sklearn.linear model.LogisticRegression

class sklearn.linear_model.LogisticRegression(penalty='l2', *, dual=False, tol=0.0001, C=1.0, fit_intercept=True, intercept_scaling=1, class_weight=None, random_state=None, solver='lbfgs', max_iter=100, multi_class='auto', verbose=0, warm_start=False, n_jobs=None, l1_ratio=None)

[source]

Logistic Regression (aka logit, MaxEnt) classifier.

In the multiclass case, the training algorithm uses the one-vs-rest (OvR) scheme if the 'multi_class' option is set to 'ovr', and uses the cross-entropy loss if the 'multi_class' option is set to 'multinomial'. (Currently the 'multinomial' option is supported only by the 'lbfgs', 'sag', 'saga' and 'newton-cg' solvers.)

This class implements regularized logistic regression using the 'liblinear' library, 'newton-cg', 'sag', 'saga' and 'lbfgs' solvers. **Note that regularization is applied by default**. It can handle both dense and sparse input. Use C-ordered arrays or CSR matrices containing 64-bit floats for optimal performance; any other input format will be converted (and copied).

The 'newton-cg', 'sag', and 'lbfgs' solvers support only L2 regularization with primal formulation, or no regularization. The 'liblinear' solver supports both L1 and L2 regularization, with a dual formulation only for the L2 penalty. The Elastic-Net regularization is only supported by the 'saga' solver.

Read more in the **User Guide**.

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Parameters:

penalty: {'l1', 'l2', 'elasticnet', None}, default='l2'

Specify the norm of the penalty:

- None: no penalty is added;
- '12': add a L2 penalty term and it is the default choice;
- '11': add a L1 penalty term;
- 'elasticnet': both L1 and L2 penalty terms are added.

Warning: Some penalties may not work with some solvers. See the parameter solver below, to know the compatibility between the penalty and solver.

New in version 0.19: I1 penalty with SAGA solver (allowing 'multinomial' + L1)

dual: bool, default=False

Dual (constrained) or primal (regularized, see also <u>this equation</u>) formulation. Dual formulation is only implemented for I2 penalty with liblinear solver. Prefer dual=False when n_samples > n_features.

tol: float, default=1e-4

Tolerance for stopping criteria.

C: float, default=1.0

Inverse of regularization strength; must be a positive float. Like in support vector machines, smaller values specify stronger regularization.

fit_intercept : bool, default=True

Specifies if a constant (a.k.a. bias or intercept) should be added to the decision function.

intercept_scaling : float, default=1

Useful only when the solver 'liblinear' is used and self.fit_intercept is set to True. In this case, x becomes [x, self.intercept_scaling], i.e. a "synthetic" feature with constant value equal to intercept_scaling is appended to the instance vector. The intercept becomes intercept_scaling * synthetic_feature_weight.

Note! the synthetic feature weight is subject to I1/I2 regularization as all other features. To lessen the effect of regularization on synthetic feature weight (and therefore on the intercept) intercept_scaling has to be increased.

class_weight: dict or 'balanced', default=None

Weights associated with classes in the form {class_label: weight}. If not given, all classes are supposed to have weight one.

The "balanced" mode uses the values of y to automatically adjust weights inversely proportional to class frequencies in the input data as $n_samples / (n_classes * np.bincount(y))$.

Note that these weights will be multiplied with sample_weight (passed through the fit method) if sample_weight is specified.

New in version 0.17: class_weight='balanced'

random state: int, RandomState instance, default=None

Used when solver == 'sag', 'saga' or 'liblinear' to shuffle the data. See Glossary for details.

solver: {'lbfgs', 'liblinear', 'newton-cg', 'newton-cholesky', 'sag', 'saga'}, default='lbfgs'

Algorithm to use in the optimization problem. Default is 'lbfgs'. To choose a solver, you might want to consider the following aspects:

- For small datasets, 'liblinear' is a good choice, whereas 'sag' and 'saga' are faster for large ones;
- For multiclass problems, only 'newton-cg', 'sag', 'saga' and 'lbfgs' handle multinomial loss;
- 'liblinear' is limited to one-versus-rest schemes.
- 'newton-cholesky' is a good choice for n_samples >> n_features, especially with one-hot encoded categorical features with rare categories. Note that it is limited to binary classification and the one-versus-rest reduction for multiclass classification. Be aware that the memory usage of this solver has a quadratic dependency on n_features because it explicitly computes the Hessian matrix.

