Package ‘LPCM’

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Author Jochen Einbeck and Ludger Evers
Maintainer Jochen Einbeck <jochen.einbeck@durham.ac.uk>
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**LPCM-package**

Fitting multivariate data patterns with local principal curves; including simple tools for data compression (projection), bandwidth selection, and measuring goodness-of-fit.

This package implements the techniques introduced in Einbeck, Tutz & Evers (2005), and successive related papers.

The main functions to be called by the user are

- **lpc**, for the estimation of the local centers of mass which make up the principal curve;
- **lpc.spline**, which is a smooth and fully parametrized cubic spline representation of the latter;
- **lpc.project**, which enables to compress data by projecting them orthogonally onto the curve;
- **lpc.coverage** and **Rc** for assessing goodness-of-fit;
- **lpc.self.coverage** for bandwidth selection;
- the generic **plot** and **print** methods for objects of class **lpc** and **lpc.spline**.

This package also contains some code for density mode detection (‘local principal points’) and mean shift clustering (as well as bandwidth selection in this context), which implements the methods presented in Einbeck (2011). See the help file for **ms**.

A second R package which will implement the extension of local principal curves to local principal surfaces and manifolds, as proposed in Einbeck, Evers & Powell (2010), is in preparation.

**Details**

- **Package:** LPCM
- **Type:** Package
- **License:** GPL (>=2)
- **LazyLoad:** yes
Acknowledgements
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Author(s)
Jochen Einbeck and Ludger Evers
Maintainer: Jochen Einbeck <jochen.einbeck@durham.ac.uk>

References

See Also
pcurve, princurve

calspeedflow

Speed-flow data from California.

Description
A ‘fundamental diagram’ with observations of speed and flow recorded from 9th of July 2007, 9am, to 10th of July 2007, 10pm, on Line 5 of the Californian Freeway SR57-N, VDS number 1202263. The data were originally measured in intervals of thirty seconds, and then aggregated over intervals of 5 minutes length.

Usage
data(calspeedflow)

Format
A data frame with 444 observations on the following 4 variables.

Date a factor with levels 07/09/2007... 07/10/2007.
Timestamp a factor with a timestamps in intervals of five minutes.
Lane5Flow a numeric vector of vehicle flow in vehicles per 5 minutes.
Lane5Speed a numeric vector of vehicle speed in miles per hour.
Source
Retrieved from PeMS.

References

Examples

```r
data(calspeedflow)
plot(calspeedflow[,3:4])
```

`coverage`  Coverage and self-coverage plots.

Description
These functions compute coverages (for any principal object), and self-coverages (only for local principal curves, these may be used for bandwidth selection).

Usage

```r
coverage.raw(X, vec, tau, weights=1, plot.type="p", print=FALSE,
label=NULL,...)
```

```r
coverage(X, vec, taumin=0.02, taumax, gridsize=25, weights=1,
plot.type="o", print=FALSE,...)
```

```r
lpc.coverage(object, taumin=0.02, taumax, gridsize=25, quick=TRUE,
plot.type="o", print=FALSE,...)
```

```r
lpc.self.coverage(X, taumin=0.02, taumax=0.5, gridsize=25, x0=1,
way = "two", scaled=TRUE, weights=1, pen=2, depth=1,
control=lpc.control(boundary=0, cross=FALSE), quick=TRUE,
plot.type="o", print=FALSE,...)
```

```r
select.self.coverage(self, smin, plot.type="o", plot.segments=NULL)
```

Arguments

- `X` a $N \times d$ data matrix.
- `object` An object of type `lpc` or `lpc.spline`.
- `vec` A matrix with $d$ columns. The rows contain the points which make up the fitted object.
- `tau` tube size.
coverage

- **taumin**: Minimal tube size.
- **taumax**: Maximal tube size.
- **weights**: An optional vector of weights. If weights are specified, then the coverage is the weighted mean of the indicator functions for falling within the tube. The function `lpc::coverage` does not have a `weights` argument, as it extracts the weights from the `$weights` component of the fitted object.
- **label**: Experimental option; don’t use.
- **gridsize**: The number of different tube sizes to consider.
- **quick**: If TRUE, an approximate coverage curve is provided by computing distances between data points and the curve through the closest local centers or mass; whereas with FALSE we use the distances of the points when projected orthogonally onto the spline representation of the local principal curve. The latter takes considerably more computing time. The resulting coverage curves are generally very similar, but the quick version may deliver little spurious peaks occasionally.
- **self**: An object of class `self`, or a matrix with two columns providing a self-coverage curve.
- **smin**: Minimum coverage for bandwidth selection. Default: 1/3 for clustering, 2/3 for principal curves.
- **plot.type**: If set to 0, no plotted output is given. Otherwise, an appropriate plot is provided, using the plotting type as specified.
- **plot.segments**: A list with default `lty=c(1,2,3), lwd=c(2,1,1), lcol=c(3,3,3))` which specifies how (and how many) bandwidth candidates, in order of decreasing negative second derivative of self-coverage, are to be highlighted.
- **print**: If TRUE, coverage values are printed on the screen as soon as computed. This is quite helpful especially if `gridsize` is large.
- **x0, way, scaled, pen, depth, control**: LPC parameters as outlined in `lpc` and `lpc.control`.

