API Reference

This is the class and function reference of scikit-learn. Please refer to the <u>full user guide</u> for further details, as the class and function raw specifications may not be enough to give full guidelines on their uses. For reference on concepts repeated across the API, see <u>Glossary of Common Terms and API Elements</u>.

sklearn.base: Base classes and utility functions

Base classes for all estimators.

Base classes

<u>base.BaseEstimator</u>	Base class for all estimators in scikit-learn.
base.BiclusterMixin	Mixin class for all bicluster estimators in scikit-learn.
base.ClassifierMixin	Mixin class for all classifiers in scikit-learn.
base.ClusterMixin	Mixin class for all cluster estimators in scikit-learn.
base.DensityMixin	Mixin class for all density estimators in scikit-learn.
base.RegressorMixin	Mixin class for all regression estimators in scikit-learn.
base.TransformerMixin	Mixin class for all transformers in scikit-learn.
base.MetaEstimatorMixin	
base.OneToOneFeatureMixin	Provides get_feature_names_out for simple transformers.
base.OutlierMixin	Mixin class for all outlier detection estimators in scikit-learn.
$\underline{\texttt{base.ClassNamePrefixFeaturesOutMixin}}$	Mixin class for transformers that generate their own names by prefixing.
<u>feature</u> <u>selection</u> . <u>SelectorMixin</u>	Transformer mixin that performs feature selection given a support mask
4	

Functions

<pre>base.clone(estimator, *[, safe])</pre>	Construct a new unfitted estimator with the same parameters.
<pre>base.is classifier(estimator)</pre>	Return True if the given estimator is (probably) a classifier.
<pre>base.is regressor(estimator)</pre>	Return True if the given estimator is (probably) a regressor.
<pre>config_context(*[, assume_finite,])</pre>	Context manager for global scikit-learn configuration.
<pre>get config()</pre>	Retrieve current values for configuration set by <u>set_config</u> .
<pre>set config([assume_finite, working_memory,])</pre>	Set global scikit-learn configuration.
<pre>show versions()</pre>	Print useful debugging information"
4	

sklearn.calibration: Probability Calibration

Calibration of predicted probabilities.

User guide: See the <u>Probability calibration</u> section for further details.

<u>calibration.CalibratedClassifierCV([...])</u> Probability calibration with isotonic regression or logistic regression.

<u>calibration.calibration curve(y_true, y_prob, *)</u> Compute true and predicted probabilities for a calibration curve.

sklearn.cluster: Clustering

The **sklearn.cluster** module gathers popular unsupervised clustering algorithms.

User guide: See the <u>Clustering</u> and <u>Biclustering</u> sections for further details.

Classes

		Perform Affinity Propagation Clustering of data.
Toggle Menu	omerativeClustering([])	Agglomerative Clustering.

<pre>cluster.Birch(*[, threshold,])</pre>	Implements the BIRCH clustering algorithm.
<pre>cluster.DBSCAN([eps, min_samples, metric,])</pre>	Perform DBSCAN clustering from vector array or distance matrix.
<pre>cluster.HDBSCAN([min_cluster_size,])</pre>	Cluster data using hierarchical density-based clustering.
<pre>cluster.FeatureAgglomeration([n_clusters,])</pre>	Agglomerate features.
<pre>cluster.KMeans([n_clusters, init, n_init,])</pre>	K-Means clustering.
<pre>cluster.BisectingKMeans([n_clusters, init,])</pre>	Bisecting K-Means clustering.
<pre>cluster.MiniBatchKMeans([n_clusters, init,])</pre>	Mini-Batch K-Means clustering.
<pre>cluster.MeanShift(*[, bandwidth, seeds,])</pre>	Mean shift clustering using a flat kernel.
<pre>cluster.OPTICS(*[, min_samples, max_eps,])</pre>	Estimate clustering structure from vector array.
<pre>cluster.SpectralClustering([n_clusters,])</pre>	Apply clustering to a projection of the normalized Laplacian.
${\color{red} \underline{\textbf{cluster}.\textbf{SpectralBiclustering}}([\textbf{n}_\textbf{clusters}, \ldots])}$	Spectral biclustering (Kluger, 2003).
cluster.SpectralCoclustering([n_clusters])	Spectral Co-Clustering algorithm (Dhillon, 2001)

Functions

Toggle Menu

<pre>cluster.affinity propagation(S, *[,])</pre>	Perform Affinity Propagation Clustering of data.
<pre>cluster.cluster optics dbscan(*,)</pre>	Perform DBSCAN extraction for an arbitrary epsilon.
<pre>cluster.cluster optics xi(*, reachability,)</pre>	Automatically extract clusters according to the Xi-steep method.
<pre>cluster.compute optics graph(X, *,)</pre>	Compute the OPTICS reachability graph.
<pre>cluster.dbscan(X[, eps, min_samples,])</pre>	Perform DBSCAN clustering from vector array or distance matrix.
<pre>cluster.estimate bandwidth(X, *[, quantile,])</pre>	Estimate the bandwidth to use with the mean-shift algorithm.
<pre>cluster.k means(X, n_clusters, *[,])</pre>	Perform K-means clustering algorithm.
<pre>cluster.kmeans_plusplus(X, n_clusters, *[,])</pre>	Init n_clusters seeds according to k-means++.
$\underline{\textbf{cluster.mean shift}}(X, *[, bandwidth, seeds,])$	Perform mean shift clustering of data using a flat kernel.
<pre>cluster.spectral clustering(affinity, *[,])</pre>	Apply clustering to a projection of the normalized Laplacian.
<pre>cluster.ward tree(X, *[, connectivity,])</pre>	Ward clustering based on a Feature matrix.
4	

sklearn.compose: Composite Estimators

Meta-estimators for building composite models with transformers

In addition to its current contents, this module will eventually be home to refurbished versions of Pipeline and FeatureUnion.

User guide: See the <u>Pipelines and composite estimators</u> section for further details.

```
<u>compose.ColumnTransformer</u>(transformers, *[, ...]) Applies transformers to columns of an array or pandas DataFrame.
compose.TransformedTargetRegressor([...])
                                                   Meta-estimator to regress on a transformed target.
<u>compose.make column transformer</u>(*transformers) Construct a ColumnTransformer from the given transformers.
                                                    Create a callable to select columns to be used with ColumnTransformer.
compose.make column selector([pattern, ...])
```

sklearn.covariance: Covariance Estimators

The sklearn.covariance module includes methods and algorithms to robustly estimate the covariance of features given a set of points. The precision matrix defined as the inverse of the covariance is also estimated. Covariance estimation is closely related to the theory of Gaussian Graphical Models.

User guide: See the <u>Covariance estimation</u> section for further details.

as(X, *[, assume_centered])

```
covariance.EmpiricalCovariance(*[, ...])
                                                 Maximum likelihood covariance estimator.
covariance.EllipticEnvelope(*[, ...])
                                                 An object for detecting outliers in a Gaussian distributed dataset.
covariance.GraphicalLasso([alpha, mode, ...])
                                                 Sparse inverse covariance estimation with an l1-penalized estimator.
covariance.GraphicalLassoCV(*[, alphas, ...])
                                                 Sparse inverse covariance w/ cross-validated choice of the l1 penalty.
covariance.LedoitWolf(*[, store_precision, ...])
                                                LedoitWolf Estimator.
covariance.MinCovDet(*[, store precision, ...])
                                                 Minimum Covariance Determinant (MCD): robust estimator of covariance.
covariance.OAS(*[, store_precision, ...])
                                                 Oracle Approximating Shrinkage Estimator as proposed in [R69773891e6a6-1].
covariance.ShrunkCovariance(*[, ...])
                                                 Covariance estimator with shrinkage.
covariance.empirical covariance(X, *[, ...])
                                                   Compute the Maximum likelihood covariance estimator.
covariance.graphical lasso(emp_cov, alpha, *)
                                                  L1-penalized covariance estimator.
covariance.ledoit_wolf(X, *[, ...])
                                                   Estimate the shrunk Ledoit-Wolf covariance matrix.
             <u>ledoit wolf shrinkage(</u>X[, ...])
```

Estimate the shrunk Ledoit-Wolf covariance matrix.

Estimate covariance with the Oracle Approximating Shrinkage as proposed in [Rca3a42e5ec35-1]

<u>covariance.shrunk covariance</u>(emp_cov[, ...]) Calculate a covariance matrix shrunk on the diagonal.

sklearn.cross_decomposition: Cross decomposition

User guide: See the <u>Cross decomposition</u> section for further details.

<pre>cross decomposition.CCA([n_components,])</pre>	Canonical Correlation Analysis, also known as "Mode B" PLS.
<pre>cross decomposition.PLSCanonical([])</pre>	Partial Least Squares transformer and regressor.
$\underline{\texttt{cross}} \ \underline{\texttt{decomposition}}. \underline{\texttt{PLSRegression}}([])$	PLS regression.
<pre>cross decomposition.PLSSVD([n_components,])</pre>	Partial Least Square SVD.

sklearn.datasets: Datasets

The <u>sklearn.datasets</u> module includes utilities to load datasets, including methods to load and fetch popular reference datasets. It also features some artificial data generators.

User guide: See the <u>Dataset loading utilities</u> section for further details.

