.

Returns Graph object that can be directly displayed in Jupyter notebooks.

Return type graphviz.Digraph

evalml.pipelines.MeanBaselineRegressionPipeline.graph_feature_importance

MeanBaselineRegressionPipeline.**graph_feature_importance** (*show_all_features=False*)

Generate a bar graph of the pipeline's feature importance

Parameters **show_all_features** (*bool, optional*) – If true, graph features with an importance value of zero. Defaults to false.

Returns plotly.Figure, a bar graph showing features and their corresponding importance

evalml.pipelines.MeanBaselineRegressionPipeline.load

static MeanBaselineRegressionPipeline.**load** (*file_path*)

Loads pipeline at file path

Parameters **file_path** (*str*) – location to load file

Returns PipelineBase object

evalml.pipelines.MeanBaselineRegressionPipeline.predict

MeanBaselineRegressionPipeline.**predict** (*X, objective=None*)

Make predictions using selected features.

Parameters

- **X** (*pd.DataFrame or np.array*) – data of shape [n_samples, n_features]
- **objective** (*Object or string*) – the objective to use to make predictions

Returns estimated labels

Return type pd.Series

evalml.pipelines.MeanBaselineRegressionPipeline.save

MeanBaselineRegressionPipeline.**save** (*file_path*)

Saves pipeline at file path

Parameters **file_path** (*str*) – location to save file

Returns None

evalml.pipelines.MeanBaselineRegressionPipeline.score

MeanBaselineRegressionPipeline.**score** (*X*, *y*, *objectives*)

Evaluate model performance on current and additional objectives

Parameters

- **X** (*pd.DataFrame* or *np.array*) – data of shape [n_samples, n_features]
- **y** (*pd.Series*) – true labels of length [n_samples]
- **objectives** (*list*) – Non-empty list of objectives to score on

Returns ordered dictionary of objective scores

Return type dict

Pipeline Utils

<i>all_pipelines</i>	Returns a complete list of all supported pipeline classes.
<i>get_pipelines</i>	Returns the pipelines allowed for a particular problem type.
<i>get_estimators</i>	Returns the estimators allowed for a particular problem type.
<i>make_pipeline</i>	Given input data, target data, an estimator class and the problem type,
<i>list_model_families</i>	List model type for a particular problem type.

evalml.pipelines.all_pipelines

evalml.pipelines.**all_pipelines** ()

Returns a complete list of all supported pipeline classes.

Returns a list of pipeline classes

Return type list[*PipelineBase*]

evalml.pipelines.get_pipelines

evalml.pipelines.**get_pipelines** (*problem_type*, *model_families=None*)

Returns the pipelines allowed for a particular problem type.

Can also optionally filter by a list of model types.

Parameters

- **problem_type** (*ProblemTypes* or *str*) – problem type to filter for

- **model_families** (*list[ModelFamily]* or *list[str]*) – model families to filter for

Returns a list of pipeline classes

Return type *list[PipelineBase]*

evalml.pipelines.get_estimators

`evalml.pipelines.get_estimators(problem_type, model_families=None)`

Returns the estimators allowed for a particular problem type.

Can also optionally filter by a list of model types.

Parameters

- **problem_type** (*ProblemTypes* or *str*) – problem type to filter for
- **model_families** (*list[ModelFamily]* or *list[str]*) – model families to filter for

Returns a list of estimator subclasses

Return type *list[class]*

evalml.pipelines.make_pipeline

`evalml.pipelines.make_pipeline(X, y, estimator, problem_type)`

Given input data, target data, an estimator class and the problem type, generates a pipeline class with a preprocessing chain which was recommended based on the inputs. The pipeline will be a subclass of the appropriate pipeline base class for the specified *problem_type*.

Parameters

- **X** (*pd.DataFrame*) – the input data of shape *[n_samples, n_features]*
- **y** (*pd.Series*) – the target labels of length *[n_samples]*
- **estimator** (*Estimator*) – estimator for pipeline
- **problem_type** – problem type for pipeline to generate

evalml.pipelines.list_model_families

`evalml.pipelines.list_model_families(problem_type)`

List model type for a particular problem type.

Parameters **problem_types** (*ProblemTypes* or *str*) – binary, multiclass, or regression

Returns a list of model families

Return type *list[ModelFamily]*

Pipeline Graph Utils

<code>precision_recall_curve</code>	Given labels and binary classifier predicted probabilities, compute and return the data representing a precision-recall curve.
<code>graph_precision_recall_curve</code>	Generate and display a precision-recall plot.
<code>roc_curve</code>	Given labels and classifier predicted probabilities, compute and return the data representing a Receiver Operating Characteristic (ROC) curve.
<code>graph_roc_curve</code>	Generate and display a Receiver Operating Characteristic (ROC) plot.
<code>confusion_matrix</code>	Confusion matrix for binary and multiclass classification.
<code>normalize_confusion_matrix</code>	Normalizes a confusion matrix.
<code>graph_confusion_matrix</code>	Generate and display a confusion matrix plot.
<code>calculate_permutation_importance</code>	Calculates permutation importance for features.
<code>graph_permutation_importance</code>	Generate a bar graph of the pipeline's permutation importance.

evalml.pipelines.precision_recall_curve

`evalml.pipelines.precision_recall_curve(y_true, y_pred_proba)`

Given labels and binary classifier predicted probabilities, compute and return the data representing a precision-recall curve.

Parameters

- **y_true** (*pd.Series* or *np.array*) – true binary labels.
- **y_pred_proba** (*pd.Series* or *np.array*) – predictions from a binary classifier, before thresholding has been applied. Note this should be the predicted probability for the “true” label.

Returns

Dictionary containing metrics used to generate a precision-recall plot, with the following keys:

- *precision*: Precision values.
- *recall*: Recall values.
- *thresholds*: Threshold values used to produce the precision and recall.
- *auc_score*: The area under the ROC curve.

Return type list

evalml.pipelines.graph_precision_recall_curve

`evalml.pipelines.graph_precision_recall_curve(y_true, y_pred_proba, title_addition=None)`

Generate and display a precision-recall plot.

Parameters

- **y_true** (*pd.Series* or *np.array*) – true binary labels.
- **y_pred_proba** (*pd.Series* or *np.array*) – predictions from a binary classifier, before thresholding has been applied. Note this should be the predicted probability for the “true” label.

- **title_addition** (*str* or *None*) – if not *None*, append to plot title. Default *None*.

Returns `plotly.Figure` representing the precision-recall plot generated

`evalml.pipelines.roc_curve`

`evalml.pipelines.roc_curve(y_true, y_pred_proba)`

Given labels and classifier predicted probabilities, compute and return the data representing a Receiver Operating Characteristic (ROC) curve.

Parameters

- **y_true** (*pd.Series* or *np.array*) – true labels.
- **y_pred_proba** (*pd.Series* or *np.array*) – predictions from a classifier, before thresholding has been applied. Note that 1 dimensional input is expected.

Returns

Dictionary containing metrics used to generate an ROC plot, with the following keys:

- *fpr_rate*: False positive rate.
- *tpr_rate*: True positive rate.
- *threshold*: Threshold values used to produce each pair of true/false positive rates.
- *auc_score*: The area under the ROC curve.

Return type `dict`

`evalml.pipelines.graph_roc_curve`

`evalml.pipelines.graph_roc_curve(y_true, y_pred_proba, custom_class_names=None, title_addition=None)`

Generate and display a Receiver Operating Characteristic (ROC) plot.

Parameters

- **y_true** (*pd.Series* or *np.array*) – true labels.
- **y_pred_proba** (*pd.Series* or *np.array*) – predictions from a classifier, before thresholding has been applied. Note this should be a one dimensional array with the predicted probability for the “true” label in the binary case.
- **custom_class_labels** (*list* or *None*) – if not *None*, custom labels for classes. Default *None*.
- **title_addition** (*str* or *None*) – if not *None*, append to plot title. Default *None*.

Returns `plotly.Figure` representing the ROC plot generated

`evalml.pipelines.confusion_matrix`

`evalml.pipelines.confusion_matrix(y_true, y_predicted, normalize_method='true')`

Confusion matrix for binary and multiclass classification.

Parameters

- **y_true** (*pd.Series* or *np.array*) – true binary labels.

- **y_pred** (*pd.Series* or *np.array*) – predictions from a binary classifier.
- **normalize_method** (*{'true', 'pred', 'all'}*) – Normalization method. Supported options are: ‘true’ to normalize by row, ‘pred’ to normalize by column, or ‘all’ to normalize by all values. Defaults to ‘true’.

Returns confusion matrix

Return type *np.array*

evalml.pipelines.normalize_confusion_matrix

`evalml.pipelines.normalize_confusion_matrix(conf_mat, normalize_method='true')`

Normalizes a confusion matrix.

Parameters

- **conf_mat** (*pd.DataFrame* or *np.array*) – confusion matrix to normalize.
- **normalize_method** (*{'true', 'pred', 'all'}*) – Normalization method. Supported options are: ‘true’ to normalize by row, ‘pred’ to normalize by column, or ‘all’ to normalize by all values. Defaults to ‘true’.

Returns A normalized version of the input confusion matrix.

evalml.pipelines.graph_confusion_matrix

`evalml.pipelines.graph_confusion_matrix(y_true, y_pred, normalize_method='true', title_addition=None)`

Generate and display a confusion matrix plot.

If *normalize_method* is set, hover text will show raw count, otherwise hover text will show count normalized with method ‘true’.

Parameters

- **y_true** (*pd.Series* or *np.array*) – true binary labels.
- **y_pred** (*pd.Series* or *np.array*) – predictions from a binary classifier.
- **normalize_method** (*{'true', 'pred', 'all'}*) – Normalization method. Supported options are: ‘true’ to normalize by row, ‘pred’ to normalize by column, or ‘all’ to normalize by all values. Defaults to ‘true’.
- **title_addition** (*str* or *None*) – if not None, append to plot title. Default None.

Returns *plotly.Figure* representing the confusion matrix plot generated

evalml.pipelines.calculate_permutation_importance

`evalml.pipelines.calculate_permutation_importance(pipeline, X, y, objective, n_repeats=5, n_jobs=None, random_state=0)`

Calculates permutation importance for features.

Parameters

- **pipeline** (*PipelineBase* or *subclass*) – fitted pipeline
- **X** (*pd.DataFrame*) – the input data used to score and compute permutation importance

- **y** (*pd.Series*) – the target labels
- **objective** (*str*, *ObjectiveBase*) – objective to score on
- **n_repeats** (*int*) – Number of times to permute a feature. Defaults to 5.
- **n_jobs** (*int or None*) – Non-negative integer describing level of parallelism used for pipelines. None and 1 are equivalent. If set to -1, all CPUs are used. For n_jobs below -1, (n_cpus + 1 + n_jobs) are used.
- **random_state** (*int, np.random.RandomState*) – The random seed/state. Defaults to 0.

Returns Mean feature importance scores over 5 shuffles.

evalml.pipelines.graph_permutation_importance

`evalml.pipelines.graph_permutation_importance(pipeline, X, y, objective, show_all_features=False)`

Generate a bar graph of the pipeline's permutation importance.

Parameters

- **pipeline** (*PipelineBase or subclass*) – Fitted pipeline
- **X** (*pd.DataFrame*) – The input data used to score and compute permutation importance
- **y** (*pd.Series*) – The target labels
- **objective** (*str, ObjectiveBase*) – Objective to score on
- **show_all_features** (*bool, optional*) – If True, graph features with a permutation importance value of zero. Defaults to False.

Returns `plotly.Figure`, a bar graph showing features and their respective permutation importance.

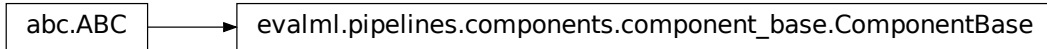
Components

Component Base Classes

Components represent a step in a pipeline.

<i>ComponentBase</i>	Base class for all components.
<i>Transformer</i>	A component that may or may not need fitting that transforms data.
<i>Estimator</i>	A component that fits and predicts given data.

evalml.pipelines.components.ComponentBase



```
class evalml.pipelines.components.ComponentBase (parameters=None, component_obj=None, random_state=0, **kwargs)
```

Base class for all components.

Methods

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data

evalml.pipelines.components.ComponentBase.__init__

```
ComponentBase.__init__ (parameters=None, component_obj=None, random_state=0, **kwargs)
```

Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.ComponentBase.clone

```
ComponentBase.clone (random_state=0)
```

Constructs a new component with the same parameters

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a RandomState instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.ComponentBase.describe

```
ComponentBase.describe (print_name=False, return_dict=False)
```

Describe a component and its parameters

Parameters

- **print_name** (*bool, optional*) – whether to print name of component
- **return_dict** (*bool, optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.ComponentBase.fit

ComponentBase.**fit** (*X*, *y=None*)

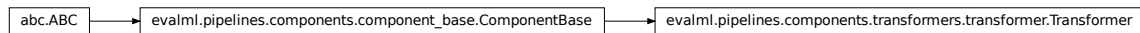
Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.Transformer



```

class evalml.pipelines.components.Transformer(parameters=None, component_obj=None, random_state=0,
                                              **kwargs)
  
```

A component that may or may not need fitting that transforms data. These components are used before an estimator.

To implement a new Transformer, define your own class which is a subclass of Transformer, including a name and a list of acceptable ranges for any parameters to be tuned during the automl search (hyperparameters). Define an `__init__` method which sets up any necessary state and objects. Make sure your `__init__` only uses standard keyword arguments and calls `super().__init__()` with a parameters dict. You may also override the `fit`, `transform`, `fit_transform` and other methods in this class if appropriate.

To see some examples, check out the definitions of any Transformer component.

Methods

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>fit_transform</code>	Fits on X and transforms X
<code>transform</code>	Transforms data X

evalml.pipelines.components.Transformer.__init__

`Transformer.__init__ (parameters=None, component_obj=None, random_state=0, **kwargs)`
Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.Transformer.clone

`Transformer.clone (random_state=0)`
Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.Transformer.describe

`Transformer.describe (print_name=False, return_dict=False)`
Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.Transformer.fit

`Transformer.fit (X, y=None)`
Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.Transformer.fit_transform

`Transformer.fit_transform (X, y=None)`
Fits on X and transforms X

Parameters

- **X** (*pd.DataFrame*) – Data to fit and transform
- **y** (*pd.DataFrame*) – Labels to fit and transform

Returns Transformed X

Return type `pd.DataFrame`

`evalml.pipelines.components.Transformer.transform`

`Transformer.transform(X, y=None)`

Transforms data X

Parameters

- **x** (`pd.DataFrame`) – Data to transform
- **y** (`pd.Series`, *optional*) – Input Labels

Returns Transformed X

Return type `pd.DataFrame`

`evalml.pipelines.components.Estimator`



class `evalml.pipelines.components.Estimator` (*parameters=None*, *component_obj=None*, *random_state=0*, ***kwargs*)

A component that fits and predicts given data.

To implement a new Transformer, define your own class which is a subclass of Transformer, including a name and a list of acceptable ranges for any parameters to be tuned during the automl search (hyperparameters). Define an `__init__` method which sets up any necessary state and objects. Make sure your `__init__` only uses standard keyword arguments and calls `super().__init__()` with a parameters dict. You may also override the `fit`, `transform`, `fit_transform` and other methods in this class if appropriate.

To see some examples, check out the definitions of any Estimator component.

Methods

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

`evalml.pipelines.components.Estimator.__init__`

`Estimator.__init__(parameters=None, component_obj=None, random_state=0, **kwargs)`

Initialize self. See `help(type(self))` for accurate signature.

evalml.pipelines.components.Estimator.clone

`Estimator.clone (random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.Estimator.describe

`Estimator.describe (print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.Estimator.fit

`Estimator.fit (X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.Estimator.predict

`Estimator.predict (X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.Estimator.predict_proba

`Estimator.predict_proba (X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

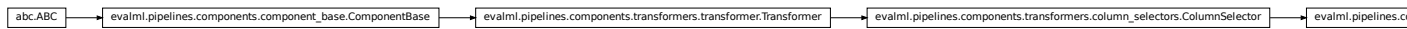
Return type pd.DataFrame

Transformers

Transformers are components that take in data as input and output transformed data.

