

# Package ‘localScore’

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**Type** Package

**Title** Package for Sequence Analysis by Local Score

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**Description** Functionalities for calculating the local score and calculating statistical relevance (p-value) to find a local Score in a sequence of given distribution (S. Mercier and J.-J. Daudin (2001) <<https://hal.archives-ouvertes.fr/hal-00714174>>) ; S. Karlin and S. Altschul (1990) <<https://www.ncbi.nlm.nih.gov/pmc/articles/PMC53667/>> ; S. Mercier, D. Cellier and F. Charlot (2003) <<https://hal.archives-ouvertes.fr/hal-00937529v1>> ; A. Lagnoux, S. Mercier and P. Valois (2017) <[doi:10.1093/bioinformatics/btw699](https://doi.org/10.1093/bioinformatics/btw699)> ).

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localScore-package      *Package for sequence analysis by local score*

---

### Description

Provides functionalities for:

- calculating the local score
- calculating statistical relevance (p-value) to find a local Score in a sequence of given distribution.

**Details**

Please refer to the vignette of this package or the manual for details on how to use this package.

**Author(s)**

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**References**

p-value:

- An Improved Approximation For Assessing The Statistical Significance of molecular Sequence Features, Mercier and al 2003
- Exact distribution for the local score of one i.i.d. random sequence, Sabine Mercier and JJ Daudin, 2001
- Limit Distributions of Maximal Segmental Score among Markov-Dependent Partial Sums, Karlin and Dembo 1992
- Methods for assessing the statistical significance of molecular sequence features by using general scoring schemes, Karlin and al 1990

local Score:

- Detection de courts segments inverses dans les genomes: methodes et applications, David Robelin 2005
- A Linear Time Algorithm for Finding All Maximal Scoring Subsequences, Constable and Bates 1985

---

automatic\_analysis      *Automatic analysis*

---

**Description**

Calculates local score and p-value for sequence(s) with integer scores.

**Usage**

```
automatic_analysis(  
  sequences,  
  model,  
  scores,  
  transition_matrix,  
  distribution,  
  method_limit = 2000,  
  score_extremes,  
  modelFunc,  
  simulated_sequence_length = 1000,  
  ...  
)
```

**Arguments**

sequences	sequences to be analysed (named list)
model	the underlying model of the sequence (either "iid" for identically independently distributed variable or "markov" for Markov chains)
scores	vector of minimum and maximum score range
transition_matrix	if the sequences are markov chains, this is their transition matrix
distribution	vector of probabilities in ascending score order (iid sequences). Note that names of the vector must be the associated scores.
method_limit	limit length from which on computation-intensive exact calculation methods for p-value are replaced by approximative methods
score_extremes	a vector with two elements: minimal score value, maximal score value
modelFunc	function to create similar sequences. In this case, Monte Carlo is used to calculate p-value
simulated_sequence_length	if a modelFunc is provided and the sequence happens to be longer than method_limit, the method karlinMonteCarlo is used. This method requires the length of the sequences that will be created by the modelFunc for estimation of Gumble parameters.
...	parameters for modelFunc

**Details**

This method picks the adequate p-value method for your input.

If no sequences are passed to this function, it will let you pick a FASTA file.

If this is the case, and if you haven't provided any score system (as you can do by passing a named list with the appropriate scores for each character), the second file dialog which will pop up is for choosing a file containing the score (and if you provide an extra column for the probabilities, they will be used, too - see section File Formats in the vignette for details).

The function then either uses empirical distribution based on your input - or if you provided a distribution, then yours - to calculate the p-value based on the length of each of the sequences given as input.

You can influence the choice of the method by providing the modelFunc argument. In this case, the function uses exclusively simulation methods (monte\_carlo, monte\_carlo\_karlin).

