

IIT Madras

ONLINE DEGREE

Machine Learning Practice
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Voting, Bagging and Random Forest

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Bagging and Boosting

Machine Learning Practice

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Namaste! Welcome to the next video of Machine Learning Practice Course. In this video we will discuss about bagging and boosting.

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Part 1: Voting, bagging and random forest

Part 2: Boosting and gradient boosting

Part 3: XGBoost

We will be recording this slide deck in three parts, in the first part, we will record voting bagging and random forest. In the second part, we will record boosting and gradient boosting. And in the third part, we will record the XGBoost functionality in sklearn.

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Voting estimators

Class: `sklearn.ensemble.VotingClassifier`

Class: `sklearn.ensemble.VotingRegressor`

Both these estimators take the following **common parameters**:

`base_estimator`

`weights`

Both these estimators implement the following **functions**:

`fit`

`predict`

`fit_transform`

`score`

`VotingClassifier` takes an **additional argument**:

`voting`

`hard`

`soft`

Let us first discuss about voting estimators. There are two voting estimators, voting classifier, and voting regressor. They are implemented as part of a `sklearn.ensemble` module. Both these estimators take the following common parameters, `base_estimator` and `weights`. And implement the following functions: `fit`, `predict`, `fit_transform` and `score`.

`Fit` is used to train the estimators. `Predict` is used to obtain the prediction for new samples. `Fit_transform` performs transformation as well as the fitment. And a `score` is used to obtain the performance of the estimator. Voting classifier takes an additional argument which is `voting`, and there are two types of voting either a hard voting or a soft voting.

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Bagging estimators

Class: `sklearn.ensemble.BaggingClassifier`

Class: `sklearn.ensemble.BaggingRegressor`

Then there are bagging estimators, two bagging estimators, again, bagging classifier and bagging regressor. They are also implemented as part of `sklearn.ensemble` module.

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Common parameters

base_estimator	None	DecisionTreeClassifier DecisionTreeRegressor
n_estimators	10	
max_samples	1.0	
max_features	1.0	
bootstrap	True	Whether samples are drawn with replacement
bootstrap_features	False	Whether features are drawn with replacement
oob_score	False	Whether to use out-of-bag samples to estimate generalization error

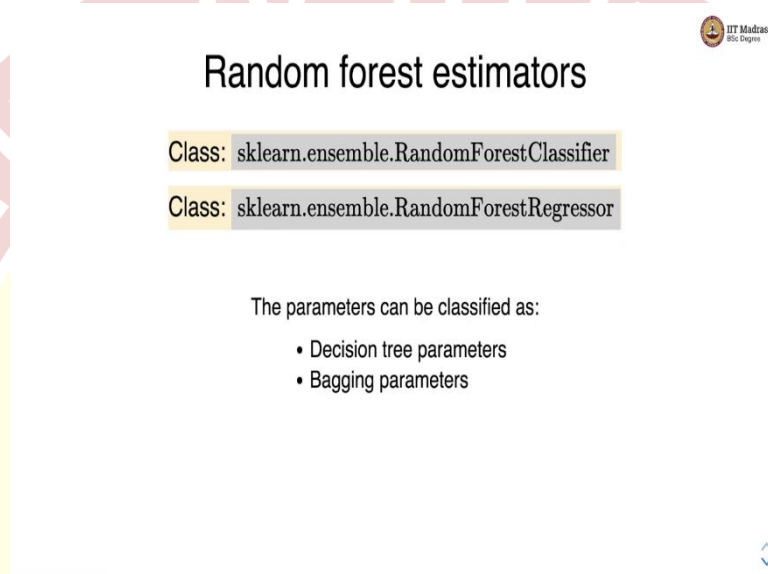
They take the following common parameters `base_estimator`, by default it is none. In case of classification, we use `DecisionTreeClassifier` as the default `base_estimator`, whereas in case of bagging regressor we use `DecisionTreeRegressor` as a default `base_estimator`. Default number of estimators is 10. `Max_samples` can be integer or float, whenever it is integer, we take those many samples.

And when is float, we basically multiply the `max_samples` with the total number of samples and take the resulting samples as the `max_samples` for sampling. `Max_features` is also either

an integer or float, and if it is integer, we select those many features. And if it is float, then we select number of features = the total number of features into max _features.

Bootstrap is true by default, and it specifies whether samples are drawn with replacement. Bootstrap _features and oob _score, both of them are false by default. And Bootstrap _features denotes or indicates whether features are drawn with the placement and oob _score denotes whether to use out of bag samples to estimate generalization error or test error.

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Random forest estimators

Class: `sklearn.ensemble.RandomForestClassifier`

Class: `sklearn.ensemble.RandomForestRegressor`

The parameters can be classified as:

- Decision tree parameters
- Bagging parameters

Then we have random forests estimators, again two estimators one for classification, and second for regression, RandomForestClassifier and RandomForestRegressor respectively. They are also implemented as part of sklearn.ensemble module. The parameters of random forests can be classified as decision tree parameters and bagging parameters.