<i>DropColumns</i>	Drops specified columns in input data.
<i>SelectColumns</i>	Selects specified columns in input data.
<i>OneHotEncoder</i>	One-hot encoder to encode non-numeric data.
<i>PerColumnImputer</i>	Imputes missing data according to a specified imputation strategy per column
<i>SimpleImputer</i>	Imputes missing data according to a specified imputation strategy.
<i>StandardScaler</i>	Standardize features: removes mean and scales to unit variance.
<i>RFRegressorSelectFromModel</i>	Selects top features based on importance weights using a Random Forest regressor.
<i>RFClassifierSelectFromModel</i>	Selects top features based on importance weights using a Random Forest classifier.

evalml.pipelines.components.DropColumns



```
class evalml.pipelines.components.DropColumns(columns=None, random_state=0,
                                              **kwargs)
```

Drops specified columns in input data.

```
name = 'Drop Columns Transformer'
```

```
model_family = 'none'
```

```
hyperparameter_ranges = {}
```

```
default_parameters = {'columns': None}
```

Instance attributes

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initializes an transformer that drops specified columns in input data.
-----------------------	--

Continued on next page

Table 75 – continued from previous page

<i>clone</i>	Constructs a new component with the same parameters
<i>describe</i>	Describe a component and its parameters
<i>fit</i>	‘Fits’ the transformer by checking if the column names are present in the dataset.
<i>fit_transform</i>	Fit transformer to data, then transform data.
<i>transform</i>	Transforms data X by dropping columns.

evalml.pipelines.components.DropColumns.__init__

`DropColumns.__init__(columns=None, random_state=0, **kwargs)`

Initializes an transformer that drops specified columns in input data.

Parameters `columns` (*list* (*string*)) – List of column names, used to determine which columns to drop.

evalml.pipelines.components.DropColumns.clone

`DropColumns.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.DropColumns.describe

`DropColumns.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- `print_name` (*bool*, *optional*) – whether to print name of component
- `return_dict` (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.DropColumns.fit

`DropColumns.fit(X, y=None)`

‘Fits’ the transformer by checking if the column names are present in the dataset.

Parameters

- `X` (*pd.DataFrame*) – Data to check.
- `y` (*pd.Series*, *optional*) – Targets.

Returns None.

evalml.pipelines.components.DropColumns.fit_transform

`DropColumns.fit_transform(X, y=None)`

Fit transformer to data, then transform data.

Parameters

- **X** (`pd.DataFrame`) – Data to transform.
- **y** (`pd.Series`, *optional*) – Targets.

Returns Transformed X.

Return type `pd.DataFrame`

evalml.pipelines.components.DropColumns.transform

`DropColumns.transform(X, y=None)`

Transforms data X by dropping columns.

Parameters

- **X** (`pd.DataFrame`) – Data to transform.
- **y** (`pd.Series`, *optional*) – Targets.

Returns Transformed X.

Return type `pd.DataFrame`

evalml.pipelines.components.SelectColumns

```
class evalml.pipelines.components.SelectColumns (columns=None, random_state=0,
                                              **kwargs)
```

Selects specified columns in input data.

```
name = 'Select Columns Transformer'
```

```
model_family = 'none'
```

```
hyperparameter_ranges = {}
```

```
default_parameters = {'columns': None}
```

Instance attributes

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initializes an transformer that drops specified columns in input data.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	'Fits' the transformer by checking if the column names are present in the dataset.
<code>fit_transform</code>	Fit transformer to data, then transform data.
<code>transform</code>	Transforms data X by selecting columns.

`evalml.pipelines.components.SelectColumns.__init__`

`SelectColumns.__init__` (*columns=None, random_state=0, **kwargs*)

Initializes an transformer that drops specified columns in input data.

Parameters `columns` (*list (string)*) – List of column names, used to determine which columns to drop.

`evalml.pipelines.components.SelectColumns.clone`

`SelectColumns.clone` (*random_state=0*)

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

`evalml.pipelines.components.SelectColumns.describe`

`SelectColumns.describe` (*print_name=False, return_dict=False*)

Describe a component and its parameters

Parameters

- `print_name` (*bool, optional*) – whether to print name of component
- `return_dict` (*bool, optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

`evalml.pipelines.components.SelectColumns.fit`

`SelectColumns.fit` (*X, y=None*)

'Fits' the transformer by checking if the column names are present in the dataset.

Parameters

- `X` (*pd.DataFrame*) – Data to check.
- `y` (*pd.Series, optional*) – Targets.

Returns None.

evalml.pipelines.components.SelectColumns.fit_transform`SelectColumns.fit_transform(X, y=None)`

Fit transformer to data, then transform data.

Parameters

- **X** (`pd.DataFrame`) – Data to transform.
- **y** (`pd.Series`, *optional*) – Targets.

Returns Transformed X.**Return type** `pd.DataFrame`**evalml.pipelines.components.SelectColumns.transform**`SelectColumns.transform(X, y=None)`

Transforms data X by selecting columns.

Parameters

- **X** (`pd.DataFrame`) – Data to transform.
- **y** (`pd.Series`, *optional*) – Targets.

Returns Transformed X.**Return type** `pd.DataFrame`**evalml.pipelines.components.OneHotEncoder**

```

abcABC → evalml.pipelines.components.component_base.ComponentBase → evalml.pipelines.components.transformers.transformer.Transformer → evalml.pipelines.components.transformers.encoders.encoder.CategoricalEncoder → evalml.pipelines

```

```

class evalml.pipelines.components.OneHotEncoder (top_n=10, categories=None,
                                                  drop=None, handle_unknown='ignore',
                                                  handle_missing='error', random_state=0,
                                                  **kwargs)

```

One-hot encoder to encode non-numeric data.

`name = 'One Hot Encoder'``model_family = 'none'``hyperparameter_ranges = {}``default_parameters = {'categories': None, 'drop': None, 'handle_missing': 'error',`**Instance attributes**

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initializes an transformer that encodes categorical features in a one-hot numeric array.”
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>fit_transform</code>	Fits on X and transforms X
<code>get_feature_names</code>	Returns names of transformed and added columns
<code>transform</code>	One-hot encode the input DataFrame.

evalml.pipelines.components.OneHotEncoder.__init__

`OneHotEncoder.__init__(top_n=10, categories=None, drop=None, handle_unknown='ignore', handle_missing='error', random_state=0, **kwargs)`
Initializes an transformer that encodes categorical features in a one-hot numeric array.”

Parameters

- **top_n** (*int*) – Number of categories per column to encode. If *None*, all categories will be encoded. Otherwise, the *n* most frequent will be encoded and all others will be dropped. Defaults to 10.
- **categories** (*list*) – A two dimensional list of categories, where *categories[i]* is a list of the categories for the column at index *i*. This can also be *None*, or “auto” if *top_n* is not *None*. Defaults to *None*.
- **drop** (*string*) – Method (“first” or “if_binary”) to use to drop one category per feature. Can also be a list specifying which method to use for each feature. Defaults to *None*.
- **handle_unknown** (*string*) – Whether to ignore or error for unknown categories for a feature encountered during *fit* or *transform*. If either *top_n* or *categories* is used to limit the number of categories per column, this must be “ignore”. Defaults to “ignore”.
- **handle_missing** (*string*) – Options for how to handle missing (NaN) values encountered during *fit* or *transform*. If this is set to “as_category” and NaN values are within the *n* most frequent, “nan” values will be encoded as their own column. If this is set to “error”, any missing values encountered will raise an error. Defaults to “error”.

evalml.pipelines.components.OneHotEncoder.clone

`OneHotEncoder.clone(random_state=0)`
Constructs a new component with the same parameters

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a *RandomState* instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.OneHotEncoder.describe

`OneHotEncoder.describe(print_name=False, return_dict=False)`
Describe a component and its parameters

Parameters

- **print_name** (*bool, optional*) – whether to print name of component
- **return_dict** (*bool, optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.OneHotEncoder.fit

`OneHotEncoder.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame or np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series, optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.OneHotEncoder.fit_transform

`OneHotEncoder.fit_transform(X, y=None)`

Fits on X and transforms X

Parameters

- **X** (*pd.DataFrame*) – Data to fit and transform
- **y** (*pd. DataFrame*) – Labels to fit and transform

Returns Transformed X

Return type pd.DataFrame

evalml.pipelines.components.OneHotEncoder.get_feature_names

`OneHotEncoder.get_feature_names()`

Returns names of transformed and added columns

Returns list of feature names not including dropped features

Return type list

evalml.pipelines.components.OneHotEncoder.transform

`OneHotEncoder.transform(X, y=None)`

One-hot encode the input DataFrame.

Parameters

- **X** (*pd.DataFrame*) – Dataframe of features.
- **y** (*pd.Series*) – Ignored.

Returns Transformed dataframe, where each categorical feature has been encoded into numerical columns using one-hot encoding.

evalml.pipelines.components.PerColumnImputer



```

class evalml.pipelines.components.PerColumnImputer(impute_strategies=None,
                                                    default_impute_strategy='most_frequent',
                                                    random_state=0, **kwargs)
    Imputes missing data according to a specified imputation strategy per column
    name = 'Per Column Imputer'
    model_family = 'none'
    hyperparameter_ranges = {}
    default_parameters = {'default_impute_strategy': 'most_frequent', 'impute_strategies'

```

Instance attributes

parameters	Returns the parameters which were used to initialize the component
------------	--

Methods:

<code>__init__</code>	Initializes a transformer that imputes missing data according to the specified imputation strategy per column.”
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits imputers on data X
<code>fit_transform</code>	Fits imputer on data X then imputes missing values in X
<code>transform</code>	Transforms data X by imputing missing values

evalml.pipelines.components.PerColumnImputer.__init__

```

PerColumnImputer.__init__(impute_strategies=None, default_impute_strategy='most_frequent',
                           random_state=0, **kwargs)

```

Initializes a transformer that imputes missing data according to the specified imputation strategy per column.”

Parameters

- **impute_strategies** (*dict*) – Column and {“impute_strategy”: strategy, “fill_value”:value} pairings. Valid values for impute strategy include “mean”, “median”,

“most_frequent”, “constant” for numerical data, and “most_frequent”, “constant” for object data types. Defaults to “most_frequent” for all columns.

When `impute_strategy == “constant”`, `fill_value` is used to replace missing data. Defaults to 0 when imputing numerical data and “missing_value” for strings or object data types.

- **default_impute_strategy** (*str*) – Impute strategy to fall back on when none is provided for a certain column. Valid values include “mean”, “median”, “most_frequent”, “constant” for numerical data, and “most_frequent”, “constant” for object data types. Defaults to “most_frequent”

`evalml.pipelines.components.PerColumnImputer.clone`

`PerColumnImputer.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

`evalml.pipelines.components.PerColumnImputer.describe`

`PerColumnImputer.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {“name”: name, “parameters”: parameters}

Returns prints and returns dictionary

Return type None or dict

`evalml.pipelines.components.PerColumnImputer.fit`

`PerColumnImputer.fit(X, y=None)`

Fits imputers on data X

Parameters

- **X** (*pd.DataFrame*) – Data to fit
- **y** (*pd.Series*, *optional*) – Input Labels

Returns self

`evalml.pipelines.components.PerColumnImputer.fit_transform`

`PerColumnImputer.fit_transform(X, y=None)`

Fits imputer on data X then imputes missing values in X

Parameters

- **X** (*pd.DataFrame*) – Data to fit and transform
- **y** (*pd.Series*) – Labels to fit and transform

Returns Transformed X

Return type *pd.DataFrame*

evalml.pipelines.components.PerColumnImputer.transform

PerColumnImputer.**transform**(X, y=None)
Transforms data X by imputing missing values

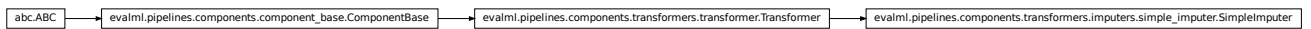
Parameters

- **X** (*pd.DataFrame*) – Data to transform
- **y** (*pd.Series*, *optional*) – Input Labels

Returns Transformed X

Return type *pd.DataFrame*

evalml.pipelines.components.SimpleImputer



```

class evalml.pipelines.components.SimpleImputer(impute_strategy='most_frequent',
                                                fill_value=None,    random_state=0,
                                                **kwargs)

    Imputes missing data according to a specified imputation strategy.

    name = 'Simple Imputer'
    model_family = 'none'
    hyperparameter_ranges = {'impute_strategy': ['mean', 'median', 'most_frequent']}
    default_parameters = {'fill_value': None, 'impute_strategy': 'most_frequent'}
  
```

Instance attributes

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initializes an transformer that imputes missing data according to the specified imputation strategy.”
<code>clone</code>	Constructs a new component with the same parameters

Continued on next page

Table 83 – continued from previous page

<i>describe</i>	Describe a component and its parameters
<i>fit</i>	Fits imputer to data
<i>fit_transform</i>	Fits on X and transforms X
<i>transform</i>	Transforms data X by imputing missing values

evalml.pipelines.components.SimpleImputer.__init__

`SimpleImputer.__init__(impute_strategy='most_frequent', fill_value=None, random_state=0, **kwargs)`

Initializes an transformer that imputes missing data according to the specified imputation strategy.”

Parameters

- **impute_strategy** (*string*) – Impute strategy to use. Valid values include “mean”, “median”, “most_frequent”, “constant” for numerical data, and “most_frequent”, “constant” for object data types.
- **fill_value** (*string*) – When `impute_strategy == “constant”`, `fill_value` is used to replace missing data. Defaults to 0 when imputing numerical data and “missing_value” for strings or object data types.

evalml.pipelines.components.SimpleImputer.clone

`SimpleImputer.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.SimpleImputer.describe

`SimpleImputer.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {“name”: name, “parameters”: parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.SimpleImputer.fit

`SimpleImputer.fit(X, y=None)`

Fits imputer to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]

- **y** (*pd.Series, optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.SimpleImputer.fit_transform

SimpleImputer.**fit_transform**(X, y=None)

Fits on X and transforms X

Parameters

- **x** (*pd.DataFrame*) – Data to fit and transform
- **y** (*pd.DataFrame*) – Labels to fit and transform

Returns Transformed X

Return type pd.DataFrame

evalml.pipelines.components.SimpleImputer.transform

SimpleImputer.**transform**(X, y=None)

Transforms data X by imputing missing values

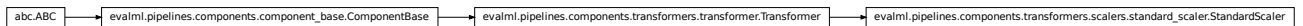
Parameters

- **x** (*pd.DataFrame*) – Data to transform
- **y** (*pd.Series, optional*) – Input Labels

Returns Transformed X

Return type pd.DataFrame

evalml.pipelines.components.StandardScaler



class evalml.pipelines.components.**StandardScaler**(*random_state=0, **kwargs*)

Standardize features: removes mean and scales to unit variance.

name = 'Standard Scaler'

model_family = 'none'

hyperparameter_ranges = {}

default_parameters = {}

Instance attributes

parameters	Returns the parameters which were used to initialize the component
------------	--

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>fit_transform</code>	Fits on X and transforms X
<code>transform</code>	Transforms data X

evalml.pipelines.components.StandardScaler.__init__

`StandardScaler.__init__(random_state=0, **kwargs)`
 Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.StandardScaler.clone

`StandardScaler.clone(random_state=0)`
 Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.StandardScaler.describe

`StandardScaler.describe(print_name=False, return_dict=False)`
 Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.StandardScaler.fit

`StandardScaler.fit(X, y=None)`
 Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.StandardScaler.fit_transform

`StandardScaler.fit_transform(X, y=None)`

Fits on X and transforms X

Parameters

- **X** (`pd.DataFrame`) – Data to fit and transform
- **y** (`pd.DataFrame`) – Labels to fit and transform

Returns Transformed X

Return type `pd.DataFrame`

evalml.pipelines.components.StandardScaler.transform

`StandardScaler.transform(X, y=None)`

Transforms data X

Parameters

- **X** (`pd.DataFrame`) – Data to transform
- **y** (`pd.Series, optional`) – Input Labels

Returns Transformed X

Return type `pd.DataFrame`

evalml.pipelines.components.RFRegressorSelectFromModel

```
class evalml.pipelines.components.RFRegressorSelectFromModel (number_features=None,
                                                             n_estimators=10,
                                                             max_depth=None,
                                                             per-
                                                             cent_features=0.5,
                                                             threshold=-inf,
                                                             n_jobs=-1,    ran-
                                                             dom_state=0,
                                                             **kwargs)
```

Selects top features based on importance weights using a Random Forest regressor.

```
name = 'RF Regressor Select From Model'
```

```
model_family = 'none'
```

```
hyperparameter_ranges = {'percent_features': Real(low=0.01, high=1, prior='uniform',
```

```
default_parameters = {'max_depth': None, 'n_estimators': 10, 'n_jobs': -1, 'number_
```

Instance attributes

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>fit_transform</code>	Fits feature selector on data X then transforms X by selecting features
<code>get_indices</code>	Get integer index of features selected
<code>get_names</code>	Get names of selected features.
<code>transform</code>	Transforms data X by selecting features

evalml.pipelines.components.RFRegressorSelectFromModel.__init__

`RFRegressorSelectFromModel.__init__(number_features=None, n_estimators=10, max_depth=None, percent_features=0.5, threshold=-inf, n_jobs=-1, random_state=0, **kwargs)`

Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.RFRegressorSelectFromModel.clone

`RFRegressorSelectFromModel.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.RFRegressorSelectFromModel.describe

`RFRegressorSelectFromModel.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.RFRegressorSelectFromModel.fit

`RFRegressorSelectFromModel.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, optional) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.RFRegressorSelectFromModel.fit_transform

`RFRegressorSelectFromModel.fit_transform(X, y=None)`

Fits feature selector on data X then transforms X by selecting features

Parameters

- **X** (*pd.DataFrame*) – Data to fit and transform
- **y** (*pd.Series*) – Labels to fit and transform

Returns Transformed X

Return type *pd.DataFrame*

evalml.pipelines.components.RFRegressorSelectFromModel.get_indices

`RFRegressorSelectFromModel.get_indices()`

Get integer index of features selected

Returns list of indices

Return type list

evalml.pipelines.components.RFRegressorSelectFromModel.get_names

`RFRegressorSelectFromModel.get_names()`

Get names of selected features.

