

Package ‘GPoM’

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Description Platform dedicated to the Global Modelling technique. Its aim is to obtain ordinary differential equations of polynomial form directly from time series. It can be applied to single or multiple time series under various conditions of noise, time series lengths, sampling, etc. This platform is developed at the Centre d'Etudes Spatiales de la Biosphere (CESBIO), UMR 5126 UPS/CNRS/CNES/IRD, 18 av. Edouard Belin, 31401 TOULOUSE, FRANCE. The developments were funded by the French program Les Enveloppes Fluides et l'Environnement (LEFE, MANU, projets GloMo, SpatioGloMo and MoMu). The French program Defi InFiNiTi (CNRS) and PNTS are also acknowledged (projects Crops'IChaos and Musc & SlowFast).

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Description

GPoM is a platform dedicated to the Global Modelling technique. Its aim is to obtain deterministic models of Ordinary Differential Equations from observational time series. It applies to single and to multiple time series. With single time series, it can be used: to detect low-dimensional determinism and low-dimensional (deterministic) chaos. It can also be used to characterize the observed behavior, using the obtained models as a proxy of the original dynamics, as far as the model validation could be checked. With multiple time series, it can be used: to detect couplings between observed variables, to infer causal networks, and to reformulate the original equations of the observed system (retro-modelling). The present package focuses on models in Ordinary Differential Equations of polynomial form. The package was designed to model weakly predictable dynamical behaviors (such as chaotic behaviors). Of course, it can also apply to more of fully predictable behavior, either linear or nonlinear. Several vignettes are associated to the package which can be used as a tutorial, and it also provides an overlook of the diversity of applications and at the performances of the tools. Users are kindly asked to quote the corresponding references when using the package (see hereafter).

Note

FOR USERS

This package was developed at Centre d'Etudes Spatiales de la Biosphere (Cesbio, UMR 5126, UPS-CNRS-CNES-IRD, <http://www.cesbio.ups-tlse.fr>). An important part of the developments were funded by the French program Les Enveloppes Fluides et l'Environnement (LEFE, MANU, projets GloMo, SpatioGloMo and MoMu). The French program Défi InFiNiTi (CNRS) and PNTS are also acknowledged (projects Crops'IChaos and Musc & SlowFast).

If you apply this package to single time series, please quote [6]. If you apply it to multivariate time series, please quote [10]. If you apply it to infer couplings among time series, please quote [8]. If you apply it to classification, please quote [11].

HISTORICAL BACKGROUND

The global modelling technique was initiated during the early 1990s [1-3]. It takes its background from the Theory of Nonlinear Dynamical Systems. Earlier investigations can also be found in the fields of Electrical Engineering and Statistics but these mainly focused on linear problems [4]. The approach became applicable to the analysis of real world environmental behaviours by the end of the 2000s [5-7]. Recent works have shown that the approach could be applied to numerous other dynamical behaviors [8-10]. Global modelling aims to obtain deterministic models directly from observed time series.

Author(s)

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References

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- [5] J. Maquet, C. Letellier, and L. A. Aguirre, 2007. Global models from the Canadian Lynx cycles as a first evidence for chaos in real ecosystems, *Juornal of Mathematical Biology*. 55(1), 21-39.
- [6] Mangiarotti S., Coudret R., Drapeau L., & Jarlan L., 2012. Polynomial search and global modeling : Two algorithms for modeling chaos, *Physical Review E*, 86, 046205.
- [7] Mangiarotti S., Drapeau L. & Letellier C., 2014. Two chaotic models for cereal crops observed from satellite in northern Morocco. *Chaos*, 24(2), 023130.
- [8] Mangiarotti S., 2015. Low dimensional chaotic models for the plague epidemic in Bombay (1896-1911). *Chaos, Solitons and Fractals*, 81A, 184-186.
- [9] Mangiarotti S., Peyre M. & Huc M., A chaotic model for the epidemic of Ebola Virus Disease in West Africa (2013-2016). *Chaos*, 26, 113112, 2016.
- [10] Mangiarotti S., 2014. Modelisation globale et Caracterisation Topologique de dynamiques environnementales - de l'analyse des enveloppes fluides et du couvert de surface de la Terre a la caracterisation topodynamique du chaos. Habilitation to Direct Research, University of Toulouse 3, France.
- [11] Mangiarotti S., Sharma A.K., Corgne S., Hubert-Moy L., Ruiz L., Sekhar M., Kerr Y., Can the global modelling technique be used for crop classification? *Chaos, Solitons & Fractals*, in press.

allMod_nVar3_dMax2 data set

Numerical description of a list of eighteen three-dimensional chaotic systems (see vignette VII_Retro-Modelling)

Description

A list named allMod_nVar3_dMax2 of matrix providing the numerical description of eighteen three-dimensional chaotic systems:

Lorenz-1963 (\$L63), Rössler-1976 (\$R76), Burke & shaw 1981 (\$BS81), Lorenz-1984 (\$L84), Nosé & Hooer 1986 (\$NH86), Genesio & Tosi 1992 (\$GT92), Spott systems 1994 (\$SprF, \$SprH, \$SprK, \$SprO, \$SprP, \$SprG, \$SprM, \$SprQ, \$SprS), Chlouverakis & Sprott 2004 (\$CS2004), Li 2007 (\$Li2007) and the Cord system by Aguirre & Letellier 2012 (\$Cord2012). Each dynamical system is provided as a matrix: each column corresponds to one equation, each lines to the polynomial coefficients which order is following the convection defined by function polabs(nVar = 3, dMax = 2).

Usage

```
allMod_nVar3_dMax2
```

Format

An object of class `list` of length 18.

Author(s)

Sylvain Mangiarotti, Mireille Huc.

References

All the references are provided in vignette VII_retro-modelling.

| | |
|-----------|---|
| allToTest | <i>A list providing the description of six models tested by the function autoGPoMoTest.</i> |
|-----------|---|

Description

List of 6 models available for tests (by autoGPoMoTest). Each model ($\$mToTest1$, $\$mToTest2$, etc.) is provided as a matrix of dimension $10 * 3$. Each column corresponds to one equation. The order of the coefficients follows the conventions defined by `poLabs(nVar = 3, dMax = 2)`.

Usage

```
allToTest
```

Format

An object of class `list` of length 6.

Author(s)

Sylvain Mangiarotti, Mireille Huc

Examples

```
#####
# example #
#####
data("allToTest")
# 6 models are available in this list:
names(allToTest)
# The parameter of their formulation (nVar and dMax)
# can be retrieved:
nVar <- dim(allToTest$mToTest6)[2]
dMax <- p2dMax(nVar = 3, pMaxKnown = dim(allToTest$mToTest6)[1])
# Their equation can be edited as follows:
visuEq(nVar, dMax, allToTest$mToTest6, approx = 2)
```

autoGPoMoSearch *Automatic search of polynomial Equations*

Description

This algorithm aims to get an ensemble of possible models which integrability will be tested later with function `autoGPoMoTest`. By default, all the terms are considered available (Some of the terms can be excluded intentionally using the option `filterReg`). The maximum size of the equation depends on the model dimension `nVar`, and on the maximum polynomial degree `dMax`. The algorithm removes polynomial terms one by one using a leave-one-out method.

