Package: ScatterDensity (via r-universe)

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Type Package

Title Density Estimation and Visualization of 2D Scatter Plots

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Description The user has the option to utilize the two-dimensional density estimation techniques called smoothed density published by Eilers and Goeman (2004) <doi:10.1093/bioinformatics/btg454>, and pareto density which was evaluated for univariate data by Thrun, Gehlert and Ultsch, 2020 <doi:10.1371/journal.pone.0238835>. Moreover, it provides visualizations of the density estimation in the form of two-dimensional scatter plots in which the points are color-coded based on increasing density. Colors are defined by the one-dimensional clustering technique called 1D distribution cluster algorithm (DDCAL) published by Lux and Rinderle-Ma (2023) <doi:10.1007/s00357-022-09428-6>.

LazyLoad yes

Imports Rcpp, RcppParallel (>= 5.1.4), pracma

Suggests DataVisualizations, ggplot2, ggExtra, plotly, FCPS, parallelDist, secr, ClusterR, geometry

Depends methods, R (>= 2.10)

LinkingTo Rcpp, RcppArmadillo, RcppParallel

NeedsCompilation yes

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Encoding UTF-8

URL https://www.deepbionics.org/

BugReports https://github.com/Mthrun/ScatterDensity/issues

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`ScatterDensity-package`

Density Estimation and Visualization of 2D Scatter Plots

Description

The user has the option to utilize the two-dimensional density estimation techniques called smoothed density published by Eilers and Goeman (2004) <doi:10.1093/bioinformatics/btg454>, and pareto density which was evaluated for univariate data by Thrun, Gehlert and Ultsch, 2020 <doi:10.1371/journal.pone.0238835>. Moreover, it provides visualizations of the density estimation in the form of two-dimensional scatter plots in which the points are color-coded based on increasing density. Colors are defined by the one-dimensional clustering technique called 1D distribution cluster algorithm (DDCAL) published by Lux and Rinderle-Ma (2023) <doi:10.1007/s00357-022-09428-6>.

Details

The DESCRIPTION file:

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Version:	0.1.1
Date:	2025-08-20
Authors@R:	c(person("Michael", "Thrun", email= "m.thrun@gmx.net",role=c("aut","cre","cph"), comment = c(ORC
Maintainer:	Michael Thrun <m.thrun@gmx.net>

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LazyLoad:	yes
Imports:	Rcpp, RcppParallel (>= 5.1.4), pracma
Suggests:	DataVisualizations, ggplot2, ggExtra, plotly, FCPS, parallelDist, secr, ClusterR, geometry
Depends:	methods, R (>= 2.10)
LinkingTo:	Rcpp, RcppArmadillo, RcppParallel
NeedsCompilation:	yes
License:	GPL-3
Encoding:	UTF-8
URL:	https://www.deepbionics.org/
BugReports:	https://github.com/Mthrun/ScatterDensity/issues
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Archs:	x64

Index of help topics:

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DensityScatter.DDCAL	Scatter density plot [Brinkmann et al., 2023]
PDEscatter	Scatter Density Plot
PointsInPolygon	PointsInPolygon
PolygonGate	PolygonGate
SampleScatter	takes a sample for a scatter plot
ScatterDensity-package	Density Estimation and Visualization of 2D Scatter Plots
SmoothedDensitiesXY	Smoothed Densities X with Y
fast_table_num	fast_table_num
inPSphere2D	2D data points in Pareto Sphere

Author(s)

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Maintainer: Michael Thrun <m.thrun@gmx.net>

Examples

```
#Todo
```

DDCAL *Density Distribution Cluster Algorithm of [Lux and Rinderle-Ma, 2023].*

Description

DDCAL is a clustering-algorithm for one-dimensional data, which heuristically finds clusters to evenly distribute the data points in low variance clusters.