By setting the method\_limit you can further decide to which extent computation-intensive methods (daudin, exact\_mcc) should be used to calculate the p-value. Remark that the warnings of the localScoreC() function have been deleted when called by automatic\_analysis() function

**Value**

A list object containing

Local score	local score...
p-value	p-value ...
Method	the method used for the calculus of the p-value

**Examples**

```

# Minimal example
l = list()
seq1 = sample(-2:1, size = 3000, replace = TRUE)
seq2 = sample(-3:1, size = 150, replace = TRUE)
l[["hello"]] = seq1
l[["world"]] = seq2
automatic_analysis(l, "iid")
# Example with a given distribution
automatic_analysis(l,"iid",scores=c(-3,1),distribution=c(0.3,0.3,0.1,0.1,0.2))
# forcing the exact method for the longest sequence
aa1=automatic_analysis(l,"iid")
aa1$hello$`method applied`
aa1$hello$`p-value`
aa2=automatic_analysis(l,"iid",method_limit=3000)
aa2$hello$`method applied`
aa2$hello$`p-value`
# Markovian example
MyTransMat <-
matrix(c(0.3,0.1,0.1,0.1,0.4, 0.3,0.2,0.2,0.2,0.1, 0.3,0.4,0.1,0.1,0.1, 0.3,0.3,0.3,0.0,0.1,
        0.1,0.3,0.2,0.3,0.1), ncol = 5, byrow=TRUE)
MySeq.CM=transmatrix2sequence(matrix = MyTransMat,length=150, score =-2:2)
MySeq.CM2=transmatrix2sequence(matrix = MyTransMat,length=110, score =-2:2)
automatic_analysis(sequences = list("x1" = MySeq.CM, "x2" = MySeq.CM2), model = "markov")

```

---

CharSequence2ScoreSequence

*Convert a character sequence into a score sequence*

---

**Description**

Convert a character sequence into a score sequence. See CharSequences2ScoreSequences() fonction for several sequences

**Usage**

```
CharSequence2ScoreSequence(sequence, dictionary)
```

**Arguments**

sequence	a character sequence
dictionary	a dictionary

**Value**

a vector of a score sequence

**Examples**

```
data(ShortSeq)
ShortSeq
data(dico)
CharSequence2ScoreSequence(ShortSeq,dico)
```

---

CharSequences2ScoreSequences

*Convert several character sequences into score sequences*

---

**Description**

Convert several character sequence into score sequences. For only one sequence see CharSequence2ScoreSequence() function.

**Usage**

```
CharSequences2ScoreSequences(sequences, dictionary)
```

**Arguments**

sequences	a list of character sequences
dictionary	a dictionary

**Value**

a list of score sequences

**Examples**

```
data(ShortSeq)
ShortSeq
data(MidSeq)
MidSeq
data(dico)
MySequences=list("A1"=ShortSeq,"A2"=MidSeq)
CharSequences2ScoreSequences(MySequences,dico)
```

---

daudin	<i>Daudin [p-value] [iid]</i>
--------	-------------------------------

---

### Description

Calculates the exact p-value in the identically and independantly distributed of a given local score, a sequence length that 'must not be too large' and for a given score distribution

### Usage

```
daudin(  
  localScore,  
  sequence_length,  
  score_probabilities,  
  sequence_min,  
  sequence_max  
)
```

### Arguments

localScore	the observed local score
sequence_length	length of the sequence
score_probabilities	the probabilities for each score from lowest to greatest
sequence_min	minimum score
sequence_max	maximum score

### Details

Small in this context depends heavily on your machine. On a 3,7GHZ machine this means for `daudin(1000, 5000, c(0.2, 0.2, 0.2, 0.1, 0.2, 0.1), -2, 3)` an execution time of ~2 seconds. This is due to the calculation method using matrix exponentiation which becomes very fast very slow. The size of the matrix of the exponentiation is equal to  $a+1$  with a the local score value. The matrix must be put at the power  $n$ , with  $n$  the sequence length. Moreover, it is known that the local score value is expected to be in mean of order  $\log(n)$ .

### Value

A double representing the probability of a local score as high as the one given as argument

### Examples

```
daudin(localScore = 4, sequence_length = 50,  
score_probabilities = c(0.2, 0.3, 0.1, 0.2, 0.1, 0.1), sequence_min = -3, sequence_max = 2)
```

---

dico

*Dictionnaire*

---

### Description

Provides the score related to each base of the sequences.

### Usage

```
dico
```

### Format

A score function for the 20 amino acid

### Source

Kyte & Doolittle (1982) J. Mol. Biol. 157, 105-132

### Examples

```
data(dico)
dico
data(MidSeq)
MidSeq
MidSeqScore=CharSequence2ScoreSequence(MidSeq,dico)
MidSeqScore[1:30]
localScoreC(MidSeqScore)$localScore
```

---

exact\_mc

*Exact method for p-value [Markov chains]*

---

### Description

Calculates the exact p-value for short numerical Markov chains. Time computation can be too large for a sequence length of several thousands, specially for a data set.