Loaders

<pre>datasets.clear data home([data_home])</pre>	Delete all the content of the data home cache.
<pre>datasets.dump svmlight file(X, y, f, *[,])</pre>	Dump the dataset in symlight / libsym file format.
<pre>datasets.fetch 20newsgroups(*[, data_home,])</pre>	Load the filenames and data from the 20 newsgroups dataset (classification).
<pre>datasets.fetch 20newsgroups vectorized(*[,])</pre>	Load and vectorize the 20 newsgroups dataset (classification).
<pre>datasets.fetch california housing(*[,])</pre>	Load the California housing dataset (regression).
<pre>datasets.fetch covtype(*[, data_home,])</pre>	Load the covertype dataset (classification).
<pre>datasets.fetch kddcup99(*[, subset,])</pre>	Load the kddcup99 dataset (classification).
<pre>datasets.fetch lfw pairs(*[, subset,])</pre>	Load the Labeled Faces in the Wild (LFW) pairs dataset (classification).
<pre>datasets.fetch lfw people(*[, data_home,])</pre>	Load the Labeled Faces in the Wild (LFW) people dataset (classification).
<pre>datasets.fetch olivetti faces(*[,])</pre>	Load the Olivetti faces data-set from AT&T (classification).
<pre>datasets.fetch openml([name, version,])</pre>	Fetch dataset from openml by name or dataset id.
<pre>datasets.fetch rcv1(*[, data_home, subset,])</pre>	Load the RCV1 multilabel dataset (classification).
<pre>datasets.fetch species distributions(*[,])</pre>	Loader for species distribution dataset from Phillips et.
<pre>datasets.get_data_home([data_home])</pre>	Return the path of the scikit-learn data directory.
<pre>datasets.load breast_cancer(*[, return_X_y,])</pre>	Load and return the breast cancer wisconsin dataset (classification).
<pre>datasets.load diabetes(*[, return_X_y,])</pre>	Load and return the diabetes dataset (regression).
<pre>datasets.load digits(*[, n_class,])</pre>	Load and return the digits dataset (classification).
<pre>datasets.load files(container_path, *[,])</pre>	Load text files with categories as subfolder names.
<pre>datasets.load iris(*[, return_X_y, as_frame])</pre>	Load and return the iris dataset (classification).
<pre>datasets.load linnerud(*[, return_X_y, as_frame])</pre>	Load and return the physical exercise Linnerud dataset.
<pre>datasets.load sample image(image_name)</pre>	Load the numpy array of a single sample image.
datasets.load sample images()	Load sample images for image manipulation.
<pre>datasets.load svmlight file(f, *[,])</pre>	Load datasets in the symlight / libsym format into sparse CSR matrix.
<pre>datasets.load svmlight files(files, *[,])</pre>	Load dataset from multiple files in SVMlight format.
<pre>datasets.load wine(*[, return_X_y, as_frame])</pre>	Load and return the wine dataset (classification).
←	

Samples generator

<pre>datasets.make_biclusters(shape, n_clusters, *)</pre>	Generate a constant block diagonal structure array for biclustering.
<pre>datasets.make blobs([n_samples, n_features,])</pre>	Generate isotropic Gaussian blobs for clustering.
<pre>datasets.make_checkerboard(shape, n_clusters, *)</pre>	Generate an array with block checkerboard structure for biclustering.
<pre>datasets.make circles([n_samples, shuffle,])</pre>	Make a large circle containing a smaller circle in 2d.
<pre>datasets.make_classification([n_samples,])</pre>	Generate a random n-class classification problem.
<pre>datasets.make friedman1([n_samples,])</pre>	Generate the "Friedman #1" regression problem.
<pre>datasets.make friedman2([n_samples, noise,])</pre>	Generate the "Friedman #2" regression problem.
<pre>datasets.make friedman3([n_samples, noise,])</pre>	Generate the "Friedman #3" regression problem.
<pre>datasets.make gaussian quantiles(*[, mean,])</pre>	Generate isotropic Gaussian and label samples by quantile.
<pre>datasets.make hastie 10 2([n_samples,])</pre>	Generate data for binary classification used in Hastie et al. 2009, Example 10.2.
<pre>datasets.make low rank matrix([n_samples,])</pre>	Generate a mostly low rank matrix with bell-shaped singular values.
Toggle Menu te moons([n_samples, shuffle,])	Make two interleaving half circles.

<pre>datasets.make multilabel classification([])</pre>	Generate a random multilabel classification problem.
<pre>datasets.make regression([n_samples,])</pre>	Generate a random regression problem.
<pre>datasets.make s curve([n_samples, noise,])</pre>	Generate an S curve dataset.
<pre>datasets.make sparse coded signal(n_samples,)</pre>	Generate a signal as a sparse combination of dictionary elements.
<pre>datasets.make sparse spd matrix([dim,])</pre>	Generate a sparse symmetric definite positive matrix.
<pre>datasets.make sparse uncorrelated([])</pre>	Generate a random regression problem with sparse uncorrelated design.
<pre>datasets.make spd matrix(n_dim, *[,])</pre>	Generate a random symmetric, positive-definite matrix.
datasets.make swiss roll([n samples, noise,])	Generate a swiss roll dataset.

sklearn.decomposition: Matrix Decomposition

The <u>sklearn.decomposition</u> module includes matrix decomposition algorithms, including among others PCA, NMF or ICA. Most of the algorithms of this module can be regarded as dimensionality reduction techniques.

User guide: See the <u>Decomposing signals in components (matrix factorization problems)</u> section for further details.

$\underline{\text{decomposition.} \textbf{DictionaryLearning}([])}$	Dictionary learning.
<pre>decomposition.FactorAnalysis([n_components,])</pre>	Factor Analysis (FA).
<pre>decomposition.FastICA([n_components,])</pre>	FastICA: a fast algorithm for Independent Component Analysis.
$\underline{\text{decomposition.IncrementalPCA}}([n_components, \ldots])$	Incremental principal components analysis (IPCA).
<pre>decomposition.KernelPCA([n_components,])</pre>	Kernel Principal component analysis (KPCA) [R396fc7d924b8-1]
$\underline{\texttt{decomposition.LatentDirichletAllocation}([])}$	Latent Dirichlet Allocation with online variational Bayes algorithm
$\underline{\texttt{decomposition.MiniBatchDictionaryLearning}([])}$	Mini-batch dictionary learning.
decomposition.MiniBatchSparsePCA([])	Mini-batch Sparse Principal Components Analysis.
<pre>decomposition.NMF([n_components, init,])</pre>	Non-Negative Matrix Factorization (NMF).
<pre>decomposition.MiniBatchNMF([n_components,])</pre>	Mini-Batch Non-Negative Matrix Factorization (NMF).
<pre>decomposition.PCA([n_components, copy,])</pre>	Principal component analysis (PCA).
<pre>decomposition.SparsePCA([n_components,])</pre>	Sparse Principal Components Analysis (SparsePCA).
<pre>decomposition.SparseCoder(dictionary, *[,])</pre>	Sparse coding.
<pre>decomposition.TruncatedSVD([n_components,])</pre>	Dimensionality reduction using truncated SVD (aka LSA).
4	
<pre>decomposition.dict learning(X, n_components,)</pre>	Solve a dictionary learning matrix factorization problem.
<pre>decomposition.dict learning online(X[,])</pre>	Solve a dictionary learning matrix factorization problem online.
<pre>decomposition.fastica(X[, n_components,])</pre>	Perform Fast Independent Component Analysis.
decomposition.non negative factorization(X)	Compute Non-negative Matrix Factorization (NMF).
<pre>decomposition.sparse encode(X, dictionary, *)</pre>	Sparse coding.
4	

sklearn.discriminant analysis: Discriminant Analysis

Linear Discriminant Analysis and Quadratic Discriminant Analysis

User guide: See the Linear and Quadratic Discriminant Analysis section for further details.

```
discriminant analysis.LinearDiscriminantAnalysis([...]) Linear Discriminant Analysis.

discriminant analysis.QuadraticDiscriminantAnalysis(*) Quadratic Discriminant Analysis.
```

sklearn.dummy: Dummy estimators

User guide: See the <u>Metrics and scoring: quantifying the quality of predictions</u> section for further details.

```
dummy.DummyClassifier(*[, strategy, ...]) DummyClassifier makes predictions that ignore the input features.
dummy.DummyRegressor(*[, strategy, ...]) Regressor that makes predictions using simple rules.
```

sklearn.ensemble: Ensemble Methods

The sklearn ensemble module includes ensemble-based methods for classification, regression and anomaly detection.

Toggle Menu

User guide: See the Ensembles: Gradient boosting, random forests, bagging, voting, stacking section for further details.