Usage

```
DDCAL(data, nClusters, minBoundary = 0.1, maxBoundary = 0.45,

numSimulations = 20, csTolerance = 0.45, csToleranceIncrease = 0.5)
```

Arguments

`data`	[1:n] Numeric vector, with the data values
`nClusters`	Scalar, number of clusters to be found
`minBoundary`	Scalar, in the range (0,1), gives the lower boundary (in percent), for the simulation. Default is 0.1
`maxBoundary`	Scalar, in the range (0,1), gives the upper boundary (in percent), for the simulation. Default is 0.45
`numSimulations`	Scalar, number of simulations/iterations of the algorithm
`csTolerance`	Scalar, in the range (0,1). Gives cluster size tolerance factor. The necessary cluster size is defined by (dataSize/nClusters - dataSize/nClusters * csTolerance). Default is 0.45
`csToleranceIncrease`	Scalar, in the range (0,1), gives the procentual increase of the csTolerance-factor, if some clusters did not reach the necessary size. Default is 0.5

Details

DDCAL creates a evenly spaced division of the min-max-normalized data from minBoundary to maxBoundary. Those divisions will be used as boundaries. The first initial clusters will be the data from min(data) to minBoundary and maxBoundary to max(data). The clusters will be extended to neighboring points, as long as the standard deviations of the clusters will be reduced. A potential clusters will be used, if they have the necessary size, given as (dataSize/nClusters - dataSize/nClusters * csTolerance). If both clusters can be used, the left cluster (which is the cluster from min(data) to minBoundary or above) is preferred. If no clusters can be found with the necessary size, then the csTolerance-factor and with it the necessary cluster size will be lowered. If a clusters is used, the next boundaries are found, which are not in the already existing clusters and the procedure is repeated with the not already clustered data, until all points are assigned to clusters.

If a matrix is given as input data, the first column of the matrix will be used as data for the clustering

Non-finite values will not be clustered, but instead will get the cluster label `NaN`.

The algorithm is not garantueed to produce the given number of clusters, given in nClusters. The found number of clusters can be lower, depending on the data and input parameters.

Value

`labels`	[1:n] Numeric vector, containing the labels for the input data points

Author(s)

Luca Brinkmann

References

[Lux and Rinderle-Ma, 2023] Lux, M., Rinderle-Ma, S.: DDCAL: Evenly Distributing Data into Low Variance Clusters Based on Iterative Feature Scaling; Springer Journal of Classification, Vol. 40, pp. 106-144, DOI: doi:10.1007/s00357022094286, 2023.

Examples

```
# Load data
if(requireNamespace("FCPS")){
data(EngyTime, package = "FCPS")
engyTimeData = EngyTime$Data
c1 = engyTimeData[,1]
c2 = engyTimeData[,2]
}else{
c1 = rnorm(n=4000)
c2 = rnorm(n=4000,1,2)
}
# Calculate Densities

densities = SmoothedDensitiesXY(c1,c2)$Densities
# Use DDCAL to cluster the densities
labels = DDCAL(densities, 9)

# Plot Densities according to labels
my_colors = c("#000066", "#3333CC", "#9999FF", "#00FFFF", "#66FF33",
                      "#FFFF00", "#FF9900", "#FF0000", "#990000")
labels = as.factor(labels)
df = data.frame(c1, c2, labels)

if(requireNamespace("ggplot2")){
ggplot2::ggplot(df, ggplot2::aes(c1, c2, color = labels)) +
  ggplot2::geom_point() +
  ggplot2::scale_color_manual(values = my_colors)
}
```

`DensityScatter.DDCAL` *Scatter density plot [Brinkmann et al., 2023]*

Description

Density estimation (PDE) [Ultsch, 2005] or "SDH" [Eilers/Goeman, 2004] used for a scatter density plot, with clustering of densities with DDCAL [Lux/Rinderle-Ma, 2023] proposed by [Brinkmann et al., 2023].

Usage

```
DensityScatter.DDCAL(X, Y, nClusters = 12, Plotter = "native",
SDHorPDE = TRUE, LimitShownPoints = FALSE,
Marginals = FALSE, na.rm=TRUE, pch, Size,
xlab="x", ylab="y", main = "",lwd = 2,
xlim=NULL,ylim=NULL,Polygon,BW = TRUE,Silent = FALSE, ...)
```