Warning: The choice of the algorithm depends on the penalty chosen. Supported penalties by solver:

- 'lbfgs' ['l2', None]
- 'liblinear' ['l1', 'l2']
- 'newton-cg' ['l2', None]
- 'newton-cholesky' ['l2', None]

Toggle Menu - ['l2', None]

' - ['elasticnet', 'l1', 'l2', None]

Note: 'sag' and 'saga' fast convergence is only guaranteed on features with approximately the same scale. You can preprocess the data with a scaler from sklearn.preprocessing.

See also: Refer to the User Guide for more information regarding <u>LogisticRegression</u> and more specifically the <u>Table</u> summarizing solver/penalty supports.

New in version 0.17: Stochastic Average Gradient descent solver.

New in version 0.19: SAGA solver.

Changed in version 0.22: The default solver changed from 'liblinear' to 'lbfgs' in 0.22.

New in version 1.2: newton-cholesky solver.

max_iter : int, default=100

Maximum number of iterations taken for the solvers to converge.

multi_class : {'auto', 'ovr', 'multinomial'}, default='auto'

If the option chosen is 'ovr', then a binary problem is fit for each label. For 'multinomial' the loss minimised is the multinomial loss fit across the entire probability distribution, *even when the data is binary*. 'multinomial' is unavailable when solver='liblinear'. 'auto' selects 'ovr' if the data is binary, or if solver='liblinear', and otherwise selects 'multinomial'.

New in version 0.18: Stochastic Average Gradient descent solver for 'multinomial' case.

Changed in version 0.22: Default changed from 'ovr' to 'auto' in 0.22.

verbose: int, default=0

For the liblinear and lbfgs solvers set verbose to any positive number for verbosity.

warm_start : bool, default=False

When set to True, reuse the solution of the previous call to fit as initialization, otherwise, just erase the previous solution. Useless for liblinear solver. See <u>the Glossary</u>.

New in version 0.17: warm_start to support lbfgs, newton-cg, sag, saga solvers.

n_jobs : int, default=None

Number of CPU cores used when parallelizing over classes if multi_class='ovr'". This parameter is ignored when the solver is set to 'liblinear' regardless of whether 'multi_class' is specified or not. None means 1 unless in a joblib.parallel_backend context. -1 means using all processors. See Glossary for more details.

I1_ratio : float, default=None

The Elastic-Net mixing parameter, with 0 <= l1_ratio <= 1. Only used if penalty='elasticnet'. Setting l1_ratio=0 is equivalent to using penalty='l2', while setting l1_ratio=1 is equivalent to using penalty='l1'. For 0 < l1_ratio <1, the penalty is a combination of L1 and L2.

Attributes:

classes_ : ndarray of shape (n_classes,)

A list of class labels known to the classifier.

coef_: ndarray of shape (1, n_features) or (n_classes, n_features)

Coefficient of the features in the decision function.

coef_ is of shape (1, n_features) when the given problem is binary. In particular, when multi_class='multinomial', coef_ corresponds to outcome 1 (True) and -coef_ corresponds to outcome 0 (False).

intercept_ : ndarray of shape (1,) or (n_classes,)

Intercept (a.k.a. bias) added to the decision function.

If fit_intercept is set to False, the intercept is set to zero. intercept_ is of shape (1,) when the given problem is binary. In particular, when multi_class='multinomial', intercept_ corresponds to outcome 1 (True) and -intercept_ corresponds to outcome 0 (False).

n_features_in_ : int

f features seen during <u>fit</u>.

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.

n_iter_: ndarray of shape (n_classes,) or (1,)

Actual number of iterations for all classes. If binary or multinomial, it returns only 1 element. For liblinear solver, only the maximum number of iteration across all classes is given.

Changed in version 0.20: In SciPy <= 1.0.0 the number of lbfgs iterations may exceed max_iter. n_iter_ will now report at most max_iter.

See also:

SGDClassifier

Incrementally trained logistic regression (when given the parameter loss="log_loss").

LogisticRegressionCV

Logistic regression with built-in cross validation.