<pre>ensemble.AdaBoostClassifier([estimator,])</pre>	An AdaBoost classifier.
<pre>ensemble.AdaBoostRegressor([estimator,])</pre>	An AdaBoost regressor.
<pre>ensemble.BaggingClassifier([estimator,])</pre>	A Bagging classifier.
<pre>ensemble.BaggingRegressor([estimator,])</pre>	A Bagging regressor.
<pre>ensemble.ExtraTreesClassifier([])</pre>	An extra-trees classifier.
<pre>ensemble.ExtraTreesRegressor([n_estimators,])</pre>	An extra-trees regressor.
<pre>ensemble.GradientBoostingClassifier(*[,])</pre>	Gradient Boosting for classification.
$\underline{\textbf{ensemble}.\textbf{GradientBoostingRegressor}(*[,])}$	Gradient Boosting for regression.
<pre>ensemble.IsolationForest(*[, n_estimators,])</pre>	Isolation Forest Algorithm.
<pre>ensemble.RandomForestClassifier([])</pre>	A random forest classifier.
$\underline{ensemble.RandomForestRegressor}([])$	A random forest regressor.
<pre>ensemble.RandomTreesEmbedding([])</pre>	An ensemble of totally random trees.
<pre>ensemble.StackingClassifier(estimators[,])</pre>	Stack of estimators with a final classifier.
<pre>ensemble.StackingRegressor(estimators[,])</pre>	Stack of estimators with a final regressor.
<pre>ensemble.VotingClassifier(estimators, *[,])</pre>	Soft Voting/Majority Rule classifier for unfitted estimators.
<pre>ensemble.VotingRegressor(estimators, *[,])</pre>	Prediction voting regressor for unfitted estimators.
$\underline{ensemble. HistGradientBoostingRegressor([])}$	Histogram-based Gradient Boosting Regression Tree.
$\underline{ensemble.HistGradientBoostingClassifier}([])$	Histogram-based Gradient Boosting Classification Tree.

sklearn.exceptions: Exceptions and warnings

The **sklearn.exceptions** module includes all custom warnings and error classes used across scikit-learn.

<u>exceptions.ConvergenceWarning</u>	Custom warning to capture convergence problems
exceptions.DataConversionWarning	Warning used to notify implicit data conversions happening in the code.
exceptions.DataDimensionalityWarning	Custom warning to notify potential issues with data dimensionality.
exceptions.EfficiencyWarning	Warning used to notify the user of inefficient computation.
exceptions.FitFailedWarning	Warning class used if there is an error while fitting the estimator.
<pre>exceptions.InconsistentVersionWarning(*,)</pre>	Warning raised when an estimator is unpickled with a inconsistent version.
<u>exceptions.NotFittedError</u>	Exception class to raise if estimator is used before fitting.
exceptions.UndefinedMetricWarning	Warning used when the metric is invalid
4	

sklearn.experimental: Experimental

The **sklearn.experimental** module provides importable modules that enable the use of experimental features or estimators.

The features and estimators that are experimental aren't subject to deprecation cycles. Use them at your own risks!

```
<u>experimental.enable iterative imputer</u> Enables IterativeImputer

<u>experimental.enable halving search cv</u> Enables Successive Halving search-estimators
```

sklearn.feature extraction: Feature Extraction

The <u>sklearn.feature extraction</u> module deals with feature extraction from raw data. It currently includes methods to extract features from text and images.

User guide: See the <u>Feature extraction</u> section for further details.

```
<u>feature extraction.DictVectorizer</u>(*[, ...]) Transforms lists of feature-value mappings to vectors. 
<u>feature extraction.FeatureHasher</u>([...]) Implements feature hashing, aka the hashing trick.
```

From images

The **sklearn.feature** extraction.image submodule gathers utilities to extract features from images.

feature exti	raction.image.extract_patches_2d()	Reshape a 2D image into a collection of patches.
Toggle Menu	action.image.grid to graph(n_x, n_y)	Graph of the pixel-to-pixel connections.

<pre>feature extraction.image.img to graph(img, *)</pre>	Graph of the pixel-to-pixel gradient connections.
feature extraction.image.reconstruct from patches 2d()	Reconstruct the image from all of its patches.
<pre>feature extraction.image.PatchExtractor(*[,])</pre>	Extracts patches from a collection of images.

From text

The sklearn.feature extraction.text submodule gathers utilities to build feature vectors from text documents.

```
feature extraction.text.CountVectorizer(*[, ...]) Convert a collection of text documents to a matrix of token counts.

feature extraction.text.HashingVectorizer(*) Convert a collection of text documents to a matrix of token occurrences.

feature extraction.text.TfidfTransformer(*) Transform a count matrix to a normalized tf or tf-idf representation.

feature extraction.text.TfidfVectorizer(*[, ...]) Convert a collection of raw documents to a matrix of TF-IDF features.
```

sklearn.feature selection: Feature Selection

The <u>sklearn.feature selection</u> module implements feature selection algorithms. It currently includes univariate filter selection methods and the recursive feature elimination algorithm.

User guide: See the <u>Feature selection</u> section for further details.

$\underline{ \textbf{feature selection}. \textbf{GenericUnivariateSelect}([])}$	Univariate feature selector with configurable strategy.
<pre>feature selection.SelectPercentile([])</pre>	Select features according to a percentile of the highest scores.
<pre>feature selection.SelectKBest([score_func, k])</pre>	Select features according to the k highest scores.
<pre>feature selection.SelectFpr([score_func, alpha])</pre>	Filter: Select the pvalues below alpha based on a FPR test.
<pre>feature selection.SelectFdr([score_func, alpha])</pre>	Filter: Select the p-values for an estimated false discovery rate.
<pre>feature selection.SelectFromModel(estimator, *)</pre>	Meta-transformer for selecting features based on importance weights.
<pre>feature selection.SelectFwe([score_func, alpha])</pre>	Filter: Select the p-values corresponding to Family-wise error rate.
<u>feature selection.SequentialFeatureSelector()</u>	Transformer that performs Sequential Feature Selection.
<pre>feature selection.RFE(estimator, *[,])</pre>	Feature ranking with recursive feature elimination.
feature selection.RFECV(estimator, *[,])	Recursive feature elimination with cross-validation to select features.
<u>feature selection.VarianceThreshold([threshold])</u>	Feature selector that removes all low-variance features.
4	
<pre>feature selection.chi2(X, y)</pre>	Compute chi-squared stats between each non-negative feature and cla

<pre>feature selection.chi2(X, y)</pre>	Compute chi-squared stats between each non-negative feature and class.
<pre>feature selection.f classif(X, y)</pre>	Compute the ANOVA F-value for the provided sample.
<pre>feature selection.f regression(X, y, *[,])</pre>	Univariate linear regression tests returning F-statistic and p-values.
<pre>feature selection.r regression(X, y, *[,])</pre>	Compute Pearson's r for each features and the target.
<pre>feature selection.mutual info classif(X, y, *)</pre>	Estimate mutual information for a discrete target variable.
$\underline{\text{feature selection.mutual info regression}}(X,y,{}^{\star})$	Estimate mutual information for a continuous target variable.

sklearn.gaussian process: Gaussian Processes

The sklearn.gaussian process module implements Gaussian Process based regression and classification.

User guide: See the <u>Gaussian Processes</u> section for further details.

```
<u>gaussian process.GaussianProcessClassifier([...])</u> Gaussian process classification (GPC) based on Laplace approximation. 
<u>gaussian process.GaussianProcessRegressor([...])</u> Gaussian process regression (GPR).
```

Kernels:

$\underline{\texttt{gaussian_process.kernels.CompoundKernel}}(kernels)$	Kernel which is composed of a set of other kernels.
<pre>gaussian_process.kernels.ConstantKernel([])</pre>	Constant kernel.
$\underline{\texttt{gaussian_process.kernels.DotProduct}}([])$	Dot-Product kernel.
$\underline{gaussian_process.kernels.ExpSineSquared([])}$	Exp-Sine-Squared kernel (aka periodic kernel).
<pre>gaussian_process.kernels.Exponentiation()</pre>	The Exponentiation kernel takes one base kernel and a scalar parameter \boldsymbol{p} and combines them via
<pre>gaussian_process.kernels.Hyperparameter()</pre>	A kernel hyperparameter's specification in form of a namedtuple.
<pre>gaussian_process.kernels.Kernel()</pre>	Base class for all kernels.
gaussian_process.kernels.Matern([])	Matern kernel.
<pre>gaussian_process.kernels.PairwiseKernel([])</pre>	Wrapper for kernels in sklearn.metrics.pairwise.
<pre>gaussian_process.kernels.Product(k1, k2)</pre>	The Product kernel takes two kernels k_1 and k_2 and combines them via
Toggle Menu cess.kernels.RBF([length_scale,])	Radial basis function kernel (aka squared-exponential kernel).

 $\begin{array}{ll} \textbf{gaussian process.kernels.RationalQuadratic}([...]) & \textbf{Rational Quadratic kernel.} \\ \textbf{gaussian process.kernels.Sum}(\textbf{k1}, \textbf{k2}) & \textbf{The Sum kernel takes two kernels } k_1 \text{ and } k_2 \text{ and combines them via} \\ \textbf{gaussian process.kernels.WhiteKernel}([...]) & \textbf{White kernel.} \\ \end{array}$

sklearn.impute: Impute

Transformers for missing value imputation

<u>impute.SimpleImputer</u>(*[, missing_values, ...])

User guide: See the <u>Imputation of missing values</u> section for further details.

Univariate imputer for completing missing values with simple strategies.

sklearn.inspection: Inspection

The **sklearn.inspection** module includes tools for model inspection.

```
inspection.partial dependence(estimator, X, ...) Partial dependence of features.
inspection.permutation importance(estimator, ...) Permutation importance for feature evaluation [Rd9e56ef97513-BRE].
```

Plotting

```
<u>inspection.DecisionBoundaryDisplay(*, xx0, ...)</u> Decisions boundary visualization.

<u>inspection.PartialDependenceDisplay(...[, ...])</u> Partial Dependence Plot (PDP).
```

sklearn.isotonic: Isotonic regression

User guide: See the <u>Isotonic regression</u> section for further details.

```
isotonic.IsotonicRegression(*[, y_min, ...]) Isotonic regression model.

isotonic.check increasing(x, y) Determine whether y is monotonically correlated with x.
isotonic.isotonic regression(y, *[, ...]) Solve the isotonic regression model.
```

sklearn.kernel approximation: Kernel Approximation

The **sklearn.kernel** approximation module implements several approximate kernel feature maps based on Fourier transforms and Count Sketches.