Returns list of the names of features selected

evalml.pipelines.components.RFRegressorSelectFromModel.transform

`RFRegressorSelectFromModel.transform(X, y=None)`

Transforms data X by selecting features

Parameters

- **X** (*pd.DataFrame*) – Data to transform
- **y** (*pd.Series*, optional) – Input Labels

Returns Transformed X

Return type `pd.DataFrame`

evalml.pipelines.components.RFClassifierSelectFromModel



```

class evalml.pipelines.components.RFClassifierSelectFromModel(number_features=None,
                                                             n_estimators=10,
                                                             max_depth=None,
                                                             per-
                                                             cent_features=0.5,
                                                             threshold=-inf,
                                                             n_jobs=-1, ran-
                                                             dom_state=0,
                                                             **kwargs)

```

Selects top features based on importance weights using a Random Forest classifier.

```
name = 'RF Classifier Select From Model'
```

```
model_family = 'none'
```

```
hyperparameter_ranges = {'percent_features': Real(low=0.01, high=1, prior='uniform',
```

```
default_parameters = {'max_depth': None, 'n_estimators': 10, 'n_jobs': -1, 'number_
```

Instance attributes

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>fit_transform</code>	Fits feature selector on data X then transforms X by selecting features
<code>get_indices</code>	Get integer index of features selected
<code>get_names</code>	Get names of selected features.
<code>transform</code>	Transforms data X by selecting features

`evalml.pipelines.components.RFClassifierSelectFromModel.__init__`

```
RFClassifierSelectFromModel.__init__(number_features=None,      n_estimators=10,
                                     max_depth=None,           percent_features=0.5,
                                     threshold=-inf,  n_jobs=-1,  random_state=0,
                                     **kwargs)
```

Initialize self. See `help(type(self))` for accurate signature.

`evalml.pipelines.components.RFClassifierSelectFromModel.clone`

```
RFClassifierSelectFromModel.clone(random_state=0)
```

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

`evalml.pipelines.components.RFClassifierSelectFromModel.describe`

```
RFClassifierSelectFromModel.describe(print_name=False, return_dict=False)
```

Describe a component and its parameters

Parameters

- `print_name` (*bool*, *optional*) – whether to print name of component
- `return_dict` (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

`evalml.pipelines.components.RFClassifierSelectFromModel.fit`

```
RFClassifierSelectFromModel.fit(X, y=None)
```

Fits component to data

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- `y` (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

`evalml.pipelines.components.RFClassifierSelectFromModel.fit_transform`

```
RFClassifierSelectFromModel.fit_transform(X, y=None)
```

Fits feature selector on data X then transforms X by selecting features

Parameters

- `X` (*pd.DataFrame*) – Data to fit and transform

- **y** (*pd.Series*) – Labels to fit and transform

Returns Transformed X

Return type *pd.DataFrame*

`evalml.pipelines.components.RFClassifierSelectFromModel.get_indices`

`RFClassifierSelectFromModel.get_indices()`

Get integer index of features selected

Returns list of indices

Return type list

`evalml.pipelines.components.RFClassifierSelectFromModel.get_names`

`RFClassifierSelectFromModel.get_names()`

Get names of selected features.

Returns list of the names of features selected

`evalml.pipelines.components.RFClassifierSelectFromModel.transform`

`RFClassifierSelectFromModel.transform(X, y=None)`

Transforms data X by selecting features

Parameters

- **x** (*pd.DataFrame*) – Data to transform
- **y** (*pd.Series*, *optional*) – Input Labels

Returns Transformed X

Return type *pd.DataFrame*

Estimators

Classifiers

Classifiers are components that output a predicted class label.

<i>CatBoostClassifier</i>	CatBoost Classifier, a classifier that uses gradient-boosting on decision trees.
<i>ElasticNetClassifier</i>	Elastic Net Classifier.
<i>ExtraTreesClassifier</i>	Extra Trees Classifier.
<i>RandomForestClassifier</i>	Random Forest Classifier.
<i>LogisticRegressionClassifier</i>	Logistic Regression Classifier.
<i>XGBoostClassifier</i>	XGBoost Classifier.
<i>BaselineClassifier</i>	Classifier that predicts using the specified strategy.

evalml.pipelines.components.CatBoostClassifier



```

class evalml.pipelines.components.CatBoostClassifier(n_estimators=1000, eta=0.03,
                                                    max_depth=6,          boot-
                                                    strap_type=None,       ran-
                                                    dom_state=0, **kwargs)
  
```

CatBoost Classifier, a classifier that uses gradient-boosting on decision trees. CatBoost is an open-source library and natively supports categorical features.

For more information, check out <https://catboost.ai/>

```
name = 'CatBoost Classifier'
```

```
model_family = 'catboost'
```

```
supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS: 'multiclass'>]
```

```
hyperparameter_ranges = {'eta': Real(low=1e-06, high=1, prior='uniform', transform='log')}
```

```
default_parameters = {'bootstrap_type': None, 'eta': 0.03, 'max_depth': 6, 'n_estimators': 1000}
```

Instance attributes

SEED_MAX	
SEED_MIN	
feature_importance	Returns importance associated with each feature.
parameters	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.CatBoostClassifier.__init__

```

CatBoostClassifier.__init__(n_estimators=1000, eta=0.03, max_depth=6, boot-
                           strap_type=None, random_state=0, **kwargs)
  
```

Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.CatBoostClassifier.clone

`CatBoostClassifier.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.CatBoostClassifier.describe

`CatBoostClassifier.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.CatBoostClassifier.fit

`CatBoostClassifier.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.CatBoostClassifier.predict

`CatBoostClassifier.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.CatBoostClassifier.predict_proba

`CatBoostClassifier.predict_proba(X)`

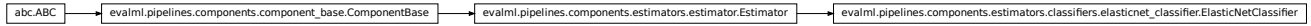
Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.components.ElasticNetClassifier



```

class evalml.pipelines.components.ElasticNetClassifier(alpha=0.5, l1_ratio=0.5,
                                                    n_jobs=-1, max_iter=1000,
                                                    random_state=0,
                                                    **kwargs)

```

Elastic Net Classifier.

```
name = 'Elastic Net Classifier'
```

```
model_family = 'linear_model'
```

```
supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS:
```

```
hyperparameter_ranges = {'alpha': Real(low=0, high=1, prior='uniform', transform='iden
```

```
default_parameters = {'alpha': 0.5, 'l1_ratio': 0.5, 'max_iter': 1000, 'n_jobs': -
```

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.ElasticNetClassifier.__init__

```
ElasticNetClassifier.__init__(alpha=0.5, l1_ratio=0.5, n_jobs=-1, max_iter=1000, random_state=0, **kwargs)
```

Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.ElasticNetClassifier.clone

`ElasticNetClassifier.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.ElasticNetClassifier.describe

`ElasticNetClassifier.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.ElasticNetClassifier.fit

`ElasticNetClassifier.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.ElasticNetClassifier.predict

`ElasticNetClassifier.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.ElasticNetClassifier.predict_proba

`ElasticNetClassifier.predict_proba(X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.components.ExtraTreesClassifier



```

class evalml.pipelines.components.ExtraTreesClassifier(n_estimators=100,
                                                       max_features='auto',
                                                       max_depth=6,
                                                       min_samples_split=2,
                                                       min_weight_fraction_leaf=0.0,
                                                       n_jobs=-1, random_state=0, **kwargs)

```

Extra Trees Classifier.

```

name = 'Extra Trees Classifier'
model_family = 'extra_trees'
supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS:
hyperparameter_ranges = {'max_depth': Integer(low=4, high=10, prior='uniform', transfo
default_parameters = {'max_depth': 6, 'max_features': 'auto', 'min_samples_split':

```

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.ExtraTreesClassifier.__init__

```

ExtraTreesClassifier.__init__(n_estimators=100, max_features='auto', max_depth=6,
                              min_samples_split=2, min_weight_fraction_leaf=0.0,
                              n_jobs=-1, random_state=0, **kwargs)

```

Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.ExtraTreesClassifier.clone

`ExtraTreesClassifier.clone (random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.ExtraTreesClassifier.describe

`ExtraTreesClassifier.describe (print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- `print_name` (*bool*, *optional*) – whether to print name of component
- `return_dict` (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.ExtraTreesClassifier.fit

`ExtraTreesClassifier.fit (X, y=None)`

Fits component to data

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- `y` (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.ExtraTreesClassifier.predict

`ExtraTreesClassifier.predict (X)`

Make predictions using selected features.

Parameters `X` (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.ExtraTreesClassifier.predict_proba

`ExtraTreesClassifier.predict_proba (X)`

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame*) – features

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.components.RandomForestClassifier



```

class evalml.pipelines.components.RandomForestClassifier(n_estimators=100,
                                                         max_depth=6, n_jobs=-
1, random_state=0,
                                                         **kwargs)

    Random Forest Classifier.

    name = 'Random Forest Classifier'
    model_family = 'random_forest'
    supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS:
    hyperparameter_ranges = {'max_depth': Integer(low=1, high=10, prior='uniform', transf
    default_parameters = {'max_depth': 6, 'n_estimators': 100, 'n_jobs': -1}
  
```

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.RandomForestClassifier.__init__

```

RandomForestClassifier.__init__(n_estimators=100, max_depth=6, n_jobs=-1, ran-
                                dom_state=0, **kwargs)
    Initialize self. See help(type(self)) for accurate signature.
  
```


evalml.pipelines.components.RandomForestClassifier.clone

`RandomForestClassifier.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.RandomForestClassifier.describe

`RandomForestClassifier.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool, optional*) – whether to print name of component
- **return_dict** (*bool, optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.RandomForestClassifier.fit

`RandomForestClassifier.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame or np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series, optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.RandomForestClassifier.predict

`RandomForestClassifier.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.RandomForestClassifier.predict_proba

`RandomForestClassifier.predict_proba(X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.components.LogisticRegressionClassifier



```
class evalml.pipelines.components.LogisticRegressionClassifier(penalty='l2',  
                                                             C=1.0,  
                                                             n_jobs=-1, ran-  
                                                             dom_state=0,  
                                                             **kwargs)  
  
    Logistic Regression Classifier.  
  
    name = 'Logistic Regression Classifier'  
    model_family = 'linear_model'  
    supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS:  
    hyperparameter_ranges = {'C': Real(low=0.01, high=10, prior='uniform', transform='iden  
    default_parameters = {'C': 1.0, 'n_jobs': -1, 'penalty': 'l2'}
```

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.LogisticRegressionClassifier.__init__

```
LogisticRegressionClassifier.__init__(penalty='l2',    C=1.0,    n_jobs=-1,    ran-  
                                     dom_state=0, **kwargs)  
    Initialize self. See help(type(self)) for accurate signature.
```

evalml.pipelines.components.LogisticRegressionClassifier.clone

`LogisticRegressionClassifier.clone (random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.LogisticRegressionClassifier.describe

`LogisticRegressionClassifier.describe (print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- `print_name` (*bool*, *optional*) – whether to print name of component
- `return_dict` (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.LogisticRegressionClassifier.fit

`LogisticRegressionClassifier.fit (X, y=None)`

Fits component to data

Parameters

- `X` (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- `y` (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.LogisticRegressionClassifier.predict

`LogisticRegressionClassifier.predict (X)`

Make predictions using selected features.

Parameters `X` (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.LogisticRegressionClassifier.predict_proba

`LogisticRegressionClassifier.predict_proba (X)`

Make probability estimates for labels.

Parameters `X` (*pd.DataFrame*) – features

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.components.XGBoostClassifier



```

class evalml.pipelines.components.XGBoostClassifier(eta=0.1,          max_depth=6,
                                                    min_child_weight=1,
                                                    n_estimators=100,      ran-
                                                    dom_state=0, **kwargs)

```

XGBoost Classifier.

```
name = 'XGBoost Classifier'
```

```
model_family = 'xgboost'
```

```
supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS:
```

```
hyperparameter_ranges = {'eta': Real(low=1e-06, high=1, prior='uniform', transform='i
```

```
default_parameters = {'eta': 0.1, 'max_depth': 6, 'min_child_weight': 1, 'n_estimat
```

Instance attributes

SEED_MAX	
SEED_MIN	
feature_importance	Returns importance associated with each feature.
parameters	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.XGBoostClassifier.__init__

```

XGBoostClassifier.__init__(eta=0.1, max_depth=6, min_child_weight=1, n_estimators=100,
                           random_state=0, **kwargs)

```

Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.XGBoostClassifier.clone

`XGBoostClassifier.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.XGBoostClassifier.describe

`XGBoostClassifier.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.XGBoostClassifier.fit

`XGBoostClassifier.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.XGBoostClassifier.predict

`XGBoostClassifier.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.XGBoostClassifier.predict_proba

`XGBoostClassifier.predict_proba(X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type pd.DataFrame

evalml.pipelines.components.BaselineClassifier



```

class evalml.pipelines.components.BaselineClassifier(strategy='mode',          ran-
                                                    dom_state=0, **kwargs)
    Classifier that predicts using the specified strategy.
    This is useful as a simple baseline classifier to compare with other classifiers.
    name = 'Baseline Classifier'
    model_family = 'baseline'
    supported_problem_types = [<ProblemTypes.BINARY: 'binary'>, <ProblemTypes.MULTICLASS:
    hyperparameter_ranges = {}
    default_parameters = {'strategy': 'mode'}
  
```

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Baseline classifier that uses a simple strategy to make predictions.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.BaselineClassifier.__init__

`BaselineClassifier.__init__(strategy='mode', random_state=0, **kwargs)`

Baseline classifier that uses a simple strategy to make predictions.

Parameters

- **strategy** (*str*) – method used to predict. Valid options are “mode”, “random” and “random_weighted”. Defaults to “mode”.
- **random_state** (*int*, *np.random.RandomState*) – seed for the random number

generator

evalml.pipelines.components.BaselineClassifier.clone

`BaselineClassifier.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.BaselineClassifier.describe

`BaselineClassifier.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool, optional*) – whether to print name of component
- **return_dict** (*bool, optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.BaselineClassifier.fit

`BaselineClassifier.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame or np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series, optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.BaselineClassifier.predict

`BaselineClassifier.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type `pd.Series`

evalml.pipelines.components.BaselineClassifier.predict_proba

`BaselineClassifier.predict_proba(X)`

Make probability estimates for labels.

Parameters *X* (`pd.DataFrame`) – features

Returns probability estimates

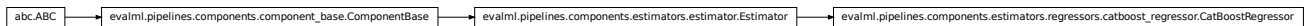
Return type `pd.DataFrame`

Regressors

Regressors are components that output a predicted target value.

<i>CatBoostRegressor</i>	CatBoost Regressor, a regressor that uses gradient-boosting on decision trees.
<i>ElasticNetRegressor</i>	Elastic Net Regressor.
<i>LinearRegressor</i>	Linear Regressor.
<i>ExtraTreesRegressor</i>	Extra Trees Regressor.
<i>RandomForestRegressor</i>	Random Forest Regressor.
<i>XGBoostRegressor</i>	XGBoost Regressor.
<i>BaselineRegressor</i>	Regressor that predicts using the specified strategy.

evalml.pipelines.components.CatBoostRegressor



```
class evalml.pipelines.components.CatBoostRegressor(n_estimators=1000, eta=0.03,
                                                    max_depth=6, boot-
                                                    strap_type=None, ran-
                                                    dom_state=0, **kwargs)
```

CatBoost Regressor, a regressor that uses gradient-boosting on decision trees. CatBoost is an open-source library and natively supports categorical features.