Usage

```
autoGPoMoSearch(data, dt, nVar = nVar, dMax = dMax, weight = NULL,
  show = 0, underSamp = NULL, filterReg = NULL)
```

Arguments

| | |
|------------------------|---|
| <code>data</code> | Input Time series: Each column is one time series that corresponds to one variable. |
| <code>dt</code> | Time sampling of the input series. |
| <code>nVar</code> | Number of variables considered in the polynomial formulation. |
| <code>dMax</code> | Maximum degree of the polynomial formulation. |
| <code>weight</code> | A vector providing the binary weighting function of the input data series (0 or 1). By default, all the values are set to 1. |
| <code>show</code> | Provide (2) or not (0-1) visual output during the running process. |
| <code>underSamp</code> | Number of points used for undersampling the data. For <code>undersamp = 1</code> the complete time series is used. For <code>undersamp = 2</code> , only one data out of two is kept, etc. |
| <code>filterReg</code> | A vector that specifies the template for the equation structure (for one single equation). The convention defined by <code>poLabs</code> is used. Value is 1 if the regressor is available, 0 if it is not. |

Value

A list of two matrices:

`$filtMemo` describes the selected terms (1 if the term is used, 0 if not)

`$KMemo` provides the corresponding coefficients

Author(s)

Sylvain Mangiarotti, Flavie Le Jean

See Also

[autoGPoMoTest](#), [gPoMo](#), [findAllSets](#), [poLabs](#)

Examples

```
# Load data
data('RosYco')
# Search for potential models
filt = autoGPoMoSearch(RosYco[,2], nVar = 3, dMax = 2,
                      dt = 1/125, show = 1)
# As an example, the equations of the fourth line has the following terms:
poLabs(nVar = 3, dMax = 2)[filt$filtMemo[5,] != 0]
# which coefficients correspond to
cbind(filt$KMemo[5,], poLabs(nVar = 3, dMax = 2))[filt$filtMemo[5,] != 0,]
```

| | |
|---------------|--|
| autoGPoMoTest | <i>Tests the numerical integrability of models and classify their dynamical regime</i> |
|---------------|--|

Description

Tests the numerical integrability of provided models (these may have been obtained with function `autoGPoMoSearch`), and classify these models as Divergent, Fixed Points, Periodic or not Unclassified (potentially chaotic).

Usage

```
autoGPoMoTest(data, tin = NULL, dt = NULL, nVar = nVar, dMax = dMax,
              show = 1, verbose = 1, allKL = allKL, IstepMin = 10,
              IstepMax = 10000, tooFarThreshold = 4, LimCyclThreshold = 0,
              fixedPtThreshold = 1e-08, method = "rk4")
```

Arguments

| | |
|----------------------|---|
| <code>data</code> | Input Time series: Each column is one time series that corresponds to one variable. |
| <code>tin</code> | Input date vector which length should correspond to the input time series. |
| <code>dt</code> | Sampling time of the input time series. |
| <code>nVar</code> | Number of variables considered in the polynomial formulation. |
| <code>dMax</code> | Maximum degree of the polynomial formulation. |
| <code>show</code> | Provide (2) or not (0-1) visual output during the running process. |
| <code>verbose</code> | Gives information (if set to 1) about the algorithm progress and keeps silent if set to 0. |
| <code>allKL</code> | A list of all the models <code>\$mToTest1</code> , <code>\$mToTest2</code> , etc. to be tested. Each model is provided as a matrix. |

| | |
|----------|-------------------------------------|
| bDrvFilt | <i>Builds the derivative filter</i> |
|----------|-------------------------------------|

Description

Build the Savitzky-Golay derivative filter (Savitzky-Golay, 1964).

Usage

```
bDrvFilt(nDrv, tstep, winL = 9)
```

Arguments

| | |
|-------|---|
| nDrv | The number of derivatives to be computed. |
| tstep | Sampling time. |
| winL | The local window length to be used for computing the derivatives [1]. |

Author(s)

Sylvain Mangiarotti

References

[1] Savitzky, A.; Golay, M.J.E., Smoothing and Differentiation of Data by Simplified Least Squares Procedures. *Analytical Chemistry* 36 (8), 1627-1639, 1964.

| | |
|--------|---|
| cano2M | <i>cano2M : Converts a model in canonical form into a matrix form</i> |
|--------|---|

Description

Converts the vectorial formulation of canonical models into a matrix formulation (that is, including explicitly all the equations). For both input, the list of terms follows the convention defined by poLabs.

Usage

```
cano2M(nVar, dMax, poly)
```

Arguments

| | |
|------|---|
| nVar | The number of variables |
| dMax | The maximum degree allowed in the formulation |
| poly | A vector of coefficients corresponding to the regressor of the canonical function |

Author(s)

Sylvain Mangiarotti, Mireille Huc

See Also

[drvSucc](#), [gPoMo](#), [poLabs](#)

Examples

```
# A vector of polynomial terms corresponding to a canonical form:
polyTerms <- c(0.2,0,-1,0.5,0,0,0,0,0)
# Convert this vector into a matrix formulation with all the equations:
K <- cano2M(3,2,polyTerms)
# Visualize the equations:
visuEq(3,2,K)
```

compDeriv

Computes the successive derivatives of a time series

Description

Computes the successive derivatives from one single time series, with the Savitzky-Golay approach (1964).

Usage

```
compDeriv(TS, nDrv, tstep, winL = 9)
```

Arguments

| | |
|-------|---|
| TS | A single time series provided as a single vector. |
| nDrv | The number of derivatives to be computed from the input series. The resulting number of output time series will thus be $nVar = nDrv + 1$. |
| tstep | Sampling Time of the input time series TS. |
| winL | The local window length used for computing the derivatives [1-2]. |

Value

A matrix containing the original variable (smoothed by the filtering process) and its nDrv first derivatives (note that winL values of the original time series will be lost both at the beginning and the end of the time series due to boundary effect).

Author(s)

Sylvain Mangiarotti

References

- [1] Savitzky, A.; Golay, M.J.E., Smoothing and Differentiation of Data by Simplified Least Squares Procedures. *Analytical Chemistry* 36 (8), 1627-1639, 1964.
- [2] Steinier J., Termonia Y., Deltour, J. Comments on smoothing and differentiation of data by simplified least square procedure. *Analytical Chemistry* 44 (11): 1906-1909, 1972.

See Also

[g1oMoId](#), [gPoMo](#), [poLabs](#)

| | |
|--------|---|
| d2pMax | <i>Provides the number of polynomial terms pMax given dMax and nVar</i> |
|--------|---|

Description

Computes the number of polynomial terms pMax used to formulate an equation given the maximal polynomial degree dMax and the number of variables nVar following the conventions as defined by fuction poLabs.

Usage

```
d2pMax(nVar, dMaxKnown)
```

Arguments

| | |
|-----------|---|
| nVar | Number of variables considered in the polynomial formulation. |
| dMaxKnown | The maximum polynomial degree dMax |

Value

The number pMax of polynomial terms used to code a polynomial equation

Author(s)

Sylvain Mangiarotti

See Also

[g1oMoId](#), [gPoMo](#), [poLabs](#)

Examples

```
#####  
# Example 1 #  
#####  
# Maximum polynomial degree ?  
# number of variables:  
nVar <- 3  
# polynomial degree:  
dMax <- 3  
# The maximal polynomial degree used for coding the polynomial is:  
d2pMax(nVar,dMax)
```

data_vignetteIII data set

Output of the vignette III_Modelling

Description

To reduce the computation time, the outputs of the simulations presented in vignette VI have been run beforehand and saved in this file.

Usage

```
data_vignetteIII
```

Format

An object of class list of length 12.

Author(s)

Sylvain Mangiarotti, Mireille Huc.

data_vignetteVI data set

Output of the vignette VI_Sensitivity

Description

To reduce the computation time, the outputs of the simulations presented in vignette VI have been run beforehand and saved in this file.

Usage

```
data_vignetteVI
```

Format

An object of class list of length 6.

Author(s)

Sylvain Mangiarotti, Mireille Huc.

data_vignetteVII data set

Output of the vignette VII_Retro-Modelling

Description

To reduce the computation time, the outputs of the simulations presented in vignette VII have been run beforehand and saved in this file.

Usage

data_vignetteVII

Format

An object of class list of length 29.

Author(s)

Sylvain Mangiarotti, Mireille Huc.

derivODE2

A subfonction for the numerical integration of polynomial equations provided in a generic form following the convention defined by function poLabs.

Description

This function provides the one step integration of polynomial Ordinary Differential Equations (ODE). This function requires the function ode (deSolve package).

