MANUAL

This manual provides reference documentation to TREMOL_Singlets from a user's and developer's perspective.

Coding style

- 4 spaces per indentation level.
- Float and Integers are lower case meanwhile vectors and arrays are upper case.
- Comments convention for functions is as follows:

```
out1 = function(Arg1, arg2):
    This is a function.
    Parameters
    - `Arg1:: Int`: array of dimensions ...
    - `arg2:: String`: string that ...
    return
    - `out1::Float`: float that ...
    """"
```

• The use of inline comments is sparingly and indicated with "#" symbol.

Requirements

- TREMOL has been tested in Linux operating system
- Julia (version 0.6.4 have been tested) https://julialang.org/downloads/oldreleases.html
- PyPlot package to generate the plots https://github.com/JuliaPy/PyPlot.jl

TREMOL_Singlets directory structure

This subsection is dedicated to list and describe the TREMOL directory structure.

Name	Description
doc/	Source files for TREMOL_Singlets documentation
examples/	Templates of basic scripts for TREMOL_Singlets modelling
TREMOL_singlets/	Source code
Description.pdf	Summary of TREMOL features and requirements
License.pdf	License file
README.pdf	Readme file

The source code is `TREMOL_singlets/`, which has the following contents:

Name	Description
main/	Main scripts to execute the simulation.
preprocessing/	Definition of the input parameters by the user
postprocessing/	Analysis of the raw data coming from main/

TREMOL_Singlets workflow

TREMOL_Singlets work flow is described in Figure 1. Pre-processing (yellow) processing (gray) and post-processing (green) phases are included.

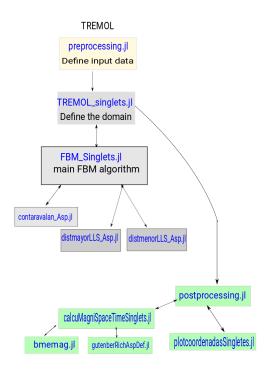


Figure 1. TREMOL_Singlets: workflow diagram

TREMOL_Singlets preprocessing

The first step to use TREMOL_Singlets code is the *preprocessing.jl* file. Here we define the input parameters that will be used in the main algorithm to model a seismic asperity rupture.

This file is seen as follows:

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Define you path

YOURPATH = "path user"

Include an external function

include("TREMOL_Singlets.jl")

Arguments

Parameters to assigned the initial conditions to the simulated domain

- `VecID ::String: Name to identify the simulated event.
- `fhi_asp::Float`: percentage of load to be transferred for any ruptured cell in the asperity domain
- `fhi_bkg::Float or Vector`: percentage of load to be transferred for any ruptured cell in the background
- `strength_asp::Integer`: strength value to define the asperity cells.
- `nbox::Integer`: Number of cells assigned to each Seismic Source.
- `nk::Integer`: Number of times that will execute a same experiment to obtain statistics *Parameters coming from the finite fault source method*
- `DurTeo::Float Vector`: rupture duration given by finite fault computation [seconds]
- `VelAsp :: Float Vector`: rupture velocity given by finite fault computation [km/s]
- `Leff::Float Vector`: Effective length size of the effective rupture area [km]
- `Weff::Float Vector`: Effective wide size of the effective rupture area [km]
- `VectorSeff::Float Vector`: Effective area [km²]
- `SaOri::Float Vector`: Ratio of the asperity size
- `YsizeOri::Float Vector`: Asperity area [km²]

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YOURPATH = "path user" include("TREMOL_Singlets.jl")

TREMOL_Singlets(VecID,fhi_asp,fhi_bkg,strength_asp,DurTeo,VelAsp,Leff,Weff,VectorSeff,SaOri,YsizeOri,nbox,nk)

Running a simulation

This section introduces the basics of running TREMOL on the command line.