Details

The function `coverage.raw` computes the coverage, i.e. the proportion of data points lying inside a circle or band with radius \( \tau \), for a fixed value `tau`. The whole coverage curve \( C(\tau) \) is constructed through function `coverage`.

Functions `coverage.raw` and `coverage` can be used for any object fitted by an unsupervised learning technique (for instance, HS principal curves, or even clustering algorithms), while the functions prefixing with `lpc` can only be used for local principal curves. The function `lpc::coverage` is a wrapper around `coverage` which takes directly a fitted `lpc` object, rather than a data matrix.

Function `select.self.coverage` extracts suitable bandwidths from the self-coverage curve, and produces a plot. The function is called from within `lpc::self.coverage`, but can also be called directly by the user (for instance, if the graphical output is to be reproduced, or if the minimum coverage `smin` is to be modified). The component `$select` contains the selected candidate bandwidths, in the order of strength of evidence provided by the self-coverage criterion (the best bandwidth comes first, etc.). A plot is produced as a by-product, which symbolizes the best bandwidth by a thick solid line.
line, the second-best by a dashed line, and the third-best by a dotted line. It is recommended to run
the self-coverage functions with fixed starting points, as in the examples below.

See Einbeck (2011) for details. Note that the original publication by Einbeck, Tutz, and Evers
(2005) uses ‘quick’ coverage curves.

Value

A list of items, and a plot (unless plot.type=0).

The function lpc.self.coverage produces an object of class self. The component $select
recommends suitable bandwidths for the use in lpc, in the order of strength of evidence. These
correspond to points of strong negative curvature (implemented via second differences) of the self-
coverage curve.

Author(s)

J. Einbeck

References

301-313.


See Also

lpc

Examples

data(gvessel)
gvessel.self <- lpc.self.coverage(gvessel[,c(2,4,5)], x0=c(35, 1870, 6.3), print=FALSE, plot.type=0)
h <- select.self.coverage(gvessel.self)$select
gvessel.lpc <- lpc(gvessel[,c(2,4,5)], h=h[1], x0=c(35, 1870, 6.3))
lpc.coverage(gvessel.lpc, gridsize=10, print=FALSE)

data(calspeedflow)
fitms <- ms(calspeedflow[,3:4])
coverage(fitms$data, fitms$cluster.center)
**followx**

*Fit an individual branch of a local principal curve.*

**Description**

Internal function of package **LPCM** called by lpc. Do not use!

**Usage**

```r
followx(Xi, x0, h, t0, iter, way, weights, pen = 2, phi = 1,
    lasteigenvector = 0, rho0 = 0.4, boundary=0.005,
    convergence.at= 0.000001, cross=FALSE)
```

**Arguments**

- `Xi`
- `x0`
- `h`
- `t0`
- `iter`
- `way`
- `weights`
- `pen`
- `phi`
- `lasteigenvector`
- `rho0`
- `boundary`
- `convergence.at`
- `cross`

**Author(s)**

JE

**See Also**

lpc
Description

(Simulated) spectral decomposition of stellar objects, generated in the framework of the Gaia project.

Usage

data(gaia)

Format

A data frame with 8286 observations on the following 22 variables.

ID  ID of the object
metallicity  metallicity (abundance); that is proportion of matter other than hydrogen and helium relative to that of the sun.
gravity  the surface gravity; that is acceleration due to gravity at the surface of the star.
temperature  the ‘effective’ temperature (K); that is the temperature of the observable part of the stellar atmosphere.
band1  photon counts in band 1
band2  photon counts in band 2
band3  photon counts in band 3
band4  photon counts in band 4
band5  photon counts in band 5
band6  photon counts in band 6
band7  photon counts in band 7
band8  photon counts in band 8
band9  photon counts in band 9
band10  photon counts in band 10
band11  photon counts in band 11
band12  photon counts in band 12
band13  photon counts in band 13
band14  photon counts in band 14
band15  photon counts in band 15
band16  photon counts in band 16
Details

Gaia is an astrophysics mission of the European Space Agency (ESA) which will undertake a detailed survey of over $10^9$ stars in our Galaxy and extragalactic objects. An important part of the scientific analysis of these data is the classification of all the objects as well as the estimation of stellar astrophysical parameters (effective stellar temperature, surface gravity, metallicity). This will be done on the basis of high-dimensional spectroscopic and astrometric data such as those ones given here.