### Usage

```
exact_mc(m, localScore, sequence_length, sequence_min, sequence_max)
```



**Arguments**

m                    Transition matrix [matrix object]  
 localScore        score for which the p-value should be calculated  
 sequence\_length   length of the sequence  
 sequence\_min     minimum score  
 sequence\_max     maximum score

**Value**

A double representing the probability of a localScore as high as the one given as argument

**Examples**

```

matrix = t(matrix(c(0.2, 0.3, 0.5, 0.3, 0.4, 0.3, 0.2, 0.4, 0.4), nrow = 3))
exact_mc(localScore = 12, m = matrix, sequence_length = 100, sequence_min = -1, sequence_max = 1)
exact_mc(localScore = 150, m = matrix, sequence_length = 1000, sequence_min = -1, sequence_max = 1)

```

---

karlin                    *Karlin [p-value] [iid]*

---

**Description**

Calculates an approximated p-value of a given local score value and a long sequence length in the identically and independantly distributed model for the sequence. See also mccc() function for another approximated method in the i.i.d. model

**Usage**

```

karlin(
  localScore,
  sequence_length,
  score_probabilities,
  sequence_min,
  sequence_max
)

```

**Arguments**

localScore        the observed local score  
 sequence\_length   length of the sequence (at least several hundreds)  
 score\_probabilities   the probabilities for each unique score from lowest to greatest  
 sequence\_min     minimum score  
 sequence\_max     maximum score

**Details**

This method works the better the longer the sequence is.

**Value**

A double representing the probability of a localScore as high as the one given as argument

**Examples**

```
karlin(150, 10000, c(0.08, 0.32, 0.08, 0.00, 0.08, 0.00, 0.00, 0.08, 0.02, 0.32, 0.02), -5, 5)
```

---

```
karlinMonteCarlo      Monte Carlo - Karlin [p-value]
```

---

**Description**

Estimates p-value, for integer scores, based on a Monte Carlo estimation of Gumble parameters from simulations of smaller sequences with same distribution. Appropriate for great sequences with length > 10<sup>3</sup>, for i.i.d and markovian sequence models.

**Usage**

```
karlinMonteCarlo(
  local_score,
  sequence_length,
  simulated_sequence_length,
  FUN,
  ...,
  numSim = 1000,
  plot = TRUE
)
```

**Arguments**

```
local_score      local score observed in a segment.
sequence_length  length of the sequence
simulated_sequence_length
                  length of simulated sequences produced by FUN
FUN              function to simulate similar sequences with.
...             parameters for FUN
numSim          number of sequences to create for estimation
plot            boolean value if to display plots for cumulated function and density
```

**Details**

The length of the simulated sequences is an argument specific to the function provided for simulation. Thus, it has to be provided also in the parameter `simulated_sequence_length` in the arguments of the "Monte Carlo - Karlin" function. It is a crucial detail as it influences precision and computation time of the result. Note that to get an appropriate estimation, the given average score must be non-positive.

**Value**

Floating value corresponding to the probability to obtain a local score with a value greater or equal to the parameter `local_score`

**Examples**

```
new = sample(-7:6, replace = TRUE, size = 1000)
#MonteCarlo taking random sample from the input sequence itself

karlinMonteCarlo(local_score = 66, sequence_length = 1000,
                 FUN = function(x, simulated_sequence_length) {return(sample(x = x,
                                     size = simulated_sequence_length, replace = TRUE))},
                 x=new, simulated_sequence_length = 1000, numSim = 1000)
```

---

karlinMonteCarlo\_double

*Monte Carlo - Karlin for real scores[p-value]*

---

**Description**

Estimates p-value, for integer scores, based on a Monte Carlo estimation of Gumble parameters from simulations of smaller sequences with same distribution. Appropriate for great sequences with length  $> 10^3$ , for i.i.d and markovian sequence models.

**Usage**

```
karlinMonteCarlo_double(
  local_score,
  sequence_length,
  simulated_sequence_length,
  FUN,
  ...,
  numSim = 1000,
  plot = TRUE
)
```

**Arguments**

local_score	local score observed in a segment.
sequence_length	length of the sequence
simulated_sequence_length	length of simulated sequences produced by FUN
FUN	function to simulate similar sequences with.
...	parameters for FUN
numSim	number of sequences to create for estimation
plot	boolean value if to display plots for cumulated function and density

**Details**

The length of the simulated sequences is an argument specific to the function provided for simulation. Thus, it has to be provided also in the parameter `simulated_sequence_length` in the arguments of the "Monte Carlo - Karlin" function. It is a crucial detail as it influences precision and computation time of the result. Note that to get an appropriate estimation, the given average score must be non-positive.

