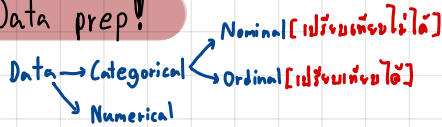


## Data prep!



→ `df.info()` : ดูว่า type

→ `df.describe()` : ดูว่าสถิติ

→ ดูว่าข้อมูล valid มั้ย เช่น อาจไม่มีค่า 0, ไม่เกิน 100 ...

\* drop ตัวแปรที่ไม่เกี่ยวข้องกัน เช่น National ID ไม่เกี่ยวข้องกับการทำนาย \* ต้องทำ

\* ตัวที่มี missing value > 50% [Unquality data]

\* ตัวแปรประเภท Categorical → ไม่ควร unique เกินไป  
↳ ไม่ควร flat เกินไป

## Preparing feature for ML

① Impute missing values (ใช้ impute target ⇒ ต้อง remove ทั้ง)

- numerical variable ⇒ Mean, Median
- Categorical variable ⇒ Mode (most-freq)

② Categorical to numeric variable

→ One-hot vector [dummy code] \* n-1 แล้ว drop

```
pd.get_dummies(df, column=['Sex', 'Embarked'], drop_first=True)
```

③ เอา Outlier ออก

→ ตัดที่ percentile ขึ้น-ลง อยู่ที่ 1%

→  $\mu \pm 3 \text{ S.D.}$

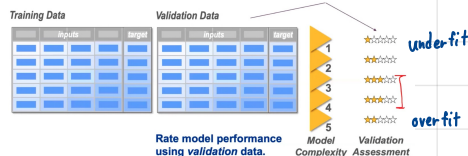
④ ถ้าการแบ่งช่วงข้อมูล ⇒ log, binning

## How Test/Train/Validate

- training data ⇒ model construction
- testing data ⇒ model evaluation
- Validating data ⇒ model tuning

\* ต้อง stratify ด้วย

### Model Performance Assessment



## Feature Selection based-on Statistics

### 1.13.2. Univariate feature selection ¶ ดูแล้ว target ⇔ สัมพันธ์กับหรือไม่

Univariate feature selection works by selecting the best features based on univariate statistical tests. It can be seen as a preprocessing step to an estimator. Scikit-learn exposes feature selection routines as objects that implement the `transform` method:

- `SelectKBest` removes all but the  $k$  highest scoring features
- `SelectPercentile` removes all but a user-specified highest scoring percentage of features
- using common univariate statistical tests for each feature: false positive rate `SelectFpr`, false discovery rate `SelectFdr`, or family wise error `SelectFwe`.
- `GenericUnivariateSelect` allows to perform univariate feature selection with a configurable strategy. This allows to select the best univariate selection strategy with hyper-parameter search estimator.

For instance, we can perform a  $\chi^2$  test to the samples to retrieve only the two best features as follows:

```
>>> from sklearn.datasets import load_iris
>>> from sklearn.feature_selection import SelectKBest
>>> from sklearn.feature_selection import chi2
>>> X, y = load_iris(return_X_y=True)
>>> X.shape
(150, 4)
>>> X_new = SelectKBest(chi2, k=2).fit_transform(X, y) # ใน X_new มีแค่ 2 feature ที่สัมพันธ์กับ target
>>> X_new.shape
(150, 2)
```

[https://scikit-learn.org/stable/modules/feature\\_selection.html](https://scikit-learn.org/stable/modules/feature_selection.html)

These objects take as input a scoring function that returns univariate scores and p-values (or only scores for `SelectKBest` and `SelectPercentile`):

- For regression: `f_regression`, `mutual_info_regression`, `r_regression` ⇔ ดูว่าความสัมพันธ์กัน
- For classification: `chi2`, `f_classif`, `mutual_info_classif`

## based-on Model

### 1.13.4. Feature selection using `SelectFromModel`

`SelectFromModel` is a meta-transformer that can be used alongside any estimator that assigns importance to each feature through a specific attribute (such as `coef_`, `feature_importances_`) or via an `importance_getter` callable after fitting. The features are considered unimportant and removed if the corresponding importance of the feature values are below the provided `threshold` parameter. Apart from specifying the threshold numerically, there are built-in heuristics for finding a threshold using a string argument. Available heuristics are "mean", "median" and float multiples of these like "0.1\*mean". In combination with the `threshold` criteria, one can use the `max_features` parameter to set a limit on the number of features to select.