Usage

derivODE2(t, x, K, regS = NULL)

Arguments

| | |
|------|--|
| t | All the dates for which the result of the numerical integration of the model must be provided |
| x | Current state vector (input from which the next state will be estimated) |
| K | A matrix providing the model description: each column corresponds to one equation which polynomial organisation is following the convention defined by function poLabs. |
| regS | Current states of each polynomial terms used in poLabs. These states can be deduced from the current state vector x (using the function regSeries). When available, it can be provided as an input to avoid unnecessary computation. |

Author(s)

Sylvain Mangiarotti

See Also

[numicano](#), [numinousy](#)

detectP1limCycl

Detection of limit cycles of period-1

Description

This algorithm aim to detect period-1 limit cycles from trajectories in the phase sapce considered in a bidimensional projection.

Usage

```
detectP1limCycl(data, LimCyclThreshold = 0.01, show = 2)
```

Arguments

| | |
|------------------|---|
| data | A matrix of the trajectory in a 2D space (if more than two columns are provided, only the two first columns are considered) |
| LimCyclThreshold | The detection threshold |
| show | Indicates the deepness of the feedback (from 0 to 2) |

Value

Indicates if a limit cycle is detected (1) or not (0)

Author(s)

Sylvain Mangiarotti

See Also[autoGPoMoTest](#)

| | |
|---------|---|
| drvSucc | <i>drvSucc : Computes the successive derivatives of a time series</i> |
|---------|---|

Description

Computes the successive derivatives from one single time series, using the Savitzky-Golay algorithm (1964).

Usage

```
drvSucc(tin = NULL, serie, nDeriv, weight = NULL, tstep = NULL,
        winL = 9)
```

Arguments

| | |
|--------|---|
| tin | Input date vector which length should correspond to the input time series. |
| serie | A single time series provided as a single vector. |
| nDeriv | The number of derivatives to be computed from the input time series. The resulting number of time series obtained in output will be $nDeriv + 1$. |
| weight | A vector providing the binary weighting function of the input data series (0 or 1). By default, all the values are set to 1. |
| tstep | Sampling time of the input time series. Used only if time vector tin is not provided. |
| winL | Number (exclusively odd number) of points of the local window used for computing the derivatives along the input time series. The Savitzky-Golay filter is used for this purpose [1,2]. |

Value

A list containing:

\$serie The original time serie

\$tin The time vector containing the dates corresponding to the original time series

\$tstep The time step (assumed to be regular)

\$tout The time vector of the output series

seriesDeriv A matrix containing the original time series (smoothed by the filtering process) in the first column and its $nDeriv + 1$ successive derivatives in the next ones. Note that winL values of the original time series will be lost, that is $(winL - 1)/2$ at the beginning and $(winL - 1)/2$ at the end of the time series due to a computation boundary effect).

Author(s)

Sylvain Mangiarotti, Mireille Huc

References

- [1] Savitzky, A.; Golay, M.J.E., Smoothing and Differentiation of Data by Simplified Least Squares Procedures. *Analytical Chemistry* 36 (8), 1627-1639, 1964.
 [2] Steinier J., Termonia Y., Deltour, J. Comments on smoothing and differentiation of data by simplified least square procedure. *Analytical Chemistry* 44 (11): 1906-1909, 1972.

See Also

[gloMoId](#), [gPoMo](#), [poLabs](#), [compDeriv](#)

Examples

```
#####
# Example 1 #
#####
# Generate a time series:
tin <- seq(0, 5, by = 0.01)
data <- 2 * sin(5*tin)
dev.new()
par(mfrow = c(3, 1))
# Compute its derivatives:
drv <- drvSucc(tin = tin, nDeriv = 2, serie = data, winL = 5)
#
# plot original and filtered series
plot(tin, data, type='l', col = 'black', xlab = 't', ylab = 'x(t)')
lines(drv$out, drv$seriesDeriv[,1], lty = 3, lwd = 3, col = 'green')
#
# analytic 1st derivative
firstD <- 10 * cos(5 * tin)
# plot both
plot(tin, firstD, type = 'l', col = 'black', xlab = 't', ylab = 'dx/dt')
lines(drv$out, drv$seriesDeriv[,2], lty = 3, lwd = 3, col = 'green')
#
# analytic 2nd derivative
scdD <- -50 * sin(5 * tin)
# plot both
plot(tin, scdD, type = 'l', col = 'black', xlab = 't', ylab = 'd2x/dt2')
lines(drv$out, drv$seriesDeriv[,3], lty=3, lwd = 3, col = 'green')

#####
# Example 2 #
#####
# load data:
data("Ross76")
tin <- Ross76[,1]
data <- Ross76[,2]
```

```

# Compute the derivatives
drvOut <- drvSucc(tin, data, nDeriv=4)
dev.new()
par(mfrow = c(3, 1))
# original and smoothed variable:
plot(drvOut$tin, drvOut$serie,
      type='p', cex = 1, xlab = 'time', ylab = 'x(t)')
lines(drvOut$tout, drvOut$seriesDeriv[,1], type='p', col='red')
lines(drvOut$tout, drvOut$seriesDeriv[,1], type='l', col='red')
# 1st derivative:
plot(drvOut$tout, drvOut$seriesDeriv[,2],
      type='p', col='red', xlab = 'time', ylab = 'dx(t)/dt')
lines(drvOut$tout, drvOut$seriesDeriv[,2], type='l', col='red')
# 2nd derivative:
plot(drvOut$tout, drvOut$seriesDeriv[,3],
      type='p', col='red', xlab = 'time', ylab = 'd2x(t)/dt2')
lines(drvOut$tout, drvOut$seriesDeriv[,3], type='l', col='red')

```

| | |
|-------------|--|
| findAllSets | <i>Find all possible sets of equation combinations considering an ensemble of possible equation.</i> |
|-------------|--|

Description

For each equation to be retrieved, an ensemble of potential formulation is given. For instance, if three possible formulations are provided for equation (1), one for equation (2) and two for equation (3). In this case, six (i.e. $3*1*2$) possible sets of equations can be obtained from these potential formulations. The aim of this program is to formulate all the potential systems from the individual formulations provided of the individual equations.

Usage

```
findAllSets(allFilt, nS = c(3), nPmin = 1, nPmax = 14)
```

Arguments

| | |
|---------|--|
| allFilt | A list with: (1) A matrix allFilt\$Xi of possible formulations for each equation (corresponding to variable Xi); And (2) a vector allFilt\$Npi providing the number of polynomial terms contained in each formulation. |
| nS | A vector providing the number of dimensions used for each input variables (see Examples 1 and 2). The dimension of the resulting model will be $nVar = \text{sum}(nS)$. |
| nPmin | Corresponds to the minimum number of parameters (and thus of polynomial term) allowed. |
| nPmax | Corresponds to the maximum number of parameters (and thus of polynomial) allowed. |

Author(s)

Sylvain Mangiarotti

See Also[autoGPoMoSearch](#)**Examples**

```
#####
# Example 1 #
#####
# We build an example
allFilt <- list()
# For equation 1 (variable X1)
allFilt$Np1 <- 1 # only one formulation with one single parameter
# For equation 2 (variable X2)
allFilt$Np2 <- c(3, 4) # two potential formulations, with respectively three and four parameters
# For equation 3 (variable X3)
allFilt$Np3 <- c(2, 4) # two potential formulations, with respectively two and four parameters
# Formulations for variables Xi:
# For X1:
allFilt$X1 <- t(as.matrix(c(0,0,0,1,0,0,0,0,0,0)))
# For X2:
allFilt$X2 <- t(matrix(c(0,-0.85,0,-0.27,0,0,0,0.46,0,0,
                        0,-0.64,0,0,0,0,0.43,0,0),
                      ncol=2, nrow=10))
# For X3:
allFilt$X3 <- t(matrix(c(0, 0.52, 0, -1.22e-05, 0, 0, 0.99, 5.38e-05, 0, 0,
                        0, 0.52, 0, 0, 0, 0, 0.99, 0, 0, 0),
                      ncol=2, nrow=10))
# From these individual we can retrieve all possible formulations
findAllSets(allFilt, nS=c(3), nPmin=1, nPmax=14)
# if only formulations with seven maximum number of terms are expected:
findAllSets(allFilt, nS=c(3), nPmin=1, nPmax=7)
```

gloMoId

*Global Model Identification***Description**

Algorithm for global modelling in polynomial and canonical formulation of Ordinary Differential Equations. Univariate Global modeling aims to obtain multidimensional models from single time series (Gouesbet & Letellier 1994, Mangiarotti et al. 2012). An example of such application can be found in Mangiarotti et al. (2014) For a multivariate application, see GPoMo (Mangiarotti 2015, Mangiarotti et al. 2016).