**Value**

Floating value corresponding to the probability to obtain a local score with value greater or equal the parameter

**Examples**

```
new = sample(-7:6, replace = TRUE, size = 1000)
#MonteCarlo taking random sample from the input sequence itself

karlinMonteCarlo_double(local_score = 66, sequence_length = 1000,
  FUN = function(x, simulated_sequence_length) {return(sample(x = x,
    size = simulated_sequence_length, replace = TRUE))},
  x=new, simulated_sequence_length = 1000, numSim = 1000)

#Markovian example (longer computation)
MyTransMat_reels <- matrix(c(0.3,0.1,0.1,0.1,0.4, 0.2,0.2,0.1,0.2,0.3, 0.3,0.4,0.1,0.1,0.1,
0.3,0.3,0.1,0.2,0.1, 0.2,0.1,0.2,0.4,0.1),ncol = 5, byrow=TRUE)

karlinMonteCarlo(local_score = 10.5,sequence_length=200,FUN = transmatrix2sequence,
matrix = MyTransMat_reels, score =c(-1,-0.5,0,0.5,1),length=1500,
plot=FALSE, numSim = 1500, simulated_sequence_length =1500)
```

---

lindley	<i>Lindley process</i>
---------	------------------------

---

**Description**

Creates a sequence of a Lindley process, also called CUSUM process, on a given sequence. For a sequence  $(X_k)_k$ , the Lindley process is defined as follows:  $W_0:=0$  and  $W_{(k+1)}=\max(0, W_k+X_{(k+1)})$ . It defines positive excursions above 0.

**Usage**

```
lindley(sequence)
```

**Arguments**

sequence            numeric sequence of a Lindley process, eg service time per customer

**Value**

a vector with the Lindley process steps

**Examples**

```
seq=c(1,2,3,-4,1,-3,-1,2,3,-4,1)
lindley(seq)
plot(1:length(seq),lindley(seq),type='b')
```

---

loadMatrixFromFile	<i>Loads matrix from csv-File</i>
--------------------	-----------------------------------

---

**Description**

Reads a csv file without header and returns the matrix. For file formats please see section "File Formats" in vignette.

**Usage**

```
loadMatrixFromFile(filepath)
```

**Arguments**

filepath            optional: Location of file on disk. If not provided, a file picker dialog will be opened.

**Value**

A Matrix Object

---

loadScoreFromFile	<i>Load score from file</i>
-------------------	-----------------------------

---

### Description

Reads a csv file with 2-3 columns and returns it as a list object of vectors, with names corresponding to the first column of the file. For details view section "File Formats" in vignette.

### Usage

```
loadScoreFromFile(filepath, ...)
```

### Arguments

filepath	optional: location of file on disk. If not provided, a file picker dialog will be opened.
...	optional: use arguments from read.csv

### Value

A List Object - Names correspond to the first column, usually Letters. If probabilities are provided, they will be loaded too and presumed to be in the third column

---

localScoreC	<i>Local score</i>
-------------	--------------------

---

### Description

Calculates the local score for a sequence of integer scores. Only provides the first occurrence of the local score. Use function suboptimalSegment() or Lindley() to obtain the others localizations of the different realizations of the local score.

### Usage

```
localScoreC(v, supressWarnings = FALSE)
```

### Arguments

v	: a sequence of integer values as vector.
supressWarnings	: if warnings should not be displayed

**Value**

A structure containing: the local score value and the begin and end index of the segment realizing this optimal score ; all the local maxima of the Lindley process (non negative excursion) and their begin and end index ; the record times of the Lindley process but only the ones corresponding to the begin index of non negative excursions