Arguments

X	Numeric vector [1:n], first feature (for x axis values)
Y	Numeric vector [1:n], second feature (for y axis values)
nClusters	(Optional) Integer defining the number of clusters (colors) used for finding a hard color transition, default is 12.
Plotter	(Optional) String, name of the plotting backend to use. Possible values are: "native", "plotly", or "ggplot2"
SDHorPDE	(Optional) Boolean, if TRUE SDH is used to calculate density, if FALSE PDE is used
LimitShownPoints	(Optional) FALSE: does nothing, TRUE: samples the number of optimal points for visualization using SampleScatter
Marginals	(Optional) Boolean, if TRUE the marginal distributions of X and Y will be plotted together with the 2D density of X and Y. Default is FALSE
na.rm	(Optional) Boolean, if TRUE non finite values will be removed
pch	(Optional) Scalar or character. Indicates the shape of data points, see plot function, symbol argument in **plotly** package, or the shape argument in **ggplot2** package, default is 20 for native and for ggplot2, and 0 for plotly
Size	(Optional) Scalar, size of data points in plot, default is 1 for native, 6 for plotly, and 3 for ggplot2
xlab	(Optional) String, title of the x axis. Default: "X", see plot() function, or similar functonality in **plotly** or **ggplot2**
ylab	(Optional) String, title of the y axis. Default: "Y", see plot() function, or similar functonality in **plotly** or **ggplot2**
main	(Optional) Character, title of the plot.
lwd	(Optional) Scalar, thickness of the lines used for the marginal distributions (only needed if Marginals=TRUE), see plot(). Default = 2
xlim	(Optional) numerical vector, min and max of x values to be plottet
ylim	(Optional) numerical vector, min and max of y values to be plottet
Polygon	(Optional) [1:p,1:2] numeric matrix that defines for x and y coordinates a polygon in magenta
BW	(Optional) Boolean, if TRUE and Plotter="ggplot2" will use a white background, if FALSE and Plotter="ggplot2", the typical ggplot2 background is used. Not needed if "Plotter="native". Default is TRUE
Silent	(Optional) Boolean, if TRUE no messages will be printed, default is FALSE
...	Further plot arguments

Details

The DensityScatter.DDCAL function generates the density of the xy data as a z coordinate. Afterwards xyz will be plotted as a contour plot. It assumens that the cases of x and y are mapped to each other meaning that a cbind(x,y) operation is allowed. The colors for the densities in the contour plot are calculated with DDCAL, which produces clusters to evenly distribute the densities in low variance clusters.

In the case of "native" as Plotter, the handle returns NULL because the basic R functon plot() is used.

For the returned density values see SmoothedDensitiesXY or PDEscatter depending on input parameter SDHorPDE for details.

Value

returns a invisible list with

DF	[1:m,1:5] of Density values, x values, y values, colors, and classification vector Cls. m=n if LimitShownPoints=FALSE, otherwise LimitShownPoints=TRUE m<n meaning that subsample is taken
PlotHandle	the plotting handle, either an object of **plotly**, **ggplot2** or NULL depending on input parameter Plotter

Author(s)

Luca Brinkmann, Michael Thrun

References

[Ultsch, 2005] Ultsch, A.: Pareto density estimation: A density estimation for knowledge discovery, In Baier, D. & Werrnecke, K. D. (Eds.), Innovations in classification, data science, and information systems, (Vol. 27, pp. 91-100), Berlin, Germany, Springer, 2005.

[Eilers/Goeman, 2004] Eilers, P. H., & Goeman, J. J.: Enhancing scatterplots with smoothed densities, Bioinformatics, Vol. 20(5), pp. 623-628. 2004.

[Lux/Rinderle-Ma, 2023] Lux, M. & Rinderle-Ma, S.: DDCAL: Evenly Distributing Data into Low Variance Clusters Based on Iterative Feature Scaling, Journal of Classification vol. 40, pp. 106-144, 2023.

[Brinkmann et al., 2023] Brinkmann, L., Stier, Q., & Thrun, M. C.: Computing Sensitive Color Transitions for the Identification of Two-Dimensional Structures, Proc. Data Science, Statistics & Visualisation (DSSV) and the European Conference on Data Analysis (ECDA), p.109, Antwerp, Belgium, July 5-7, 2023.

Examples

```
# Create two bimodial distributions
x1=rnorm(n = 7500,mean = 0,sd = 1)
y1=rnorm(n = 7500,mean = 0,sd = 1)
x2=rnorm(n = 7500,mean = 2.5,sd = 1)
```

```
y2=rnorm(n = 7500,mean = 2.5,sd = 1)
x=c(x1,x2)
y=c(y1,y2)

DensityScatter.DDCAL(x, y, Marginals = TRUE)
```

`fast_table_num` *fast_table_num*

Description

Faster table computation in R compared to table() by omitting as.character and as.factor