Example:

For a model dimension nVar=3, the global model will read:

$$\begin{aligned}dX1/dt &= X2 \\dX2/dt &= X3 \\dX3/dt &= P(X1, X2, X3).\end{aligned}$$

Usage

```
gloMoId(series, tin = NULL, dt = NULL, nVar = NULL, dMax = 1,
        weight = NULL, show = 1, filterReg = NULL, winL = 9)
```

Arguments

| | |
|-----------|---|
| series | The original data set: either a single vector corresponding to the original variable; Or a matrix containing the original variable in the first column and its successive derivatives in the next columns. In the latter case, for the construction of n-dimensional model, series should have $nVar + 1$ columns since one more derivative will be necessary to identify the model parameters. Variable nVar will be set equal to n. In the former case, that is when only a single vector is provided, the derivatives will be automatically recomputed. Therefore, the dimension nVar expected for the model has to be provided. |
| tin | Input date vector which length should correspond to the input time series. |
| dt | Sampling time of the input time series. |
| nVar | Number of variables considered in the polynomial formulation. |
| dMax | Maximum degree of the polynomial formulation. |
| weight | A vector providing the binary weighting function of the input data series (0 or 1). By default, all the values are set to 1. |
| show | Provide (2) or not (0-1) visual output during the running process. |
| filterReg | A vector that specifies the template for the equation structure (for one single equation). The convention defined by poLabs is used. Value is 1 if the regressor is available, 0 if it is not. |
| winL | Total number of points used for computing the derivatives of the input time series. This parameter will be used as an input in function drvSucc to compute the derivatives. |

Value

A list of five elements :

\$init The original time series and the successive derivatives used for the modeling.

\$filterReg The structure of the output model. Value is 1 if the regressor is available, 0 if it is not. The terms order is given by function poLabs.

\$K Values of the identified coefficients corresponding to the regressors defined in filterReg.

`$resTot` The variance of the residual signal of the model.

`$resSsMod` The variance of the residual signal of the closer submodels.

`$finalWeight` Weighting series after boundary values were removed.

Author(s)

Sylvain Mangiarotti, Laurent Drapeau, Mireille Huc

References

- [1] Gouesbet G., Letellier C., Global vector-field reconstruction by using a multivariate polynomial L2 approximation on nets, *Physical Review E*, 49 (6), 4955-4972, 1994.
- [2] Mangiarotti S., Coudret R., Drapeau L., & Jarlan L., Polynomial search and global modeling : Two algorithms for modeling chaos, *Physical Review E*, 86, 046205, 2012.
- [3] Mangiarotti S., Drapeau L. & Letellier C., Two chaotic models for cereal crops observed from satellite in northern Morocco. *Chaos*, 24(2), 023130, 2014.
- [4] Mangiarotti S., Low dimensional chaotic models for the plague epidemic in Bombay (1896-1911), *Chaos, Solitons & Fractals*, 81(A), 184-196, 2015.
- [5] Mangiarotti S., Peyre M. & Huc M., A chaotic model for the epidemic of Ebola Virus Disease in West Africa (2013-2016). *Chaos*, 26, 113112, 2016.

See Also

[gPoMo](#), [autoGPoMoSearch](#), [autoGPoMoTest](#), [poLabs](#)

Examples

```
#####
# Example 1 #
#####
# load data
data("Ross76")
tin <- Ross76[,1]
data <- Ross76[,2:3]

# Polynomial identification
reg <- gloMoId(data[0:500,2], dt=1/100, nVar=2, dMax=2, show=0)

#####
# Example 2 #
#####
# load data
data(NDVI)

# Definition of the Model structure
```

```

terms <- c(1, 0, 0, 0, 1, 1, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1, 1, 1, 1)
poLabs(3,3,terms==1)
reg <- gloMoId(NDVI [,1:1], dt=1/125, nVar=3, dMax=3,
              show=0, filterReg=terms==1)

## Not run:
#####
# Example 3 #
#####
# load data
data("Ross76")
# time vector
tin <- Ross76[1:500,1]
# single time series
series <- Ross76[1:500,3]
# some noise is added
series[1:100] <- series[1:100] + 0.01 * runif(1:100, min = -1, max = 1)
series[301:320] <- series[301:320] + 0.05 * runif(1:20, min = -1, max = 1)
# weighting function
W <- tin * 0 + 1
W[1:100] <- 0 # the first hundred values will not be considered
W[301:320] <- 0 # twenty other values will not be considered either
reg <- gloMoId(series, dt=1/100, weight = W, nVar=3, dMax=2, show=1)
visuEq(3, 2, reg$K, approx = 4)
# first weight which value not equal to zero:
i1 = which(reg$finalWeight == 1)[1]
v0 <- reg$init[i1,1:3]

reconstr <- numicano(nVar=3, dMax=2, Istep=5000, onestep=1/250, PolyTerms=reg$K,
                  v0=v0, method="ode45")
plot(reconstr$reconstr[,2], reconstr$reconstr[,3], type='l', lwd = 3,
     main='phase portrait', xlab='time t', ylab = 'x(t)', col='orange')

# original data:
lines(reg$init[,1], reg$init[,2], type='l',
      main='phase portrait', xlab='x', ylab = 'dx/dt', col='black')
# initial condition
lines(v0[1], v0[2], type = 'p', col = 'red')

## End(Not run)

```

Description

Algorithm for a Generalized Polynomial formulation of multivariate Global Modeling. Global modeling aims to obtain multidimensional models from single time series [1-2]. In the generalized

(polynomial) formulation provided in this function, it can also be applied to multivariate time series [3-4].

Example:

Note that nS provides the number of dimensions used from each variable

case I

For nS=c(2, 3) means that 2 dimensions are reconstructed from variable 1: the original variable X1 and its first derivative X2), and 3 dimensions are reconstructed from variable 2: the original variable X3 and its first and second derivatives X4 and X5. The generalized model will thus be such as:

$$dX1/dt = X2$$

$$dX2/dt = P1(X1, X2, X3, X4, X5)$$

$$dX3/dt = X4$$

$$dX4/dt = X5$$

$$dX5/dt = P2(X1, X2, X3, X4, X5).$$

case II

For nS=c(1, 1, 1, 1) means that only the original variables X1, X2, X3 and X4 will be used. The generalized model will thus be such as:

$$dX1/dt = P1(X1, X2, X3, X4)$$

$$dX2/dt = P2(X1, X2, X3, X4)$$

$$dX3/dt = P3(X1, X2, X3, X4)$$

$$dX4/dt = P4(X1, X2, X3, X4).$$

Usage

```
gPoMo(data, tin = NULL, dtFixe = NULL, dMax = 2, nS = c(3), winL = 9,
       weight = NULL, show = 1, verbose = 1, underSamp = NULL, EqS = NULL,
       IstepMin = 2, IstepMax = 2000, nPmin = 1, nPmax = 14,
       method = "lsoda")
```