**Examples**

```
seq.OneSegment=c(1,-2,3,1,-1,2)
# one segment realizing the local score value
localScoreC(seq.OneSegment)
seq.TwoSegments=c(1,-2,3,1,2,-2,-2,-1,1,-2,3,1,2,-1,-2,-2,-1,1)
# two segments realizing the local score value
localScoreC(seq.TwoSegments)
# only the first realization
localScoreC(seq.TwoSegments)$localScore
# all the realization of the local together with the suboptimal ones
localScoreC(seq.TwoSegments)$suboptimalSegmentScores
# for small sequences, you can also use lindley() fonction to check if
# several segments achieve the local Score
lindley(seq.TwoSegments)
plot(1:length(seq.TwoSegments),lindley(seq.TwoSegments),type='b')
seq.TwoSegments.InSameExcursion=c(1,-2,3,2,-1,0,1,-2,-2,-4,1)
localScoreC(seq.TwoSegments.InSameExcursion)
# lindley() shows two realizations in the same excursion (no 0 value between the two LS values)
lindley(seq.TwoSegments.InSameExcursion)
# same beginning index but two possible ending indexes
# only one excursion realizes the local score even in there is two possible length of segment
localScoreC(seq.TwoSegments.InSameExcursion)$suboptimalSegmentScores
plot(1:length(seq.TwoSegments.InSameExcursion),lindley(seq.TwoSegments.InSameExcursion),type='b')
```

---

localScoreC\_double      *Local score for sequences of floating values*

---

**Description**

Calculates the local score for a sequence of doubles. Only provides the first occurrence. Use function suboptimalSegment() or Lindley() to obtain the others localizations of the different realizations of the local score.

**Usage**

```
localScoreC_double(v, supressWarnings = FALSE)
```

**Arguments**

v                      A sequence of values as vector.  
supressWarnings      if warnings should be displayed

**Value**

A structure containing: the local score value and the begin and end index of the segment realizing this optimal score ; all the local maxima of the Lindley process (non negative excursion) and their begin and end index ; the record times of the Lindley process but only the ones corresponding to the begin index of non negative excursions

**Examples**

```

localScoreC_double(c(1.2,-2.1,3.5,1.7,-1.1,2.3))
seq.TwoSegments=c(1.2,-2.1,3.5,1.7,2,-2,-2,-3.5,1,3.5,1.7,1,-2,-2)
# two segments realizing the local score value
localScoreC(seq.TwoSegments)
# only the first realization
localScoreC(seq.TwoSegments)$localScore
# all the realization of the local together with the suboptimal ones
localScoreC(seq.TwoSegments)$suboptimalSegmentScores
# for small sequences, you can also use lindley() fonction to check if
# several segments achieve the local score
lindley(seq.TwoSegments)
plot(1:length(seq.TwoSegments),lindley(seq.TwoSegments),type='b')
seq.TwoSegments.InSameExcursion=c(1,-2,3,2,-1,0,1,-2,-2)
localScoreC(seq.TwoSegments.InSameExcursion)
# lindley() shows two realizations in the same excursion (no 0 value between the two LS values)
lindley(seq.TwoSegments.InSameExcursion)
plot(1:length(seq.TwoSegments.InSameExcursion),lindley(seq.TwoSegments.InSameExcursion),type='b')
# same beginning index but two possible ending indexes
# only one excursion realizes the local score even in there is two possible length of segment
localScoreC(seq.TwoSegments.InSameExcursion)$suboptimalSegmentScores

```

---

LongSeq

*Long protein sequence*

---

**Description**

A long protein sequence.

**Usage**

LongSeq

**Format**

A character string with 1093 characters corresponding to Q60519.fasta in UniProt Data base.

**Source**

<https://www.uniprot.org/>



**Examples**

```

data(LongSeq)
LongSeq
nchar(LongSeq)
data(dico)
LongSeqScore=CharSequence2ScoreSequence(LongSeq,dico)
LongSeqScore[1:50]
localScoreC(LongSeqScore)$localScore
LS=localScoreC(LongSeqScore)$localScore[1]
prob1 = scoreSequences2probabilityVector(list(LongSeqScore))
daudin(localScore = LS, sequence_length = nchar(LongSeq),
        score_probabilities = prob1,
        sequence_min = min(LongSeqScore),
        sequence_max = max(LongSeqScore))
karlin(localScore = LS, sequence_length = nchar(LongSeq),
        score_probabilities = prob1,
        sequence_min = min(LongSeqScore),
        sequence_max = max(LongSeqScore))

```

---

mcc

*MCC [p-value] [iid]*


---

**Description**

Calculates an approximated p-value for a given local score value and a medium to long sequence length in the identically and independantly distributed model

**Usage**

```

mcc(
  localScore,
  sequence_length,
  score_probabilities,
  sequence_min,
  sequence_max
)

```

**Arguments**

localScore      the observed local score  
sequence\_length      length of the sequence (up to one hundred)  
score\_probabilities      the probabilities for each unique score from lowest to greatest  
sequence\_min      minimum score  
sequence\_max      maximum score

**Details**

This method is actually an improved method of Karlin and produces more precise results. It should be privileged whenever possible.