More precisely, the spectral data come in form of photon counts ("fluxes") observed in (originally) 96 wavelength intervals ("bands"), see Bailer-Jones (2010) for more details. The data given here are a 16-dimensional subset created by binning/selecting from the 96 bands. The counts given here are standardized, i.e. they are divided by the total number of incoming photons over all filters (in other words, they add up to 1). Note that these data are simulated using computer models. The satellite which will collect the actual data will be launched in 2012.

The 16-d spectral data have been used in Einbeck, Evers and Bailer-Jones (2008) as well as Einbeck, Evers and Powell (2010) in order to predict the stellar temperature.

Source

Coryn Bailer-Jones (MPIA Heidelberg).

References


Examples

data(gaia)
s <- sample(nrow(gaia),200)
library(lattice)
splom(gaia[s,5:20], cex=0.3,pscales=0)

gaia.pc <- princomp(gaia[s,5:20])
temp <- gaia$temperature
tempcol <- (temp[s]- min(temp[s]))/max(temp[s]- min(temp[s]))
library(scatterplot3d)
scatterplot3d(gaia.pc$scores[,c(2,1,3)], pch="+",
color=rgb(sqrt(tempcol),0,1-sqrt(tempcol)))

# This is a 3D scatterplot of the first three principal component scores;
# with higher stellar temperatures shaded in red colour.
**gvessel**  

*North Atlantic Water Temperature Data.*

**Description**

These are observations taken over nine days in May 2000 by the German vessel Gauss in the North Atlantic.

**Usage**

```r
data(gvessel)
```

**Format**

A data frame with 643 observations on the following 7 variables.

- `dayRg` an integer for the day at which the measurement was taken.
- `salg` a numeric vector with measurements of salinity according to the PSS (Practical Salinity Scale).
- `tempg` a numeric vector with measurements of water temperature in degrees Celsius.
- `depthg` a numeric vector with the water depths (in meters) at which the measurements were taken.
- `oxyg` a numeric vector with measurements of oxygen content (mm per litre of water).
- `longg` longitude
- `latg` latitude

**Source**


**References**


**Examples**

```r
data(gvessel)
pairs(gvessel[,c(3,2,4,5)])
tcol <- (gvessel$tempg- min(gvessel$tempg))/(max(gvessel$tempg)- min(gvessel$tempg))
require(scatterplot3d)
scatterplot3d(gvessel[,2],gvessel[,4],gvessel[,5], color=rgb(tcol,0,1-tcol))
```
Auxiliary kernel and distance functions.

Description

Internal \texttt{LPCM} functions which are normally not to be called by the user.

Usage

\begin{verbatim}
  kern(y, x = 0, h = 1)
  kernd(X, x, h)
  kdex(X, x, h)
  distancevector(X, y, d = "euclid", na.rm = TRUE)
  vecdist(X,Y)
  mindist(X,y)
  enorm(x)
\end{verbatim}

Arguments

- \texttt{x} a number or vector.
- \texttt{y} a vector.
- \texttt{h} a bandwidth.
- \texttt{X} a matrix.
- \texttt{Y} a matrix.
- \texttt{d} type of distance measure (only ‘euclid’).
- \texttt{na.rm} ...

Details

\texttt{kern} specifies the base kernel (by default Gaussian) used in \texttt{lpc}; \texttt{kernd} is the corresponding multivariate product kernel. \texttt{kdex} is a pointwise multivariate kernel density estimator.

\texttt{distancevector} makes use of function \texttt{vdisseuclid} from \texttt{R} package \texttt{hopach} (but that package does not need to be loaded).

Author(s)

JE

References

Description

This is the main function which computes the actual local principal curve, i.e. a sequence of local centers of mass.