Usage

```
fast_table_num(x, y, edges_x, edges_y, redefine = TRUE,

byrow = FALSE, all.inside = FALSE, rightmost.closed = FALSE,

sort = FALSE, na.rm = FALSE, names = FALSE,extendOutput=FALSE)
```

Arguments

`x`	[1:n] numerical vector
`y`	[1:n] numerical vector
`edges_x`	Optional, [1:(k_1+1)] numerical vector defining the specific borders in `x` default `unique(x)` for categorical scale
`edges_y`	Optional, [1:(k_2+1)] numerical vector defining the specific borders in `y`, default `unique(y)` for categorical scale
`redefine`	Optional, boolean TRUE: resets counts in `y` direction in order from 1:k_2 to k_2:1
`byrow`	Optional, boolean, If FALSE (the default) the count matrix is filled by columns, otherwise the matrix is filled by rows.
`all.inside`	Optional, boolean, if TRUE, the returned indices are coerced into 1,...,N-1, i.e., 0 is mapped to 1 and N to N-1
`rightmost.closed`	Optional, boolean, if TRUE, the rightmost interval, vec[N-1] .. vec[N] is treated as closed
`sort`	Optional, boolean, if TRUE, `edges_x`, `edges_y` are sorted non-decreasingly, NA/NaN are gthen ignored
`na.rm`	Optional, boolean, if TRUE, only complete observations are taken into account
`names`	Optional, boolean, if TRUE, output matrix is named by `edges_x`[1:k_1] and `edges_y`[1:k_2] (left-sided)
`extendOutput`	Optional, boolean, default FALSE, if TRUE, list is the output, otherwise numerical matrix, see below.

Details

edges_x and edges_y must be sorted non-decreasingly. Beware that kernels are centers of bins, `edges_x`, `edges_y` are borders of bins. If edges are given, `edges_x`, `edges_y` can contain Inf,-Inf borders. In that case, edges always define n-1 bins lying within the edges. data outside first edge or last edge are ignored. Edges have either to be sorted non-decreasingly or set sort=T.

If edges are not given, set sort=T. In this case, they define the unique number of points. Then the number of edges internally sets the number of bins.

Beware that in matrix notation, `count matrix` would be expected be ordered [1:k_1,1:k_2] instead of [1:k_2,1:k_1]. Here we use the ordering that intuitively is given in plot(x,y), i.e. x are columns and y are rows.

Value

`extendOutput==FALSE`: [1:k_2,1:k_1] numerical matrix of counts

If `extendOutput==TRUE`, then list of named elements

`count matrix`: [1:k_2,1:k_1] numerical matrix of counts

`x_idx`: [1:k_1] numerical vector of counts for `x` based on `edges_x`

`y_idx`: [1:k_2] numerical vector of counts for `y` based on `edges_y`

Author(s)

Michael Thrun

See Also

`table`, `findInterval`

Examples

```
if(requireNamespace("FCPS")){
data(Hepta,package ="FCPS")
Cls=Hepta$Cls
Cls1=Cls+1
#k unqiue points define k bins
fast_table_num(Cls,Cls1,

redefine = FALSE,names=TRUE)==as.matrix(table(Cls,Cls1))
}
#k unqiue points define k bins
tab=fast_table_num(rnorm(100),rnorm(100),redefine=FALSE,sort=TRUE)

#set k+1 edges to get k bins
x=rnorm(100)
y=rnorm(100)
binsxy=5
edgex=seq(from=min(x),to=max(x),length.out=binsxy+1)
edgesy=seq(from=min(y),to=max(y),length.out=binsxy+1)
fast_table_num(x,y,edgex,edgesy,
```

```
redefine=FALSE,names=TRUE,rightmost.closed =TRUE)

#definition of counts analog to plotting
x = c(rnorm(1000, mean=-5), rnorm(1000, mean=5))
y = rnorm(2000)
edgesx = seq(min(x), max(x), length.out=512+1)
edgesy = seq(min(y), max(y), length.out=256+1)
joint_table = fast_table_num(x, y, edgesx, edgesy)

plot(x,y)
plot(colSums(joint_table),xlab="x marginal",

ylab="sum of counts",main="x-values are stored in columns")

plot(rowSums(joint_table),xlab="y marginal",

ylab="sum of counts",main="y-values are stored in rows")
```

`inPSphere2D` *2D data points in Pareto Sphere*

Description

This function determines the 2D data points inside a ParetoSphere with ParetoRadius.