User guide: See the <u>Kernel Approximation</u> section for further details.

```
kernel approximation. AdditiveChi2Sampler(*) Approximate feature map for additive chi2 kernel.

kernel approximation. Nystroem([kernel, ...]) Approximate a kernel map using a subset of the training data.

kernel approximation. PolynomialCountSketch(*) Polynomial kernel approximation via Tensor Sketch.

kernel approximation. RBFSampler(*[, gamma, ...]) Approximate a RBF kernel feature map using random Fourier features.

kernel approximation. SkewedChi2Sampler(*[, ...]) Approximate feature map for "skewed chi-squared" kernel.
```

sklearn.kernel_ridge: Kernel Ridge Regression

Module **sklearn.kernel ridge** implements kernel ridge regression.

User guide: See the <u>Kernel ridge regression</u> section for further details.

kernel ridge.KernelRidge([alpha, kernel, ...]) Kernel ridge regression.

Toggle Menu

sklearn.linear_model: Linear Models

The **sklearn.linear** model module implements a variety of linear models.

User guide: See the <u>Linear Models</u> section for further details.

The following subsections are only rough guidelines: the same estimator can fall into multiple categories, depending on its parameters.

Linear classifiers

<u>linear model.LogisticRegression([penalty,])</u>	Logistic Regression (aka logit, MaxEnt) classifier.
<pre>linear model.LogisticRegressionCV(*[, Cs,])</pre>	Logistic Regression CV (aka logit, MaxEnt) classifier.
<pre>linear model.PassiveAggressiveClassifier(*)</pre>	Passive Aggressive Classifier.
<pre>linear model.Perceptron(*[, penalty, alpha,])</pre>	Linear perceptron classifier.
<u>linear model.RidgeClassifier([alpha,])</u>	Classifier using Ridge regression.
<pre>linear model.RidgeClassifierCV([alphas,])</pre>	Ridge classifier with built-in cross-validation.
<pre>linear model.SGDClassifier([loss, penalty,])</pre>	Linear classifiers (SVM, logistic regression, etc.) with SGD training.
<pre>linear model.SGDOneClassSVM([nu,])</pre>	Solves linear One-Class SVM using Stochastic Gradient Descent.
←	

Classical linear regressors

<u>linear model.Ridge</u> ([alpha, fit_intercept,]) Linear least squares with I2 regularization. Ridge regression with built-in cross-validation.	<pre>linear model.LinearRegression(*[,])</pre>	Ordinary least squares Linear Regression.
linear model.RidgeCV([alphas,]) Ridge regression with built-in cross-validation.	<pre>linear model.Ridge([alpha, fit_intercept,])</pre>	Linear least squares with I2 regularization.
	<pre>linear model.RidgeCV([alphas,])</pre>	Ridge regression with built-in cross-validation.
<u>linear model.SGDRegressor</u> ([loss, penalty,]) Linear model fitted by minimizing a regularized empirical loss with S	<u>linear model.SGDRegressor</u> ([loss, penalty,])	Linear model fitted by minimizing a regularized empirical loss with SGD.

Regressors with variable selection

The following estimators have built-in variable selection fitting procedures, but any estimator using a L1 or elastic-net penalty also performs variable selection: typically SGDClassifier with an appropriate penalty.

<pre>linear model.ElasticNet([alpha, l1_ratio,])</pre>	Linear regression with combined L1 and L2 priors as regularizer.
<u>linear model.ElasticNetCV</u> (*[, l1_ratio,])	Elastic Net model with iterative fitting along a regularization path.
<pre>linear model.Lars(*[, fit_intercept,])</pre>	Least Angle Regression model a.k.a.
<pre>linear model.LarsCV(*[, fit_intercept,])</pre>	Cross-validated Least Angle Regression model.
<pre>linear model.Lasso([alpha, fit_intercept,])</pre>	Linear Model trained with L1 prior as regularizer (aka the Lasso).
<pre>linear model.LassoCV(*[, eps, n_alphas,])</pre>	Lasso linear model with iterative fitting along a regularization path.
<pre>linear model.LassoLars([alpha,])</pre>	Lasso model fit with Least Angle Regression a.k.a.
<pre>linear model.LassoLarsCV(*[, fit_intercept,])</pre>	Cross-validated Lasso, using the LARS algorithm.
<pre>linear model.LassoLarsIC([criterion,])</pre>	Lasso model fit with Lars using BIC or AIC for model selection.
<pre>linear model.OrthogonalMatchingPursuit(*[,])</pre>	Orthogonal Matching Pursuit model (OMP).
<pre>linear model.OrthogonalMatchingPursuitCV(*)</pre>	Cross-validated Orthogonal Matching Pursuit model (OMP).

Bayesian regressors

```
<u>linear model.ARDRegression</u>(*[, max_iter, ...]) Bayesian ARD regression.

<u>linear model.BayesianRidge</u>(*[, max_iter, ...]) Bayesian ridge regression.
```

Multi-task linear regressors with variable selection

These estimators fit multiple regression problems (or tasks) jointly, while inducing sparse coefficients. While the inferred coefficients may differ between the tasks, they are constrained to agree on the features that are selected (non-zero coefficients).

<u>linear model.MultiTaskElasticNet([alpha,])</u>	Multi-task ElasticNet model trained with L1/L2 mixed-norm as regularizer.
<u>linear model.MultiTaskElasticNetCV(*[,])</u>	Multi-task L1/L2 ElasticNet with built-in cross-validation.
<u>linear model.MultiTaskLasso([alpha,])</u>	Multi-task Lasso model trained with L1/L2 mixed-norm as regularizer.
<pre>linear model.MultiTaskLassoCV(*[, eps,])</pre>	Multi-task Lasso model trained with L1/L2 mixed-norm as regularizer.
→	

Outlier-robust regressors

Any estimator using the Huber loss would also be robust to outliers, e.g. **SGDRegressor** with loss='huber'.

```
linear model.HuberRegressor(*[, epsilon, ...])L2-regularized linear regression model that is robust to outliers.linear model.QuantileRegressor(*[, ...])Linear regression model that predicts conditional quantiles.linear model.RANSACRegressor([estimator, ...])RANSAC (RANdom SAmple Consensus) algorithm.linear model.TheilSenRegressor(*[, ...])Theil-Sen Estimator: robust multivariate regression model.
```

Generalized linear models (GLM) for regression

These models allow for response variables to have error distributions other than a normal distribution:

```
      linear model.PoissonRegressor(*[, alpha, ...])
      Generalized Linear Model with a Poisson distribution.

      linear model.TweedieRegressor(*[, power, ...])
      Generalized Linear Model with a Tweedie distribution.

      linear model.GammaRegressor(*[, alpha, ...])
      Generalized Linear Model with a Gamma distribution.
```

Miscellaneous

<u>linear model.PassiveAggressiveRegressor</u> (*[,])	Passive Aggressive Regressor.
<pre>linear model.enet path(X, y, *[, l1_ratio,])</pre>	Compute elastic net path with coordinate descent.
<pre>linear model.lars path(X, y[, Xy, Gram,])</pre>	Compute Least Angle Regression or Lasso path using the LARS algorithm [1].
<pre>linear model.lars path gram(Xy, Gram, *,)</pre>	The lars_path in the sufficient stats mode [1].
<pre>linear model.lasso path(X, y, *[, eps,])</pre>	Compute Lasso path with coordinate descent.
<pre>linear model.orthogonal mp(X, y, *[,])</pre>	Orthogonal Matching Pursuit (OMP).
<pre>linear model.orthogonal mp gram(Gram, Xy, *)</pre>	Gram Orthogonal Matching Pursuit (OMP).
<pre>linear model.ridge regression(X, y, alpha, *)</pre>	Solve the ridge equation by the method of normal equations.

sklearn.manifold: Manifold Learning

The **sklearn.manifold** module implements data embedding techniques.

User guide: See the <u>Manifold learning</u> section for further details.

<pre>manifold.Isomap(*[, n_neighbors, radius,])</pre>	Isomap Embedding.
<pre>manifold.LocallyLinearEmbedding(*[,])</pre>	Locally Linear Embedding.
<pre>manifold.MDS([n_components, metric, n_init,])</pre>	Multidimensional scaling.
$\underline{\texttt{manifold.SpectralEmbedding}}([n_components, \ldots])$	Spectral embedding for non-linear dimensionality reduction.
<pre>manifold.TSNE([n_components, perplexity,])</pre>	T-distributed Stochastic Neighbor Embedding.
4	
<pre>manifold.locally linear embedding(X, *,)</pre>	Perform a Locally Linear Embedding analysis on the data.
<pre>manifold.smacof(dissimilarities, *[,])</pre>	Compute multidimensional scaling using the SMACOF algori
<pre>manifold.spectral embedding(adjacency, *[,])</pre>	Project the sample on the first eigenvectors of the graph Lap
<pre>manifold.trustworthiness(X, X_embedded, *[,])</pre>	Indicate to what extent the local structure is retained.
4	

sklearn.metrics: Metrics

See the <u>Metrics and scoring: quantifying the quality of predictions</u> section and the <u>Pairwise metrics</u>, <u>Affinities and Kernels</u> section of the user guide for further details.