Usage

\[
lpc(X, \ h, \ t\theta = \text{mean}(h), \ x\theta, \ \text{way = "two"}, \ \text{scaled = TRUE}, \\
\ \text{weights=1}, \ \text{pen = 2}, \ \text{depth = 1}, \ \text{control=lpc.control()})
\]

Arguments

- \(X\) data matrix with \(N\) rows (observations) and \(d\) columns (variables).
- \(h\) bandwidth. May be either specified as a single number, then the same bandwidth is used in all dimensions, or as a \(d\)-dimensional bandwidth vector. The default setting is 10 percent of the range in each direction. If \(\text{scaled = TRUE}\) then the bandwidth has to be specified in fractions of the data range, e.g. \(h= c(0.2,0.1)\), rather than absolute values.
- \(t\theta\) scalar step length. Default setting is \(t\theta=h\), if \(h\) is a scalar, and \(t\theta=\text{mean}(h)\), if \(h\) is a vector.
- \(x\theta\) specifies the choice of starting points. The default choice \(x\theta=1\) will select one suitable starting point automatically (in form of a local density mode). The second built-in option \(x\theta=0\) will use all local density modes as starting points, hence produce as many branches as modes. Optionally, one can also set one or more starting points manually here. This can be done in form of a matrix, where each row corresponds to a starting point, or in form of a vector, where starting points are read in consecutive order from the entries of the vector. The starting point has always to be specified on the original data scale, even if \(\text{scaled = TRUE}\). A fixed number of starting points can be enforced through option \(\text{mult}\) in \(\text{lpc.control}\).
- \(\text{way}\) "one": go only in direction of the first local eigenvector, "back": go only in opposite direction, "two": go from starting point in both directions.
- \(\text{scaled}\) if \(\text{TRUE}\), scales each variable by dividing through its range (see also the Notes section below).
- \(\text{weights}\) a vector of observation weights (can also be used to exclude individual observations from the computation by setting their weight to zero.)
- \(\text{pen}\) power used for angle penalization (see [1]). If set to 0, the angle penalization is switched off.
- \(\text{depth}\) maximum depth of branches (\(\phi_{max}\) in [2]), restricted to the values 1,2 or 3 (The original LPC branch has depth 1. If, along this curve, a point features a high second local PC, this launches a new starting point, and the resulting branch has depth 2. If, along this branch, a point features a high second local PC, this launches a new starting point, and the resulting branch has depth 3.)
lpc

control Additional parameters steering particularly the starting-, boundary-, and convergence behavior of the fitted curve. See lpc.control.

Value
A list of items:

LPC The coordinates of the local centers of mass of the fitted principal curve.
Parametrization Curve parameters and branch labels for each local center of mass.
h The bandwidth used for the curve estimation.
to The constant $t_0$ used for the curve estimation.
starting.points The coordinates of the starting point(s) used.
data The data frame used for curve estimation.
scaled Logical.
weights The vector of weights used for curve estimation.
control The settings used in lpc.control()
Misc Miscellanea.

Note
All values provided in the output refer to the scaled data, if scaled=TRUE. Use unscale to convert the results back to the original data scale.

The option scaled=TRUE scales the data by dividing each variable through their range. This differs from the usual scaling through the standard deviation as common for PCA, but we found the algorithm and the default bandwidth selection to work more reliably this way. If you wish to scale by the standard deviation, please do that by feeding the scaled data directly into the lpc function, i.e. lpc(sweep(data, 2, sd(data), "/"), h, t0, ..., scaled=FALSE, ...).

Author(s)
J. Einbeck and L. Evers. See LPCM-package for further acknowledgements.

References
Examples

```r
data(calspeedflow)
lpc1 <- lpc(calspeedflow[,3:4])
plot(lpc1)

data(mussels, package="dr")
lpc2 <- lpc(mussels[,3], x0=as.numeric(mussels[49,4]), scaled=FALSE)
plot(lpc2, curvecol=2)

data(gaia)
s <- sample(nrow(gaia), 200)
gaia.pc <- princomp(gaia[s,5:20])
lpc3 <- lpc(gaia.pc$scores[,c(2,1,3)], scaled=FALSE)
plot(lpc3, curvecol=2, type=c("curve","mass"))

# Simulated letter 'E' with branched LPC
ex <- c(rep(0,40), seq(0,1,length=20), seq(0,1,length=20), seq(0,1,length=20))
ey <- c(seq(0,2,length=40), rep(0,20), rep(1,20), rep(2,20))
sex <- rnorm(100,0,0.01); sey <- rnorm(100,0,0.01)
exy <- sex+ex; eyy <- sey+ey
ex2 <- ex+ex; ey2 <- ey+ey
e1 <- cbind(ex,e1); e2 <- cbind(ex2,ey2)
lpc.e1 <- lpc(e1, h=c(0.1,0.1), depth=2, scaled=FALSE)
plot(lpc.e1, type=c("curve","mass", "start"))
```

---

### lpc.control

**Auxiliary parameters for controlling local principal curves.**

#### Description

This function bundles parameters controlling mainly the starting-, convergence-, boundary-, and stopping-behaviour of the local principal curve. It will be used only inside the `lpc()` function argument.