Usage

```
inPSphere2D(data, paretoRadius=NULL,Compute="Cpp")
```

Arguments

`data`	numeric matrix of data.
`paretoRadius`	numeric value. radius of P-spheres. If not given, calculate by the function 'paretoRad'
`Compute`	Either `Cpp` or `Parallel`, defines if multicore computation should be performed

Value

numeric vector with the number of data points inside a P-sphere with ParetoRadius.

Author(s)

Felix Pape

PDEscatter *Scatter Density Plot*

Description

Concept of Pareto density estimation (PDE) proposed for univsariate data by [Ultsch, 2005] and comparet to varius density estimation techniques by [Thrun et al., 2020] for univariate data is here applied for a scatter density plot. It was also applied in [Thrun and Ultsch, 2018] to bivariate data, but is not yet compared to other techniques.

Usage

```
PDEscatter(x,y,SampleSize,

na.rm=FALSE,PlotIt=TRUE,ParetoRadius,Compute="Cpp",

sampleParetoRadius, NrOfContourLines=20,Plotter='native',

DrawTopView = TRUE, xlab="X", ylab="Y", main="PDEscatter",

xlim, ylim, Legendlab_ggplot="value", NormalizeAsDensity=FALSE)
```

Arguments

x	Numeric vector [1:n], first feature (for x axis values)
y	Numeric vector [1:n], second feature (for y axis values)
SampleSize	Numeric m, positiv scalar, maximum size of the sample used for calculation. High values increase runtime significantly. The default is that no sample is drawn
na.rm	Function may not work with non finite values. If these cases should be automatically removed, set parameter TRUE
ParetoRadius	Numeric, positiv scalar, the Pareto Radius. If omitted (or 0), calculate by paretoRad.
Compute	Either Cpp or Parallel, defines if multicore computation should be performed
sampleParetoRadius	Numeric, positiv scalar, maximum size of the sample used for estimation of "kernel", should be significantly lower than SampleSize because requires distance computations which is memory expensive
PlotIt	TRUE: plots with function call FALSE: Does not plot, plotting can be done using the list element Handle -1: Computes density only, does not perfom any preperation for plotting meaning that Handle=NULL
NrOfContourLines	Numeric, number of contour lines to be drawn. 20 by default.

Plotter	String, name of the plotting backend to use. Possible values are: "native", "ggplot", "plotly"
DrawTopView	Boolean, True means contur is drawn, otherwise a 3D plot is drawn. Default: TRUE
xlab	String, title of the x axis. Default: "X", see plot() function
ylab	String, title of the y axis. Default: "Y", see plot() function
main	string, the same as "main" in plot() function
xlim	see plot() function
ylim	see plot() function
Legendlab_ggplot	String, in case of Plotter="ggplot" label for the legend. Default: "value"
NormalizeAsDensity	Boolean, TRUE: normalizes densities to integral equal to one

Details

The PDEscatter function generates the density of the xy data as a z coordinate. Afterwards xyz will be plotted either as a contour plot or a 3d plot. It assumens that the cases of x and y are mapped to each other meaning that a cbind(x,y) operation is allowed. This function plots the PDE on top of a scatterplot. Variances of x and y should not differ by extreme numbers, otherwise calculate the percentiles on both first. If DrawTopView=FALSE only the plotly option is currently available. If another option is chosen, the method switches automatically there.

The method was succesfully used in [Thrun, 2018; Thrun/Ultsch 2018].

PlotIt=FALSE is usefull if one likes to perform adjustements like axis scaling prior to plotting with **ggplot2** or **plotly**. In the case of "native"" the handle returns NULL because the basic R functon plot() is used

Value

List of:

X	Numeric vector [1:m],m<=n, first feature used in the plot or the kernels used
Y	Numeric vector [1:m],m<=n, second feature used in the plot or the kernels used
Densities	Numeric vector [1:m],m<=n, Number of points within the ParetoRadius of each point, i.e. density information
Matrix3D	1:n,1:3] marix of x,y and density information
ParetoRadius	ParetoRadius used for PDEscatter
Handle	Handle of the plot object. Information-string if native R plot is used.

Note

MT contributed with several adjustments

Author(s)

Felix Pape

References

[Thrun/Ultsch, 2018] Thrun, M. C., & Ultsch, A. : Effects of the payout system of income taxes to municipalities in Germany, in Papiez, M. & Smiech,, S. (eds.), Proc. 12th Professor Aleksander Zelias International Conference on Modelling and Forecasting of Socio-Economic Phenomena, pp. 533-542, Cracow: Foundation of the Cracow University of Economics, Cracow, Poland, 2018.