For more information, check out <https://catboost.ai/>

```
name = 'CatBoost Regressor'
```

```
model_family = 'catboost'
```

```
supported_problem_types = [<ProblemTypes.REGRESSION: 'regression'>]
```

```
hyperparameter_ranges = {'eta': Real(low=1e-06, high=1, prior='uniform', transform='i
```

```
default_parameters = {'bootstrap_type': None, 'eta': 0.03, 'max_depth': 6, 'n_estim
```

Instance attributes

SEED_MAX

Continued on next page

Table 106 – continued from previous page

SEED_MIN	
feature_importance	Returns importance associated with each feature.
parameters	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Build a model
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.CatBoostRegressor.__init__

`CatBoostRegressor.__init__(n_estimators=1000, eta=0.03, max_depth=6, bootstrap_type=None, random_state=0, **kwargs)`
 Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.CatBoostRegressor.clone

`CatBoostRegressor.clone(random_state=0)`
 Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.CatBoostRegressor.describe

`CatBoostRegressor.describe(print_name=False, return_dict=False)`
 Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.CatBoostRegressor.fit

`CatBoostRegressor.fit(X, y=None)`
 Build a model

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.CatBoostRegressor.predict

`CatBoostRegressor.predict(X)`
Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.components.CatBoostRegressor.predict_proba

`CatBoostRegressor.predict_proba(X)`
Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.components.ElasticNetRegressor



```

class evalml.pipelines.components.ElasticNetRegressor(alpha=0.5, l1_ratio=0.5,
                                                       max_iter=1000, normalize=False,
                                                       random_state=0, **kwargs)

```

Elastic Net Regressor.

name = 'Elastic Net Regressor'

model_family = 'linear_model'

supported_problem_types = [<ProblemTypes.REGRESSION: 'regression'>]

hyperparameter_ranges = {'alpha': Real(low=0, high=1, prior='uniform', transform='identity')}

default_parameters = {'alpha': 0.5, 'l1_ratio': 0.5, 'max_iter': 1000, 'normalize': False}

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.ElasticNetRegressor.__init__

`ElasticNetRegressor.__init__(alpha=0.5, l1_ratio=0.5, max_iter=1000, normalize=False, random_state=0, **kwargs)`
 Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.ElasticNetRegressor.clone

`ElasticNetRegressor.clone(random_state=0)`
 Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.ElasticNetRegressor.describe

`ElasticNetRegressor.describe(print_name=False, return_dict=False)`
 Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.ElasticNetRegressor.fit

`ElasticNetRegressor.fit(X, y=None)`
 Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.ElasticNetRegressor.predict

`ElasticNetRegressor.predict(X)`
Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.components.ElasticNetRegressor.predict_proba

`ElasticNetRegressor.predict_proba(X)`
Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.components.LinearRegressor



```
class evalml.pipelines.components.LinearRegressor(fit_intercept=True, normalize=False, n_jobs=-1, random_state=0, **kwargs)
```

Linear Regressor.

name = 'Linear Regressor'

model_family = 'linear_model'

supported_problem_types = [*<ProblemTypes.REGRESSION: 'regression'>*]

hyperparameter_ranges = {'fit_intercept': [True, False], 'normalize': [True, False]}

default_parameters = {'fit_intercept': True, 'n_jobs': -1, 'normalize': False}

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
---------------------------------	--

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Table 110 – continued from previous page

<code>parameters</code>	Returns the parameters which were used to initialize the component
-------------------------	--

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.LinearRegressor.__init__

`LinearRegressor.__init__` (*fit_intercept=True, normalize=False, n_jobs=-1, random_state=0, **kwargs*)

Initialize self. See `help(type(self))` for accurate signature.

evalml.pipelines.components.LinearRegressor.clone

`LinearRegressor.clone` (*random_state=0*)

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.LinearRegressor.describe

`LinearRegressor.describe` (*print_name=False, return_dict=False*)

Describe a component and its parameters

Parameters

- **print_name** (*bool, optional*) – whether to print name of component
- **return_dict** (*bool, optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.LinearRegressor.fit

`LinearRegressor.fit` (*X, y=None*)

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, optional) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.LinearRegressor.predict

`LinearRegressor.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.components.LinearRegressor.predict_proba

`LinearRegressor.predict_proba(X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.components.ExtraTreesRegressor



```

class evalml.pipelines.components.ExtraTreesRegressor(n_estimators=100,
                                                       max_features='auto',
                                                       max_depth=6,
                                                       min_samples_split=2,
                                                       min_weight_fraction_leaf=0.0,
                                                       n_jobs=-1, random_state=0,
                                                       **kwargs)

```

Extra Trees Regressor.

name = 'Extra Trees Regressor'

model_family = 'extra_trees'

supported_problem_types = [`<ProblemTypes.REGRESSION: 'regression'>`]

hyperparameter_ranges = {'max_depth': `Integer(low=4, high=10, prior='uniform', transfo`

default_parameters = {'max_depth': 6, 'max_features': 'auto', 'min_samples_split':

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

`evalml.pipelines.components.ExtraTreesRegressor.__init__`

`ExtraTreesRegressor.__init__(n_estimators=100, max_features='auto', max_depth=6, min_samples_split=2, min_weight_fraction_leaf=0.0, n_jobs=-1, random_state=0, **kwargs)`

Initialize self. See `help(type(self))` for accurate signature.

`evalml.pipelines.components.ExtraTreesRegressor.clone`

`ExtraTreesRegressor.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

`evalml.pipelines.components.ExtraTreesRegressor.describe`

`ExtraTreesRegressor.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.ExtraTreesRegressor.fit

ExtraTreesRegressor.**fit** (*X*, *y=None*)

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [*n_samples*, *n_features*]
- **y** (*pd.Series*, *optional*) – the target training labels of length [*n_samples*]

Returns *self*

evalml.pipelines.components.ExtraTreesRegressor.predict

ExtraTreesRegressor.**predict** (*X*)

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.components.ExtraTreesRegressor.predict_proba

ExtraTreesRegressor.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.components.RandomForestRegressor

```
abc.ABC → evalml.pipelines.components.component_base.ComponentBase → evalml.pipelines.components.estimators.estimator.Estimator → evalml.pipelines.components.estimators.regressors.rf_regressor.RandomForestRegressor
```

```
class evalml.pipelines.components.RandomForestRegressor (n_estimators=100,  
                                                    max_depth=6, n_jobs=-1,  
                                                    random_state=0,  
                                                    **kwargs)
```

Random Forest Regressor.

name = 'Random Forest Regressor'

model_family = 'random_forest'

supported_problem_types = [*<ProblemTypes.REGRESSION: 'regression'>*]

hyperparameter_ranges = {'max_depth': *Integer(low=1, high=32, prior='uniform', transfo*

default_parameters = {'max_depth': 6, 'n_estimators': 100, 'n_jobs': -1}

Instance attributes

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

`evalml.pipelines.components.RandomForestRegressor.__init__`

`RandomForestRegressor.__init__(n_estimators=100, max_depth=6, n_jobs=-1, random_state=0, **kwargs)`
 Initialize self. See `help(type(self))` for accurate signature.

`evalml.pipelines.components.RandomForestRegressor.clone`

`RandomForestRegressor.clone(random_state=0)`
 Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

`evalml.pipelines.components.RandomForestRegressor.describe`

`RandomForestRegressor.describe(print_name=False, return_dict=False)`
 Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

`evalml.pipelines.components.RandomForestRegressor.fit`

`RandomForestRegressor.fit(X, y=None)`
 Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, optional) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.RandomForestRegressor.predict

RandomForestRegressor.**predict** (*X*)

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.components.RandomForestRegressor.predict_proba

RandomForestRegressor.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type *pd.DataFrame*

evalml.pipelines.components.XGBoostRegressor



```

class evalml.pipelines.components.XGBoostRegressor(eta=0.1, max_depth=6,
                                                    min_child_weight=1,
                                                    n_estimators=100, random_state=0, **kwargs)

```

XGBoost Regressor.

name = 'XGBoost Regressor'

model_family = 'xgboost'

supported_problem_types = [<ProblemTypes.REGRESSION: 'regression'>]

hyperparameter_ranges = {'eta': Real(low=1e-06, high=1, prior='uniform', transform='i

default_parameters = {'eta': 0.1, 'max_depth': 6, 'min_child_weight': 1, 'n_estimat

Instance attributes

<code>SEED_MAX</code>	
<code>SEED_MIN</code>	
<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Initialize self.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.XGBoostRegressor.__init__

`XGBoostRegressor.__init__(eta=0.1, max_depth=6, min_child_weight=1, n_estimators=100, random_state=0, **kwargs)`
 Initialize self. See help(type(self)) for accurate signature.

evalml.pipelines.components.XGBoostRegressor.clone

`XGBoostRegressor.clone(random_state=0)`
 Constructs a new component with the same parameters

Parameters `random_state` (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.XGBoostRegressor.describe

`XGBoostRegressor.describe(print_name=False, return_dict=False)`
 Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {"name": name, "parameters": parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.XGBoostRegressor.fit

`XGBoostRegressor.fit(X, y=None)`
 Fits component to data

Parameters

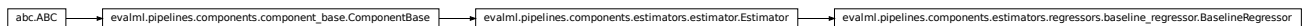
- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, optional) – the target training labels of length [n_samples]

Returns self**evalml.pipelines.components.XGBoostRegressor.predict**XGBoostRegressor.**predict** (*X*)

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features**Returns** estimated labels**Return type** *pd.Series***evalml.pipelines.components.XGBoostRegressor.predict_proba**XGBoostRegressor.**predict_proba** (*X*)

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features**Returns** probability estimates**Return type** *pd.DataFrame***evalml.pipelines.components.BaselineRegressor**

```
class evalml.pipelines.components.BaselineRegressor (strategy='mean', random_state=0, **kwargs)
```

Regressor that predicts using the specified strategy.

This is useful as a simple baseline regressor to compare with other regressors.

name = 'Baseline Regressor'**model_family** = 'baseline'**supported_problem_types** = [*<ProblemTypes.REGRESSION: 'regression'>*]**hyperparameter_ranges** = {}**default_parameters** = {'strategy': 'mean'}**Instance attributes**

<code>feature_importance</code>	Returns importance associated with each feature.
<code>parameters</code>	Returns the parameters which were used to initialize the component

Methods:

<code>__init__</code>	Baseline regressor that uses a simple strategy to make predictions.
<code>clone</code>	Constructs a new component with the same parameters
<code>describe</code>	Describe a component and its parameters
<code>fit</code>	Fits component to data
<code>predict</code>	Make predictions using selected features.
<code>predict_proba</code>	Make probability estimates for labels.

evalml.pipelines.components.BaselineRegressor.__init__

`BaselineRegressor.__init__(strategy='mean', random_state=0, **kwargs)`

Baseline regressor that uses a simple strategy to make predictions.

Parameters

- **strategy** (*str*) – method used to predict. Valid options are “mean”, “median”. Defaults to “mean”.
- **random_state** (*int*, *np.random.RandomState*) – seed for the random number generator

evalml.pipelines.components.BaselineRegressor.clone

`BaselineRegressor.clone(random_state=0)`

Constructs a new component with the same parameters

Parameters **random_state** (*int*) – the value to seed the random state with. Can also be a `RandomState` instance. Defaults to 0.

Returns A new instance of this component with identical parameters

evalml.pipelines.components.BaselineRegressor.describe

`BaselineRegressor.describe(print_name=False, return_dict=False)`

Describe a component and its parameters

Parameters

- **print_name** (*bool*, *optional*) – whether to print name of component
- **return_dict** (*bool*, *optional*) – whether to return description as dictionary in the format {“name”: name, “parameters”: parameters}

Returns prints and returns dictionary

Return type None or dict

evalml.pipelines.components.BaselineRegressor.fit

`BaselineRegressor.fit(X, y=None)`

Fits component to data

Parameters

- **X** (*pd.DataFrame* or *np.array*) – the input training data of shape [n_samples, n_features]
- **y** (*pd.Series*, optional) – the target training labels of length [n_samples]

Returns self

evalml.pipelines.components.BaselineRegressor.predict

`BaselineRegressor.predict(X)`

Make predictions using selected features.

Parameters **X** (*pd.DataFrame*) – features

Returns estimated labels

Return type *pd.Series*

evalml.pipelines.components.BaselineRegressor.predict_proba

`BaselineRegressor.predict_proba(X)`

Make probability estimates for labels.

Parameters **X** (*pd.DataFrame*) – features

Returns probability estimates

Return type *pd.DataFrame*

Objective Functions**Objective Base Classes**

<i>ObjectiveBase</i>	Base class for all objectives.
<i>BinaryClassificationObjective</i>	Base class for all binary classification objectives.
<i>MulticlassClassificationObjective</i>	Base class for all multiclass classification objectives.
<i>RegressionObjective</i>	Base class for all regression objectives.

evalml.objectives.ObjectiveBase

class evalml.objectives.ObjectiveBase
Base class for all objectives.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.ObjectiveBase.objective_function

classmethod ObjectiveBase.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.ObjectiveBase.score

ObjectiveBase.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [*n_samples*]
- **y_true** (*pd.Series*) – actual class labels of length [*n_samples*]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.ObjectiveBase.validate_inputs`

`ObjectiveBase.validate_inputs(y_true, y_predicted)`

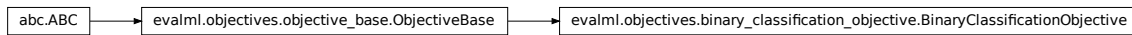
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

`evalml.objectives.BinaryClassificationObjective`



class `evalml.objectives.BinaryClassificationObjective`

Base class for all binary classification objectives.

problem_type (*ProblemTypes*): Type of problem this objective is. Set to *ProblemTypes.BINARY*.
can_optimize_threshold (*bool*): Determines if threshold used by objective can be optimized or not.

Methods

<code>decision_function</code>	Apply a learned threshold to predicted probabilities to get predicted classes.
<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>optimize_threshold</code>	Learn a binary classification threshold which optimizes the current objective.
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

`evalml.objectives.BinaryClassificationObjective.decision_function`

`BinaryClassificationObjective.decision_function(ypred_proba, threshold=0.5, X=None)`

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float, optional*) – Threshold used to make a prediction. Defaults to

0.5.

- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns predictions

`evalml.objectives.BinaryClassificationObjective.objective_function`

classmethod `BinaryClassificationObjective.objective_function` (*y_true*,
y_predicted,
X=None)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.BinaryClassificationObjective.optimize_threshold`

`BinaryClassificationObjective.optimize_threshold` (*ypred_proba*, *y_true*, *X=None*)

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

`evalml.objectives.BinaryClassificationObjective.score`

`BinaryClassificationObjective.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.BinaryClassificationObjective.validate_inputs

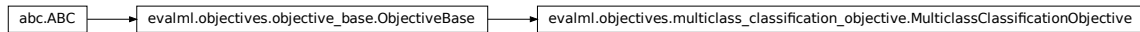
`BinaryClassificationObjective.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.MulticlassClassificationObjective

class evalml.objectives.**MulticlassClassificationObjective**

Base class for all multiclass classification objectives.

problem_type (*ProblemTypes*): Type of problem this objective is. Set to *ProblemTypes.MULTICLASS*.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.MulticlassClassificationObjective.objective_function

classmethod `MulticlassClassificationObjective.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: **y_predicted** (*pd.Series*) : predicted values of length [n_samples] **y_true** (*pd.Series*) : actual class labels of length [n_samples] **X** (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.MulticlassClassificationObjective.score

`MulticlassClassificationObjective.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length [n_samples]
- **y_true** (`pd.Series`) – actual class labels of length [n_samples]
- **X** (`pd.DataFrame` or `np.array`) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.MulticlassClassificationObjective.validate_inputs

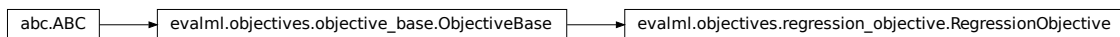
`MulticlassClassificationObjective.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length [n_samples]
- **y_true** (`pd.Series`) – actual class labels of length [n_samples]

Returns None

evalml.objectives.RegressionObjective**class evalml.objectives.RegressionObjective**

Base class for all regression objectives.

problem_type (`ProblemTypes`): Type of problem this objective is. Set to `ProblemTypes.REGRESSION`.