The **sklearn.metrics** module includes score functions, performance metrics and pairwise metrics and distance computations.

Model Selection Interface

See the The scoring parameter: defining model evaluation rules section of the user guide for further details.

metrics.check scoring(estimator[, scoring,]) Determine scorer from user options.
metrics.get scorer(scoring)	Get a scorer from string.
<pre>metrics.get scorer names()</pre>	Get the names of all available scorers.
<pre>metrics.make scorer(score func, *[,])</pre>	Make a scorer from a performance metric or loss function.

Classification metrics

See the <u>Classification metrics</u> section of the user guide for further details.

<pre>metrics.accuracy_score(y_true, y_pred, *[,])</pre>	Accuracy classification score.	
<pre>metrics.auc(X, y)</pre>	Compute Area Under the Curve (AUC) using the trapezoidal rule.	
<pre>metrics.average_precision_score(y_true,)</pre>	Compute average precision (AP) from prediction scores.	
<pre>metrics.balanced accuracy score(y_true,)</pre>	Compute the balanced accuracy.	
<pre>metrics.brier_score_loss(y_true, y_prob, *)</pre>	Compute the Brier score loss.	
<pre>metrics.class likelihood ratios(y_true,)</pre>	Compute binary classification positive and negative likelihood ratios.	
<pre>metrics.classification report(y_true, y_pred, *)</pre>	Build a text report showing the main classification metrics.	
metrics.cohen kappa score(y1, y2, *[,])	Compute Cohen's kappa: a statistic that measures inter-annotator agreement.	
<pre>metrics.confusion matrix(y_true, y_pred, *)</pre>	Compute confusion matrix to evaluate the accuracy of a classification.	
<pre>metrics.dcg score(y_true, y_score, *[, k,])</pre>	Compute Discounted Cumulative Gain.	
<pre>metrics.det curve(y_true, y_score[,])</pre>	Compute error rates for different probability thresholds.	
<pre>metrics.f1 score(y_true, y_pred, *[,])</pre>	Compute the F1 score, also known as balanced F-score or F-measure.	
<pre>metrics.fbeta score(y_true, y_pred, *, beta)</pre>	Compute the F-beta score.	
<pre>metrics.hamming loss(y_true, y_pred, *[,])</pre>	Compute the average Hamming loss.	
<pre>metrics.hinge loss(y_true, pred_decision, *)</pre>	Average hinge loss (non-regularized).	
<pre>metrics.jaccard score(y_true, y_pred, *[,])</pre>	Jaccard similarity coefficient score.	
<pre>metrics.log_loss(y_true, y_pred, *[, eps,])</pre>	Log loss, aka logistic loss or cross-entropy loss.	
<pre>metrics.matthews corrcoef(y_true, y_pred, *)</pre>	Compute the Matthews correlation coefficient (MCC).	
<pre>metrics.multilabel confusion matrix(y_true,)</pre>	Compute a confusion matrix for each class or sample.	
<pre>metrics.ndcg score(y_true, y_score, *[, k,])</pre>	Compute Normalized Discounted Cumulative Gain.	
<pre>metrics.precision recall curve(y_true,)</pre>	Compute precision-recall pairs for different probability thresholds.	
metrics.precision recall fscore support()	Compute precision, recall, F-measure and support for each class.	
<pre>metrics.precision score(y_true, y_pred, *[,])</pre>	Compute the precision.	
<pre>metrics.recall score(y_true, y_pred, *[,])</pre>	Compute the recall.	
<pre>metrics.roc auc score(y_true, y_score, *[,])</pre>	Compute Area Under the Receiver Operating Characteristic Curve (ROC AUC) scores.	from prediction
<pre>metrics.roc curve(y_true, y_score, *[,])</pre>	Compute Receiver operating characteristic (ROC).	
<pre>metrics.top k accuracy_score(y_true, y_score, *)</pre>	Top-k Accuracy classification score.	
<pre>metrics.zero one loss(y_true, y_pred, *[,])</pre>	Zero-one classification loss.	
4		

Regression metrics

See the $\underline{\text{Regression metrics}}$ section of the user guide for further details.

<pre>metrics.explained variance score(y_true,)</pre>	Explained variance regression score function.
<pre>metrics.max error(y_true, y_pred)</pre>	The max_error metric calculates the maximum residual error.
<pre>metrics.mean absolute error(y_true, y_pred, *)</pre>	Mean absolute error regression loss.
<pre>metrics.mean squared error(y_true, y_pred, *)</pre>	Mean squared error regression loss.
<pre>metrics.mean squared log_error(y_true, y_pred, *)</pre>	Mean squared logarithmic error regression loss.
<pre>metrics.median absolute error(y_true, y_pred, *)</pre>	Median absolute error regression loss.
<pre>metrics.mean absolute percentage error()</pre>	Mean absolute percentage error (MAPE) regression loss.
<pre>metrics.r2 score(y_true, y_pred, *[,])</pre>	R^2 (coefficient of determination) regression score function.
<pre>metrics.mean_poisson_deviance(y_true, y_pred, *)</pre>	Mean Poisson deviance regression loss.
<pre>metrics.mean_gamma_deviance(y_true, y_pred, *)</pre>	Mean Gamma deviance regression loss.
<pre>metrics.mean tweedie deviance(y_true, y_pred, *)</pre>	Mean Tweedie deviance regression loss.
<pre>metrics.d2 tweedie score(y_true, y_pred, *)</pre>	D^2 regression score function, fraction of Tweedie deviance explained.
<pre>metrics.mean_pinball loss(y_true, y_pred, *)</pre>	Pinball loss for quantile regression.
<pre>metrics.d2 pinball score(y_true, y_pred, *)</pre>	D^2 regression score function, fraction of pinball loss explained.
<pre>metrics.d2 absolute error score(y_true,)</pre>	D^2 regression score function, fraction of absolute error explained.
4	

Multilabel ranking metrics

See the $\underline{\text{Multilabel ranking metrics}}$ section of the user guide for further details.

<pre>metrics.coverage error(y_true, y_score, *[,])</pre>	Coverage error measure.
<pre>metrics.label ranking average precision score()</pre>	Compute ranking-based average precision.
<pre>metrics.label ranking_loss(y_true, y_score, *)</pre>	Compute Ranking loss measure.

Clustering metrics

See the <u>Clustering performance evaluation</u> section of the user guide for further details.

The sklearn.metrics.cluster submodule contains evaluation metrics for cluster analysis results. There are two forms of evaluation:

- supervised, which uses a ground truth class values for each sample.
- unsupervised, which does not and measures the 'quality' of the model itself.

<pre>metrics.adjusted mutual info score([,])</pre>	Adjusted Mutual Information between two clusterings.
<pre>metrics.adjusted rand score(labels_true,)</pre>	Rand index adjusted for chance.
metrics.calinski harabasz score(X, labels)	Compute the Calinski and Harabasz score.
<pre>metrics.davies bouldin score(X, labels)</pre>	Compute the Davies-Bouldin score.
<pre>metrics.completeness_score(labels_true,)</pre>	Compute completeness metric of a cluster labeling given a ground truth.
<pre>metrics.cluster.contingency matrix([,])</pre>	Build a contingency matrix describing the relationship between labels.
<pre>metrics.cluster.pair confusion matrix()</pre>	Pair confusion matrix arising from two clusterings [R9ca8fd06d29a-1].
<pre>metrics.fowlkes mallows score(labels_true,)</pre>	Measure the similarity of two clusterings of a set of points.
$\underline{\texttt{metrics.homogeneity}}\ \underline{\texttt{completeness}}\ \underline{\texttt{v}}\ \underline{\texttt{measure}}()$	Compute the homogeneity and completeness and V-Measure scores at once.
<pre>metrics.homogeneity_score(labels_true,)</pre>	Homogeneity metric of a cluster labeling given a ground truth.
<pre>metrics.mutual info score(labels_true,)</pre>	Mutual Information between two clusterings.
<pre>metrics.normalized mutual info score([,])</pre>	Normalized Mutual Information between two clusterings.
<pre>metrics.rand score(labels_true, labels_pred)</pre>	Rand index.
<pre>metrics.silhouette score(X, labels, *[,])</pre>	Compute the mean Silhouette Coefficient of all samples.
<pre>metrics.silhouette samples(X, labels, *[,])</pre>	Compute the Silhouette Coefficient for each sample.
<pre>metrics.v measure score(labels_true,[, beta])</pre>	V-measure cluster labeling given a ground truth.
4	

Biclustering metrics

See the <u>Biclustering evaluation</u> section of the user guide for further details.

metrics.consensus score(a, b, *[, similarity]) The similarity of two sets of biclusters.

Distance metrics

 $\underline{\texttt{metrics.DistanceMetric}} \quad \text{Uniform interface for fast distance metric functions}.$

Pairwise metrics

See the <u>Pairwise metrics</u>, <u>Affinities and Kernels</u> section of the user guide for further details.