For examples on how it is to be used refer to the sections below.

#### Examples

- Model-based and sequential feature selection

#### 1.13.4.1. L1-based feature selection

Linear models penalized with the L1 norm have sparse solutions: many of their estimated coefficients are zero. When the goal is to reduce the dimensionality of the data to use with another classifier, they can be used along with `SelectFromModel` to select the non-zero coefficients. In particular, sparse estimators useful for this purpose are the Lasso for regression, and of `LogisticRegression` and `LinearSVC` for classification:

```
>>> from sklearn.svm import LinearSVC
>>> from sklearn.datasets import load_iris
>>> from sklearn.feature_selection import SelectFromModel
>>> X, y = load_iris(return_X_y=True)
>>> X.shape
(150, 4)
>>> lsvc = LinearSVC(C=0.01, penalty="l1", dual=False).fit(X, y)
>>> model = SelectFromModel(lsvc, prefit=True)
>>> X_new = model.transform(X)
>>> X_new.shape
(150, 3)
```

## scikit-learn

Machine learning in Python

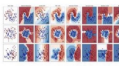
- Simple and efficient tools for predictive data analysis
- Accessible to everybody, and reusable in various contexts
- Built on NumPy, SciPy, and matplotlib
- Open source, commercially usable - BSD license

## Classification Category

Identifying which category an object belongs to.

**Applications:** Spam detection, image recognition.

**Algorithms:** SVM, nearest neighbors, random forest, and more...



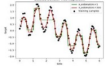
Examples

## Regression Numerical

Predicting a continuous valued attribute associated with an object.

**Applications:** Drug response, stock prices.

**Algorithms:** SVR, nearest neighbors, random forest, and more...



Examples

## Clustering

Automatic grouping of similar objects into sets.

**Applications:** Customer segmentation, grouping coherent outcomes.

**Algorithms:** K-Means, spectral clustering, mean-shift, and more...



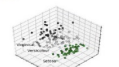
Examples

## Dimensionality reduction

Reducing the number of random variables to consider.

**Applications:** Visualization, increased efficiency.

**Algorithms:** PCA, feature selection, non-negative matrix factorization, and more...



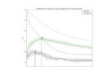
Examples

## Model selection

Comparing, validating and choosing parameters and models.

**Applications:** Improved accuracy via parameter tuning.

**Algorithms:** grid search, cross validation, metrics, and more...



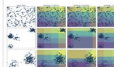
Examples

## Preprocessing

Feature extraction and normalization.

**Applications:** Transforming input data such as text for use with machine learning algorithms.

**Algorithms:** preprocessing, feature extraction, and more...



Examples

- Supervised  $\Rightarrow$  with target
- Unsupervised  $\Rightarrow$  without target
- RL

## CHULA ENGINEERING

## Estimator Interface

## Decision Trees

We'll start just by training a single decision tree.

```

In [1]: from sklearn.tree import DecisionTreeClassifier

In [11]: from sklearn.datasets import load_breast_cancer
In [12]: X_train, X_test, y_train, y_test = train_test_split(
    cancer.data, cancer.target,
    stratify=cancer.target, random_state=1337)

tree = DecisionTreeClassifier(random_state=7331)
tree.fit(X_train, y_train)

```

## Prediction and Evaluation

Let's evaluate our decision tree.