Methods

<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

evalml.objectives.RegressionObjective.objective_function

```
classmethod RegressionObjective.objective_function(y_true, y_predicted,  
                                                    X=None)
```

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (pd.Series) : predicted values of length [n_samples] *y_true* (pd.Series) : actual class labels of length [n_samples] *X* (pd.DataFrame or np.array) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.RegressionObjective.score

```
RegressionObjective.score(y_true, y_predicted, X=None)
```

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (pd.Series) – predicted values of length [n_samples]
- **y_true** (pd.Series) – actual class labels of length [n_samples]
- **X** (pd.DataFrame or np.array) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.RegressionObjective.validate_inputs

```
RegressionObjective.validate_inputs(y_true, y_predicted)
```

Validates the input based on a few simple checks.

Parameters

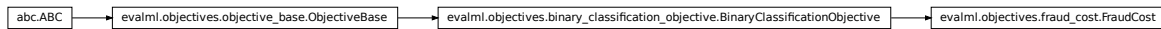
- **y_predicted** (pd.Series) – predicted values of length [n_samples]
- **y_true** (pd.Series) – actual class labels of length [n_samples]

Returns None

Domain-Specific Objectives

<i>FraudCost</i>	Score the percentage of money lost of the total transaction amount process due to fraud.
<i>LeadScoring</i>	Lead scoring.

evalml.objectives.FraudCost



class evalml.objectives.**FraudCost** (*retry_percentage=0.5*, *interchange_fee=0.02*,
fraud_payout_percentage=1.0, *amount_col='amount'*)
 Score the percentage of money lost of the total transaction amount process due to fraud.

Methods

<code>__init__</code>	Create instance of FraudCost
<code>decision_function</code>	Determine if a transaction is fraud given predicted probabilities, threshold, and dataframe with transaction amount.
<code>objective_function</code>	Calculate amount lost to fraud per transaction given predictions, true values, and dataframe with transaction amount.
<code>optimize_threshold</code>	Learn a binary classification threshold which optimizes the current objective.
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

evalml.objectives.FraudCost.__init__

`FraudCost.__init__` (*retry_percentage=0.5*, *interchange_fee=0.02*, *fraud_payout_percentage=1.0*,
amount_col='amount')
 Create instance of FraudCost

Parameters

- **retry_percentage** (*float*) – What percentage of customers that will retry a transaction if it is declined. Between 0 and 1. Defaults to .5
- **interchange_fee** (*float*) – How much of each successful transaction you can collect. Between 0 and 1. Defaults to .02
- **fraud_payout_percentage** (*float*) – Percentage of fraud you will not be able to collect. Between 0 and 1. Defaults to 1.0
- **amount_col** (*str*) – Name of column in data that contains the amount. Defaults to “amount”

evalml.objectives.FraudCost.decision_function

`FraudCost.decision_function` (*ypred_proba*, *threshold=0.0*, *X=None*)
 Determine if a transaction is fraud given predicted probabilities, threshold, and dataframe with transaction amount.

Parameters

- **ypred_proba** (*pd.Series*) – Predicted probabilities
- **x** (*pd.DataFrame*) – Dataframe containing transaction amount
- **threshold** (*float*) – Dollar threshold to determine if transaction is fraud

Returns Series of predicted fraud labels using X and threshold

Return type *pd.Series*

evalml.objectives.FraudCost.objective_function

`FraudCost.objective_function(y_true, y_predicted, X)`

Calculate amount lost to fraud per transaction given predictions, true values, and dataframe with transaction amount.

Parameters

- **y_predicted** (*pd.Series*) – predicted fraud labels
- **y_true** (*pd.Series*) – true fraud labels
- **x** (*pd.DataFrame*) – dataframe with transaction amounts

Returns amount lost to fraud per transaction

Return type *float*

evalml.objectives.FraudCost.optimize_threshold

`FraudCost.optimize_threshold(ypred_proba, y_true, X=None)`

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **x** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.FraudCost.score

`FraudCost.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **x** (*pd.DataFrame or np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.FraudCost.validate_inputs

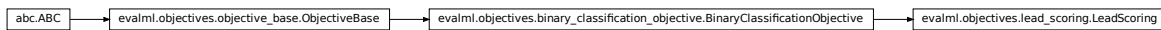
`FraudCost.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.LeadScoring

class evalml.objectives.LeadScoring(*true_positives=1, false_positives=-1*)

Lead scoring.

Methods

<code>__init__</code>	Create instance.
<code>decision_function</code>	Apply a learned threshold to predicted probabilities to get predicted classes.
<code>objective_function</code>	Calculate the profit per lead.
<code>optimize_threshold</code>	Learn a binary classification threshold which optimizes the current objective.
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

evalml.objectives.LeadScoring.__init__

`LeadScoring.__init__(true_positives=1, false_positives=-1)`

Create instance.

Parameters

- **true_positives** (*int*) – reward for a true positive
- **false_positives** (*int*) – cost for a false positive. Should be negative.

evalml.objectives.LeadScoring.decision_function

`LeadScoring.decision_function(ypred_proba, threshold=0.5, X=None)`

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float, optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.LeadScoring.objective_function

LeadScoring.objective_function (*y_true, y_predicted, X=None*)

Calculate the profit per lead.

Parameters

- **y_predicted** (*pd.Series*) – predicted labels
- **y_true** (*pd.Series*) – true labels
- **X** (*pd.DataFrame*) – None, not used.

Returns profit per lead

Return type float

evalml.objectives.LeadScoring.optimize_threshold

LeadScoring.optimize_threshold (*ypred_proba, y_true, X=None*)

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.LeadScoring.score

LeadScoring.score (*y_true, y_predicted, X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame or np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.LeadScoring.validate_inputs

LeadScoring.**validate_inputs**(*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

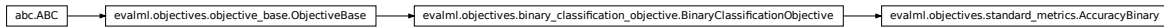
Classification Objectives

<i>AccuracyBinary</i>	Accuracy score for binary classification.
<i>AccuracyMulticlass</i>	Accuracy score for multiclass classification.
<i>AUC</i>	AUC score for binary classification.
<i>AUCMacro</i>	AUC score for multiclass classification using macro averaging.
<i>AUCMicro</i>	AUC score for multiclass classification using micro averaging.
<i>AUCWeighted</i>	AUC Score for multiclass classification using weighted averaging.
<i>BalancedAccuracyBinary</i>	Balanced accuracy score for binary classification.
<i>BalancedAccuracyMulticlass</i>	Balanced accuracy score for multiclass classification.
<i>F1</i>	F1 score for binary classification.
<i>F1Micro</i>	F1 score for multiclass classification using micro averaging.
<i>F1Macro</i>	F1 score for multiclass classification using macro averaging.
<i>F1Weighted</i>	F1 score for multiclass classification using weighted averaging.
<i>LogLossBinary</i>	Log Loss for binary classification.
<i>LogLossMulticlass</i>	Log Loss for multiclass classification.
<i>MCCBinary</i>	Matthews correlation coefficient for binary classification.
<i>MCCMulticlass</i>	Matthews correlation coefficient for multiclass classification.
<i>Precision</i>	Precision score for binary classification.
<i>PrecisionMicro</i>	Precision score for multiclass classification using micro averaging.
<i>PrecisionMacro</i>	Precision score for multiclass classification using macro averaging.
<i>PrecisionWeighted</i>	Precision score for multiclass classification using weighted averaging.
<i>Recall</i>	Recall score for binary classification.
<i>RecallMicro</i>	Recall score for multiclass classification using micro averaging.
<i>RecallMacro</i>	Recall score for multiclass classification using macro averaging.

Continued on next page

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<i>RecallWeighted</i>	Recall score for multiclass classification using weighted averaging.
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evalml.objectives.AccuracyBinary

class evalml.objectives.**AccuracyBinary**

Accuracy score for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.AccuracyBinary.decision_function

AccuracyBinary.**decision_function** (*ypred_proba*, *threshold=0.5*, *X=None*)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float*, *optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.AccuracyBinary.objective_function

AccuracyBinary.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: `y_predicted` (`pd.Series`) : predicted values of length `[n_samples]` `y_true` (`pd.Series`) : actual class labels of length `[n_samples]` `X` (`pd.DataFrame` or `np.array`) : extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.AccuracyBinary.optimize_threshold`

`AccuracyBinary.optimize_threshold(ypred_proba, y_true, X=None)`

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **`ypred_proba`** (`list`) – The classifier’s predicted probabilities
- **`y_true`** (`list`) – The ground truth for the predictions.
- **`X`** (`pd.DataFrame`, *optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

`evalml.objectives.AccuracyBinary.score`

`AccuracyBinary.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **`y_predicted`** (`pd.Series`) – predicted values of length `[n_samples]`
- **`y_true`** (`pd.Series`) – actual class labels of length `[n_samples]`
- **`X`** (`pd.DataFrame` or `np.array`) – extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns score

`evalml.objectives.AccuracyBinary.validate_inputs`

`AccuracyBinary.validate_inputs(y_true, y_predicted)`

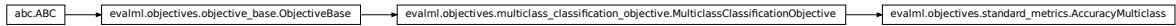
Validates the input based on a few simple checks.

Parameters

- **`y_predicted`** (`pd.Series`) – predicted values of length `[n_samples]`
- **`y_true`** (`pd.Series`) – actual class labels of length `[n_samples]`

Returns None

evalml.objectives.AccuracyMulticlass



class evalml.objectives.**AccuracyMulticlass**
Accuracy score for multiclass classification.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.AccuracyMulticlass.objective_function

AccuracyMulticlass.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.AccuracyMulticlass.score

AccuracyMulticlass.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [*n_samples*]
- **y_true** (*pd.Series*) – actual class labels of length [*n_samples*]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

evalml.objectives.AccuracyMulticlass.validate_inputs

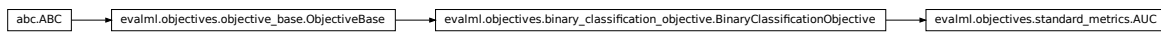
AccuracyMulticlass.**validate_inputs** (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.AUC

class evalml.objectives.**AUC**
AUC score for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.AUC.decision_function

AUC.**decision_function** (*ypred_proba*, *threshold=0.5*, *X=None*)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float*, *optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.AUC.objective_function

`AUC.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: `y_predicted` (`pd.Series`) : predicted values of length `[n_samples]` `y_true` (`pd.Series`) : actual class labels of length `[n_samples]` `X` (`pd.DataFrame` or `np.array`) : extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.AUC.optimize_threshold

`AUC.optimize_threshold(ypred_proba, y_true, X=None)`

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (`list`) – The classifier’s predicted probabilities
- **y_true** (`list`) – The ground truth for the predictions.
- **X** (`pd.DataFrame`, *optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.AUC.score

`AUC.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length `[n_samples]`
- **y_true** (`pd.Series`) – actual class labels of length `[n_samples]`
- **X** (`pd.DataFrame` or `np.array`) – extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns score

evalml.objectives.AUC.validate_inputs

`AUC.validate_inputs(y_true, y_predicted)`

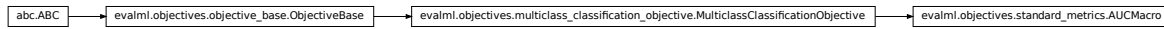
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length `[n_samples]`
- **y_true** (`pd.Series`) – actual class labels of length `[n_samples]`

Returns None

evalml.objectives.AUCMacro



class evalml.objectives.AUCMacro

AUC score for multiclass classification using macro averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.AUCMacro.objective_function

AUCMacro.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.AUCMacro.score

AUCMacro.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [*n_samples*]
- **y_true** (*pd.Series*) – actual class labels of length [*n_samples*]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

evalml.objectives.AUCMacro.validate_inputs

AUCMacro.**validate_inputs**(*y_true*, *y_predicted*)

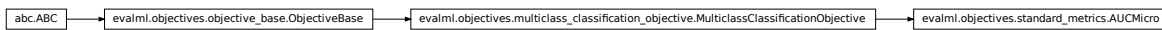
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.AUCMicro



class evalml.objectives.AUCMicro

AUC score for multiclass classification using micro averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.AUCMicro.objective_function

AUCMicro.**objective_function**(*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.AUCMicro.score

AUCMicro.**score**(*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.AUCMicro.validate_inputs

`AUCMicro.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.AUCWeighted



class evalml.objectives.AUCWeighted

AUC Score for multiclass classification using weighted averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.AUCWeighted.objective_function

`AUCWeighted.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: y_predicted (*pd.Series*) : predicted values of length [n_samples] y_true (*pd.Series*) : actual class labels of length [n_samples] X (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.AUCWeighted.score

AUCWeighted.**score**(*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.AUCWeighted.validate_inputs

AUCWeighted.**validate_inputs**(*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.BalancedAccuracyBinary



class evalml.objectives.BalancedAccuracyBinary

Balanced accuracy score for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.BalancedAccuracyBinary.decision_function

BalancedAccuracyBinary.**decision_function**(ypred_proba, threshold=0.5, X=None)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float, optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.BalancedAccuracyBinary.objective_function

BalancedAccuracyBinary.**objective_function**(y_true, y_predicted, X=None)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: y_predicted (pd.Series) : predicted values of length [n_samples] y_true (pd.Series) : actual class labels of length [n_samples] X (pd.DataFrame or np.array) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.BalancedAccuracyBinary.optimize_threshold

BalancedAccuracyBinary.**optimize_threshold**(ypred_proba, y_true, X=None)

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.BalancedAccuracyBinary.score

BalancedAccuracyBinary.**score**(y_true, y_predicted, X=None)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.BalancedAccuracyBinary.validate_inputs

BalancedAccuracyBinary.**validate_inputs** (*y_true*, *y_predicted*)

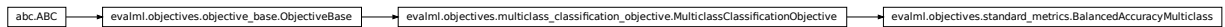
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.BalancedAccuracyMulticlass



class evalml.objectives.BalancedAccuracyMulticlass

Balanced accuracy score for multiclass classification.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.BalancedAccuracyMulticlass.objective_function

BalancedAccuracyMulticlass.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.BalancedAccuracyMulticlass.score

BalancedAccuracyMulticlass.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.BalancedAccuracyMulticlass.validate_inputs

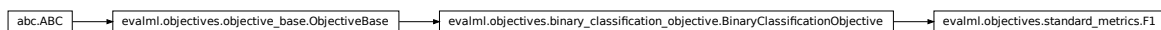
BalancedAccuracyMulticlass.**validate_inputs** (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.F1

class evalml.objectives.**F1**
F1 score for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.F1.decision_function

F1.decision_function (*ypred_proba*, *threshold=0.5*, *X=None*)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float*, *optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.F1.objective_function

F1.objective_function (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.F1.optimize_threshold

F1.optimize_threshold (*ypred_proba*, *y_true*, *X=None*)

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.F1.score

F1.score (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [*n_samples*]
- **y_true** (*pd.Series*) – actual class labels of length [*n_samples*]

- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.F1.validate_inputs

F1.validate_inputs (*y_true*, *y_predicted*)

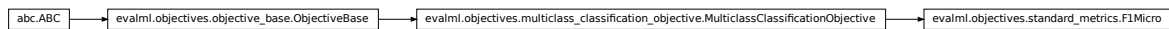
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.F1Micro



class evalml.objectives.F1Micro

F1 score for multiclass classification using micro averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.F1Micro.objective_function

F1Micro.objective_function (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.F1Micro.score

`F1Micro.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length [n_samples]
- **y_true** (`pd.Series`) – actual class labels of length [n_samples]
- **X** (`pd.DataFrame` or `np.array`) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.F1Micro.validate_inputs

`F1Micro.validate_inputs(y_true, y_predicted)`

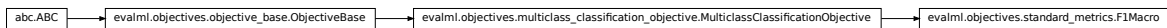
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length [n_samples]
- **y_true** (`pd.Series`) – actual class labels of length [n_samples]

Returns None

evalml.objectives.F1Macro



class evalml.objectives.F1Macro

F1 score for multiclass classification using macro averaging.

Methods

<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

evalml.objectives.F1Macro.objective_function

`F1Macro.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: `y_predicted` (`pd.Series`) : predicted values of length `[n_samples]` `y_true` (`pd.Series`) : actual class labels of length `[n_samples]` `X` (`pd.DataFrame` or `np.array`) : extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.F1Macro.score`

`F1Macro.score` (`y_true`, `y_predicted`, `X=None`)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **`y_predicted`** (`pd.Series`) – predicted values of length `[n_samples]`
- **`y_true`** (`pd.Series`) – actual class labels of length `[n_samples]`
- **`X`** (`pd.DataFrame` or `np.array`) – extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns score

`evalml.objectives.F1Macro.validate_inputs`

`F1Macro.validate_inputs` (`y_true`, `y_predicted`)

Validates the input based on a few simple checks.