<pre>metrics.pairwise.additive chi2 kernel(X[, Y])</pre>	Compute the additive chi-squared kernel between observations in X and Y.
<pre>metrics.pairwise.chi2 kernel(X[, Y, gamma])</pre>	Compute the exponential chi-squared kernel between X and Y.
<pre>metrics.pairwise.cosine similarity(X[, Y,])</pre>	Compute cosine similarity between samples in X and Y.
<pre>metrics.pairwise.cosine distances(X[, Y])</pre>	Compute cosine distance between samples in X and Y.
<pre>metrics.pairwise.distance metrics()</pre>	Valid metrics for pairwise_distances.
metrics.pairwise.euclidean distances(X[, Y,])	Compute the distance matrix between each pair from a vector array X and Y.
metrics.pairwise.haversine distances(X[, Y])	Compute the Haversine distance between samples in X and Y.
<pre>metrics.pairwise.kernel metrics()</pre>	Valid metrics for pairwise_kernels.
<pre>metrics.pairwise.laplacian kernel(X[, Y, gamma])</pre>	Compute the laplacian kernel between X and Y.
metrics.pairwise.linear kernel(X[, Y,])	Compute the linear kernel between X and Y.
metrics.pairwise.manhattan distances(X[, Y,])	Compute the L1 distances between the vectors in X and Y.
metrics.pairwise.nan euclidean distances(X)	Calculate the euclidean distances in the presence of missing values.
<pre>metrics.pairwise.pairwise kernels(X[, Y,])</pre>	Compute the kernel between arrays X and optional array Y.
<pre>metrics.pairwise.polynomial kernel(X[, Y,])</pre>	Compute the polynomial kernel between X and Y.
<pre>metrics.pairwise.rbf kernel(X[, Y, gamma])</pre>	Compute the rbf (gaussian) kernel between X and Y.
<pre>metrics.pairwise.sigmoid kernel(X[, Y,])</pre>	Compute the sigmoid kernel between X and Y.
<pre>metrics.pairwise.paired euclidean distances(X, Y)</pre>	Compute the paired euclidean distances between X and Y.
metrics.pairwise.paired manhattan distances(X, Y)	Compute the paired L1 distances between X and Y.
metrics.pairwise.paired cosine distances(X, Y)	Compute the paired cosine distances between X and Y.
<pre>metrics.pairwise.paired distances(X, Y, *[,])</pre>	Compute the paired distances between X and Y.
<pre>metrics.pairwise distances(X[, Y, metric,])</pre>	Compute the distance matrix from a vector array X and optional Y.
<pre>metrics.pairwise distances argmin(X, Y, *[,])</pre>	Compute minimum distances between one point and a set of points.
<pre>metrics.pairwise distances argmin min(X, Y, *)</pre>	Compute minimum distances between one point and a set of points.

metrics.pairwise distances chunked(X[, Y, ...]) Generate a distance matrix chunk by chunk with optional reduction.

Plotting

See the Visualizations section of the user guide for further details.

<pre>metrics.ConfusionMatrixDisplay([,])</pre>	Confusion Matrix visualization.
<pre>metrics.DetCurveDisplay(*, fpr, fnr[,])</pre>	DET curve visualization.
<pre>metrics.PrecisionRecallDisplay(precision,)</pre>	Precision Recall visualization.
<pre>metrics.PredictionErrorDisplay(*, y_true, y_pred)</pre>	Visualization of the prediction error of a regression model.
<pre>metrics.RocCurveDisplay(*, fpr, tpr[,])</pre>	ROC Curve visualization.
<pre>calibration.CalibrationDisplay(prob_true,)</pre>	Calibration curve (also known as reliability diagram) visualization.

sklearn.mixture: Gaussian Mixture Models

The **sklearn.mixture** module implements mixture modeling algorithms.

User guide: See the <u>Gaussian mixture models</u> section for further details.

<u>mixture.BayesianGaussianMixture</u>(*[, ...]) Variational Bayesian estimation of a Gaussian mixture. <u>mixture.GaussianMixture</u>([n_components, ...]) Gaussian Mixture.

sklearn.model_selection: Model Selection

User guide: See the <u>Cross-validation: evaluating estimator performance</u>, <u>Tuning the hyper-parameters of an estimator</u> and <u>Learning curve</u> sections for further details.

Splitter Classes

<pre>model selection.GroupKFold([n_splits])</pre>	K-fold iterator variant with non-overlapping groups.
<pre>model selection.GroupShuffleSplit([])</pre>	Shuffle-Group(s)-Out cross-validation iterator
<pre>model selection.KFold([n_splits, shuffle,])</pre>	K-Folds cross-validator
<pre>model selection.LeaveOneGroupOut()</pre>	Leave One Group Out cross-validator
<pre>model selection.LeavePGroupsOut(n_groups)</pre>	Leave P Group(s) Out cross-validator
<pre>model selection.LeaveOneOut()</pre>	Leave-One-Out cross-validator
<pre>model selection.LeavePOut(p)</pre>	Leave-P-Out cross-validator
<pre>model selection.PredefinedSplit(test_fold)</pre>	Predefined split cross-validator
<pre>model selection.RepeatedKFold(*[, n_splits,])</pre>	Repeated K-Fold cross validator.
$\underline{model\ selection.RepeatedStratifiedKFold}(*[,])$	Repeated Stratified K-Fold cross validator.
<pre>model selection.ShuffleSplit([n_splits,])</pre>	Random permutation cross-validator
<pre>model selection.StratifiedKFold([n_splits,])</pre>	Stratified K-Folds cross-validator.
<pre>model selection.StratifiedShuffleSplit([])</pre>	Stratified ShuffleSplit cross-validator
<pre>model selection.StratifiedGroupKFold([])</pre>	Stratified K-Folds iterator variant with non-overlapping groups.
<pre>model selection.TimeSeriesSplit([n_splits,])</pre>	Time Series cross-validator
4	

Splitter Functions

model selection.check cv([cv, y, classifier]) Input checker utility for building a cross-validator.

model selection.train test split(*arrays[, ...]) Split arrays or matrices into random train and test subsets.

Hyper-parameter optimizers

<pre>model selection.GridSearchCV(estimator,)</pre>	Exhaustive search over specified parameter values for an estimator.
$\underline{model\ selection.HalvingGridSearchCV}([,\])$	Search over specified parameter values with successive halving.
<pre>model selection.ParameterGrid(param_grid)</pre>	Grid of parameters with a discrete number of values for each.
<pre>model selection.ParameterSampler([,])</pre>	Generator on parameters sampled from given distributions.
$\underline{model\ selection.RandomizedSearchCV}([,\])$	Randomized search on hyper parameters.
model_selection.HalvingRandomSearchCV([,])	Randomized search on hyper parameters.

model selection.cross validate(estimator, X) Evaluate metric(s) by cross-validation and also record fit/score times. model selection.cross val predict(estimator, X) Generate cross-validated estimates for each input data point. model selection.cross val score(estimator, X) Evaluate a score by cross-validation. <u>model selection.learning curve</u>(estimator, X, ...) Learning curve. <u>model selection.permutation test score(...)</u> Evaluate the significance of a cross-validated score with permutations. <u>model selection.validation curve</u>(estimator, ...) Validation curve.

Visualization

Model validation

```
<u>model_selection.LearningCurveDisplay(*, ...)</u>
                                                      Learning Curve visualization.
<u>model selection.ValidationCurveDisplay(*, ...)</u> Validation Curve visualization.
```

sklearn.multiclass: Multiclass classification

Multiclass classification strategies

This module implements multiclass learning algorithms:

- one-vs-the-rest / one-vs-all
- one-vs-one
- error correcting output codes

The estimators provided in this module are meta-estimators: they require a base estimator to be provided in their constructor. For example, it is possible to use these estimators to turn a binary classifier or a regressor into a multiclass classifier. It is also possible to use these estimators with multiclass estimators in the hope that their accuracy or runtime performance improves.

All classifiers in scikit-learn implement multiclass classification; you only need to use this module if you want to experiment with custom multiclass strategies.

The one-vs-the-rest meta-classifier also implements a predict_proba method, so long as such a method is implemented by the base classifier. This method returns probabilities of class membership in both the single label and multilabel case. Note that in the multilabel case, probabilities are the marginal probability that a given sample falls in the given class. As such, in the multilabel case the sum of these probabilities over all possible labels for a given sample *will not* sum to unity, as they do in the single label case.

User guide: See the <u>Multiclass classification</u> section for further details.

```
multiclass.OneVsRestClassifier(estimator, *)
                                                One-vs-the-rest (OvR) multiclass strategy.
multiclass.OneVsOneClassifier(estimator, *)
                                                 One-vs-one multiclass strategy.
multiclass.OutputCodeClassifier(estimator, *) (Error-Correcting) Output-Code multiclass strategy.
```

sklearn.multioutput: Multioutput regression and classification

This module implements multioutput regression and classification.

The estimators provided in this module are meta-estimators: they require a base estimator to be provided in their constructor. The meta-estimator extends single output estimators to multioutput estimators.

User guide: See the <u>Multilabel classification</u>, <u>Multiclass-multioutput classification</u>, and <u>Multioutput regression</u> sections for further details.

```
multioutput.ClassifierChain(base_estimator, *)
                                                    A multi-label model that arranges binary classifiers into a chain.
multioutput.MultiOutputRegressor(estimator, *)
                                                    Multi target regression.
multioutput.MultiOutputClassifier(estimator, *)
                                                   Multi target classification.
multioutput.RegressorChain(base_estimator, *)
                                                    A multi-label model that arranges regressions into a chain.
```

sklearn.naive_bayes: Naive Bayes

The sklearn.naive bayes module implements Naive Bayes algorithms. These are supervised learning methods based on applying Bayes' theorem with strong (naive) feature independence assumptions.

See the Naive Bayes section for further details.