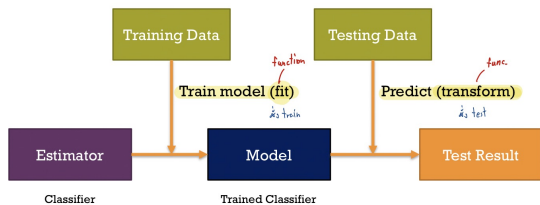
```

In [13]: predictions = tree.predict(X_test)

In [14]: from sklearn.metrics import classification_report, confusion_matrix

In [15]: print(classification_report(y_test, predictions))

```



## Training Phase

```

from sklearn.tree import DecisionTreeClassifier
from sklearn.datasets import load_breast_cancer
from sklearn.model_selection import train_test_split

cancer = load_breast_cancer() # Get some data
X_train, X_test, y_train, y_test = train_test_split(
    cancer.data, cancer.target,
    stratify=cancer.target, random_state=1337)

tree = DecisionTreeClassifier(random_state=7331)
tree.fit(X_train, y_train) # Learn a Decision Function

```

Ref: C. Travis Johnson, Mines Linux Users Group on April 27, 2017

## CHULA ENGINEERING

## Testing Phase (Prediction)

```

>>> from sklearn.metrics import classification_report
>>> y_true = [0, 1, 2, 2, 2]
>>> y_pred = [0, 0, 2, 2, 1]
>>> target_names = ['class 0', 'class 1', 'class 2']
>>> print(classification_report(y_true, y_pred, target_names=target_names))

```

	precision	recall	f1-score	support
class 0	0.50	1.00	0.67	1
class 1	0.00	0.00	0.00	1
class 2	1.00	0.67	0.80	3
micro avg	0.60	0.60	0.60	5
macro avg	0.50	0.56	0.49	5
weighted avg	0.70	0.60	0.61	5

Decision Tree Classifier  $\Rightarrow$  ใช้กับโครง Classification

```

select_features = 'age' @param ['age']
target_column = 'SeriousDlqinYrs' @param ['SeriousDlqinYrs']

select_features = [select_features]
# Setup Features & Target
features = train_df[select_features]
target = train_df[target_col]

# Setup Parameters
criterion = 'Entropy' @param ['Entropy', 'Gini']
max_depth = 1
dtree = tree.DecisionTreeClassifier(
    criterion=criterion.lower(),
    max_depth=max_depth,
)

# Training your tree
dtree.fit(features, target)

```

```

select_features: age
target_column: SeriousDlqinYrs
Criterion: Entropy
Gini

```

```

DecisionTreeClassifier(ccp_alpha=0.0, class_weight=None, criterion='entropy',
    max_depth=1, max_features=None, max_leaf_nodes=None,
    min_impurity_decrease=0.0, min_impurity_split=None,
    min_samples_leaf=1, min_samples_split=2,
    min_weight_fraction_leaf=0.0, presort='deprecated',
    random_state=None, splitter='best')

```

## skLearn.tree.DecisionTreeClassifier

```

class sklearn.tree.DecisionTreeClassifier(criterion='gini', splitter='best', max_depth=None, min_samples_split=2,
    min_samples_leaf=1, min_weight_fraction_leaf=0.0, max_features=None, random_state=None, max_leaf_nodes=None,
    min_impurity_decrease=0.0, min_impurity_split=None, class_weight=None, presort='deprecated', ccp_alpha=0.0)

```

A decision tree classifier.

Read more in the User Guide.

Parameters:

- criterion** : {'gini', 'entropy'}, default='gini'
 The function to measure the quality of a split. Supported criteria are "gini" for the Gini impurity and "entropy" for the information gain.
- splitter** : {'best', 'random'}, default='best'
 The strategy used to choose the split at each node. Supported strategies are "best" to choose the best split and "random" to choose the best random split.
- max\_depth** : int, default=None
 The maximum depth of the tree. If None, then nodes are expanded until all leaves are pure or until all leaf contain less than min\_samples\_split samples.
- min\_samples\_split** : int or float, default=2
 The minimum number of samples required to split an internal node.