Parameters

- **`y_predicted`** (`pd.Series`) – predicted values of length `[n_samples]`
- **`y_true`** (`pd.Series`) – actual class labels of length `[n_samples]`

Returns None

`evalml.objectives.F1Weighted`



class `evalml.objectives.F1Weighted`

F1 score for multiclass classification using weighted averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.F1Weighted.objective_function

`F1Weighted.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: `y_predicted` (`pd.Series`) : predicted values of length `[n_samples]` `y_true` (`pd.Series`) : actual class labels of length `[n_samples]` `X` (`pd.DataFrame` or `np.array`) : extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.F1Weighted.score

`F1Weighted.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length `[n_samples]`
- **y_true** (`pd.Series`) – actual class labels of length `[n_samples]`
- **X** (`pd.DataFrame` or `np.array`) – extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns score

evalml.objectives.F1Weighted.validate_inputs

`F1Weighted.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (`pd.Series`) – predicted values of length `[n_samples]`
- **y_true** (`pd.Series`) – actual class labels of length `[n_samples]`

Returns None

evalml.objectives.LogLossBinary



class evalml.objectives.LogLossBinary
Log Loss for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.LogLossBinary.decision_function

LogLossBinary.**decision_function** (*ypred_proba*, *threshold=0.5*, *X=None*)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float*, *optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.LogLossBinary.objective_function

LogLossBinary.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.LogLossBinary.optimize_threshold

`LogLossBinary.optimize_threshold(y_pred_proba, y_true, X=None)`
Learn a binary classification threshold which optimizes the current objective.

Parameters

- **y_pred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.LogLossBinary.score

`LogLossBinary.score(y_true, y_predicted, X=None)`
Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame or np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.LogLossBinary.validate_inputs

`LogLossBinary.validate_inputs(y_true, y_predicted)`
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.LogLossMulticlass



class evalml.objectives.LogLossMulticlass
Log Loss for multiclass classification.

Methods

<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

`evalml.objectives.LogLossMulticlass.objective_function`

`LogLossMulticlass.objective_function` (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (`pd.Series`) : predicted values of length [*n_samples*] *y_true* (`pd.Series`) : actual class labels of length [*n_samples*] *X* (`pd.DataFrame` or `np.array`) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.LogLossMulticlass.score`

`LogLossMulticlass.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]
- ***X*** (`pd.DataFrame` or `np.array`) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.LogLossMulticlass.validate_inputs`

`LogLossMulticlass.validate_inputs` (*y_true*, *y_predicted*)

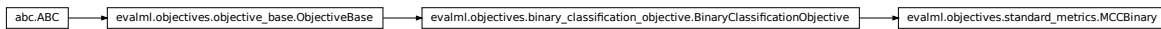
Validates the input based on a few simple checks.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]

Returns None

evalml.objectives.MCCBinary



class evalml.objectives.MCCBinary
Matthews correlation coefficient for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.MCCBinary.decision_function

MCCBinary.**decision_function**(ypred_proba, threshold=0.5, X=None)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float, optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.MCCBinary.objective_function

MCCBinary.**objective_function**(y_true, y_predicted, X=None)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: y_predicted (pd.Series) : predicted values of length [n_samples] y_true (pd.Series) : actual class labels of length [n_samples] X (pd.DataFrame or np.array) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.MCCBinary.optimize_threshold

MCCBinary.**optimize_threshold**(ypred_proba, y_true, X=None)

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.MCCBinary.score

MCCBinary.**score**(y_true, y_predicted, X=None)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame or np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.MCCBinary.validate_inputs

MCCBinary.**validate_inputs**(y_true, y_predicted)

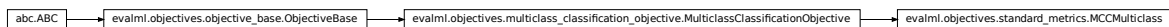
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.MCCMulticlass



class evalml.objectives.MCCMulticlass

Matthews correlation coefficient for multiclass classification.

Methods

<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

`evalml.objectives.MCCMulticlass.objective_function`

`MCCMulticlass.objective_function` (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (`pd.Series`) : predicted values of length [*n_samples*] *y_true* (`pd.Series`) : actual class labels of length [*n_samples*] *X* (`pd.DataFrame` or `np.array`) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.MCCMulticlass.score`

`MCCMulticlass.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]
- ***X*** (`pd.DataFrame` or `np.array`) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.MCCMulticlass.validate_inputs`

`MCCMulticlass.validate_inputs` (*y_true*, *y_predicted*)

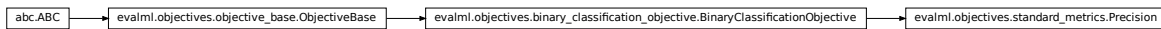
Validates the input based on a few simple checks.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]

Returns None

evalml.objectives.Precision



class evalml.objectives.Precision

Precision score for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.Precision.decision_function

Precision.**decision_function** (*ypred_proba*, *threshold=0.5*, *X=None*)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **threshold** (*float*, *optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame*, *optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.Precision.objective_function

Precision.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.Precision.optimize_threshold

`Precision.optimize_threshold(ypred_proba, y_true, X=None)`

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.Precision.score

`Precision.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame or np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.Precision.validate_inputs

`Precision.validate_inputs(y_true, y_predicted)`

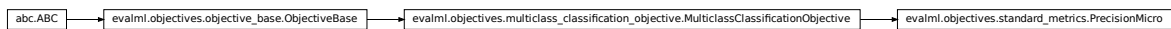
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.PrecisionMicro



class evalml.objectives.PrecisionMicro

Precision score for multiclass classification using micro averaging.

Methods

<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

`evalml.objectives.PrecisionMicro.objective_function`

`PrecisionMicro.objective_function` (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (`pd.Series`) : predicted values of length [*n_samples*] *y_true* (`pd.Series`) : actual class labels of length [*n_samples*] *X* (`pd.DataFrame` or `np.array`) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.PrecisionMicro.score`

`PrecisionMicro.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]
- ***X*** (`pd.DataFrame` or `np.array`) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.PrecisionMicro.validate_inputs`

`PrecisionMicro.validate_inputs` (*y_true*, *y_predicted*)

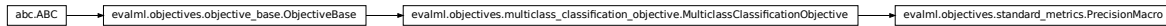
Validates the input based on a few simple checks.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]

Returns None

evalml.objectives.PrecisionMacro



class evalml.objectives.PrecisionMacro

Precision score for multiclass classification using macro averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.PrecisionMacro.objective_function

PrecisionMacro.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.PrecisionMacro.score

PrecisionMacro.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [*n_samples*]
- **y_true** (*pd.Series*) – actual class labels of length [*n_samples*]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

evalml.objectives.PrecisionMacro.validate_inputs

PrecisionMacro.**validate_inputs**(*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.PrecisionWeighted

class evalml.objectives.PrecisionWeighted

Precision score for multiclass classification using weighted averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.PrecisionWeighted.objective_function

PrecisionWeighted.**objective_function**(*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.PrecisionWeighted.score

PrecisionWeighted.**score**(*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.PrecisionWeighted.validate_inputs

PrecisionWeighted.**validate_inputs** (*y_true*, *y_predicted*)

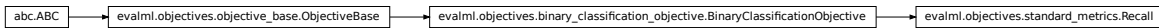
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.Recall



class evalml.objectives.Recall
Recall score for binary classification.

Methods

<i>decision_function</i>	Apply a learned threshold to predicted probabilities to get predicted classes.
<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>optimize_threshold</i>	Learn a binary classification threshold which optimizes the current objective.
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.Recall.decision_function

Recall.**decision_function** (*ypred_proba*, *threshold=0.5*, *X=None*)

Apply a learned threshold to predicted probabilities to get predicted classes.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities

- **threshold** (*float, optional*) – Threshold used to make a prediction. Defaults to 0.5.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns predictions

evalml.objectives.Recall.objective_function

`Recall.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: `y_predicted` (*pd.Series*) : predicted values of length `[n_samples]` `y_true` (*pd.Series*) : actual class labels of length `[n_samples]` `X` (*pd.DataFrame or np.array*) : extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.Recall.optimize_threshold

`Recall.optimize_threshold(ypred_proba, y_true, X=None)`

Learn a binary classification threshold which optimizes the current objective.

Parameters

- **ypred_proba** (*list*) – The classifier’s predicted probabilities
- **y_true** (*list*) – The ground truth for the predictions.
- **X** (*pd.DataFrame, optional*) – Any extra columns that are needed from training data.

Returns Optimal threshold for this objective

evalml.objectives.Recall.score

`Recall.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length `[n_samples]`
- **y_true** (*pd.Series*) – actual class labels of length `[n_samples]`
- **X** (*pd.DataFrame or np.array*) – extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns score

evalml.objectives.Recall.validate_inputs

`Recall.validate_inputs(y_true, y_predicted)`

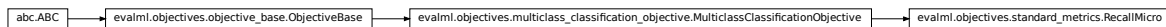
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.RecallMicro



class evalml.objectives.RecallMicro

Recall score for multiclass classification using micro averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.RecallMicro.objective_function

`RecallMicro.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: **y_predicted** (*pd.Series*) : predicted values of length [n_samples] **y_true** (*pd.Series*) : actual class labels of length [n_samples] **X** (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.RecallMicro.score

`RecallMicro.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.RecallMicro.validate_inputs

`RecallMicro.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.RecallMacro



class evalml.objectives.RecallMacro

Recall score for multiclass classification using macro averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.RecallMacro.objective_function

`RecallMacro.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: y_predicted (*pd.Series*) : predicted values of length [n_samples] y_true (*pd.Series*) : actual class labels of length [n_samples] X (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.RecallMacro.score

`RecallMacro.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.RecallMacro.validate_inputs

`RecallMacro.validate_inputs(y_true, y_predicted)`

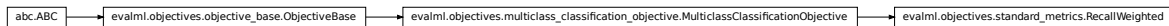
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.RecallWeighted



class evalml.objectives.RecallWeighted

Recall score for multiclass classification using weighted averaging.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.RecallWeighted.objective_function

`RecallWeighted.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: `y_predicted` (`pd.Series`) : predicted values of length `[n_samples]` `y_true` (`pd.Series`) : actual class labels of length `[n_samples]` `X` (`pd.DataFrame` or `np.array`) : extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.RecallWeighted.score`

`RecallWeighted.score(y_true, y_predicted, X=None)`

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **`y_predicted`** (`pd.Series`) – predicted values of length `[n_samples]`
- **`y_true`** (`pd.Series`) – actual class labels of length `[n_samples]`
- **`X`** (`pd.DataFrame` or `np.array`) – extra data of shape `[n_samples, n_features]` necessary to calculate score

Returns score

`evalml.objectives.RecallWeighted.validate_inputs`

`RecallWeighted.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

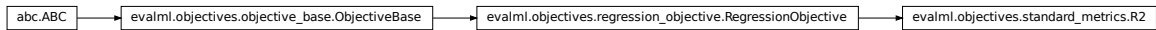
- **`y_predicted`** (`pd.Series`) – predicted values of length `[n_samples]`
- **`y_true`** (`pd.Series`) – actual class labels of length `[n_samples]`

Returns None

Regression Objectives

<code>R2</code>	Coefficient of determination for regression.
<code>MAE</code>	Mean absolute error for regression.
<code>MSE</code>	Mean squared error for regression.
<code>MeanSquaredLogError</code>	Mean squared log error for regression.
<code>MedianAE</code>	Median absolute error for regression.
<code>MaxError</code>	Maximum residual error for regression.
<code>ExpVariance</code>	Explained variance score for regression.
<code>RootMeanSquaredError</code>	Root mean squared error for regression.
<code>RootMeanSquaredLogError</code>	Root mean squared log error for regression.

evalml.objectives.R2



class evalml.objectives.R2
Coefficient of determination for regression.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.R2.objective_function

R2.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.R2.score

R2.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.R2.validate_inputs

R2.**validate_inputs** (*y_true*, *y_predicted*)

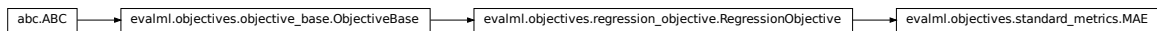
Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.MAE



class evalml.objectives.**MAE**
 Mean absolute error for regression.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.MAE.objective_function

MAE.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.MAE.score

MAE.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

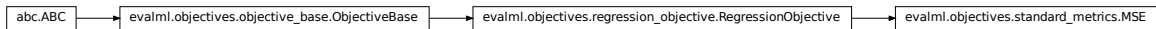
- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score**evalml.objectives.MAE.validate_inputs****MAE.validate_inputs** (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None**evalml.objectives.MSE****class** evalml.objectives.**MSE**

Mean squared error for regression.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.MSE.objective_function**MSE.objective_function** (*y_true*, *y_predicted*, *X=None*)**Computes the relative value of the provided predictions compared to the actual labels, according a specified metric**

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.MSE.score

MSE.score (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.MSE.validate_inputs

MSE.validate_inputs (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.MeanSquaredLogError



class evalml.objectives.MeanSquaredLogError

Mean squared log error for regression.

Only valid for nonnegative inputs. Otherwise, will throw a ValueError

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

`evalml.objectives.MeanSquaredLogError.objective_function`

`MeanSquaredLogError.objective_function` (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (`pd.Series`) : predicted values of length [*n_samples*] *y_true* (`pd.Series`) : actual class labels of length [*n_samples*] *X* (`pd.DataFrame` or `np.array`) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.MeanSquaredLogError.score`

`MeanSquaredLogError.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]
- ***X*** (`pd.DataFrame` or `np.array`) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.MeanSquaredLogError.validate_inputs`

`MeanSquaredLogError.validate_inputs` (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]

Returns None

`evalml.objectives.MedianAE`



class `evalml.objectives.MedianAE`
Median absolute error for regression.

Methods

<code>objective_function</code>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<code>score</code>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<code>validate_inputs</code>	Validates the input based on a few simple checks.

`evalml.objectives.MedianAE.objective_function`

`MedianAE.objective_function` (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (`pd.Series`) : predicted values of length [*n_samples*] *y_true* (`pd.Series`) : actual class labels of length [*n_samples*] *X* (`pd.DataFrame` or `np.array`) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.MedianAE.score`

`MedianAE.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]
- ***X*** (`pd.DataFrame` or `np.array`) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.MedianAE.validate_inputs`

`MedianAE.validate_inputs` (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]

Returns None

evalml.objectives.MaxError



class evalml.objectives.**MaxError**
Maximum residual error for regression.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.MaxError.objective_function

MaxError.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [*n_samples*] *y_true* (*pd.Series*) : actual class labels of length [*n_samples*] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.MaxError.score

MaxError.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [*n_samples*]
- **y_true** (*pd.Series*) – actual class labels of length [*n_samples*]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

evalml.objectives.MaxError.validate_inputs

MaxError.**validate_inputs** (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.ExpVariance

class evalml.objectives.**ExpVariance**

Explained variance score for regression.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.ExpVariance.objective_function

ExpVariance.**objective_function** (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (*pd.Series*) : predicted values of length [n_samples] *y_true* (*pd.Series*) : actual class labels of length [n_samples] *X* (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.ExpVariance.score

ExpVariance.**score** (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.ExpVariance.validate_inputs

`ExpVariance.validate_inputs(y_true, y_predicted)`

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.RootMeanSquaredError



class evalml.objectives.RootMeanSquaredError

Root mean squared error for regression.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

evalml.objectives.RootMeanSquaredError.objective_function

`RootMeanSquaredError.objective_function(y_true, y_predicted, X=None)`

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: **y_predicted** (*pd.Series*) : predicted values of length [n_samples] **y_true** (*pd.Series*) : actual class labels of length [n_samples] **X** (*pd.DataFrame* or *np.array*) : extra data of shape [n_samples, n_features] necessary to calculate score

Returns numerical value used to calculate score

evalml.objectives.RootMeanSquaredError.score

RootMeanSquaredError.**score**(*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]
- **X** (*pd.DataFrame* or *np.array*) – extra data of shape [n_samples, n_features] necessary to calculate score

Returns score

evalml.objectives.RootMeanSquaredError.validate_inputs

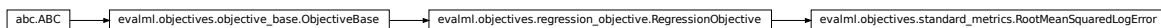
RootMeanSquaredError.**validate_inputs**(*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

- **y_predicted** (*pd.Series*) – predicted values of length [n_samples]
- **y_true** (*pd.Series*) – actual class labels of length [n_samples]

Returns None

evalml.objectives.RootMeanSquaredLogError**class evalml.objectives.RootMeanSquaredLogError**

Root mean squared log error for regression.

Only valid for nonnegative inputs. Otherwise, will throw a ValueError.