As with Karlin, the method works the better the longer the sequence.

**Value**

A double representing the probability of a local score as high as the one given as argument

**Examples**

```
mcc(40, 100, c(0.08, 0.32, 0.08, 0.00, 0.08, 0.00, 0.00, 0.08, 0.02, 0.32, 0.02), -6, 4)
mcc(40, 10000, c(0.08, 0.32, 0.08, 0.00, 0.08, 0.00, 0.00, 0.08, 0.02, 0.32, 0.02), -6, 4)
```

---

MidSeq

*Protein sequence*

---

**Description**

A protein sequence.

**Usage**

```
MidSeq
```

**Format**

A character string with 219 characters corresponding to P49755.fasta query in UniProt Data base.

**Source**

<https://www.uniprot.org/>

**Examples**

```
data(MidSeq)
MidSeq
nchar(MidSeq)
data(dico)
MidSeqScore=CharSequence2ScoreSequence(MidSeq,dico)
MidSeqScore[1:30]
localScoreC(MidSeqScore)$localScore
prob1 = scoreSequences2probabilityVector(list(MidSeqScore))
daudin(localScore = 52, sequence_length = nchar(MidSeq),
        score_probabilities = prob1,
        sequence_min = min(MidSeqScore),
        sequence_max = max(MidSeqScore))
score=-5:5
prob2=c(0.15,0.15,0.1,0.1,0.0,0.05,0.15,0.05,0.2,0.0,0.05)
```

```

daudin(localScore = 52, sequence_length = nchar(MidSeq),
       score_probabilities = prob2,
       sequence_min = min(MidSeqScore),
       sequence_max = max(MidSeqScore))

```

---

monteCarlo	<i>Monte Carlo method [p-value]</i>
------------	-------------------------------------

---

### Description

Calculates an empirical p-value based on simulations of similar integer sequences of the same length. Perfect for small sequences (both markov chains and identically and independantly distributed) with length  $\sim 10^3$ . See function monteCarlo\_double() for possible real scores.

### Usage

```
monteCarlo(local_score, FUN, ..., plot = TRUE, numSim = 1000)
```

### Arguments

local_score	local score observed in a segment.
FUN	function to simulate similar sequences with.
...	parameters for FUN
plot	boolean value if to display plots for cumulated function and density
numSim	number of sequences to generate during simulation

### Value

Floating value corresponding to the probability to obtain a local score with value greater or equal to the parameter

### Examples

```

monteCarlo(120, FUN = rbinom, n = 100, size = 5, prob=0.2)

new = sample(-7:6, replace = TRUE, size = 1000)
#MonteCarlo taking random sample from the input sequence itself

monteCarlo(local_score = 20, FUN = function(x) {return(sample(x = x,
size = length(x), replace = TRUE))}, x=new)

# Markovian example
MyTransMat <-
+ matrix(c(0.3,0.1,0.1,0.1,0.4, 0.2,0.2,0.1,0.2,0.3, 0.3,0.4,0.1,0.1,0.1, 0.3,0.3,0.1,0.0,0.3,
+         0.1,0.1,0.2,0.3,0.3), ncol = 5, byrow=TRUE)

monteCarlo(local_score = 50,

```

```
FUN = transmatrix2sequence, matrix = MyTransMat,
length=150, score = c(-2,-1,0,2,3), plot=FALSE, numSim = 5000)
```

---

monteCarlo\_double      *Monte Carlo method for real score case [p-value]*

---

### Description

Calculates an empirical p-value based on simulations of similar sequences of the same length. Perfect for small sequences (both markov chains and identically and independantly distributed) with length  $\sim 10^3$ . Function dedicated for real score case.