[Ultsch, 2005] Ultsch, A.: Pareto density estimation: A density estimation for knowledge discovery, In Baier, D. & Werrnecke, K. D. (Eds.), Innovations in classification, data science, and information systems, (Vol. 27, pp. 91-100), Berlin, Germany, Springer, 2005.

[Thrun et al., 2020] Thrun, M. C., Gehlert, T. & Ultsch, A.: Analyzing the Fine Structure of Distributions, PLoS ONE, Vol. 15(10), pp. 1-66, DOI doi:10.1371/journal.pone.0238835, 2020.

Examples

```
#taken from [Thrun/Ultsch, 2018]
if(requireNamespace("DataVisualizations")){
data("ITS",package = "DataVisualizations")
data("MTY",package = "DataVisualizations")
Inds=which(ITS<900&MTY<8000)
plot(ITS[Inds],MTY[Inds],main='Bimodality is not visible in normal scatter plot')


PDEscatter(ITS[Inds],MTY[Inds],xlab = 'ITS in EUR',

ylab ='MTY in EUR' ,main='Pareto Density Estimation indicates Bimodality' )


}
```

`PointsInPolygon` *PointsInPolygon*

Description

Defines a Cls based on points in a given polygon.

Usage

```
PointsInPolygon(Points, Polygon, PlotIt = FALSE, ...)
```

Arguments

`Points`	[1:n,1:2] xy cartesian coordinates of a projection
`Polygon`	Numerical matrix of 2 columns defining a closed polygon
`PlotIt`	TRUE: Plots marked points
`...`	`BMUorProjected`: Default == FALSE, If TRUE assuming BestMatches of ESOM instead of Projected Points `main`: title of plot Further Plotting Arguments,xlab etc used in `Classplot`

Details

We assume that polygon is closed, i.e., that the last point connects to the fist point

Value

Numerical classification vector Cls with 1 = outside polygon and 2 = inside polygon

Author(s)

Michael Thrun

See Also

Classplot

Examples

```
XY=cbind(runif(80,min = -1,max = 1),rnorm(80))
#closed polygon
polymat <- cbind(x = c(0,1,1,0), y = c(0,0,1,1))
#takes sometimes more than 5 sec

Cls=PointsInPolygon(XY,polymat,PlotIt = TRUE)
```

PolygonGate *PolygonGate*

Description

A specific Gate defined by xy coordinates that result in a closed polygon is applied to the flowcytometry data.

Usage

```
PolygonGate(Data, Polygon, GateVars,  PlotIt = FALSE, PlotSampleSize = 1000)
```

Arguments

Data	numerical matrix n x d
Polygon	numerical marix of two columns defining the coordiantes of the polygon. polygon assumed to be closed, i.e.,last coordinate connects to first coordinate.
GateVars	vector, either column index in Data of X and Y coordinates of gate or its variable names as string
PlotIt	if TRUE: plots a sample of data in the two selected variables and marks point inside the gate as yellow and outside as magenta
PlotSampleSize	size pof the plottet sample

Details

Gates are alwaxs two dimensional, i.e., require two filters, although all dimensions of data are filted by the gates. Only high-dimensional points inside the polygon (gate) are given back

Value

list of

DataInGate	m x d numerical matrix with m<=n of data points inside the gate
InGateInd	index of length m for the datapoints in original matrix

Note

if GateVars is not found a text is given back which will state this issue

Author(s)

Michael Thrun

See Also

PointsInPolygon

Examples

```
Data <- matrix(runif(1000), ncol = 10)
colnames(Data)=paste0("GateVar",1:ncol(Data))
poly <- cbind(x = c(0.2,0.5,0.8), y = c(0.2,0.8,0.2))
#set PlotIt TRUE for understanding the example


#Select index
V=PolygonGate(Data,poly,c(5,8),PlotIt=FALSE,100)

#select var name
V=PolygonGate(Data,poly,c("GateVar5","GateVar8"),PlotIt=FALSE,100)
```

SampleScatter *takes a sample for a scatter plot*

Description

Given 2D points having X and Y coordinates takes a sample, such that these points are is optimally visualized if a plot function is called.