## เลือกไม่รู้แปลว่าโง่

```

from sklearn.feature_selection import SelectFromModel
from sklearn.ensemble import AdaBoostRegressor
from sklearn.datasets import load_boston
from numpy import array

boston = load_boston()
X = boston.data
y = boston.target

print("Feature data dimension: ", X.shape)

selector = SelectFromModel(AdaBoostRegressor(
    random_state=0, n_estimators=50))
selector = selector.fit(X, y)

status = selector.get_support()
print("Selection status: ", status)

features = array(boston.feature_names)
print("All features")
print(features)

print("Selected features")
print(features[status])
selector.transform(X)

```

## Hyper param.

## Important Parameters in Decision Tree

- Splitting measure (criterion)** : gini / entropy
- Maximum depth** : ~5-10 (depend on number of feature)
- Maximum leaf nodes** : depend on number of class (target) and feature
- Minimum sample split** : 5 ~ 20% (depend on number of data)
- Minimum impurity decrease** : (default 0)
- Pruning adjustment (ccp\_alpha)** : 0.001 - 0.01 (depend on size of tree)

\* Stop when Leaf node  
 \* Min Max depth  
 \* Minimum leaf size

และจากค่า Tree มาแล้ว  $\Rightarrow$  `tree.feature_importances_`

ดูตามลำดับของตัวแปร

# Random Forest [Strong base line!] ⇒ ง่าย Classifier, Regressor

Importance Params: # Tree

# Columns (Variables)

# Rows (examples)

\* max\_features: sqrt, default = 'sqrt'

⇒ max\_features ดีกว่า

ถ้ามี 100 features ก็ควรใช้ประมาณ 50-80

อันนี้ default = 'sqrt'

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3.2.4.3.1. sklearn.ensemble.RandomForestClassifier

class sklearn.ensemble.RandomForestClassifier(estimators=100, criterion='gini', max\_depth=None, min\_samples\_split=2, min\_samples\_leaf=1, min\_weight\_fraction=0.0, max\_features='auto', max\_leaf\_nodes=None, min\_impurity\_decrease=0.0, min\_impurity\_split=None, bootstrap=True, oob\_score=False, n\_jobs=None, random\_state=None, verbose=0, warm\_start=False, class\_weight=None, ccp\_alpha=0.0, max\_samples=None) [source]

A random forest classifier.

A random forest is a meta estimator that fits a number of decision tree classifiers on various sub-samples of the dataset and uses averaging to improve the predictive accuracy and control over-fitting. The sub-sample size is always the same as the original input sample size but the samples are drawn with replacement if `bootstrap=True` (default).

Read more in the User Guide.

Parameters: **n\_estimators**: integer, optional (default=100)  
The number of trees in the forest.

Changed in version 0.22: The default value of `n_estimators` changed from 10 to 100 in 0.22.

**criterion**: string, optional (default='gini')  
The function to measure the quality of a split. Supported criteria are "gini" for the Gini impurity and "entropy" for the information gain. Note that this parameter is tree-specific.

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3.2.4.3.2. sklearn.ensemble.RandomForestRegressor

class sklearn.ensemble.RandomForestRegressor(estimators=100, criterion='mse', max\_depth=None, min\_samples\_split=2, min\_samples\_leaf=1, min\_weight\_fraction=0.0, max\_features='auto', max\_leaf\_nodes=None, min\_impurity\_decrease=0.0, min\_impurity\_split=None, bootstrap=True, oob\_score=False, n\_jobs=None, random\_state=None, verbose=0, warm\_start=False, ccp\_alpha=0.0, max\_samples=None) [source]

A random forest regressor.

A random forest is a meta estimator that fits a number of classifying decision trees on various sub-samples of the dataset and uses averaging to improve the predictive accuracy and control over-fitting. The sub-sample size is always the same as the original input sample size but the samples are drawn with replacement if `bootstrap=True` (default).

Read more in the User Guide.

Parameters: **n\_estimators**: integer, optional (default=10)  
The number of trees in the forest.

Changed in version 0.22: The default value of `n_estimators` changed from 10 to 100 in 0.22.