#### Usage

```r
lpc.control(iter=100, cross=TRUE,
            boundary = 0.005, convergence.at = 0.00001,
            mult=NA, ms.h=NA, ms.sub=30,
            pruning.thresh=0.0, rho0=0.4)
```

#### Arguments

- **iter**
  - Maximum number of iterations on either side of the starting point within each branch.

- **cross**
  - Logical parameter. If TRUE, curves are stopped when they come too close to an existing branch. Used in the self-coverage function.
boundary
This boundary correction [2] reduces the bandwidth adaptively once the relative
difference of parameter values between two centers of mass falls below the given
threshold. This measure delays convergence and enables the curve to proceed
further into the end points. If set to 0, this boundary correction is switched off.

convergence.at
This forces the curve to stop if the relative difference of parameter values be-
tween two centers of mass falls below the given threshold. If set to 0, then the
curve will always stop after exactly iter iterations.

mult
numerical value which enforces a fixed number of starting points. If the num-
ber given here is larger than the number of starting points provided at x0, then
the missing points will be set at random (For example, if d = 2, mult=3, and
x0=c(58.5, 17.8, 80, 20), then one gets the starting points (58.5, 17.8),
(80,20), and a randomly chosen third one. Another example for such a situa-
tion is x0=NULL with mult=1, in which one random starting point is chosen). If
the number given here is smaller the number of starting points provided at x0,
then only the first mult starting points will be used.

ms.h
sets the bandwidth (vector) for the initial mean shift procedure which finds the
local density modes, and, hence, the starting points for the LPC. If unspecified,
the bandwidth h used in function lpc is used here too.

ms.sub
proportion of data points (default=30) which are used to initialize mean shift
trajectories for the mode finding. In fact, we use
\[ \min(\max(ms.sub, \floor(ms.sub*\text{N}/100)), 10*ms.sub) \]
trajectories.

pruning.thresh
Prunes branches corresponding to higher-depth starting points if their density
estimate falls below this threshold. Typically, a value between 0.0 and 1.0. The
setting 0.0 means no pruning.

rho0
A numerical value which steers the birth process of higher-depth starting points.
Usually, between 0.3 and 0.4 (see reference [1]).

Value
A list of the nine specified input parameters, which can be read by the control argument of the
lpc function.

Author(s)
JE

References
Springer, Heidelberg, pages 256-263.

Examples

```r
data(calspeedflow)
fit1 <- lpc(calspeedflow[,c(3,4)], x0=c(50,60), scaled=TRUE,
            control=lpc.control(iter=20, boundary=0))
plot(fit1, type=c("curve","start","mass"))
```

---

**Description**

Projects a new observation onto the spline representation of the local principal curve.

**Usage**

```r
lpc.project(object, newdata, ...)
```

**Arguments**

- `object`: Object of class `lpc` or `lpc.spline`.
- `newdata`: A data frame containing the new data to be projected.
- `...`: Additional arguments to be passed to `lpc.project.spline`.

**Value**

- `closest.pi`: Projection index of projected point(s) (in cubic spline parametrization).
- `closest.or.pi`: Projection index of projected point(s) (in terms of the original LPC parametrization).
- `closest.coords`: Coordinates of projected data point(s).
- `closest.dist`: Euclidean distance between data point(s) and their projected counterpart(s).
- `closest.branch`: ID of branch onto which the data point was projected (the IDs get allocated in the output component `$Parametrization` of function `lpc`).

**Note**

The parametrization of the cubic spline function is not exactly the same as that of the original LPC. The reason is that the latter uses Euclidean distances between centers of masses, while the former uses the arc length along the cubic spline. The differences are normally quite small, though.

**Author(s)**

J. Einbeck and L. Evers
References


See Also

lpc, lpc.spline

Examples

data(gvessel)
gvessel.lpc <- lpc(gvessel[,c(2,4,5)], scaled=TRUE, h=0.11, x0=c(35, 1870, 6.3))
lpc.project(gvessel.lpc, newdata=data.frame(salg=35, dephtg=2000, oxyg=6))

lpc.spline

Representing local principal curves through a cubic spline.

Description

Fis a natural cubic spline component-wise through the series of local centers of mass. This provides a continuous parametrization in terms of arc length distance, which can be used to compute a projection index for the original or new data points.

Usage

lpc.spline(lpcobject, optimize = TRUE, compute.Rc=FALSE, project=FALSE, ...)

Arguments

lpcobject Object of class lpc.
optimize Boolean. If TRUE, optimize is used to find the point on the curve with minimum distance. Otherwise, data points are only projected onto the closest knot.
compute.Rc Boolean. If TRUE, the goodness-of-fit measure $R_C$ suggested in [1] is computed and returned (using the scaled data, if scaled=TRUE in lpcobject).
project Boolean. If TRUE, projections onto curve are computed.
... Additional arguments to be passed to lpc.project.spline

Details

See reference [2].
Value

- **knots.pi**
  LPC parameters (in cubic spline parametrization) at position of the knots of the spline function (these are not identical to the LPC mass points!)