### Usage

```
monteCarlo_double(local_score, FUN, ..., plot = TRUE, numSim = 1000)
```

### Arguments

local_score	local score observed in a segment.
FUN	function to simulate similar sequences with.
...	parameters for FUN
plot	Boolean value if to display plots for cumulated function and density
numSim	number of sequences to generate during simulation

### Value

Floating value corresponding to the probability to obtain a local score with value greater or equal to the parameter

### Examples

```
score_reels=c(-1,-0.5,0,0.5,1)
proba_score_reels=c(0.2,0.3,0.1,0.2,0.2)
sample_from_model <- function(score.sple,proba.sple, length.sple){sample(score.sple,
                                size=length.sple, prob=proba.sple, replace=TRUE)}

monteCarlo_double(5.5,FUN=sample_from_model, plot = TRUE,
score.sple=score_reels,proba.sple=proba_score_reels, length.sple=100, numSim = 1000)
```

---

MySeqList	<i>Several sequences</i>
-----------	--------------------------

---

**Description**

A vector of character strings

**Usage**

```
MySeqList
```

**Format**

A list of 285 character strings with their entry codes as names

**Source**

Structural Classification Of Proteins database (SCOP). More precisely this data contain the 285 protein sequences of the data called "CF\_scop2dom\_20140205aa" with length from 31 to 404.

**Examples**

```
data(MySeqList)
head(MySeqList)
MySeqList[1]
nchar(MySeqList[1])
summary(sapply(MySeqList, nchar))
data(dico)
MySeqScoreList=lapply(MySeqList, FUN=CharSequence2ScoreSequence, dico)
head(MySeqScoreList)
AA=automatic_analysis(sequences=MySeqScoreList, model='iid')
AA[[1]]
# the p-value of the first 10 sequences
sapply(AA, function(x){x$`p-value`})[1:10]
# the 20th smallest p-values
sort(sapply(AA, function(x){x$`p-value`})) [1:20]
which(sapply(AA, function(x){x$`p-value`}) < 0.05)
table(sapply(AA, function(x){x$`method`}))
# The maximum sequence length equals 404 so it here normal that the exact method is used for
# all the 606 sequences of the data base
# Score distribution learnt on the data set
scoreSequences2probabilityVector(MySeqScoreList)
```

---

RealScores2IntegerScores

*Convert a real scores vector into an integer scores vector*

---

## Description

Convert real scores into integer scores

## Usage

```
RealScores2IntegerScores(RealScore, ProbRealScore, coef = 10)
```

## Arguments

RealScore	vector of real scores
ProbRealScore	vector of probability
coef	coefficient

## Details

Convert real scores into integer scores by multiplying real scores by a coefficient (default 10) and then assigning probability to corresponding extended (from the minimum to the maximum) integer scores

## Value

list containing `ExtendedIntegerScore` and `ProbExtendedIntegerScore`

## Examples

```
score <- c(-1,-0.5,0,0.5,1)
prob.score <- c(0.2,0,0.4,0.1,0.3)
(res1 <- RealScores2IntegerScores(score, prob.score, coef=10))
prob.score.err <- c(0.1,0,0.4,0.1,0.3)
(res2 <- RealScores2IntegerScores(score, prob.score.err, coef=10))
# When coef=1, the function can handle integer scores
ex.integer.score <- c(-3,-1,0,1, 5)
(res3 <- RealScores2IntegerScores(ex.integer.score, prob.score, coef=1))
```

---

`scoreDictionary2probabilityVector`*Check for missing scores values in the score distribution*

---

**Description**

Get extremes scores, then create a complete list and set a probability equal to zeros for not present scores

**Usage**

```
scoreDictionary2probabilityVector(list, score_extremes)
```

**Arguments**

`list`                vector of scores  
`score_extremes`    vector of probability

**Value**

vector containing all score values between extremes and the probability equal to 0 for missing score

**Examples**

```
Mylist=list("x1"=c(-2,0.1), "x2"=c(0,0.7), "x3"=c(1,0.2))  
scoreDictionary2probabilityVector(list=Mylist, score_extremes=c(-2,1))
```

---

`scoreSequences2probabilityVector`*Empirical distribution from sequences*

---

**Description**

Builds empirical distribution from a list of numerical sequences

**Usage**

```
scoreSequences2probabilityVector(sequences)
```

**Arguments**

`sequences`        list of numerical sequences

**Details**

By determining the extreme scores in the sequences, this function creates a vector of probabilities including values that do not occur at all. In this it differs from `table()`. For example, two sequences containing values from 1:2 and 5:6 will produce a vector of size 6.

**Value**

empirical distribution from minimum score to maximum score as a vector of floating numbers.