Arguments

| | |
|--------|--|
| data | Input Time series: Each column is one time series that corresponds to one variable. |
| tin | Input date vector which length should correspond to the input time series. |
| dtFixe | Time step used for the analysis. It should correspond to the sampling time of the input data. Note that for very large and very small time steps, alternative units may be used in order to stabilize the numerical computation. |
| dMax | Maximum degree of the polynomial formulation. |
| nS | A vector providing the number of dimensions used for each input variables (see Examples 1 and 2). The dimension of the resulting model will be nVar = sum(nS). |
| winL | Total number of points used for computing the derivatives of the input time series. This parameter will be used as an input in function drvSucc to compute the derivatives. |
| weight | A vector providing the binary weighting function of the input data series (0 or 1). By default, all the values are set to 1. |

| | |
|-----------|---|
| show | Provide (2) or not (0-1) visual output during the running process. |
| verbose | Gives information (if set to 1) about the algorithm progress and keeps silent if set to 0. |
| underSamp | Number of points used for undersampling the data. For <code>undersamp = 1</code> the complete time series is used. For <code>undersamp = 2</code> , only one data out of two is kept, etc. |
| EqS | Model template including all allowed regressors. Each column corresponds to one equation. Each line corresponds to one polynomial term as defined by function <code>poLabs</code> . |
| IstepMin | The minimum number of integration step to start of the analysis (by default <code>IstepMin = 10</code>). |
| IstepMax | The maximum number of integration steps for stopping the analysis (by default <code>IstepMax = 10000</code>). |
| nPmin | Corresponds to the minimum number of parameters (and thus of polynomial term) allowed. |
| nPmax | Corresponds to the maximum number of parameters (and thus of polynomial) allowed. |
| method | The integration technique used for the numerical integration. By default, the fourth-order Runge-Kutta method (<code>method = 'rk4'</code>) is used. Other methods such as <code>'ode45'</code> or <code>'lsoda'</code> may also be chosen. See package <code>deSolve</code> for details. |

Value

A list containing:

`$tin` The time vector of the input time series

`$inputdata` The input time series

`$tfiltdata` The time vector of the filtered time series (boundary removed)

`$filtdata` A matrix of the filtered time series with its derivatives

`$okMod` A vector classifying the models: diverging models (0), periodic models of period-1 (-1), unclassified models (1).

`$coeff` A matrix with the coefficients of one selected model

`$models` A list of all the models to be tested `$mToTest1`, `$mToTest2`, etc. and all selected models `$model1`, `$model2`, etc.

`$tout` The time vector of the output time series (vector length corresponding to the longest numerical integration duration)

`$stockoutreg` A list of matrices with the integrated trajectories (variable X1 in column 1, X2 in 2, etc.) of all the models `$model1`, `$model2`, etc.

Author(s)

Sylvain Mangiarotti, Flavie Le Jean, Mireille Huc

References

- [1] Gouesbet G. & Letellier C., 1994. Global vector-field reconstruction by using a multivariate polynomial L2 approximation on nets, *Physical Review E*, 49 (6), 4955-4972.
- [2] Mangiarotti S., Coudret R., Drapeau L. & Jarlan L., Polynomial search and Global modelling: two algorithms for modeling chaos. *Physical Review E*, 86(4), 046205.
- [3] Mangiarotti S., Le Jean F., Huc M. & Letellier C., Global Modeling of aggregated and associated chaotic dynamics. *Chaos, Solitons and Fractals*, 83, 82-96.
- [4] S. Mangiarotti, M. Peyre & M. Huc, 2016. A chaotic model for the epidemic of Ebola virus disease in West Africa (2013-2016). *Chaos*, 26, 113112.

See Also

[gloMoId](#), [autoGPoMoSearch](#), [autoGPoMoTest](#)
[autoGPoMoSearch](#), [autoGPoMoTest](#), [visuOutGP](#), [poLabs](#), [predictab](#), [drvSucc](#)

Examples

```
#Example 1
data("Ross76")
tin <- Ross76[,1]
data <- Ross76[,3]
dev.new()
out1 <- gPoMo(data, tin=tin, dMax = 2, nS=c(3), show = 1,
              IstepMax = 1000, nPmin = 9, nPmax = 11)
visuEq(3, 2, out1$models$model1, approx = 4)

## Not run:
#Example 2
data("Ross76")
tin <- Ross76[,1]
data <- Ross76[,2:4]
dev.new()
out2 <- gPoMo(data, tin=tin, dMax = 2, nS=c(1,1,1), show = 1,
              IstepMin = 10, IstepMax = 3000, nPmin = 7, nPmax = 8)
# the simplest model able to reproduce the observed dynamics is model #5
visuEq(3, 2, out2$models$model5, approx = 4) # the original Rossler system is thus retrieved

## End(Not run)

## Not run:
#Example 3
data("Ross76")
tin <- Ross76[,1]
data <- Ross76[,2:3]
# model template:
EqS <- matrix(1, ncol = 3, nrow = 10)
EqS[,1] <- c(0,0,0,1,0,0,0,0,0,0)
EqS[,2] <- c(1,1,0,1,0,1,1,1,1,1)
EqS[,3] <- c(0,1,0,0,0,0,1,1,0,0)
visuEq(3, 2, EqS, substit = c('X','Y','Z'))
```

```

dev.new()
out3 <- gPoMo(data, tin=tin, dMax = 2, nS=c(2,1), show = 1,
             EqS = EqS, IstepMin = 10, IstepMax = 2000,
             nPmin = 9, nPmax = 11)

## End(Not run)

## Not run:
## Example 4
# load data
data("TSallMod_nVar3_dMax2")
#multiple (six) time series
tin <- TSallMod_nVar3_dMax2$SprK$reconstr[1:400,1]
TSRo76 <- TSallMod_nVar3_dMax2$R76$reconstr[,2:4]
TSSprK <- TSallMod_nVar3_dMax2$SprK$reconstr[,2:4]
data <- cbind(TSRo76,TSSprK)[1:400,]
dev.new()
# generalized Polynomial modelling
out4 <- gPoMo(data, tin = tin, dMax = 2, nS = c(1,1,1,1,1,1),
             show = 0, method = 'rk4',
             IstepMin = 2, IstepMax = 3,
             nPmin = 13, nPmax = 13)

# the original Rossler (variables x, y and z) and Sprott (variables u, v and w)
# systems are retrieved:
visuEq(6, 2, out4$models$model347, approx = 4,
      substit = c('x', 'y', 'z', 'u', 'v', 'w'))
# to check the robustness of the model, the integration duration
# should be chosen longer (at least IstepMax = 4000)

## End(Not run)

```

GSproc

Gram-Schmidt procedure

Description

Computes regressors coefficients using the Gram-Schmidt procedure.

Usage

```
GSproc(polyK, ivec, weight = NULL)
```

Arguments

polyK One list including $\$Y$ and $\$phy$ with: $\$Y$ a matrix for which the i th column will be used to add one orthogonal vector to the $(i-1)$ th vectors of the current orthogonal base; and $\$phy$ such as the current orthogonal base is given by the $(i-1)$ th first columns of matrix $\text{polyK}\$phy$.

| | |
|--------|---|
| ivec | Defines i , the current vector of <code>polyK\$Y</code> and the current orthogonal base of <code>pParam\$phy</code> . |
| weight | The weighing vector. |

Value

`uNew` The model parameterization, that is: The residual orthogonal vector that can be included into the current orthogonal base. If the current base is empty, `uNew` is equal to the input vector of `$Y`; if the base is complete, `uNew` equals 0.

Author(s)

Sylvain Mangiarotti

NDVI

A time series of vegetation index measured from satellite

Description

A time series of 28 years of Normalized Difference Vegetation Index measured from space by the Advanced Very High Resolution Radiometer (AVHRR) sensor from 1982 to 2008 (see reference [1] for details).

Usage

NDVI

Format

An object of class `data.frame` with 9618 rows and 4 columns.

Author(s)

Sylvain Mangiarotti, Flavie Le Jean

References

[1] Mangiarotti S., Drapeau L. & Letellier C., 2014. Two chaotic models for cereal crops observed from satellite in northern Morocco.

 numicano

Numerical Integration of models in ODE of polynomial form

Description

Function for the numerical integration of Ordinary Differential Equations of polynomial form.