Notes

The underlying C implementation uses a random number generator to select features when fitting the model. It is thus not uncommon, to have slightly different results for the same input data. If that happens, try with a smaller tol parameter.

Predict output may not match that of standalone liblinear in certain cases. See differences from liblinear in the narrative documentation.

References

L-BFGS-B - Software for Large-scale Bound-constrained Optimization

Ciyou Zhu, Richard Byrd, Jorge Nocedal and Jose Luis Morales. http://users.iems.northwestern.edu/~nocedal/lbfgsb.html

LIBLINEAR - A Library for Large Linear Classification

https://www.csie.ntu.edu.tw/~cjlin/liblinear/

SAG – Mark Schmidt, Nicolas Le Roux, and Francis Bach

Minimizing Finite Sums with the Stochastic Average Gradient https://hal.inria.fr/hal-00860051/document

SAGA - Defazio, A., Bach F. & Lacoste-Julien S. (2014).

"SAGA: A Fast Incremental Gradient Method With Support for Non-Strongly Convex Composite Objectives"

Hsiang-Fu Yu, Fang-Lan Huang, Chih-Jen Lin (2011). Dual coordinate descent

methods for logistic regression and maximum entropy models. Machine Learning 85(1-2):41-75. https://www.csie.ntu.edu.tw/~cjlin/papers/maxent_dual.pdf

Examples

Methods

<pre>decision function(X)</pre>	Predict confidence scores for samples.
<pre>densify()</pre>	Convert coefficient matrix to dense array format.
<pre>fit(X, y[, sample_weight])</pre>	Fit the model according to the given training data.
<pre>get_metadata_routing()</pre>	Get metadata routing of this object.
<pre>get params([deep])</pre>	Get parameters for this estimator.
<pre>predict(X)</pre>	Predict class labels for samples in X.
Toggle Menu <u>proba(</u> X)	Predict logarithm of probability estimates.

<u>predict_proba(</u> X)	Probability estimates.
<pre>score(X, y[, sample_weight])</pre>	Return the mean accuracy on the given test data and labels.
<pre>set_fit_request(*[, sample_weight])</pre>	Request metadata passed to the fit method.
set_params(**params)	Set the parameters of this estimator.
<pre>set_score_request(*[, sample_weight])</pre>	Request metadata passed to the score method.
<pre>sparsify()</pre>	Convert coefficient matrix to sparse format.

decision_function(X)

Predict confidence scores for samples.

The confidence score for a sample is proportional to the signed distance of that sample to the hyperplane.

Parameters:

X: {array-like, sparse matrix} of shape (n_samples, n_features)

The data matrix for which we want to get the confidence scores.

Returns:

scores: ndarray of shape (n_samples,) or (n_samples, n_classes)

Confidence scores per (n_samples, n_classes) combination. In the binary case, confidence score for self.classes_[1] where >0 means this class would be predicted.

densify() [source]

Convert coefficient matrix to dense array format.

Converts the coef_ member (back) to a numpy.ndarray. This is the default format of coef_ and is required for fitting, so calling this method is only required on models that have previously been sparsified; otherwise, it is a no-op.

Returns:

self

Fitted estimator.

fit(X, y, sample_weight=None)

[source]

Fit the model according to the given training data.

Parameters:

X: {array-like, sparse matrix} of shape (n_samples, n_features)

Training vector, where n_samples is the number of samples and n_features is the number of features.

y: array-like of shape (n_samples,)

Target vector relative to X.

sample_weight : array-like of shape (n_samples,) default=None

Array of weights that are assigned to individual samples. If not provided, then each sample is given unit weight.

New in version 0.17: sample_weight support to LogisticRegression.

Returns:

self

Fitted estimator.

Notes

The SAGA solver supports both float64 and float32 bit arrays.

get_metadata_routing()

[source]

Get metadata routing of this object.

Please check <u>User Guide</u> on how the routing mechanism works.

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Returns:

routing: MetadataRequest

A <u>MetadataRequest</u> encapsulating routing information.

get_params(deep=True)

Get parameters for this estimator.

Parameters:

deep: bool, default=True

If True, will return the parameters for this estimator and contained subobjects that are estimators.

Returns:

params: dict

Parameter names mapped to their values.

predict(X)

Predict class labels for samples in X.