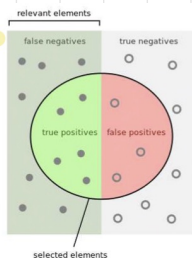
**criterion**: string, optional (default='mse')  
The function to measure the quality of a split. Supported criteria are "mse" for the mean squared error, which is equal to variance reduction as feature selection criterion, and "mae" for the mean absolute error.

New in version 0.18: Mean Absolute Error (MAE) criterion.

## Confusion Matrix

## Precision, Recall, F1 // Confusion Matrix

	Predicted: NO	Predicted: YES	
Actual: NO	TN = 50	FP = 10	60
Actual: YES	FN = 5	TP = 100	105
	55	110	



■ Precision = correctly predict =  $\frac{TP}{TP + FP}$

■ Recall = coverage =  $\frac{TP}{TP + FN}$

■ F1 =  $(2 * \text{pre} * \text{rec}) / (\text{pre} + \text{rec})$

Accuracy =  $\frac{TP + TN}{n}$



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Examples using sklearn.metrics.confusion\_matrix

sklearn.metrics.confusion\_matrix(y\_true, y\_pred, labels=None, sample\_weight=None, normalize=None) [source]

Compute confusion matrix to evaluate the accuracy of a classification.

By definition a confusion matrix  $C$  is such that  $C_{ij}$  is equal to the number of observations known to be in group  $i$  and predicted to be in group  $j$ .

Thus in binary classification, the count of true negatives is  $C_{0,0}$ , false negatives is  $C_{1,0}$ , true positives is  $C_{1,1}$  and false positives is  $C_{0,1}$ .

Read more in the User Guide.

Parameters: **y\_true**: array-like of shape  $(n_{\text{samples}})$   
Ground truth (correct) target values.

**y\_pred**: array-like of shape  $(n_{\text{samples}})$   
Predicted target values.

**labels**: array-like of shape  $(n_{\text{classes}})$ , optional  
The labels of the target classes. If None, the labels are inferred from the targets.

**sample\_weight**: array-like of shape  $(n_{\text{samples}})$ , optional  
Sample weights used in the calculation.

**normalize**: {'true', 'pred', 'all'}, optional  
Whether to normalize the matrix. If 'true', the matrix is divided by the number of observations in the target classes. If 'pred', the matrix is divided by the number of observations in the predicted classes. If 'all', the matrix is divided by the total number of observations.

[https://scikit-learn.org/stable/modules/generated/sklearn.metrics.confusion\\_matrix.html](https://scikit-learn.org/stable/modules/generated/sklearn.metrics.confusion_matrix.html)

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Examples using sklearn.metrics.classification\_report

sklearn.metrics.classification\_report(y\_true, y\_pred, labels=None, target\_names=None, sample\_weight=None, digits=2, output\_dict=False, zero\_division='warn') [source]

Build a text report showing the main classification metrics.

Read more in the User Guide.

Parameters: **y\_true**: array-like of shape  $(n_{\text{samples}})$   
Ground truth (correct) target values.

**y\_pred**: array-like of shape  $(n_{\text{samples}})$   
Predicted target values.

**labels**: array-like of shape  $(n_{\text{classes}})$ , optional  
The labels of the target classes. If None, the labels are inferred from the targets.

**target\_names**: array-like of shape  $(n_{\text{classes}})$ , optional  
The names of the target classes. If None, the names are inferred from the targets.

**sample\_weight**: array-like of shape  $(n_{\text{samples}})$ , optional  
Sample weights used in the calculation.

**digits**: int, optional  
The number of digits for the floating-point numbers.

**output\_dict**: bool, optional  
Whether to return the metrics as a dictionary.

**zero\_division**: {'warn', 'raise'}, optional  
Whether to raise an exception or warn when there is a zero division.

[https://scikit-learn.org/stable/modules/generated/sklearn.metrics.classification\\_report.html](https://scikit-learn.org/stable/modules/generated/sklearn.metrics.classification_report.html)

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Examples using sklearn.metrics.mean\_absolute\_error

sklearn.metrics.mean\_absolute\_error(y\_true, y\_pred, sample\_weight=None) [source]

Mean absolute error regression loss.