- **knots.coords**
  Coordinates of the spline knots.

- **closest.pi**
  Parameter of the projected data points.

- **closest.coords**
  Coordinates of projected data points.

- **closest.dist**
  Euclidean distance between original and projected data point.

- **closest.branch**
  ID Number of the branch on which the data point was projected (the IDs are given in the output of function `lpc`).

- **Rc**
  Value of $R_C$.

- **project**
  repeats the input value of project.

- **lpcobject**
  returns the provided object `lpcobject`.

- **splinefun**
  returns the cubic spline function (generated by `lpc.splinefun`).

Warning

Careful with options `project` and `compute.Rc` - they can take rather long if the data set is large!

Note

The parametrization of the cubic spline function is not exactly the same as that of the original LPC. The reason is that the latter uses Euclidean distances between centers of masses, while the former uses the arc length along the cubic spline. However, the differences are normally quite small.

Author(s)

J. Einbeck and L. Evers

References


See Also

- `lpc`

Examples

```r
data(gvessel)
gvessel.lpc <- lpc(gvessel[,c(2,4,5)], h=0.11, x0=c(35, 1870, 6.3))
gvessel.spline <- lpc.spline(gvessel.lpc)
plot(gvessel.spline, lwd=2)
```
Auxiliary functions for spline fitting and projection.

Description

Internal functions of package \texttt{LPCM} called by \texttt{lpc.spline} and others. These will rarely be called directly by the user.

Usage

\begin{verbatim}
\texttt{lpc.splinefun(lpcobject)}
\texttt{lpc.fit.spline(lpcsl, num.knots = 100)}
\texttt{lpc.spline.eval(lpcsl, or.pi, branch = 0)}
\texttt{lpc.project.spline(lpcsl, newdata, num.knots = 100, optimize = TRUE)}
\texttt{lpc.curve.length(lpcsl, or.pi, branch = 0, total.subdivisions = 10000, min.subdivisions = 100)}
\end{verbatim}

Arguments

- \texttt{lpcobject} Object of type \texttt{lpc}.
- \texttt{lpcsl} Object generated by \texttt{lpc.splinefun}.
- \texttt{num.knots} number of spline knots.
- \texttt{or.pi} original projection index.
- \texttt{branch} branch ID.
- \texttt{newdata} new data frame.
- \texttt{optimize} Boolean.
- \texttt{total.subdivisions} total number of subdivisions for arc length computation.
- \texttt{min.subdivisions} minimum number of subdivisions for arc length computation.

Author(s)

L. Evers and J. Einbeck

See Also

\texttt{lpc.spline}
Mean shift clustering.

Description

Functions for mean shift, iterative mean shift, mean shift clustering, and bandwidth selection for mean shift clustering (based on self-coverage). The main function is `ms` which, for a given bandwidth, detects the local modes ('local principal points') and performs the clustering.

These functions implement the techniques presented in Einbeck (2011).

Usage

```r
meanshift(X, x, h)
ms.rep(X, x, h, plotms=1, thresh= 0.00000001, iter=200)
ms(X, h, subset, thr=0.0001, scaled= TRUE, iter=200, plotms=2,
   or.labels=NULL, ...)
ms.self.coverage(X, taumin=0.02, taumax=0.5, gridsize=25,
   thr=0.0001, scaled=TRUE, cluster=FALSE, plot.type="o",
   or.labels=NULL, print=FALSE, ...)
```

Arguments

- `X` data matrix.
- `h` bandwidth (by default, 10 percent of the data range).
- `x` point from which we wish to shift to the local mean.
- `subset` vector specifying a subset of 1:n, where n is the sample size. This allows to run the iterative mean shift procedure only from a subset of points (if unspecified, 1:n is used here, i.e. each data point serves as a starting point).
- `scaled` logical (if TRUE, each variable is divided by its range).
- `taumin,taumax,gridsize` determine the grid of bandwidths to investigate.
- `thresh, iter` mean shift iterations are stopped when the mean shift length (relative to the length of x) falls below `thresh`, or after `iter` iterations (whatever event happens first).
- `thr` adjacent mean shift clusters are merged if their relative distance falls below this threshold.
- `cluster` if TRUE, distances are always measured to the cluster to which an observation is assigned, rather than to the nearest cluster.
- `plotms, plot.type, or.labels, ...` graphical parameters.
- `print` if TRUE, coverage values are printed on the screen as soon as computed. This is quite helpful especially if `gridsize` is large.
Details

The methods implemented here can be used for density mode estimation, clustering, and the selection of starting points for the LPC algorithm.