Methods

<i>objective_function</i>	Computes the relative value of the provided predictions compared to the actual labels, according a specified metric
<i>score</i>	Returns a numerical score indicating performance based on the differences between the predicted and actual values.
<i>validate_inputs</i>	Validates the input based on a few simple checks.

`evalml.objectives.RootMeanSquaredLogError.objective_function`

`RootMeanSquaredLogError.objective_function` (*y_true*, *y_predicted*, *X=None*)

Computes the relative value of the provided predictions compared to the actual labels, according a specified metric

Arguments: *y_predicted* (`pd.Series`) : predicted values of length [*n_samples*] *y_true* (`pd.Series`) : actual class labels of length [*n_samples*] *X* (`pd.DataFrame` or `np.array`) : extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns numerical value used to calculate score

`evalml.objectives.RootMeanSquaredLogError.score`

`RootMeanSquaredLogError.score` (*y_true*, *y_predicted*, *X=None*)

Returns a numerical score indicating performance based on the differences between the predicted and actual values.

Parameters

- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]
- ***X*** (`pd.DataFrame` or `np.array`) – extra data of shape [*n_samples*, *n_features*] necessary to calculate score

Returns score

`evalml.objectives.RootMeanSquaredLogError.validate_inputs`

`RootMeanSquaredLogError.validate_inputs` (*y_true*, *y_predicted*)

Validates the input based on a few simple checks.

Parameters

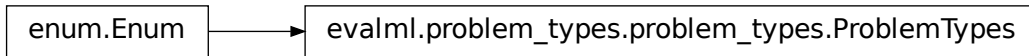
- ***y_predicted*** (`pd.Series`) – predicted values of length [*n_samples*]
- ***y_true*** (`pd.Series`) – actual class labels of length [*n_samples*]

Returns None

Problem Types

ProblemTypes

Enum for type of machine learning problem: BINARY, MULTICLASS, or REGRESSION.

evalml.problem_types.ProblemTypes

class evalml.problem_types.**ProblemTypes**

Enum for type of machine learning problem: BINARY, MULTICLASS, or REGRESSION.

handle_problem_types

Handles problem_type by either returning the ProblemTypes or converting from a str.

evalml.problem_types.handle_problem_types

evalml.problem_types.**handle_problem_types** (*problem_type*)

Handles problem_type by either returning the ProblemTypes or converting from a str.

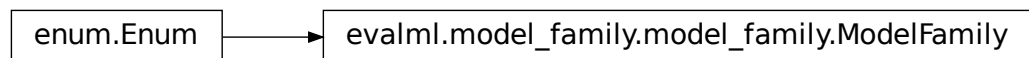
Parameters **problem_type** (*str* or *ProblemTypes*) – problem type that needs to be handled

Returns ProblemTypes

Model Family

ModelFamily

Enum for family of machine learning models.

evalml.model_family.ModelFamily

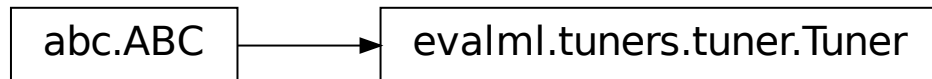
class evalml.model_family.**ModelFamily**

Enum for family of machine learning models.

Tuners

<i>Tuner</i>	Defines API for Tuners.
<i>SKOptTuner</i>	Bayesian Optimizer.
<i>GridSearchTuner</i>	Grid Search Optimizer.
<i>RandomSearchTuner</i>	Random Search Optimizer.

evalml.tuners.Tuner



class evalml.tuners.**Tuner** (*pipeline_hyperparameter_ranges*, *random_state=0*)

Defines API for Tuners.

Tuners implement different strategies for sampling from a search space. They're used in EvalML to search the space of pipeline hyperparameters.

Methods

<i>__init__</i>	Base Tuner class
<i>add</i>	Register a set of hyperparameters with the score obtained from training a pipeline with those hyperparameters.
<i>is_search_space_exhausted</i>	Optional.
<i>propose</i>	Returns a suggested set of parameters to train and score a pipeline with, based off the search space dimensions and prior samples.

evalml.tuners.Tuner.__init__

Tuner.__init__ (*pipeline_hyperparameter_ranges*, *random_state=0*)

Base Tuner class

Parameters

- **pipeline_hyperparameter_ranges** (*dict*) – a set of hyperparameter ranges corresponding to a pipeline's parameters
- **random_state** (*int*, *np.random.RandomState*) – The random state

evalml.tuners.Tuner.add**Tuner.add**(*pipeline_parameters*, *score*)

Register a set of hyperparameters with the score obtained from training a pipeline with those hyperparameters.

Parameters

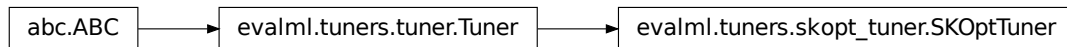
- **pipeline_parameters** (*dict*) – a dict of the parameters used to evaluate a pipeline
- **score** (*float*) – the score obtained by evaluating the pipeline with the provided parameters

Returns None**evalml.tuners.Tuner.is_search_space_exhausted****Tuner.is_search_space_exhausted**()

Optional. If possible search space for tuner is finite, this method indicates whether or not all possible parameters have been scored.

Returns Returns true if all possible parameters in a search space has been scored.**Return type** bool**evalml.tuners.Tuner.propose****Tuner.propose**()

Returns a suggested set of parameters to train and score a pipeline with, based off the search space dimensions and prior samples.

Returns proposed pipeline parameters**Return type** dict**evalml.tuners.SKOptTuner**

class evalml.tuners.**SKOptTuner**(*pipeline_hyperparameter_ranges*, *random_state=0*)
 Bayesian Optimizer.

Methods`__init__`

Init SKOptTuner

Continued on next page

Table 168 – continued from previous page

<code>add</code>	Add score to sample
<code>is_search_space_exhausted</code>	Optional.
<code>propose</code>	Returns a suggested set of parameters to train and score a pipeline with, based off the search space dimensions and prior samples.

`evalml.tuners.SKOptTuner.__init__`

`SKOptTuner.__init__(pipeline_hyperparameter_ranges, random_state=0)`

Init SKOptTuner

Parameters

- **`pipeline_hyperparameter_ranges`** (*dict*) – a set of hyperparameter ranges corresponding to a pipeline’s parameters
- **`random_state`** (*int*, *np.random.RandomState*) – The random state

`evalml.tuners.SKOptTuner.add`

`SKOptTuner.add(pipeline_parameters, score)`

Add score to sample

Parameters

- **`pipeline_parameters`** (*dict*) – a dict of the parameters used to evaluate a pipeline
- **`score`** (*float*) – the score obtained by evaluating the pipeline with the provided parameters

Returns None

`evalml.tuners.SKOptTuner.is_search_space_exhausted`

`SKOptTuner.is_search_space_exhausted()`

Optional. If possible search space for tuner is finite, this method indicates whether or not all possible parameters have been scored.

Returns Returns true if all possible parameters in a search space has been scored.

Return type bool

`evalml.tuners.SKOptTuner.propose`

`SKOptTuner.propose()`

Returns a suggested set of parameters to train and score a pipeline with, based off the search space dimensions and prior samples.

Returns proposed pipeline parameters

Return type dict

evalml.tuners.GridSearchTuner



class evalml.tuners.**GridSearchTuner** (*pipeline_hyperparameter_ranges*, *n_points=10*, *random_state=0*)
Grid Search Optimizer.

Example

```

>>> tuner = GridSearchTuner({'My Component': {'param a': [0.0, 10.0], 'param b': [
  ↳ 'a', 'b', 'c']}}, n_points=5)
>>> proposal = tuner.propose()
>>> assert proposal.keys() == {'My Component'}
>>> assert proposal['My Component'] == {'param a': 0.0, 'param b': 'a'}
  
```

Methods

<code>__init__</code>	Generate all of the possible points to search for in the grid
<code>add</code>	Not applicable to grid search tuner as generated parameters are not dependent on scores of previous parameters.
<code>is_search_space_exhausted</code>	Checks if it is possible to generate a set of valid parameters.
<code>propose</code>	Returns parameters from <code>_grid_points</code> iterations

evalml.tuners.GridSearchTuner.__init__

`GridSearchTuner.__init__` (*pipeline_hyperparameter_ranges*, *n_points=10*, *random_state=0*)
Generate all of the possible points to search for in the grid

Parameters

- **pipeline_hyperparameter_ranges** (*dict*) – a set of hyperparameter ranges corresponding to a pipeline’s parameters
- **n_points** – The number of points to sample from along each dimension defined in the *space* argument
- **random_state** – Unused in this class

`evalml.tuners.GridSearchTuner.add`

`GridSearchTuner.add(pipeline_parameters, score)`

Not applicable to grid search tuner as generated parameters are not dependent on scores of previous parameters.

Parameters

- **pipeline_parameters** (*dict*) – a dict of the parameters used to evaluate a pipeline
- **score** (*float*) – the score obtained by evaluating the pipeline with the provided parameters

`evalml.tuners.GridSearchTuner.is_search_space_exhausted`

`GridSearchTuner.is_search_space_exhausted()`

Checks if it is possible to generate a set of valid parameters. Stores generated parameters in `self.curr_params` to be returned by `propose()`.

Raises `NoParamsException` – If a search space is exhausted, then this exception is thrown.

Returns If no more valid parameters exists in the search space, return false.

Return type `bool`

`evalml.tuners.GridSearchTuner.propose`

`GridSearchTuner.propose()`

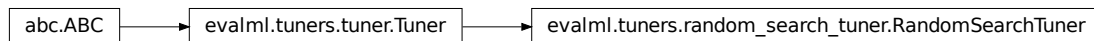
Returns parameters from `_grid_points` iterations

If all possible combinations of parameters have been scored, then `NoParamsException` is raised.

Returns proposed pipeline parameters

Return type `dict`

`evalml.tuners.RandomSearchTuner`



```
class evalml.tuners.RandomSearchTuner(pipeline_hyperparameter_ranges,          ran-  
                                     dom_state=0, with_replacement=False, replace-  
                                     ment_max_attempts=10)
```

Random Search Optimizer.

Example

```
>>> tuner = RandomSearchTuner({'My Component': {'param a': [0.0, 10.0], 'param b': ['a', 'b', 'c']}}, random_state=42)
>>> proposal = tuner.propose()
>>> assert proposal.keys() == {'My Component'}
>>> assert proposal['My Component'] == {'param a': 3.7454011884736254, 'param b': 'c'}
```

Methods

<code>__init__</code>	Sets up check for duplication if needed.
<code>add</code>	Not applicable to random search tuner as generated parameters are not dependent on scores of previous parameters.
<code>is_search_space_exhausted</code>	Checks if it is possible to generate a set of valid parameters.
<code>propose</code>	Generate a unique set of parameters.

evalml.tuners.RandomSearchTuner.__init__

`RandomSearchTuner.__init__(pipeline_hyperparameter_ranges, random_state=0, with_replacement=False, replacement_max_attempts=10)`

Sets up check for duplication if needed.

Parameters

- **pipeline_hyperparameter_ranges** (*dict*) – a set of hyperparameter ranges corresponding to a pipeline’s parameters
- **random_state** – Unused in this class
- **with_replacement** – If false, only unique hyperparameters will be shown
- **replacement_max_attempts** – The maximum number of tries to get a unique set of random parameters. Only used if tuner is initialized with `with_replacement=True`

evalml.tuners.RandomSearchTuner.add

`RandomSearchTuner.add(pipeline_parameters, score)`

Not applicable to random search tuner as generated parameters are not dependent on scores of previous parameters.

Parameters

- **pipeline_parameters** (*dict*) – a dict of the parameters used to evaluate a pipeline
- **score** (*float*) – the score obtained by evaluating the pipeline with the provided parameters

evalml.tuners.RandomSearchTuner.is_search_space_exhausted

`RandomSearchTuner.is_search_space_exhausted()`

Checks if it is possible to generate a set of valid parameters. Stores generated parameters in `self`.

`curr_params` to be returned by `propose()`.

Raises `NoParamsException` – If a search space is exhausted, then this exception is thrown.

Returns If no more valid parameters exists in the search space, return false.

Return type bool

`evalml.tuners.RandomSearchTuner.propose`

`RandomSearchTuner.propose()`

Generate a unique set of parameters.

If tuner was initialized with `with_replacement=True` and the tuner is unable to generate a unique set of parameters after `replacement_max_attempts` tries, then `NoParamsException` is raised.

Returns proposed pipeline parameters

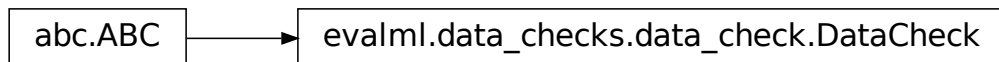
Return type dict

Data Checks

Data Check Classes

<i>DataCheck</i>	Base class for all data checks.
<i>InvalidTargetDataCheck</i>	Checks if the target labels contain missing or invalid data.
<i>HighlyNullDataCheck</i>	Checks if there are any highly-null columns in the input.
<i>IDColumnsDataCheck</i>	Check if any of the features are likely to be ID columns.
<i>LabelLeakageDataCheck</i>	Check if any of the features are highly correlated with the target.
<i>OutliersDataCheck</i>	Checks if there are any outliers in input data by using an Isolation Forest to obtain the anomaly score of each index and then using IQR to determine score anomalies.

`evalml.data_checks.DataCheck`



class `evalml.data_checks.DataCheck`

Base class for all data checks. Data checks are a set of heuristics used to determine if there are problems with input data.

name = `'DataCheck'`

Instance attributes

Methods:

<code>validate</code>	Inspects and validates the input data, runs any necessary calculations or algorithms, and returns a list of warnings and errors if applicable.
-----------------------	--

`evalml.data_checks.DataCheck.validate`

`DataCheck.validate` (*X*, *y=None*)

Inspects and validates the input data, runs any necessary calculations or algorithms, and returns a list of warnings and errors if applicable.

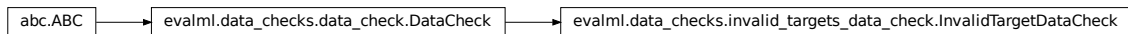
Parameters

- **X** (*pd.DataFrame*) – the input data of shape [n_samples, n_features]
- **y** (*pd.Series*, *optional*) – the target data of length [n_samples]

Returns list of `DataCheckError` and `DataCheckWarning` objects

Return type list (*DataCheckMessage*)

`evalml.data_checks.InvalidTargetDataCheck`



class `evalml.data_checks.InvalidTargetDataCheck`

Checks if the target labels contain missing or invalid data.

name = `'InvalidTargetDataCheck'`

Instance attributes

Methods:

<code>validate</code>	Checks if the target labels contain missing or invalid data.
-----------------------	--

evalml.data_checks.InvalidTargetDataCheck.validate

`InvalidTargetDataCheck.validate(X, y)`

Checks if the target labels contain missing or invalid data.

Parameters

- **X** (`pd.DataFrame`, `pd.Series`, `np.array`, `list`) – Features. Ignored.
- **y** – Target labels to check for invalid data.

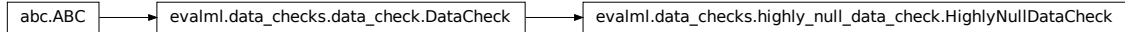
Returns list with `DataCheckErrors` if any invalid data is found in target labels.

Return type list (`DataCheckError`)

Example

```
>>> X = pd.DataFrame({})
>>> y = pd.Series([0, 1, None, None])
>>> target_check = InvalidTargetDataCheck()
>>> assert target_check.validate(X, y) == [DataCheckError("2 row(s) (50.0%)
of target values are null", "InvalidTargetDataCheck")]
```

evalml.data_checks.HighlyNullDataCheck



class `evalml.data_checks.HighlyNullDataCheck` (`pct_null_threshold=0.95`)

Checks if there are any highly-null columns in the input.

name = `'HighlyNullDataCheck'`

Instance attributes

—

Methods:

<code>__init__</code>	Checks if there are any highly-null columns in the input.
<code>validate</code>	Checks if there are any highly-null columns in the input.

evalml.data_checks.HighlyNullDataCheck.__init__

`HighlyNullDataCheck.__init__(pct_null_threshold=0.95)`

Checks if there are any highly-null columns in the input.

Parameters `pct_null_threshold` (*float*) – If the percentage of NaN values in an input feature exceeds this amount, that feature will be considered highly-null. Defaults to 0.95.

`evalml.data_checks.HighlyNullDataCheck.validate`

`HighlyNullDataCheck.validate` (*X*, *y=None*)

Checks if there are any highly-null columns in the input.

