skfda.ml.classification.NearestNeighbors

class skfda.ml.classification.NearestNeighbors(n_neighbors=5, radius=1.0, algorithm='auto', leaf_size=30, metric=<function lp_distance>, metric_params=None, n_jobs=1, sklearn_metric=False) [source]

Unsupervised learner for implementing neighbor searches.

Parameters:

- n_neighbors (int, optional (default = 5)) Number of neighbors to use by default for kneighbors() queries.
- radius (float, optional (default = 1.0)) Range of parameter space to use by default for radius_neighbors() queries.
- algorithm ({'auto', 'ball_tree', 'brute'}, optional) –
 Algorithm used to compute the nearest neighbors:
 - 'ball_tree' will use sklearn.neighbors.BallTree .
 - o 'brute' will use a brute-force search.
 - 'auto' will attempt to decide the most appropriate algorithm based on the values passed to fit() method.
- leaf_size (int, optional (default = 30)) Leaf size passed to BallTree or KDTree. This can affect the speed of the construction and query, as well as the memory required to store the tree. The optimal value depends on the nature of the problem.
- metric (string or callable, (default) lp_distance) the distance metric to use for the tree. The default metric is the Lp distance. See the documentation of the metrics module for a list of available metrics.
- **metric_params** (*dict*, *optional* (*default* = *None*)) Additional keyword arguments for the metric function.
- n_jobs (int or None, optional (default=None)) The number of parallel jobs to run for neighbors search. None means 1 unless in a joblib.parallel_backend context. -1 means using all processors.
 Doesn't affect fit() method.
- sklearn_metric (boolean, optional (default = False)) Indicates if the metric used is a sklearn distance between vectors (see
 sklearn.neighbors.DistanceMetric) or a functional metric of the module
 skfda.misc.metrics .

Examples

We will fit a Nearest Neighbors estimator

```
>>> from skfda.ml.classification import NearestNeighbors
>>> neigh = NearestNeighbors(radius=.3)
>>> neigh.fit(fd)
NearestNeighbors(algorithm='auto', leaf_size=30,...)
```

Now we can query the k-nearest neighbors.

```
>>> distances, index = neigh.kneighbors(fd[:2])
>>> index # Index of k-neighbors of samples 0 and 1
array([[ 0,  7,  6, 11,  2],...)
```

```
>>> distances.round(2) # Distances to k-neighbors array([[ 0. , 0.28, 0.29, 0.29, 0.3 ], [ 0. , 0.27, 0.28, 0.29, 0.3 ]])
```

We can query the neighbors in a given radius too.

```
>>> distances, index = neigh.radius_neighbors(fd[:2])
>>> index[0]
array([ 0,  2,  6,  7, 11]...)
```

```
>>> distances[0].round(2) # Distances to neighbors of the sample 0 array([ 0. , 0.3 , 0.29, 0.28, 0.29])
```

See also

```
KNeighborsClassifier, RadiusNeighborsClassifier, KNeighborsScalarRegressor, RadiusNeighborsScalarRegressor, NearestCentroids
```

See Nearest Neighbors in the sklearn online documentation for a discussion of the choice of algorithm and leaf_size.

This class wraps the sklearn classifier sklearn.neighbors.KNeighborsClassifier.

https://en.wikipedia.org/wiki/K-nearest_neighbor_algorithm

Initialize the nearest neighbors searcher.

Methods

init ([n_neighbors, radius, algorithm,])	Initialize the nearest neighbors searcher.
fit (X[, y])	Fit the model using X as training data.
get_params ([deep])	Get parameters for this estimator.
kneighbors ([X, n_neighbors, return_distance])	Finds the K-neighbors of a point.
kneighbors_graph ([X, n_neighbors, mode])	Computes the (weighted) graph of k-Neighbor
radius_neighbors ([X, radius, return_distance])	Finds the neighbors within a given radius of a
<pre>radius_neighbors_graph ([X, radius, mode])</pre>	Computes the (weighted) graph of Neighbors
set_params (**params)	Set the parameters of this estimator.