Usage

```
numicano(nVar, dMax, Istep = 1000, onestep = 1/125, KL = NULL,
        PolyTerms = NULL, v0 = NULL, method = "rk4")
```

Arguments

| | |
|-----------|--|
| nVar | Number of variables considered in the polynomial formulation. |
| dMax | Maximum degree of the polynomial formulation. |
| Istep | The number of integration time steps |
| onestep | Time step length |
| KL | Matrix formulation of the model to integrate numerically |
| PolyTerms | Vectorial formulation of the model (only for models of canonical form) |
| v0 | The initial conditions (a vector which length should correspond to the model dimension nVar) |
| method | The integration method (See package deSolve), by default method = 'rk4'. |

Value

A list of two variables:

`$KL` The model in its matrix formulation

`$reconstr` The integrated trajectory (first column is the time, next columns are the model variables)

Author(s)

Sylvain Mangiarotti

See Also

[derivODE2](#), [numinoisy](#)

Examples

```
#####
# Example 1 #
#####
# For a model of general form (here the rossler model)
# model dimension:
nVar = 3
# maximal polynomial degree
dMax = 2
# Number of parameter number (by default)
pMax <- d2pMax(nVar, dMax)
# convention used for the model formulation
poLabs(nVar, dMax)
# Definition of the Model Function
a = 0.520
b = 2
c = 4
Eq1 <- c(0,-1, 0,-1, 0, 0, 0, 0, 0, 0)
Eq2 <- c(0, 0, 0, a, 0, 0, 1, 0, 0, 0)
Eq3 <- c(b,-c, 0, 0, 0, 0, 0, 1, 0, 0)
K <- cbind(Eq1, Eq2, Eq3)
# Edition of the equations
visuEq(nVar, dMax, K)
# initial conditions
v0 <- c(-0.6, 0.6, 0.4)
# model integration
reconstr <- numicano(nVar, dMax, Istep=1000, onestep=1/50, KL=K,
                    v0=v0, method="ode45")
# Plot of the simulated time series obtained
dev.new()
plot(reconstr$reconstr[,2], reconstr$reconstr[,3], type='l',
     main='phase portrait', xlab='x(t)', ylab = 'y(t)')

## Not run:
#####
# Example 2 #
#####
# For a model of canonical form
# model dimension:
nVar = 4
# maximal polynomial degree
dMax = 3
# Number of parameter number (by default)
pMax <- d2pMax(nVar, dMax)
# Definition of the Model Function
PolyTerms <- c(281000, 0, 0, 0, -2275, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0,
              861, 0, 0, 0, -878300, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0)
# terms used in the model
poLabs(nVar, dMax, PolyTerms!=0)
# initial conditions
v0 <- c(0.54, 3.76, -90, -5200)
# model integration
```

```

reconstr <- numicano(nVar, dMax, Istep=500, onestep=1/250, PolyTerms=PolyTerms,
                    v0=v0, method="ode45")
# Plot of the simulated time series obtained
plot(reconstr$reconstr[,2], reconstr$reconstr[,3], type='l',
      main='phase portrait', xlab='x', ylab = 'dx/dt')
# Edition of the equations
visuEq(nVar, dMax, reconstr$KL)

## End(Not run)

```

| | |
|-----------|---|
| numinoisy | <i>Generates time series of deterministic-behavior with stochastic perturbations (measurement and/or dynamical noise)</i> |
|-----------|---|

Description

Generates time series from Ordinary Differential Equations perturbed by dynamical and/or measurement noises

Usage

```

numinoisy(x0, t, K, varData = NULL, txVarBruitA = NULL,
          txVarBruitM = NULL, varBruitA = NULL, varBruitM = NULL, taux = NULL,
          freq = NULL, variables = NULL, method = NULL)

```

Arguments

| | |
|-------------|--|
| x0 | The initial conditions. Should be a vector which size must be equal to the model dimension $\dim(K)[2]$ (the number of variables of the model defined by matrix K). |
| t | A vector providing all the dates for which the output are expected. |
| K | The Ordinary Differential Equations used to model the dynamics. The number of column should correspond to the number of variables, the number of lines to the number of parameters following the convention defined by <code>poLabs(nVar, dMax)</code> . |
| varData | A vector of size nVar providing the characteristic variances of each variable of the dynamical systems in ODE defined by matrix K. If not provided, this variance is automatically estimated. |
| txVarBruitA | A vector defining the ratio of ADDITIVE noise for each variable of the dynamical system in ODE. The additive noise is added at the end of the numerical integration process. The ratio is defined relatively to the signal variance of each variable. |
| txVarBruitM | A vector defining the ratio of DYNAMICAL noise for each variable of the dynamical system in ODE. This noise is a perturbation added at each numerical integration step. The ratio is defined relatively to the signal variance of each variable. |

| | |
|-----------|---|
| varBruitA | A vector defining the variance of ADDITIVE noise for each variable of the dynamical system in ODE. The additive noise is added at the end of the numerical integration process. |
| varBruitM | A vector defining the variance of DYNAMICAL noise for each variable of the dynamical system in ODE. This noise is a perturbation added at each numerical integration step. |
| taux | Generates random gaps in time series. Parameter <i>taux</i> defines the ratio of data to be kept (e.g. for <i>taux</i> = 0.75, 75 percents of the data are kept). |
| freq | Subsamples the time series. Parameter <i>freq</i> defines the periodicity of data kept (e.g. for <i>freq</i> = 3, 1 data out of 3 is kept). |
| variables | Defines which variables must be generated. |
| method | Defines the numerical integration method to be used. The fourth-order Runge-Kutta method is used by default (<i>method</i> = 'rk4'). Other method may be used (such as 'ode45' or 'lsoda'), see function <i>ode</i> from package <i>deSolve</i> for details. |

Value

A list of two variables:

\$donnees The integrated trajectory (first column is the time, next columns are the model variables)

\$bruitM The level of dynamical noise

\$bruitA The level of additive noise

\$vectBruitM The vector of the dynamical noise used to produce the time series

\$vectBruitA The vector of the additive noise used to produce the time series

\$cart_type The level standard deviation

Author(s)

Sylvain Mangiarotti, Malika Chassan

Examples

```
#####
# Example 1 #
#####
# Rossler Model formulation
# The model dimension
nVar = 3
# maximal polynomial degree
```

```

dMax = 2
a = 0.520
b = 2
c = 4
Eq1 <- c(0,-1, 0,-1, 0, 0, 0, 0, 0, 0)
Eq2 <- c(0, 0, 0, a, 0, 0, 1, 0, 0, 0)
Eq3 <- c(b,-c, 0, 0, 0, 0, 0, 1, 0, 0)
K <- cbind(Eq1, Eq2, Eq3)
# Edit the equations
visuEq(nVar, dMax, K)
# initial conditions
v0 <- c(-0.6, 0.6, 0.4)
# output time required
timeOut = (0:1000)/50
# variance of additive noise
varBruitA = c(0,0,0)^2
# variance of multiplitive noise
varBruitM = c(2E-2, 0, 2E-2)^2
# numerical integration with noise
integr <- numinoisy(v0, timeOut, K, varBruitA = varBruitA, varBruitM = varBruitM, freq = 1)
# Plot of the simulated time series obtained
dev.new()
plot(integr$donnees[,2], integr$donnees[,3], type='l',
     main='phase portrait', xlab='x(t)', ylab = 'y(t)')
dev.new()
par(mfrow = c(3, 1))
plot(integr$donnees[,1], integr$donnees[,2], type='l',
     main='phase portrait', xlab='x(t)', ylab = 'y(t)')
lines(integr$donnees[,1], integr$vectBruitM[,2]*10, type='l',
     main='phase portrait', xlab='x(t)', ylab = 'e(t)*10', col='red')
plot(integr$donnees[,1], integr$donnees[,3], type='l',
     main='phase portrait', xlab='x(t)', ylab = 'y(t)')
lines(integr$donnees[,1], integr$vectBruitM[,3]*10, type='l',
     main='phase portrait', xlab='x(t)', ylab = 'e(t)*10', col='red')
plot(integr$donnees[,1], integr$donnees[,4], type='l',
     main='phase portrait', xlab='x(t)', ylab = 'y(t)')
lines(integr$donnees[,1], integr$vectBruitM[,4]*10, type='l',
     main='phase portrait', xlab='x(t)', ylab = 'e(t)*10', col='red')

```

odeBruitMult2

For the numerical integration of ordinary differential equations with dynamical noise.

Description

A subfunction for the numerical integration of Ordinary Differential Equations provided in a generic polynomial form. Model formulation follows the convention defined by function `poLabs`.