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Bagging parameters

- The number of trees are specified by `n_estimators`.
 - Default #trees for classification = 10
 - Default #trees for regression = 100
- `bootstrap` specifies whether to use bootstrap samples for training.
 - `True` : bootstrapped samples are used.
 - `False` : whole dataset is used.
- `oob_score` specifies whether to use out-of-bag samples for estimating generalization error. It is only available when `bootstrap = True`.

Let us first look at bagging parameters. The number of trees are specified by `n_estimators`. Default number of trees used for classification instead, whereas default number of trees used for regression are 100. Then we have `bootstrap` flag that specifies whether to use Bootstrap samples for training when it is true bootstrap samples are used. And when Bootstrap is false whole dataset is used for training the estimators. `oob_score` specifies whether to use out of bag samples for estimating generalization error. It is only available when Bootstrap is = true.

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- `max_samples` specifies the number of samples to be drawn while bootstrapping.
 - `None` : Use all samples in the training data.
 - `int` : Use `max_samples` samples from the training data.
 - `float` : Use `max_samples` * total number of samples from the training data. The value should be between 0 and 1.
- `random_state` controls randomness of features and samples selected during bootstrap.

Then `max_samples` specifies the number of samples to be drawn while bootstrapping. If it is specified as none, we use all samples in the training data. If it is specified as integer, then we use `max_samples` from training data. If it is float, we use `max_samples` into total number of

samples from the training data. If it is float, the value should always be between 0 and 1. And `random_state` controls randomness of features and sample selected during bootstrap.

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- The number of features to be considered while splitting is specified by `max_features`.
 - `auto`, `sqrt`, `log2`, `int`, `float`

Value	max_features
int	value specified
float	value * # features
auto	$\sqrt{\text{\#features}}$
sqrt	$\sqrt{\text{\#features}}$
log2	$\log_2(\text{\#features})$
None	\#features

The number of features to be considered while splitting the decision tree is specified by `max_features`. There could be five different values either `auto`, square root, `log2`, integer and float. Whenever we specify an integer value, we use the `max_features` as the value specified in the parameter.

If it is float, then the total number of features used are value into total number of features. If it is `auto`, then we use square root of number of features. If it is square root, `sqrt`, we use square root of number of features while making the splitting decision. If it is `log2`, we take number of features, log of number of features to the base 2. And if it is none, then we take all the features for deciding on the split in the decision tree.

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Decision tree parameters

- The criteria for splitting the node is specified through `criterion`.
 - Default for classification: `gini`
 - Default for regression: `squared_error`
- The `depth of the tree` is controlled by `max_depth`. The default value is `None`, which means the tree will be grown until all leaf nodes are pure or until leaves contain less than `min_samples_split` samples.
- We will continue to split the internal node until they contain `min_samples_split` samples.
 - Whenever it is specified as an integer, then it is considered as a number.
 - Whenever it is specified as a float, and the `min_samples_split` is calculated as $\text{min_samples_split} \times n$.

Let us look at some of the decision tree parameters. We are the criteria for splitting the node and it is specified to criterion default for classification is gini and default for regression is squared_error. The depth of the tree is controlled by a max_depth parameter, the default value is none which means the tree will be grown until all leaf nodes are pure or until leaves content less than min_samples_split samples.

We will continue to split internal load until they contain min_samples_split samples. Whenever it is specified as an integer then it is considered as a number and whenever it is specified as a float then the min_samples_split is calculated as the min_samples_split into total number of samples.

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- The tree growth can also be controlled by `min_impurity_decrease` parameter.
 - A node will be split if it reduces impurity at least by the value specified in this parameter.
- The complexity of tree can also be controlled by `ccp_alpha` parameter through minimal cost complexity pruning procedure.

The tree growth can be controlled by `min_impurity_decrease` parameter. A node will be split if it reduces impurity, at least by the value specified in this parameter. The complexity of the tree can also be controlled by `ccp_alpha` parameter through minimal cost complexity pruning procedure. So, `ccp_alpha` is a parameter used in minimal cost complexity pruning.

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Trained random forest estimators

- `estimators_` member variable contains a collection of fitted estimators.
- `feature_importances_` member variable contains a list of important features.

Estimator `_member` variable contains a collection of fitted estimators. Feature `_importances_` member variable contains a list of important features.

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Training and inference for random forest

- `fit` builds forest of trees from the training dataset with the specified parameters.
- `decision_path` returns decision path in the forest.
- `predict` returns class label in classification and output value in regression.
- `predict_proba` and `predict_log_proba` returns probabilities and their logs for classification set up.

We use `fit` method for building forest of trees from training data with the specified parameters. `Decision_path` returns decision path in the forest. `Predict` returns class label in classification, and outputs a value in regression. And `predict_proba` and `predict_log_proba` are functions available in classification and the return probabilities and the log of probabilities.

In this video, we studied a sklearn functionality for Voting, Bagging and Random Forest. We looked at how to set up classification and regression with Voting, Bagging and Random Forest.