Parameters:

X: {array-like, sparse matrix} of shape (n_samples, n_features)

The data matrix for which we want to get the predictions.

Returns:

y_pred : ndarray of shape (n_samples,)

Vector containing the class labels for each sample.

predict_log_proba(X)

Predict logarithm of probability estimates.

The returned estimates for all classes are ordered by the label of classes.

Parameters:

X : array-like of shape (n_samples, n_features)

Vector to be scored, where n_samples is the number of samples and n_features is the number of features.

Returns:

T: array-like of shape (n_samples, n_classes)

Returns the log-probability of the sample for each class in the model, where classes are ordered as they are in self.classes_.

predict_proba(X)

Probability estimates.

The returned estimates for all classes are ordered by the label of classes.

For a multi_class problem, if multi_class is set to be "multinomial" the softmax function is used to find the predicted probability of each class. Else use a one-vs-rest approach, i.e. calculate the probability of each class assuming it to be positive using the logistic function. and normalize these values across all the classes.

Parameters:

X : array-like of shape (n_samples, n_features)

Vector to be scored, where n_samples is the number of samples and n_features is the number of features.

Returns:

T: array-like of shape (n_samples, n_classes)

Returns the probability of the sample for each class in the model, where classes are ordered as they are in self.classes_.

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score(X, y, sample_weight=None)

[source]

Return the mean accuracy on the given test data and labels.

In multi-label classification, this is the subset accuracy which is a harsh metric since you require for each sample that each label set be correctly predicted.

Parameters:

X : array-like of shape (n_samples, n_features)

Test samples.

y: array-like of shape (n_samples,) or (n_samples, n_outputs)

True labels for x.

sample_weight : array-like of shape (n_samples,), default=None

Sample weights.

Returns:

score: float

Mean accuracy of self.predict(X) w.r.t. y.

set_fit_request(*, sample_weight: bool | None | str = '\$UNCHANGED\$') → LogisticRegression

[source]

Request metadata passed to the fit method.

Note that this method is only relevant if enable_metadata_routing=True (see sklearn.set_config). Please see User Guide on how the routing mechanism works.

The options for each parameter are:

- True: metadata is requested, and passed to fit if provided. The request is ignored if metadata is not provided.
- False: metadata is not requested and the meta-estimator will not pass it to fit.
- None: metadata is not requested, and the meta-estimator will raise an error if the user provides it.
- str: metadata should be passed to the meta-estimator with this given alias instead of the original name.

The default (sklearn.utils.metadata_routing.UNCHANGED) retains the existing request. This allows you to change the request for some parameters and not others.

New in version 1.3.

Note: This method is only relevant if this estimator is used as a sub-estimator of a meta-estimator, e.g. used inside a <u>Pipeline</u>. Otherwise it has no effect.

Parameters:

sample_weight : str, True, False, or None, default=sklearn.utils.metadata_routing.UNCHANGED

Metadata routing for sample_weight parameter in fit.

Returns:

self : object

The updated object.

set_params(**params)

[source]

Set the parameters of this estimator.

The method works on simple estimators as well as on nested objects (such as Pipeline). The latter have parameters of the form component> component compon

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Parameters:

**params: dict

Estimator parameters.

Returns:

self: *estimator instance*Estimator instance.

set_score_request(*, sample_weight: bool | None | str = '\$UNCHANGED\$') → LogisticRegression

[source]

Request metadata passed to the score method.

Note that this method is only relevant if enable_metadata_routing=True (see sklearn.set_config). Please see User Guide on how the routing mechanism works.

The options for each parameter are:

- True: metadata is requested, and passed to score if provided. The request is ignored if metadata is not provided.
- False: metadata is not requested and the meta-estimator will not pass it to score.
- None: metadata is not requested, and the meta-estimator will raise an error if the user provides it.
- str: metadata should be passed to the meta-estimator with this given alias instead of the original name.

The default (sklearn.utils.metadata_routing.UNCHANGED) retains the existing request. This allows you to change the request for some parameters and not others.

New in version 1.3.

Note: This method is only relevant if this estimator is used as a sub-estimator of a meta-estimator, e.g. used inside a Pipeline. Otherwise it has no effect.