Toggle Menu

```
    naive bayes.BernoulliNB(*[, alpha, ...])
    Naive Bayes classifier for multivariate Bernoulli models.
    naive bayes.CategoricalNB(*[, alpha, ...])
    Naive Bayes classifier for categorical features.
    naive bayes.ComplementNB(*[, alpha, ...])
    The Complement Naive Bayes classifier described in Rennie et al. (2003).
    naive bayes.GaussianNB(*[, priors, ...])
    Gaussian Naive Bayes (GaussianNB).
    naive bayes.MultinomialNB(*[, alpha, ...])
    Naive Bayes classifier for multinomial models.
```

sklearn.neighbors: Nearest Neighbors

The **sklearn.neighbors** module implements the k-nearest neighbors algorithm.

User guide: See the Nearest Neighbors section for further details.

<pre>neighbors.BallTree(X[, leaf_size, metric])</pre>	BallTree for fast generalized N-point problems
<pre>neighbors.KDTree(X[, leaf_size, metric])</pre>	KDTree for fast generalized N-point problems
<pre>neighbors.KernelDensity(*[, bandwidth,])</pre>	Kernel Density Estimation.
<pre>neighbors.KNeighborsClassifier([])</pre>	Classifier implementing the k-nearest neighbors vote.
$\underline{neighbors.KNeighborsRegressor}([n_neighbors,])$	Regression based on k-nearest neighbors.
<pre>neighbors.KNeighborsTransformer(*[, mode,])</pre>	Transform X into a (weighted) graph of k nearest neighbors.
<pre>neighbors.LocalOutlierFactor([n_neighbors,])</pre>	Unsupervised Outlier Detection using the Local Outlier Factor (LOF).
$\underline{\texttt{neighbors.RadiusNeighborsClassifier}}([])$	Classifier implementing a vote among neighbors within a given radius.
$\underline{neighbors.RadiusNeighborsRegressor}([radius, \ldots])$	Regression based on neighbors within a fixed radius.
<u>neighbors.RadiusNeighborsTransformer(*[,])</u>	Transform X into a (weighted) graph of neighbors nearer than a radius.
neighbors.RadiusNeighborsTransformer(*[,]) neighbors.NearestCentroid([metric,])	Transform X into a (weighted) graph of neighbors nearer than a radius. Nearest centroid classifier.
-	, , , , , ,
neighbors.NearestCentroid([metric,])	Nearest centroid classifier.
<pre>neighbors.NearestCentroid([metric,]) neighbors.NearestNeighbors(*[, n_neighbors,])</pre>	Nearest centroid classifier. Unsupervised learner for implementing neighbor searches.
<pre>neighbors.NearestCentroid([metric,]) neighbors.NearestNeighbors(*[, n_neighbors,])</pre>	Nearest centroid classifier. Unsupervised learner for implementing neighbor searches.
<pre>neighbors.NearestCentroid([metric,]) neighbors.NearestNeighbors(*[, n_neighbors,]) neighbors.NeighborhoodComponentsAnalysis([])</pre>	Nearest centroid classifier. Unsupervised learner for implementing neighbor searches. Neighborhood Components Analysis.
<pre>neighbors.NearestCentroid([metric,]) neighbors.NearestNeighbors(*[, n_neighbors,]) neighbors.NeighborhoodComponentsAnalysis([]) neighbors.kneighbors_graph(X, n_neighbors, *)</pre>	Nearest centroid classifier. Unsupervised learner for implementing neighbor searches. Neighborhood Components Analysis. Compute the (weighted) graph of k-Neighbors for points in X.

sklearn.neural_network: Neural network models

The **sklearn.neural network** module includes models based on neural networks.

User guide: See the Neural network models (supervised) and Neural network models (unsupervised) sections for further details.

```
      neural network.BernoulliRBM([n_components, ...])
      Bernoulli Restricted Boltzmann Machine (RBM).

      neural network.MLPClassifier([...])
      Multi-layer Perceptron classifier.

      neural network.MLPRegressor([...])
      Multi-layer Perceptron regressor.
```

sklearn.pipeline: Pipeline

The sklearn.pipeline module implements utilities to build a composite estimator, as a chain of transforms and estimators.

User guide: See the <u>Pipelines and composite estimators</u> section for further details.

```
pipeline.FeatureUnion(transformer_list, *[, ...]) Concatenates results of multiple transformer objects.
pipeline.Pipeline(steps, *[, memory, verbose]) Pipeline of transforms with a final estimator.
pipeline.Pipeline(*steps[, memory, verbose]) Construct a Binaline from the given estimators.
```

```
pipeline.make pipeline(*steps[, memory, verbose]) Construct a Pipeline from the given estimators.
pipeline.make union(*transformers[, n_jobs, ...]) Construct a FeatureUnion from the given transformers.
```

sklearn.preprocessing: Preprocessing and Normalization

The sklearn.preprocessing module includes scaling, centering, normalization, binarization methods.

User guide: See the Preprocessing data section for further details.

Toggle Menu <u>g.Binarizer</u>(*[, threshold, copy]) Binarize data (set feature values to 0 or 1) according to a threshold.

<pre>preprocessing.FunctionTransformer([func,])</pre>	Constructs a transformer from an arbitrary callable.
<pre>preprocessing.KBinsDiscretizer([n_bins,])</pre>	Bin continuous data into intervals.
<pre>preprocessing.KernelCenterer()</pre>	Center an arbitrary kernel matrix K .
<pre>preprocessing.LabelBinarizer(*[, neg_label,])</pre>	Binarize labels in a one-vs-all fashion.
<pre>preprocessing.LabelEncoder()</pre>	Encode target labels with value between 0 and n_classes-1.
$\underline{\texttt{preprocessing.MultiLabelBinarizer}}(\texttt{*}[, \ldots])$	Transform between iterable of iterables and a multilabel format.
<pre>preprocessing.MaxAbsScaler(*[, copy])</pre>	Scale each feature by its maximum absolute value.
<pre>preprocessing.MinMaxScaler([feature_range,])</pre>	Transform features by scaling each feature to a given range.
<pre>preprocessing.Normalizer([norm, copy])</pre>	Normalize samples individually to unit norm.
$\underline{\textbf{preprocessing.0} \textbf{oneHotEncoder}}(\texttt{*[}, \texttt{categories},])$	Encode categorical features as a one-hot numeric array.
<pre>preprocessing.OrdinalEncoder(*[,])</pre>	Encode categorical features as an integer array.
$\underline{\textbf{preprocessing.PolynomialFeatures}} ([\texttt{degree}, \dots])$	Generate polynomial and interaction features.
<pre>preprocessing.PowerTransformer([method,])</pre>	Apply a power transform featurewise to make data more Gaussian-like.
$\underline{\textbf{preprocessing.QuantileTransformer}}(*[,])$	Transform features using quantiles information.
$\underline{\textbf{preprocessing.RobustScaler}}(*[, \dots])$	Scale features using statistics that are robust to outliers.
$\underline{\textbf{preprocessing.SplineTransformer}}([n_knots, \ldots])$	Generate univariate B-spline bases for features.
<pre>preprocessing.StandardScaler(*[, copy,])</pre>	Standardize features by removing the mean and scaling to unit variance.
<pre>preprocessing.TargetEncoder([categories,])</pre>	Target Encoder for regression and classification targets.
<pre>preprocessing.add dummy feature(X[, value])</pre>	Augment dataset with an additional dummy feature.
<pre>preprocessing.binarize(X, *[, threshold, copy])</pre>	Boolean thresholding of array-like or scipy.sparse matrix.
<pre>preprocessing.label binarize(y, *, classes)</pre>	Binarize labels in a one-vs-all fashion.
<pre>preprocessing.maxabs scale(X, *[, axis, copy])</pre>	Scale each feature to the [-1, 1] range without breaking the sparsity.
<pre>preprocessing.minmax scale(X[,])</pre>	Transform features by scaling each feature to a given range.
<pre>preprocessing.normalize(X[, norm, axis,])</pre>	Scale input vectors individually to unit norm (vector length).
<pre>preprocessing.quantile transform(X, *[,])</pre>	Transform features using quantiles information.
<pre>preprocessing.robust scale(X, *[, axis,])</pre>	Standardize a dataset along any axis.
<pre>preprocessing.scale(X, *[, axis, with_mean,])</pre>	Standardize a dataset along any axis.
<pre>preprocessing.power transform(X[, method,])</pre>	Parametric, monotonic transformation to make data more Gaussian-like.
4	

sklearn.random_projection: Random projection

Random Projection transformers.

Random Projections are a simple and computationally efficient way to reduce the dimensionality of the data by trading a controlled amount of accuracy (as additional variance) for faster processing times and smaller model sizes.

The dimensions and distribution of Random Projections matrices are controlled so as to preserve the pairwise distances between any two samples of the dataset.

The main theoretical result behind the efficiency of random projection is the <u>Johnson-Lindenstrauss lemma (quoting Wikipedia)</u>:

In mathematics, the Johnson-Lindenstrauss lemma is a result concerning low-distortion embeddings of points from high-dimensional into low-dimensional Euclidean space. The lemma states that a small set of points in a high-dimensional space can be embedded into a space of much lower dimension in such a way that distances between the points are nearly preserved. The map used for the embedding is at least Lipschitz, and can even be taken to be an orthogonal projection.

User guide: See the Random Projection section for further details.

<u>random projection.GaussianRandomProjection([...])</u> Reduce dimensionality through Gaussian random projection.

<u>random projection.SparseRandomProjection([...])</u> Reduce dimensionality through sparse random projection.