Chen (1995) showed that, if the mean shift is computed iteratively, the resulting sequence of local means converges to a mode of the estimated density function. By assigning each data point to the mode to which it has converged, this turns into a clustering technique.

The concepts of coverage and self-coverage, which were originally introduced in the principal curve context, adapt straightforwardly to this setting.

The goodness-of-fit measure $R_c$ can also be applied in this context. For instance, a value of $R_c = 0.8$ means that, after the clustering, the mean absolute residual length has been reduced by 80% (compared to the distances to the overall mean).

Value

The main function `ms` produces an object of class `ms`, with components:

- `cluster.center`: a matrix which gives the coordinates of the estimated density modes (i.e., of the mean-shift based cluster centers).
- `cluster.label`: assigns each data point to the cluster center to which its mean shift trajectory has converged.
- `closest.label`: assigns each data point to the closest cluster center in terms of Euclidean distance.
- `data`: the data frame (scaled if `scaled=TRUE`).
- `scaled`: boolean.
- `scaled.by`: the data were scaled by dividing each variable through the values provided in this vector.

For all other functions, use `names()`. 

Author(s)

J. Einbeck. See `LPCM-package` for further acknowledgements.

References


See Also

$R_c$, `lpc.self.coverage`
Examples

data(faithful)
  # Mean shift clustering with user-defined bandwidth (5 percent of data range)
  fit <- ms(faithful, h=0.05)

  # Goodness-of-fit
  coverage(fit$data, fit$cluster.center)
  Rc(fit)

  # Bandwidth selection via self-coverage
  ## Not run: foo <- ms.self.coverage(faithful, gridsize= 50, taumin=0.1,
  ## taumax=0.5, plot.type="o")
  h <- select.self.coverage(foo)$select
  fit <- ms(faithful, h=h[1])
  ## End(Not run)

plot.lpc

Plotting local principal curves

Description

Takes an object of class lpc or lpc.spline and plots any subset of the following components of
the local principal curve: Centers of mass; the curve connecting the local centers of mass; the cubic
spline representation of the curve; the projections onto the curve; the starting points.

Usage

## S3 method for class 'lpc'
plot(x, type, unscale = TRUE, lwd = 1, datcol = "grey60",
     datpch = 21, masscol = NULL, masspch = 15, curvecol = 1, splinecol = 3,
     projectcol = 4, startcol = NULL, startpch = NULL,...)

## S3 method for class 'lpc.spline'
plot(x, type, unscale = TRUE, lwd = 1, datcol = "grey60",
     datpch = 21, masscol = NULL, masspch = 15, curvecol = 1, splinecol = 3,
     projectcol = 4, startcol = NULL, startpch = NULL,...)

Arguments

x
  an object of class lpc or lpc.spline.

type
  a vector of type c("mass", "spline", ...) with possible entries mass, curve, spline, project, start.

unscale
  if TRUE, then data (and all fitted components) are scaled back to their original
  scale; otherwise the scaled data are plotted (only relevant if scaled=TRUE in the
  fitted object).

lwd
  width of curves.

datcol
  color of data points.
plot.lpc

datpch  
plotting symbol for data points.

masscol  
color of centers of mass (see below).

masspch  
plotting symbol for centers of mass.

curvecol  
color of the curve interpolating the local centers of mass (this is the "local principal curve").

splinect  
color of the spline representation of the local principal curve.

projectcol  
color of projections onto the spline representation of the local principal curve.

startcol  
color of the plotted starting points.

startpch  
plotting symbol for starting points; needs to be either a single symbol, or a vector of symbols of the same length as the number of starting points.

...  
further arguments passed to plot or scatterplot3d.

Value

A 2D plot, 3D plot, or a pairs plot (depending on the data dimension $d$).

The most flexible plotting option is masscol. Depending on the length of the specified vector, this will be interpreted differently. If a scalar is provided, the corresponding color will be given to all centers of mass. If the length of the vector is larger than 1, then this option will assign different colours to different depths, or different branch numbers, or to individual data points, depending on the length. The default setting is assigning colours according to depth, in the order red, blue, black.

With increasing dimension $d$, less plotting options tend to be supported. The nicest plots are obtained for $d = 2$ and $d = 3$.

Warning

This function computes all missing information (if possible), so computation will take the longer the less informative the given object is, and the more advanced aspects are asked to plot!