**Examples**

```
seq1 = sample(7:8, size = 10, replace = TRUE)
seq2 = sample(2:3, size = 15, replace = TRUE)
l = list(seq1, seq2)
scoreSequences2probabilityVector(l)
```

---

sequences2transmatrix *Transition matrix from sequence(s)*

---

**Description**

Calculates the transition matrix by counting occurrences of tuples in given vector list

**Usage**

```
sequences2transmatrix(sequences)
```

**Arguments**

sequences      Sequences to be analyzed, as list of numeric vectors

**Details**

The transition matrix will be structured so that the lowest score corresponds to the first column and row and the highest score corresponds to the last column and row. Note that the resulting matrix is not stochastic because it can occur rows filled up with only 0 for not observed score in Min Value Max Value interval.

**Value**

A list object containing

Transition Matrix

Transition Matrix

Min Value      minimal score found in supplied sequences

Max Value      maximal score found in supplied sequences



**Examples**

```
seq = sample(-1:1, size = 20, replace = TRUE)
seq2 = sample(-6:1, size = 20, replace = TRUE)
seq3 = sample(3:6, size = 50, replace = TRUE)
sequences2transmatrix(list(seq, seq2, seq3))
```

ShortSeq

*Short protein sequence***Description**

A short protein sequence of 31 amino acids corresponding to Q09FU3.fasta query in UniProt Data base.

**Usage**

```
ShortSeq
```

**Format**

A character string with 31 characters "MLTITSYFGFLLAALTITSVLFIGLNKIRLI"

**Source**

<https://www.uniprot.org/>

**Examples**

```
data(ShortSeq)
ShortSeq
nchar(ShortSeq)
data(dico)
SeqScore=CharSequence2ScoreSequence(ShortSeq,dico)
SeqScore
localScoreC(SeqScore)$localScore
LS=localScoreC(SeqScore)$localScore[1]
prob1 = scoreSequences2probabilityVector(list(SeqScore))
daudin(localScore = LS, sequence_length = nchar(ShortSeq),
        score_probabilities = prob1,
        sequence_min = min(SeqScore),
        sequence_max = max(SeqScore))

score=-5:5
prob2=c(0.15,0.15,0.1,0.1,0.0,0.05,0.15,0.05,0.2,0.0,0.05)
sum(prob2*score)
karlin(localScore = LS, sequence_length = nchar(ShortSeq),
        score_probabilities = prob2,
        sequence_min = min(SeqScore),
        sequence_max = max(SeqScore))
```

---

 stationary\_distribution

*Stationary distribution [Markov chains]*


---

**Description**

Calculates stationary distribution of markov transition matrix by use of eigenvectors of length 1

**Usage**

```
stationary_distribution(m)
```

**Arguments**

m                    Transition Matrix [matrix object]

**Value**

A vector with the probabilities

**Examples**

```
B = t(matrix (c(0.2, 0.8, 0.4, 0.6), nrow = 2))
stationary_distribution(B)
```

---

 transmatrix2sequence    *Sampling function for Markov chains*


---

**Description**

Creates Markov chains based on a transition matrix. Can be used as parameter for the Monte Carlo function.

**Usage**

```
transmatrix2sequence(matrix, length, initialIndex, score)
```

**Arguments**

matrix                transition matrix of Markov process  
 length                length of sequence to be sampled  
 initialIndex        (optional) index of matrix which should be initial value of sequence. If none supplied, a value from the stationary distribution is sampled as initial value.  
 score                (optional) a vector representing the scores (in ascending order) of the matrix index. If supplied, the result will be a vector of these values.

**Details**

The transition matrix is considered representing the transition from one score to another such that the score in the first row is the lowest and the last row are the transitions from the highest score to another. The matrix must be stochastic (no rows filled up with only '0' values).

**Value**

a Markov chain sampled from the transition matrix

**Examples**

```
B = t(matrix(c(0.2, 0.8, 0.4, 0.6), nrow = 2))
transmatrix2sequence(B, length = 10)
MyTransMat <-
  matrix(c(0.3,0.1,0.1,0.1,0.4, 0.2,0.2,0.1,0.2,0.3, 0.3,0.4,0.1,0.1,0.1, 0.3,0.3,0.1,0.0,0.3,
    0.1,0.1,0.2,0.3,0.3), ncol = 5, byrow=TRUE)
MySeq.CM=transmatrix2sequence(matrix = MyTransMat,length=90, score =c(-2,-1,0,2,3))
MySeq.CM
```

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