<u>random projection.johnson lindenstrauss min dim(...)</u> Find a 'safe' number of components to randomly project to.

sklearn.semi_supervised: Semi-Supervised Learning

The <u>sklearn.semi supervised</u> module implements semi-supervised learning algorithms. These algorithms utilize small amounts of labeled data and large amounts of unlabeled data for classification tasks. This module includes Label Propagation.

User guide: See the Semi-supervised learning section for further details.

sed.LabelPropagation([kernel, ...]) Label Propagation classifier.

semi supervised.LabelSpreading([kernel, ...]) LabelSpreading model for semi-supervised learning.
semi supervised.SelfTrainingClassifier(...) Self-training classifier.

sklearn.svm: Support Vector Machines

The **sklearn.svm** module includes Support Vector Machine algorithms.

User guide: See the Support Vector Machines section for further details.

Estimators

```
    svm.LinearSVC([penalty, loss, dual, tol, C, ...])
    svm.LinearSVR(*[, epsilon, tol, C, loss, ...])
    Linear Support Vector Regression.
    svm.NuSVC(*[, nu, kernel, degree, gamma, ...])
    Nu-Support Vector Classification.
    svm.NuSVR(*[, nu, C, kernel, degree, gamma, ...])
    Nu Support Vector Regression.
    svm.OneClassSVM(*[, kernel, degree, gamma, ...])
    Unsupervised Outlier Detection.
    svm.SVC(*[, C, kernel, degree, gamma, ...])
    C-Support Vector Classification.
    svm.SVR(*[, kernel, degree, gamma, coef0, ...])
    Epsilon-Support Vector Regression.
```

 $\underline{\text{svm.11 min c}}(X,\,y,\,{}^*[,\,loss,\,fit_intercept,\,...]) \quad \text{Return the lowest bound for C}.$

sklearn.tree: Decision Trees

The sklearn.tree module includes decision tree-based models for classification and regression.

User guide: See the <u>Decision Trees</u> section for further details.

```
tree.DecisionTreeClassifier(*[, criterion, ...]) A decision tree classifier.
tree.DecisionTreeRegressor(*[, criterion, ...]) A decision tree regressor.
tree.ExtraTreeClassifier(*[, criterion, ...]) An extremely randomized tree classifier.
tree.ExtraTreeRegressor(*[, criterion, ...]) An extremely randomized tree regressor.
```

<u>tree.export_graphviz</u>(decision_tree[, ...]) Export a decision tree in DOT format.

<u>tree.export_text</u>(decision_tree, *[, ...]) Build a text report showing the rules of a decision tree.

Plotting

<u>tree.plot_tree</u>(decision_tree, *[, ...]) Plot a decision tree.

sklearn.utils: Utilities

The sklearn.utils module includes various utilities.

Developer guide: See the <u>Utilities for Developers</u> page for further details.

<u>utils.Bunch</u>(**kwargs) Container object exposing keys as attributes.

<pre>utils.arrayfuncs.min_pos(X)</pre>	Find the minimum value of an array over positive values
utils.as float array(X, *[, copy,])	Convert an array-like to an array of floats.
<pre>utils.assert all finite(X, *[, allow_nan,])</pre>	Throw a ValueError if X contains NaN or infinity.
<pre>utils.check X y(X, y[, accept_sparse,])</pre>	Input validation for standard estimators.
<pre>utils.check array(array[, accept_sparse,])</pre>	Input validation on an array, list, sparse matrix or similar.
<pre>utils.check scalar(x, name, target_type, *)</pre>	Validate scalar parameters type and value.
utils.check consistent length(*arrays)	Check that all arrays have consistent first dimensions.
utils.check random state(Seed)	Turn seed into a np.random.RandomState instance.
utils.class weight.compute class weight()	Estimate class weights for unbalanced datasets.
utils.class weight.compute sample weight()	Estimate sample weights by class for unbalanced datasets.
Toggle Menu ated([extra])	Decorator to mark a function or class as deprecated.
Toggic McHa	

utils.estimator checks.check estimator([])	Check if estimator adheres to scikit-learn conventions.
utils.estimator checks.parametrize with checks()	Pytest specific decorator for parametrizing estimator checks.
utils.estimator html repr(estimator)	Build a HTML representation of an estimator.
<pre>utils.extmath.safe sparse dot(a, b, *[,])</pre>	Dot product that handle the sparse matrix case correctly.
<pre>utils.extmath.randomized range finder(A, *,)</pre>	Compute an orthonormal matrix whose range approximates the range of A.
<pre>utils.extmath.randomized svd(M, n_components, *)</pre>	Compute a truncated randomized SVD.
utils.extmath.fast logdet(A)	Compute logarithm of determinant of a square matrix.
<pre>utils.extmath.density(w, **kwargs)</pre>	Compute density of a sparse vector.
<pre>utils.extmath.weighted mode(a, w, *[, axis])</pre>	Return an array of the weighted modal (most common) value in the passed array.
<pre>utils.gen batches(n, batch_size, *[,])</pre>	Generator to create slices containing batch_size elements from 0 to n.
<pre>utils.gen even slices(n, n_packs, *[, n_samples])</pre>	Generator to create n_packs evenly spaced slices going up to n.
utils.graph.single source shortest path length()	Return the length of the shortest path from source to all reachable nodes.
utils.indexable(*iterables)	Make arrays indexable for cross-validation.
<pre>utils.metaestimators.available if(check)</pre>	An attribute that is available only if check returns a truthy value.
<pre>utils.multiclass.type of target(y[, input_name])</pre>	Determine the type of data indicated by the target.
<pre>utils.multiclass.is multilabel(y)</pre>	Check if y is in a multilabel format.
utils.multiclass.unique labels(*yS)	Extract an ordered array of unique labels.
utils.murmurhash3 32	Compute the 32bit murmurhash3 of key at seed.
<pre>utils.resample(*arrays[, replace,])</pre>	Resample arrays or sparse matrices in a consistent way.
<pre>utils. safe indexing(X, indices, *[, axis])</pre>	Return rows, items or columns of X using indices.
<pre>utils.safe mask(X, mask)</pre>	Return a mask which is safe to use on X.
<pre>utils.safe sqr(X, *[, copy])</pre>	Element wise squaring of array-likes and sparse matrices.
<pre>utils.shuffle(*arrays[, random_state, n_samples])</pre>	Shuffle arrays or sparse matrices in a consistent way.
<pre>utils.sparsefuncs.incr mean variance axis(X,)</pre>	Compute incremental mean and variance along an axis on a CSR or CSC matrix.
<pre>utils.sparsefuncs.inplace column scale(X, scale)</pre>	Inplace column scaling of a CSC/CSR matrix.
<pre>utils.sparsefuncs.inplace row scale(X, scale)</pre>	Inplace row scaling of a CSR or CSC matrix.
utils.sparsefuncs.inplace swap row(X, m, n)	Swap two rows of a CSC/CSR matrix in-place.
utils.sparsefuncs.inplace swap column(X, m, n)	Swap two columns of a CSC/CSR matrix in-place.
<pre>utils.sparsefuncs.mean variance axis(X, axis)</pre>	Compute mean and variance along an axis on a CSR or CSC matrix.
<pre>utils.sparsefuncs.inplace csr column scale(X,)</pre>	Inplace column scaling of a CSR matrix.
utils.sparsefuncs fast.inplace csr row normalize 11	Inplace row normalize using the I1 norm
utils.sparsefuncs fast.inplace csr row normalize 12	Inplace row normalize using the I2 norm
utils.random.sample without replacement	Sample integers without replacement.
<pre>utils.validation.check is fitted(estimator)</pre>	Perform is_fitted validation for estimator.
<u>utils.validation.check memory</u> (memory)	Check that memory is joblib.Memory-like.
<pre>utils.validation.check symmetric(array, *[,])</pre>	Make sure that array is 2D, square and symmetric.
<pre>utils.validation.column or 1d(y, *[, dtype,])</pre>	Ravel column or 1d numpy array, else raises an error.
utils.validation.has fit parameter()	Check whether the estimator's fit method supports the given parameter.
utils.metadata routing.get routing for object([Obj])	Get a Metadata{Router, Request} instance from the given object.
<pre>utils.metadata routing.MetadataRouter(OWNer)</pre>	Stores and handles metadata routing for a router object.
utils.metadata routing.MetadataRequest(OWNer)	Contains the metadata request info of a consumer.
<pre>utils.metadata routing.MethodMapping()</pre>	Stores the mapping between callee and caller methods for a router.
utils.metadata routing.process routing(Obj,)	Validate and route input parameters.

Specific utilities to list scikit-learn components:

<pre>utils.discovery.all estimators([type_filter])</pre>	Get a list of all estimators from sklearn.
utils.discovery.all displays()	Get a list of all displays from sklearn.
<pre>utils.discovery.all functions()</pre>	Get a list of all functions from sklearn.

Utilities from joblib:

```
utils.parallel.delayedDecorator used to capture the arguments of a function.utils.parallel backendDecorator used to capture the arguments of a function.Utils.register parallel backendChange the default backend used by Parallel inside a with block.Register a new Parallel backend factory.
```

 $\underline{\textbf{utils.parallel.Parallel}} \ ([\textbf{n_jobs}, \textbf{backend}, \dots]) \ \ \textbf{Tweak of } \underline{\textbf{joblib.Parallel}} \ \textbf{that propagates the scikit-learn configuration}.$

Recently deprecated