Parameters

- **X** (*pd.DataFrame*, *pd.Series*, *np.array*, *list*) – features
- **y** – Ignored.

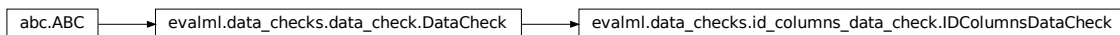
Returns list with a `DataCheckWarning` if there are any highly-null columns.

Return type list (*DataCheckWarning*)

Example

```
>>> df = pd.DataFrame({
...     'lots_of_null': [None, None, None, None, 5],
...     'no_null': [1, 2, 3, 4, 5]
... })
>>> null_check = HighlyNullDataCheck(pct_null_threshold=0.8)
>>> assert null_check.validate(df) == [DataCheckWarning("Column 'lots_of_null"
↪ ' is 80.0% or more null", "HighlyNullDataCheck")]
```

`evalml.data_checks.IDColumnsDataCheck`



class `evalml.data_checks.IDColumnsDataCheck` (*id_threshold=1.0*)

Check if any of the features are likely to be ID columns.

name = 'IDColumnsDataCheck'

Instance attributes

—

Methods:

<code>__init__</code>	Check if any of the features are likely to be ID columns.
<code>validate</code>	Check if any of the features are likely to be ID columns.

evalml.data_checks.IDColumnsDataCheck.__init__

IDColumnsDataCheck.__init__(id_threshold=1.0)

Check if any of the features are likely to be ID columns.

Parameters **id_threshold** (*float*) – the probability threshold to be considered an ID column. Defaults to 1.0.

evalml.data_checks.IDColumnsDataCheck.validate

IDColumnsDataCheck.validate(X, y=None)

Check if any of the features are likely to be ID columns. Currently performs these simple checks:

- column name is “id”
- column name ends in “_id”
- column contains all unique values (and is not float / boolean)

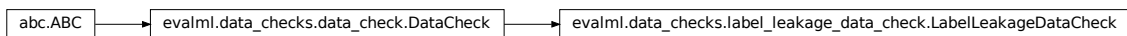
Parameters

- **X** (*pd.DataFrame*) – The input features to check
- **threshold** (*float*) – the probability threshold to be considered an ID column. Defaults to 1.0

Returns A dictionary of features with column name or index and their probability of being ID columns

Example

```
>>> df = pd.DataFrame({
...     'df_id': [0, 1, 2, 3, 4],
...     'x': [10, 42, 31, 51, 61],
...     'y': [42, 54, 12, 64, 12]
... })
>>> id_col_check = IDColumnsDataCheck()
>>> assert id_col_check.validate(df) == [DataCheckWarning("Column 'df_id' is
↳ 100.0% or more likely to be an ID column", "IDColumnsDataCheck")]
```

evalml.data_checks.LabelLeakageDataCheck

class evalml.data_checks.LabelLeakageDataCheck (pct_corr_threshold=0.95)

Check if any of the features are highly correlated with the target.

name = 'LabelLeakageDataCheck'

Instance attributes

Methods:

<code>__init__</code>	Check if any of the features are highly correlated with the target.
<code>validate</code>	Check if any of the features are highly correlated with the target.

`evalml.data_checks.LabelLeakageDataCheck.__init__`

`LabelLeakageDataCheck.__init__(pct_corr_threshold=0.95)`

Check if any of the features are highly correlated with the target.

Currently only supports binary and numeric targets and features.

Parameters `pct_corr_threshold` (*float*) – The correlation threshold to be considered leakage. Defaults to 0.95.

`evalml.data_checks.LabelLeakageDataCheck.validate`

`LabelLeakageDataCheck.validate(X, y)`

Check if any of the features are highly correlated with the target.

Currently only supports binary and numeric targets and features.

Parameters

- `X` (*pd.DataFrame*) – The input features to check
- `y` (*pd.Series*) – The labels

Returns list with a `DataCheckWarning` if there is label leakage detected.

Return type list (*DataCheckWarning*)

Example

```
>>> X = pd.DataFrame({
...     'leak': [10, 42, 31, 51, 61],
...     'x': [42, 54, 12, 64, 12],
...     'y': [12, 5, 13, 74, 24],
... })
>>> y = pd.Series([10, 42, 31, 51, 40])
>>> label_leakage_check = LabelLeakageDataCheck(pct_corr_threshold=0.8)
>>> assert label_leakage_check.validate(X, y) == [DataCheckWarning("Column
↪ 'leak' is 80.0% or more correlated with the target", "LabelLeakageDataCheck
↪ ")]
```

evalml.data_checks.OutliersDataCheck



```
class evalml.data_checks.OutliersDataCheck (random_state=0)
```

Checks if there are any outliers in input data by using an Isolation Forest to obtain the anomaly score of each index and then using IQR to determine score anomalies. Indices with score anomalies are considered outliers.

```
name = 'OutliersDataCheck'
```

Instance attributes

—

Methods:

<code>__init__</code>	Checks if there are any outliers in the input data.
<code>validate</code>	Checks if there are any outliers in a dataframe by using an Isolation Forest to obtain the anomaly score of each index and then using IQR to determine score anomalies.

evalml.data_checks.OutliersDataCheck.__init__

```
OutliersDataCheck.__init__(random_state=0)
```

Checks if there are any outliers in the input data.

Parameters `random_state` (`int`, `np.random.RandomState`) – The random seed/state. Defaults to 0.

evalml.data_checks.OutliersDataCheck.validate

```
OutliersDataCheck.validate(X, y=None)
```

Checks if there are any outliers in a dataframe by using an Isolation Forest to obtain the anomaly score of each index and then using IQR to determine score anomalies. Indices with score anomalies are considered outliers.

Parameters

- `X` (`pd.DataFrame`) – features
- `y` – Ignored.

Returns A set of indices that may have outlier data.

Example

```
>>> df = pd.DataFrame({
...     'x': [1, 2, 3, 40, 5],
...     'y': [6, 7, 8, 990, 10],
...     'z': [-1, -2, -3, -1201, -4]
... })
>>> outliers_check = OutliersDataCheck()
>>> assert outliers_check.validate(df) == [DataCheckWarning("Row '3' is_
↳ likely to have outlier data", "OutliersDataCheck")]
```

<i>DataChecks</i>	A collection of data checks.
<i>DefaultDataChecks</i>	A collection of basic data checks that is used by AutoML by default.

evalml.data_checks.DataChecks

evalml.data_checks.data_checks.DataChecks

class evalml.data_checks.**DataChecks** (*data_checks=None*)
 A collection of data checks.

Methods

<i>__init__</i>	A collection of data checks.
<i>validate</i>	Inspects and validates the input data against data checks and returns a list of warnings and errors if applicable.

evalml.data_checks.DataChecks.__init__

DataChecks.__init__ (*data_checks=None*)
 A collection of data checks.

Parameters **data_checks** (*list* (*DataCheck*)) – list of DataCheck objects

evalml.data_checks.DataChecks.validate

`DataChecks.validate` (X , $y=None$)

Inspects and validates the input data against data checks and returns a list of warnings and errors if applicable.

Parameters

- **X** (*pd.DataFrame*) – the input data of shape [n_samples, n_features]
- **y** (*pd.Series*) – the target labels of length [n_samples]

Returns list containing `DataCheckMessage` objects

Return type list (*DataCheckMessage*)

evalml.data_checks.DefaultDataChecks

evalml.data_checks.data_checks.DataChecks

evalml.data_checks.default_data_checks.DefaultDataChecks

class evalml.data_checks.DefaultDataChecks (*data_checks=None*)

A collection of basic data checks that is used by AutoML by default. Includes `HighlyNullDataCheck`, `IDColumnsDataCheck`, and `LabelLeakageDataCheck`.

Methods

<code>__init__</code>	A collection of basic data checks.
<code>validate</code>	Inspects and validates the input data against data checks and returns a list of warnings and errors if applicable.

evalml.data_checks.DefaultDataChecks.__init__

`DefaultDataChecks.__init__` (*data_checks=None*)

A collection of basic data checks.

Parameters **data_checks** (*list (DataCheck)*) – Ignored.

evalml.data_checks.DefaultDataChecks.validate

`DefaultDataChecks.validate` (X , $y=None$)

Inspects and validates the input data against data checks and returns a list of warnings and errors if applicable.

Parameters

- **X** (*pd.DataFrame*) – the input data of shape [n_samples, n_features]

- **y** (*pd.Series*) – the target labels of length [n_samples]

Returns list containing DataCheckMessage objects

Return type list (*DataCheckMessage*)

Data Check Messages

<i>DataCheckMessage</i>	Base class for all DataCheckMessages.
<i>DataCheckError</i>	DataCheckMessage subclass for errors returned by data checks.
<i>DataCheckWarning</i>	DataCheckMessage subclass for warnings returned by data checks.

evalml.data_checks.DataCheckMessage

evalml.data_checks.data_check_message.DataCheckMessage

class evalml.data_checks.DataCheckMessage (*message, data_check_name*)

Base class for all DataCheckMessages.

message_type = None

Methods:

<code>__init__</code>	Message returned by a DataCheck, tagged by name.”
<code>__str__</code>	String representation of data check message, equivalent to self.message attribute.
<code>__eq__</code>	Checks for equality.

evalml.data_checks.DataCheckMessage.__init__

DataCheckMessage.**__init__** (*message, data_check_name*)

Message returned by a DataCheck, tagged by name.”

Parameters

- **message** (*str*) – Message string
- **data_check_name** (*str*) – Name of data check

evalml.data_checks.DataCheckMessage.__str__

DataCheckMessage.__str__()

String representation of data check message, equivalent to self.message attribute.

evalml.data_checks.DataCheckMessage.__eq__

DataCheckMessage.__eq__(*other*)

Checks for equality. Two DataCheckMessage objs are considered equivalent if their message type and message are equivalent.

evalml.data_checks.DataCheckError

class evalml.data_checks.DataCheckError(*message*, *data_check_name*)

DataCheckMessage subclass for errors returned by data checks.

message_type = 'error'

Methods:

<code>__init__</code>	Message returned by a DataCheck, tagged by name.”
<code>__str__</code>	String representation of data check message, equivalent to self.message attribute.
<code>__eq__</code>	Checks for equality.

evalml.data_checks.DataCheckError.__init__

DataCheckError.__init__(*message*, *data_check_name*)

Message returned by a DataCheck, tagged by name.”

Parameters

- **message** (*str*) – Message string
- **data_check_name** (*str*) – Name of data check

evalml.data_checks.DataCheckError.__str__

DataCheckError.__str__()

String representation of data check message, equivalent to self.message attribute.

evalml.data_checks.DataCheckError.__eq__`DataCheckError.__eq__` (*other*)

Checks for equality. Two `DataCheckMessage` objs are considered equivalent if their message type and message are equivalent.

evalml.data_checks.DataCheckWarning

class `evalml.data_checks.DataCheckWarning` (*message, data_check_name*)

`DataCheckMessage` subclass for warnings returned by data checks.

message_type = 'warning'

Methods:

<code>__init__</code>	Message returned by a <code>DataCheck</code> , tagged by name.”
<code>__str__</code>	String representation of data check message, equivalent to <code>self.message</code> attribute.
<code>__eq__</code>	Checks for equality.

evalml.data_checks.DataCheckWarning.__init__

`DataCheckWarning.__init__` (*message, data_check_name*)

Message returned by a `DataCheck`, tagged by name.”

Parameters

- **message** (*str*) – Message string
- **data_check_name** (*str*) – Name of data check

evalml.data_checks.DataCheckWarning.__str__

`DataCheckWarning.__str__` ()

String representation of data check message, equivalent to `self.message` attribute.

evalml.data_checks.DataCheckWarning.__eq__

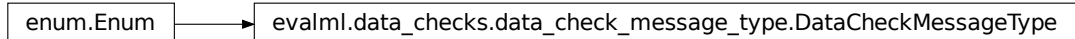
`DataCheckWarning.__eq__` (*other*)

Checks for equality. Two `DataCheckMessage` objs are considered equivalent if their message type and message are equivalent.

Data Check Message Types

<i>DataCheckMessageType</i>	Enum for type of data check message: WARNING or ERROR.
-----------------------------	--

evalml.data_checks.DataCheckMessageType



class evalml.data_checks.DataCheckMessageType
Enum for type of data check message: WARNING or ERROR.

Utils

<i>import_or_raise</i>	Attempts to import the requested library by name.
<i>convert_to_seconds</i>	Converts a string describing a length of time to its length in seconds.
<i>get_random_state</i>	Generates a numpy.random.RandomState instance using seed.
<i>get_random_seed</i>	Given a numpy.random.RandomState object, generate an int representing a seed value for another random number generator.

evalml.utils.import_or_raise

evalml.utils.import_or_raise(*library*, *error_msg=None*)
Attempts to import the requested library by name. If the import fails, raises an ImportError.

Parameters

- **library** (*str*) – the name of the library
- **error_msg** (*str*) – error message to return if the import fails

evalml.utils.convert_to_seconds

evalml.utils.convert_to_seconds(*input_str*)
Converts a string describing a length of time to its length in seconds.

evalml.utils.get_random_state

evalml.utils.get_random_state(*seed*)
Generates a numpy.random.RandomState instance using seed.

Parameters `seed` (*None, int, np.random.RandomState object*) – seed to use to generate `numpy.random.RandomState`. Must be between `SEED_BOUNDS.min_bound` and `SEED_BOUNDS.max_bound`, inclusive. Otherwise, an exception will be thrown.

`evalml.utils.get_random_seed`

`evalml.utils.get_random_seed(random_state, min_bound=0, max_bound=2147483647)`

Given a `numpy.random.RandomState` object, generate an int representing a seed value for another random number generator. Or, if given an int, return that int.

To protect against invalid input to a particular library’s random number generator, if an int value is provided, and it is outside the bounds “[`min_bound`, `max_bound`)”, the value will be projected into the range between the `min_bound` (inclusive) and `max_bound` (exclusive) using modular arithmetic.

Parameters

- **`random_state`** (*int, numpy.random.RandomState*) – random state
- **`min_bound`** (*None, int*) – if not default of `None`, will be min bound when generating seed (inclusive). Must be less than `max_bound`.
- **`max_bound`** (*None, int*) – if not default of `None`, will be max bound when generating seed (exclusive). Must be greater than `min_bound`.

Returns seed for random number generator

Return type int

1.6.17 FAQ

What is the difference between EvalML and other AutoML libraries?

EvalML optimizes machine learning pipelines on *custom practical objectives* instead of vague machine learning loss functions so that it will find the best pipelines for your specific needs. Furthermore, EvalML *pipelines* are able to take in all kinds of data (missing values, categorical, etc.) as long as the data are in a single table. EvalML also allows you to build your own pipelines with existing or custom components so you can have more control over the AutoML process. Moreover, EvalML also provides you with support in the form of *data checks* to ensure that you are aware of potential issues your data may cause with machine learning algorithms.

How does EvalML handle missing values?

EvalML contains imputation components in its pipelines so that missing values are taken care of. EvalML optimizes over different types of imputation to search for the best possible pipeline. You can find more information about components [here](#) and in the API reference [here](#).

How does EvalML handle categorical encoding?

EvalML provides a *one-hot-encoding component* in its pipelines for categorical variables. EvalML plans to support other encoders in the future.

How does EvalML handle feature selection?

EvalML currently utilizes scikit-learn’s `SelectFromModel` with a Random Forest classifier/regressor to handle feature selection. EvalML plans on supporting more feature selectors in the future. You can find more information in the API reference [here](#).

How is feature importance calculated?

Feature importance depends on the estimator used. Variable coefficients are used for regression-based estimators (Logistic Regression and Linear Regression) and Gini importance is used for tree-based estimators (Random Forest and XGBoost).

How does hyperparameter tuning work?

EvalML tunes hyperparameters for its pipelines through Bayesian optimization. In the future we plan to support more optimization techniques such as random search.

Can I create my own objective metric?

Yes you can! You can *create your own custom objective* so that EvalML optimizes the best model for your needs.

How does EvalML avoid overfitting?

EvalML provides *data checks* to combat overfitting. Such data checks include detecting label leakage, unstable pipelines, hold-out datasets and cross validation. EvalML defaults to using Stratified K-Fold cross-validation for classification problems and K-Fold cross-validation for regression problems but allows you to utilize your own cross-validation methods as well.

Can I create my own pipeline for EvalML?

Yes! EvalML allows you to create *custom pipelines* using modular components. This allows you to customize EvalML pipelines for your own needs or for AutoML.

Does EvalML work with X algorithm?

EvalML is constantly improving and adding new components and will allow your own algorithms to be used as components in our pipelines.

[]:

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