Author(s)

JE

References


See Also

lpc, lpc.spline
print.lpc

Printing output for lpc and lpc.spline objects

Examples

```r
data(calspeedflow)
lpc1 <- lpc(calspeedflow[,3:4])
plot(lpc1, type=c("spline","project"), lwd=2)
dlpc <- lpc.spline(lpc1)
print(lpc)
```

Description

Takes an object of class `lpc` or `lpc.spline` and displays some standard output.

Usage

```r
## S3 method for class 'lpc'
print(x, digits = max(3,getOption("digits") - 3), ...)
## S3 method for class 'lpc.spline'
print(x, digits = max(3,getOption("digits") - 3), ...)
```

Arguments

- `x` an object of class `lpc` or `lpc.spline`.
- `digits` not yet in use.
- `...` further arguments.

Value

Some short text.

Author(s)

JE

See Also

- `lpc`

Examples

```r
data(calspeedflow)
lpc1 <- lpc(calspeedflow[,3:4])
print(lpc1)
lpc2 <- lpc.spline(lpc1)
print(lpc2)
```
Measuring goodness-of-fit for principal objects.

Description

These functions compute the ‘coverage coefficient’ $RC$ for local principal curves, local principal points (i.e., kernel density estimates obtained through iterated mean shift), and other principal objects.

Usage

Rc(x,...)

## S3 method for class 'lpc'
Rc(x,...)

## S3 method for class 'lpc.spline'
Rc(x,...)

## S3 method for class 'ms'
Rc(x,...)

base.Rc(data, closest.coords, type="curve")

Arguments

x an object used to select a method.

... Further arguments passed to or from other methods (not needed yet).

data A data matrix.

closest.coords A matrix of coordinates of the projected data.

type For principal curves, don’t modify. For principal points, set "points".

Details

Rc computes the coverage coefficient $RC$, a quantity which estimates the goodness-of-fit of a fitted principal object. This quantity can be interpreted similar to the coefficient of determination in regression analysis: Values close to 1 indicate a good fit, while values close to 0 indicate a ‘bad’ fit (corresponding to linear PCA).

For objects of type lpc, lpc.spline, and ms, S3 methods are available which use the generic function Rc. This, in turn, calls the base function base.Rc, which can also be used manually if the fitted object is of another class. In principle, function base.Rc can be used for assessing goodness-of-fit of any principal object provided that the coordinates (closest.coords) of the projected data are available. For instance, for HS principal curves fitted via princurve, this information is contained in component $s$, and for a a k-means object, say fitk, this information can be obtained via fitk$centers[fitk$cluster,]. Set type="points" in the latter case.

The function Rc attempts to compute all missing information, so computation will take the longer the less informative the given object x is. Note also, Rc looks up the option scaled in the fitted
object, and accounts for the scaling automatically. Important: If the data were scaled, then do NOT unscale the results by hand in order to feed the unscaled version into base.Rc, this will give a wrong result.

In terms of methodology, these functions compute $R_C$ directly through the mean reduction of absolute residual length, rather than through the area above the coverage curve.

These functions do currently not account for observation weights, i.e. $R_C$ is computed through the unweighted mean reduction in absolute residual length (even if weights have been used for the curve fitting).

Acknowledgements

Contributions (in form of pieces of code, or useful suggestions for improvements) by Jo Dwyer, Mohammad Zayed, and Ben Oakley are gratefully acknowledged.

Author(s)

J. Einbeck and L. Evers.

References


See Also

lpc.spline, codems, coverage.

Examples

data(calspeedflow)
lpc1 <- lpc.spline(lpc(calspeedflow[,3:4]), project=TRUE)
Rc(lpc1)  # is the same as:
base.Rc(lpc1$lpcobject$data, lpc1$closest.coords)

ms1 <- ms(calspeedflow[,3:4], plotms=0)
Rc(ms1)  # is the same as:
base.Rc(ms1$data, ms1$cluster.center[ms1$closest.label,], type="points")
Unscaling local principal objects.

Description

unscale takes an object of type lpc, lpc.spline, or ms, which had been fitted using option scaled=TRUE, and transforms the scaled components back to the original data scale.

Usage

unscale(x, ...)

## S3 method for class 'lpc'
unscale(x, ...)

## S3 method for class 'lpc.spline'
unscale(x, ...)

## S3 method for class 'ms'
unscale(x, ...)

Arguments

x an object used to select a method.

... Further arguments passed to or from other methods (not needed yet).

Value

A list of relevant items, such as LPC, start, cluster.centers, etc., which gives the unscaled versions of these quantities (some of them may carry the value NULL, if the corresponding information was not available from x).

Author(s)

JE

See Also

lpc, lpc.spline, ms

Examples

data(gvessel)
unscale(lpc(gvessel[,c(2,4,5)], h=0.11, x0=c(35, 1870, 6.3)) )
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