Usage

```
odeBruitMult2(x0, t, K, varData = NULL, txVarBruitM = NULL,
              varBruitM = NULL, method = NULL)
```

Arguments

| | |
|-------------|--|
| x0 | Initial conditions |
| t | All the dates for which the result of the numerical integration of the model must be provided |
| K | A matrix providing the model description: each column corresponds to one equation which polynomial organisation is following the convention defined by function poLabs. |
| varData | A vector of size nVar providing the characteristic variances of each variable of the dynamical systems in ODE defined by matrix K. If not provided, this variance is automatically estimated. |
| txVarBruitM | A vector defining the ratio of DYNAMICAL noise for each variable of the dynamical system in ODE. This noise is a perturbation added at each numerical integration step. The ratio is defined relatively to the signal variance of each variable. |
| varBruitM | A vector defining the variance of DYNAMICAL noise for each variable of the dynamical system in ODE. This noise is a perturbation added at each numerical integration step. |
| method | Numerical method used in the integration process. (see ode function in deSolve package for details). |

Author(s)

Sylvain Mangiarotti, Malika Chassan

See Also

[numinoisy](#)

| | |
|--------|--|
| p2dMax | <i>p2dMax : provides the maximum polynomial degree dMax given the number of variables nVar and the number of possible polynomial terms pMax.</i> |
|--------|--|

Description

Find the maximum polynomial degree dMax given the number of polynomial terms pMax and the system dimension nVar.

Usage

```
p2dMax(nVar, pMaxKnown)
```

Arguments

nVar Number of variables considered in the polynomial formulation.
 pMaxKnown The number of polynomial terms

Value

dMax The maximum polynomial degree

Author(s)

Sylvain Mangiarotti, Laurent Drapeau

See Also

[gloMoId](#), [gPoMo](#), [poLabs](#)

Examples

```
#####
# Example 1 #
#####
# Maximum polynomial degree ?
# number of variables:
nVar <- 3
# size of the polynomial vector:
pMax <- 10
# The maximal polynomial degree used for coding the polynomial is:
p2dMax(nVar,pMax)

#####
# Example 2 #
#####
# for pMax = 462 and nVar = 6, then dMax is:
p2dMax(6,462)
# indeed:
length(poLabs(nVar=6, dMax=5))
```

 paramId

For parameter Identification

Description

Estimate the polynomial coefficients.

Usage

```
paramId(allForK, drv, weight)
```

Arguments

| | |
|---------|--|
| allFork | The list of input parameters |
| drv | The derivative (on the equation left hand) |
| weight | The weighting series |

Value

allFork The initial list completed with the model parameters.

Author(s)

Sylvain Mangiarotti

poLabs

Polynomial labels order

Description

Defines the order of the polynomial labels given the number of variables nVar and the maximum polynomial degree dMax.

Usage

```
poLabs(nVar, dMax, findIt = NULL, Xnote = "X")
```

Arguments

| | |
|--------|--|
| nVar | The number of variables |
| dMax | The maximum degree allowed in the formulation |
| findIt | A vector of selected terms. |
| Xnote | Enables to defines the notation used for the variable, by default Xnote = 'X'. |

Value

lbls A vector of characters. Each element is the expression of one polynomial term, such as $X_1^2 X_3 X_4$

Author(s)

Sylvain Mangiarotti

See Also

[visuEq](#)

Examples

```

#Regressor order for three variables \eqn{(X1,X2,X3)} (nVar = 3) for a maximum
#polynomial degree equal to 2 (dMax = 2): poLabs(3,2)
#and for two variables only : poLabs(2,2)

# For a quadratic equation of two variables,
# the polynomial \deqn{P(X1,X2) = 0.5 + 0.3 X1 -0.25 X1 X2}
# could thus be written as a vector Pvec such as:
Pvec = c(0.5, 0, 0, 0.3, -0.25, 0)
# considering the convention corresponding to
poLabs(2,2)
# Indeed:
poLabs(2, 2, findIt = Pvec!=0)
# An alternative notation can be used with parameter Xnote
poLabs(2, 2, findIt = Pvec!=0, Xnote = 'w')
# or also
poLabs(2, 2, findIt = Pvec!=0, Xnote = c('x','y'))

```

| | |
|-----------|--|
| predictab | <i>Estimate the models performance obtained with GPoMo in term of predictability</i> |
|-----------|--|

Description

The algorithm aims to estimate automatically the forecasting performances of the models obtained with gPoMo.

Usage

```

predictab(ogp, fullt = NULL, fulldata = NULL, hp = NULL, Nech = 50,
  show = 1, selecmod = NULL, id = 1, selV = 1, na.rm = FALSE)

```

Arguments

| | |
|----------|---|
| ogp | The output list obtained from function gPoMo. |
| fullt | Time vector of the data set for which predictability will be tested |
| fulldata | Data set for which predictability will be tested |
| hp | Time vector of the horizon of prediction |
| Nech | Number of simulations |
| show | Provide (2) or not (0-1) visual output during the running process. |
| selecmod | A vector of the model selected. |
| id | The type of model to identify. id = 1 corresponds to unidentified models, that is, potentially chaotic. |
| selV | Selected variable for the analysis |
| na.rm | Indicates if the NA should be removed (na.rm = TRUE) or not (na.rm = FALSE). |

Value

ErrmodAll A list of matrix \$Predmod1, \$Predmod2, etc. and \$Errmod1, \$Errmod2, etc. providing respectively the forecasting and the forecasting error of models 1, 2, etc. Each column corresponds to one simulation starting from a specific initial condition. Each line corresponds to one horizon of prediction. Vectors corresponding to the initial condition time tE and the horizon of prediction hpE are also provided in \$tE and \$hpE, respectively.

Author(s)

Sylvain Mangiarotti, Mireille Huc

Examples

```
# load data
data("Ross76")
# time vector
tin <- Ross76[seq(1, 3000, by = 8), 1]
# single time series
data <- Ross76[seq(1, 3000, by = 8), 3]
# dev.new()
# plot(tin, data, xlab = 'time', ylab = 'y(t)')

# global modelling
# results are put in list outputGPoM
outputGPoM <- gPoMo(data[1:300], tin = tin[1:300], dMax = 2, nS=c(3),
                    show = 0, method = 'rk4',
                    nPmin = 10, nPmax = 12,
                    IstepMin = 150, IstepMax = 151)
#
visuOutGP(outputGPoM)

#####
# and test predictability #
#####
outpred <- predictab(outputGPoM, hp = 15, Nech = 30)

# manual visualisation of the outputs (e.g. for model 1):
dev.new()
image(outpred$tE, outpred$hpE, t(outpred$Errmod1),
      xlab = 't', ylab = 'hp', main = 'Errmod1')
```

Description

Generate the conventional order of the polynomial terms for the polynomial description. It is formulated as a matrix of exponents: Each column of the matrix (a,b,c, ...) corresponds to a product of the nVar available variables X1, X2, X3, etc., that is, $X1^a X2^b X3^c$, etc.

Usage

```
regOrd(nVar, dMax)
```

Arguments

| | |
|------|---|
| nVar | The number of variables |
| dMax | The maximum degree allowed in the formulation |

Value

A matrix of exponents. Each column corresponds to one polynomial term. Each line correspond to the exponent of one variable. For example, a column of three exponents (0, 2, 1) corresponds to the monomial $X1^0 * X2^2 * X3^1$, that is $X2^2 X3$.

Author(s)

Sylvain Mangiarotti

See Also

[poLabs](#)

regSeries

Estimates the monomial time series

Description

Creates time series by multiplying given time series among them.