Usage

```
SampleScatter(X, Y, ThresholdPoints = 20,

DensityThreshold, nbins = 100,na.rm=TRUE, PlotIt = FALSE)
```

Arguments

X	[1:n] x coordinate
Y	[1:n] y coordinate
ThresholdPoints	how many points per grid are allowed to overlap, default 20
DensityThreshold	default is internally estimated, scalar above zero and below 1, SDH threshold for density computed with SmoothedDensitiesXY, usually good values are around 0.1
nbins	number of bins in grid, default 100x100
na.rm	Function may not work with non finite values. If these cases should be automatically removed, set parameter TRUE
PlotIt	Plots the remaining points

Details

"Optimally"" visualized in the sense that not too much point overap visually. The lower the value for ThresholdPoints, the smaller is the sample that is taken by the function.

Value

SubsampleInd[1:m] indices of m points, m<n, that will be relevant for a optimal scatter plot

Author(s)

Michael Thrun

See Also

SmoothedDensitiesXY

Examples

```
if(requireNamespace("DataVisualizations")){
data("ITS",package = "DataVisualizations")
data("MTY",package = "DataVisualizations")
sample_ind=SampleScatter(ITS,MTY,PlotIt=TRUE)
}else{
#sample random data
ITS=rnorm(10000)
```

```
MTY=rnorm(10000)
sample_ind=SampleScatter(ITS,MTY,ThresholdPoints = 5)
del_ind=setdiff(1:length(ITS),sample_ind)
plot(ITS,MTY,type="p",pch=20,col="grey",main="Grey=full data, red=overlapping data points")
points(ITS[del_ind],MTY[del_ind],type="p",pch=20,col="red")
}
```

SmoothedDensitiesXY *Smoothed Densities X with Y*

Description

Density is the smothed histogram density at [X,Y] of [Eilers/Goeman, 2004]

Usage

```
SmoothedDensitiesXY(X, Y, nbins, lambda, Xkernels, Ykernels,

isXDiscrete=FALSE, isYDiscrete=FALSE, Compute="Cpp", PlotIt = FALSE)
```

Arguments

X	Numeric vector [1:n], first feature (for x axis values)
Y	Numeric vector [1:n], second feature (for y axis values),
nbins	Optional, number of bins, default is 250, if scalar nbins=nx=ny otherwise c(nx,ny) vector of length 2
lambda	Optional, smoothing factor used by the density estimator, default is lambda = 20
Xkernels	Optional, bin kernels in x direction are given, if not set nx bins in range of data are generated with seq
Ykernels	Optional, bin kernels y direction are given, if not set ny bins in range of data are generated with seq
isXDiscrete	whether or not the first numeric vector resembles a continuous or discrete measurement
isYDiscrete	whether or not the second numeric vector resembles a continuous or discrete measurement
Compute	Either R, Cpp or Parallel, defines how the heavy-load computation is performed
PlotIt	FALSE: no plotting, TRUE: simple plot

Details

lambda has to chosen by the user and is a sensitive parameter and a lambda = 20 roughly means that the smoothing is over 20 bins around a given point.

Value

List of:

`Densities`	numeric vector [1:n] is the smothed density in 3D
`Xkernels`	numeric vector [1:nx], nx defined by `nbins`, such that mesh(Xkernels,Ykernels,F) form the (not NaN) smothed densisties
`Ykernels`	numeric vector [1:ny], nx defined by `nbins`, such that mesh(Xkernels,Ykernels,F) form the (not NaN) smothed densisties
`GridDensity`	matrix [1:nx,1:ny] beeing the smoothed 2D histogram
`Points2GridInd`	an index such that `Densities = GridDensity[Points2GridInd]`

Author(s)

Michael Thrun

References

[Eilers/Goeman, 2004] Eilers, P. H., & Goeman, J. J.: Enhancing scatterplots with smoothed densities, Bioinformatics, Vol. 20(5), pp. 623-628.DOI: doi:10.1093/bioinformatics/btg454, 2004.

Examples

```
if(requireNamespace("DataVisualizations")){
data("ITS",package = "DataVisualizations")
data("MTY",package = "DataVisualizations")
Inds=which(ITS<900&MTY<8000)
V=SmoothedDensitiesXY(ITS[Inds],MTY[Inds])
}else{
#sample random data
ITS=rnorm(1000)
MTY=rnorm(1000)
V=SmoothedDensitiesXY(ITS,MTY)
}
```

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