Parameters: **y\_true**: array-like of shape  $(n_{\text{samples}})$  or  $(n_{\text{samples}}, n_{\text{outputs}})$   
Ground truth (correct) target values.

**y\_pred**: array-like of shape  $(n_{\text{samples}})$  or  $(n_{\text{samples}}, n_{\text{outputs}})$   
Estimated target values.

**sample\_weight**: array-like of shape  $(n_{\text{samples}})$ , optional  
Sample weights used in the calculation.

**Returns**: loss: float  
A positive floating point value.

Examples

```
>>> from sklearn.metrics import mean_absolute_error
>>> y_true = [3, -0.5, 2, 7]
>>> y_pred = [2.5, 0.0, 2, 8]
>>> mean_absolute_error(y_true, y_pred)
0.5
>>> y_true = [[0.5, 1], [-1, 1], [7, -6]]
>>> y_pred = [[0, 2], [-1, 2], [8, -5]]
>>> mean_absolute_error(y_true, y_pred)
0.75
```



# Linear Regression → <sup>L1</sup> <sup>L2</sup> <sup>Feature Selection</sup> <sup>lasso</sup> <sup>Ridge</sup>

CHULA ENGINEERING [https://scikit-learn.org/stable/modules/generated/sklearn.linear\\_model.LinearRegression.html](https://scikit-learn.org/stable/modules/generated/sklearn.linear_model.LinearRegression.html)

learn

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sklearn.linear\_model.LinearRegression  
Examples using  
sklearn.linear\_model.LinearRegression

sklearn.linear\_model.LinearRegression

class sklearn.linear\_model.LinearRegression(fit\_intercept=True, normalize=False, copy\_X=True, n\_jobs=None)

Ordinary least squares Linear Regression.

LinearRegression fits a linear model with coefficients  $w = (w_1, \dots, w_p)$  to minimize the residual sum of squares between the observed targets in the dataset, and the targets predicted by the linear approximation.

Parameters:

**fit\_intercept** : bool, optional, default True  
Whether to calculate the intercept for this model. If set to False, no intercept will be used in calculations (i.e. data is expected to be centered).

**normalize** : bool, optional, default False  
This parameter is ignored when `fit_intercept` is set to False. If True, the regressors  $X$  will be normalized before regression by subtracting the mean and dividing by the l2-norm. If you wish to standardize, please use `StandardScaler` before calling `fit` on an estimator with `normalize=False`.  
Deprecated since version 1.0: normalize was deprecated in version 1.0 and will be removed in 1.2.

Attributes:

**coef\_** : array of shape (n\_features, ) or (n\_targets, n\_features)  
Estimated coefficients for the linear regression problem. If multiple targets are passed during the fit (y 2D), only one target is passed, this is a 1D array of

**intercept\_** : float or array of shape (n\_targets,)  
Independent term in the linear model. Set to 0.0 if `fit_intercept = False`.

**n\_features\_in\_** : int  
Number of features seen during fit.  
New in version 0.24.

**feature\_names\_in\_** : ndarray of shape (n\_features\_in\_,)  
Names of features seen during fit. Defined only when  $X$  has feature names that are all strings.  
New in version 1.0.

estimator = Lasso()

selector = SelectFromModel(estimator)

estimator = Ridge()

selector = SelectFromModel(estimator, threshold=?)

sklearn.linear\_model.Ridge <sup>L2</sup>

class sklearn.linear\_model.Ridge(alpha=1.0, \*, fit\_intercept=True, max\_iter=None, tol=0.001, solver='auto', positive=False, random\_state=None)

Linear least squares with l2 regularization.