Usage

```
regSeries(nVar, dMax, series, pReg = NULL)
```

Arguments

| | |
|--------|---|
| nVar | Number of variables considered in the polynomial formulation. |
| dMax | Maximum degree of the polynomial formulation. |
| series | A matrix containing the original time series from which the monomials are built. Each column corresponds to one given variable. |
| pReg | A matrix filled, for each column, with powers of time series used to create. |

Value

rpFull A matrix of time series. Each column corresponds to one regressor such as $X_1^2 X_3 X_4$

Author(s)

Sylvain Mangiarotti

Examples

```
data(TSallMod_nVar3_dMax2)
sprttK <- as.matrix(TSallMod_nVar3_dMax2$SprK$reonstr)[,2:4]
dMax <- 2
nVar <- dim(sprttK)[2]

#Example 1
polySeries2 <- regSeries(nVar, dMax, sprttK)

#Example 2
p <- c(1,3,1)
polySeries2 <- regSeries(nVar, dMax, sprttK, pReg=p)
```

Rossler-1976 data set *Time series of the Rossler-1976 system*

Description

The Rössler system is the 3-dimensional chaotic system

$$dx/dt = -y - z$$

$$dy/dt = x + ay$$

$$dz/dt = b + z(x - c),$$

discovered by Otto E. Rössler in 1976 [1]. The following parameters and initial conditions were used to produce the present data set:

$$a = 0.520, b = 2, c = 4$$

$$\text{and } (x_0, y_0, z_0) = (-0.04298734, 1.025536, 0.09057987).$$

The following four columns are provided:

(1) time t, (2) x(t), (3) y(t) and (4) z(t).

For this parameterization, the Rössler system produces a chaotic behavior characterized by a regime non-coherent in phase (oscillations duration can be very different from one oscillation to another).

Usage

Ross76

Format

An object of class deSolve (inherits from matrix) with 4000 rows and 4 columns.

Author(s)

Sylvain Mangiarotti, Flavie Le Jean, Malika Chassan, Laurent Drapeau, Mireille Huc.

References

[1] O. Rössler, 1976. An Equation for Continuous Chaos, Physics Letters, 71A, 2-3, 155-157.

RosYco

Twelve Rossler-1976 time series (exclusively variable y)

Description

Twelve independant Rossler-1976 time series (variable y). The parameters used to generate the time series correspond to a phase coherent behavior. Details can be found in [1]

Usage

RosYco

Format

An object of class `matrix` with 3000 rows and 12 columns.

Details

Another set of time series of the Rossler-1976 chaotic system

Author(s)

Sylvain Mangiarotti, Flavie Le Jean.

References

[1] Mangiarotti S., Le Jean F., Huc M. & Letellier C., Global Modeling of aggregated and associated chaotic dynamics. Chaos, Solitons and Fractals, 83, 82-96.

TSallMod_nVar3_dMax2 data set

Time series of three-dimensional chaotic systems (for vignette VII_Retro-Modelling)

Description

A list of matrix providing the time series in a list named TSallMod_nVar3_dMax2 of eighteen three-dimensional chaotic systems: Lorenz-1963 (\$L63), Rössler-1976 (\$R76), Burke & shaw 1981 (\$BS81), Lorenz-1984 (\$L84), Nosé & Hooer 1986 (\$NH86), Genesio & Tosi 1992 (\$GT92), Spott systems 1994 (\$SprF, \$SprH, \$SprK, \$SprO, \$SprP, \$SprG, \$SprM, \$SprQ, \$SprS), Chlouverakis & Sprott 2004 (\$CS2004), Li 2007 (\$Li2007) and the Cord system by Aguirre & Letellier 2012 (\$Cord2012). Time series are provided in a matrix in which each column corresponds to one variable of the dynamical systems.

Usage

TSallMod_nVar3_dMax2

Format

An object of class list of length 18.

Author(s)

Sylvain Mangiarotti, Mireille Huc.

References

References for the systems are provided in vignette ‘VII_retro-modelling‘.

visuEq

Displays the models Equations

Description

Displays the model equations for a polynomial model which description is provided as a matrix K, each column corresponding to one equation. The coefficients of the polynomial terms are given following the order defined by function polAbs.

Usage

visuEq(nVar, dMax, K, substit = 0, approx = FALSE)

Arguments

| | |
|---------|---|
| nVar | The number of variables |
| dMax | The maximum degree allowed in the formulation |
| K | A matrix providing the model description: each column corresponds to one equation which polynomial organisation is following the convention defined by function poLabs. |
| substit | Applies substitutions to the default values: for <code>substit = 0</code> (default value), variables are chosen as X1, X2, ... for <code>substit = 1</code> , variable X1, X2, ... will be replaced by x, y, z, ... for <code>substit = 2</code> , the codes provides a LaTeX-like formulation of the model. The variables name can also be defined explicitly as follows: for <code>substit = c('x', 'H', 'T1')</code> , variables X1, X2, X3 ... will be replaced by x, H and T1. |
| approx | The number of extra digits to be used: for <code>approx = FALSE</code> (default value) digits are edited with double precision; for <code>approx = TRUE</code> , only the minimum number of digits is edited (in order to have all the terms different from 0) for <code>approx = 1, 2, etc.</code> then respectively 1, 2, etc. digits are added to the minimum number of digits corresponding to <code>approx = TRUE</code> . |

Author(s)

Sylvain Mangiarotti

Examples

```
#EQUATIONS VISUALISATION
# number of variables:
nVar <- 3
# maximum polynomial degree:
dMax <- 2
# polynomial organization:
poLabs(nVar,dMax)
# model construction
KL = matrix(0, ncol = 3, nrow = 10)
KL[1,1] <- KL[2,2] <- 1
KL[4,1] <- -1
KL[5,3] <- -0.123456789
# Equations visualisation:
# (a) by default, variables names X1, X2, X3 are used
visuEq(nVar, dMax, KL)
# (b) for substit=1, variables names x, y, z are used instead
visuEq(nVar, dMax, KL, approx = TRUE, substit=1)
# (c) the name of the variables can also be chosen manually
visuEq(nVar, dMax, KL, approx = 3, substit=c('U', 'V', 'W'))

# A canonical model can be provided as a single vector
polyTerms <- c(0.2,0,-1,0.5,0,0,0,0,0,0)
visuEq(3,2,KL)
```

visuOutGP

*visuOutGP : get a quick information of gPoMo output***Description**

The algorithm aims to get a quick information about the outputs obtained with gPoMo.

Usage

```
visuOutGP(ogp, selecmod = NULL, id = 1, prioMinMax = "data",
  opt3D = "TRUE")
```

Arguments

| | |
|------------|---|
| ogp | The output list obtained from gPoMo. |
| selecmod | A vector of the selected model. Maximum 24 models can be presented at the same time. |
| id | The type of model to identify. id = 1 corresponds to the unidentified models, that is, potentially chaotic models). |
| prioMinMax | Gives the priority for the plots among: "data", "model", "dataonly" and "modelonly". |
| opt3D | Provides a 3D plot (x,y,z) when opt = 'TRUE' (the rgl library is required). |

Value

A Matrix describing the terms composing each model by row. The first row corresponds to the model detection (1 unclarified, 2 diverging, 0 is fixed point, -n with n an integer, is period-n cycle')

Author(s)

Sylvain Mangiarotti

Examples

```
# load data
data("Ross76")
# # time vector
tin <- Ross76[seq(1, 3000, by = 8), 1]
# single time series
data <- Ross76[seq(1, 3000, by = 8), 3]
dev.new()
plot(tin, data, type = 'l')
# global modelling
# results are put in list outputGPoM
outputGPoM <- gPoMo(data, tin=tin, dMax = 2, nS=c(3), show = 0,
  nPmin = 9, nPmax = 12, method = 'rk4',
  IstepMin = 200, IstepMax = 201)
visuOutGP(outputGPoM)
```

| | |
|---------|-------------------------------|
| wInProd | <i>Weighted inner product</i> |
|---------|-------------------------------|

Description

Computes weighted inner products.

Usage

```
wInProd(A1, A2, weight = NULL)
```

Arguments

| | |
|--------|-----------------------|
| A1 | The input matrix 1. |
| A2 | The input matrix 2. |
| weight | The weighting vector. |

Value

inP The weighted inner product.

Examples

```
#####  
#Example 1 #  
#####  
A1 = c(0,1,2,0,1,3)  
A2 = c(1,2,0,0,4,1)  
wInProd(A1, A2)  
  
#####  
#Example 2 #  
#####  
A1 = c(0,1,2,0,1,3)  
A2 = c(1,2,0,0,4,1)  
w = c(0,0,0,1,1,1)  
wInProd(A1, A2, weight = w)
```

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