Parameters:

sample_weight : str, True, False, or None, default=sklearn.utils.metadata_routing.UNCHANGED

Metadata routing for sample_weight parameter in score.

Returns:

self : object

The updated object.

sparsify()

Convert coefficient matrix to sparse format.

Converts the coef_ member to a scipy.sparse matrix, which for L1-regularized models can be much more memory- and storage-efficient than the usual numpy.ndarray representation.

The intercept_ member is not converted.

Returns:

self

Fitted estimator.

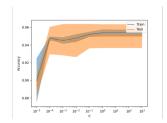
Notes

For non-sparse models, i.e. when there are not many zeros in coef_, this may actually *increase* memory usage, so use this method with care. A rule of thumb is that the number of zero elements, which can be computed with (coef_ == 0).sum(), must be more than 50% for this to provide significant benefits.

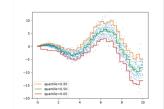
After calling this method, further fitting with the partial_fit method (if any) will not work until you call densify.

Examples using sklearn.linear_model.LogisticRegression

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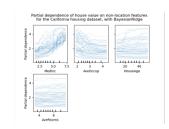
Release Highlights for scikit-learn 1.3



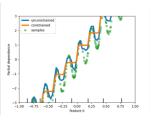
Release Highlights for scikit-learn 1.1



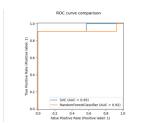
Release Highlights for scikit-learn 1.0



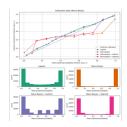
Release Highlights for scikit-learn 0.24



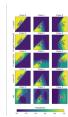
Release Highlights for scikit-learn 0.23



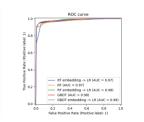
Release Highlights for scikit-learn 0.22



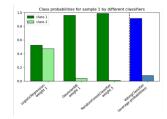
Probability Calibration curves



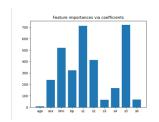
Plot classification probability



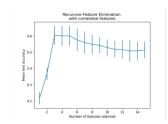
Feature transformations with ensembles of trees



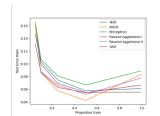
Plot class probabilities calculated by the VotingClassifier



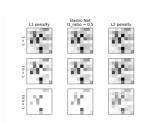
Model-based and sequential feature selection



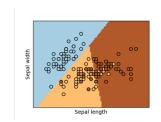
Recursive feature elimination with cross-validation



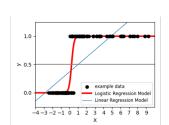
Comparing various online solvers



L1 Penalty and Sparsity in Logistic Regression



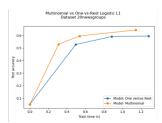
Logistic Regression 3class Classifier



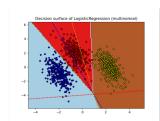
Logistic function



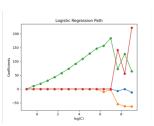
MNIST classification using multinomial logistic + L1



Multiclass sparse logistic regression on 20newgroups



Plot multinomial and One-vs-Rest Logistic Regression



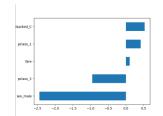
Regularization path of L1-Logistic Regression



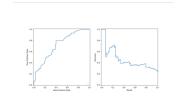
Displaying Pipelines



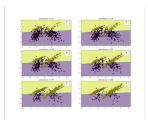
Displaying estimators and complex pipelines



Introducing the set_output API

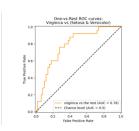


Visualizations with Display Objects



Class Likelihood Ratios to measure classification performance

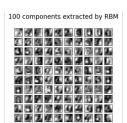
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Multiclass Receiver Operating Characteristic (ROC)



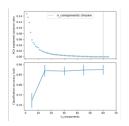
Multilabel classification using a classifier chain



Restricted Boltzmann Machine features for digit classification



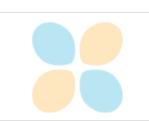
Column Transformer with Mixed Types



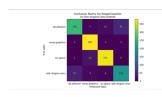
Pipelining: chaining a PCA and a logistic regression



Feature discretization



Digits Classification Exercise



Classification of text documents using sparse features

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