Minimizes the objective function:  
$$\frac{1}{2} \|y - Xw\|_2^2 + \alpha \|w\|_2^2$$

This model solves a regression model where the loss function is the l2-norm. Also known as Ridge Regression or Tikhonov regularization (i.e., when  $y$  is a 2d-array of shape (n\_samples, n\_targets)).  
Read more in the User Guide.

Parameters:

**alpha** : float, ndarray of shape (n\_targets,)  
Constant that multiplies the l2 term, controls float i.e. in  $[0, \infty)$ .  
  
When `alpha = 0`, the objective is equivalent to ordinary least squares, solved by the `LinearRegression` object.  
  
If an array is passed, penalties are assumed to be independent.  
  
**fit\_intercept** : bool, default=True  
Whether to fit the intercept for this model. If `False`, the data is expected to be centered.

sklearn.linear\_model.Lasso <sup>L1</sup>

class sklearn.linear\_model.Lasso(alpha=1.0, \*, fit\_intercept=True, normalize='deprecated', precompute='auto', max\_iter=1000, tol=0.0001, warm\_start=False, positive=False, random\_state=None, selection='cyclic')

Linear regression with l1 regularization (aka the Lasso).

The optimization objective for Lasso is:  
$$\frac{1}{2} \|y - Xw\|_2^2 + \alpha \|w\|_1$$

Technically the Lasso model is optimizing the same objective function as the Elastic Net with `l1_ratio=1`.  
Read more in the User Guide.

Parameters:

**alpha** : float, default=1.0  
Constant that multiplies the l1 term, controlling regularization strength. `alpha` must be float i.e. in  $[0, \infty)$ .  
  
When `alpha = 0`, the objective is equivalent to ordinary least squares, solved by the `LinearRegression` object.  
  
**fit\_intercept** : bool, default=True  
Whether to calculate the intercept for this model. If set to `False`, no intercept will be used (i.e. data is expected to be centered).

sklearn.linear\_model.ElasticNet <sup>L1&L2</sup>

class sklearn.linear\_model.ElasticNet(alpha=1.0, \*, l1\_ratio=0.5, fit\_intercept=True, normalize='deprecated', precompute='auto', max\_iter=1000, copy\_X=True, tol=0.0001, warm\_start=False, positive=False, random\_state=None, selection='cyclic')

Linear regression with combined l1 and l2 priors as regularizer.

Minimizes the objective function:  
$$\frac{1}{2} \|y - Xw\|_2^2 + \alpha (\text{l1\_ratio} \|w\|_1 + (1 - \text{l1\_ratio}) \|w\|_2^2)$$

If you are interested in controlling the l1 and l2 penalty separately, keep in mind that this is equivalent to:  
$$a \|w\|_1 + 0.5 \times b \|w\|_2^2$$
  
where:  
$$\alpha = a + b \text{ and } \text{l1\_ratio} = a / (a + b)$$

The parameter `l1_ratio` corresponds to `alpha` in the glmnet R package while `alpha` corresponds to the `lambda` parameter in glmnet. Specifically, `l1_ratio = 1` is the lasso penalty. Currently, `l1_ratio < 0.01` is not reliable, unless you supply your own sequence of `alpha`.  
Read more in the User Guide.

Parameters:

**alpha** : float, default=1.0  
Constant that multiplies the penalty terms. Defaults to 1.0. See the notes for the exact mathematical meaning of this parameter. `alpha = 0` is equivalent to an ordinary least square, solved by the `LinearRegression` object. For numerical reasons, using `alpha = 0` with the `Lasso` object is not advised. Given this, you should use the `LinearRegression` object.  
  
**l1\_ratio** : float, default=0.5  
The ElasticNet mixing parameter, with `0 <= l1_ratio <= 1`. For `l1_ratio = 0` the penalty is an l2 penalty. For `l1_ratio = 1` it is an l1 penalty. For `0 < l1_ratio < 1`, the penalty is a combination of l1 and l2.  
  
**fit\_intercept** : bool, default=True  
Whether the intercept should be estimated or not. If `False`, the data is assumed to be already centered.

