Arguments

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numeric matrix of data, or an object that can be coerced to such a matrix (such as a numeric vector or a data frame with all numeric columns).

centers

either the number of clusters, say k, or a set of initial (distinct) cluster centres. If a number, a random set of (distinct) rows in x is chosen as the initial centres.

the maximum number of iterations allowed. iter.max

if centers is a number, how many random sets should be chosen? nstart

algorithm character: may be abbreviated. Note that "Lloyd" and "Forgy" are alternative names for one algorithm.

an \mathbb{R} object of class "kmeans", typically the result ob of ob <- kmeans (..). object

character: may be abbreviated. "centers" causes fitted to return cluster centers (one for each method input point) and "classes" causes fitted to return a vector of class assignments.

logical or integer number, currently only used in the default method ("Hartigan-Wong"): if positive trace (or true), tracing information on the progress of the algorithm is produced. Higher values may produce more tracing information.

not used.

sklearn.cluster.KMeans

class $sklearn.cluster.KMeans(n_clusters=8, *, init='k-means++', n_init=10, max_iter=300, tol=0.0001, verbose=0, random_state=None, copy_x=True, algorithm='lloyd') [source]$

K-Means clustering.

Read more in the User Guide.

Parameters:

n_clusters : int, default=8

The number of clusters to form as well as the number of centroids to generate.

init : {'k-means++', 'random'}, callable or array-like of shape (n_clusters, n_features), default='k-means++'

Method for initialization:

'k-means++': selects initial cluster centers for k-mean clustering in a smart way to speed up convergence. See section Notes in k_init for more details.

'random': choose n_clusters observations (rows) at random from data for the initial centroids.

If an array is passed, it should be of shape (n_clusters, n_features) and gives the initial centers.