- 1. Install Julia v0.6.
- 2. Add **PyPlot package** following the instructions defined in https://github.com/JuliaPy/PyPlot.il.
- 3. Define **YOURPATH** which is the path route to code and must to be defined in two scripts manually <u>`~/TREMOL singlets/preprocessing.il` script</u>
 - `~/TREMOL singlets/main/TREMOL Singlets.il
- 4. After it, in *preprocessing.jl* script define the input arguments. To execute TREMOL_Singlets there are two options:
- a) from a Linux bash run

\$ julia YOURPATH/REMOL_singlets/TREMOL_singlets/preprocessing.jl

This option do not allow to show the plots produced in the functions TREMOL_singlets.jl and postprocessing.jl. However, the output plots and files will be save in the folder /PATH/TREMOL_singlets/Results_PlotFiles

b) from a julia prompt

 $include (join path ("/YOURPATH/TREMOL_singlets/TREMOL_singlets/preprocessing/", "preprocessing .jl"))$

Visualization of results

As outputs TREMOL_Singlets generated eigth plots as ".pdf" and one file in ASCII ".dat" Plots are generated in four different functions, such as: TREMOL_singlets.jl, calcuMagniSpaceTimeSinglets.jl, plotcoordenadasSingletes.jl and gutenberRichAspDef.jl.

- TREMOL_singlets.jl: generates the plot of the initial strength configuration
 "EventoID*AsperezaSpatial.pdf"
- calcuMagniSpaceTimeSinglets.jl: generates three plots:
- 1. **"EventID-MagnitudeTime.pdf":** the magnitude of the events as function of the computed time
- 2. "**EventID-FrequencyMagnitude.pdf**": a histogram of the frequency magnitude of the simulated quakes
- 3. "**EventID-DurationFrequency.pdf**": the rupture duration of each simulated earthquake
- gutenberRichAspDef.jl: the cumulative number of earthquakes vs magnitude are generated for four different scale relations "**EventID-Grfit.pdf**"
- plotcoordenadasSingletes.jl: this function generates three plots
- 1. "EventID-Mean_Load.pdf": mean load value vs time since the simulation starts
- 2. **"EventID-IntereventRate.pdf":** the inter-event rate of the synthetic earthquakes
- 3. **"EventID-SpatialDistribution.pdf":** Final spatial distribution of the rupture groups shown in different colors each one.

These plots are saved in the directory /YOURPATH/TREMOL_singlets/Results_PlotsFiles. The user could deactivated the plots generation isplaying a "#" symbol at the beginning of the any PyPlot line.

If the number of realizations of each experitment is larger than one (nk >1) then the plots "**EventoID*AsperezaSpatial.pdf**" and "**EventID-SpatialDistribution.pdf**" is configured to be saved only for the first realization (i.e. nk=1)

The file "**EventID-MagnitudeStatisticalResults.dat**" contains in each column the following information:

- 1. b-value computed trough the function beemag, il using Somerville relation
- 2. maximum magnitude using Somerville relation
- 3. b-value computed trough the function beamag, il using Mai relation
- 4. maximum magnitude using Mai relation
- 5. b-value computed trough the function bmemag.jl using Mai-VL relation
- 6. maximum magnitude using Mai-VL relation
- 7. b-value computed trough the function bmemag.jl using Ramirez relation
- 8. maximum magnitude using Ramirez relation
- 9. ratio of the largest simulated earthquake [cells] and the total number of cells
- 10. largest simulated earthquake in [cells]
- 11. the size of the asperity defined in the random range
- 12. the random number to define the size of the asperity size
- 13. original asperity size coming from preprocessing input data
- 14. number of steps realized in the algorithm
- 15. maximum magnitude using Somerville relation
- 16. length of the area earhquake (in km/2) divided by the rupture velocity (km/s).
- 17. equivalent rupture time in seconds, considering the longest length divided by the rupture velocity
- 18. equivalent rupture time in seconds considering the root square of the area divided by the rupture velocity
- 19. rupture velocity, velAsp
- 20.area of the largest simulated earthquake
- 21. sqrt(area of the largest simulated earthquake)/velAsp
- 22.sqrt(area of the largest simulated earthquake)/DurTeo
- 23. velAsp*DurTeo
- 24. sqrt(area of the largest simulated earthquake)/velAsp

Each row corresponds to one execution controlled by the number "**nk**" defined in preprocessing.jl script.

The user can realize the corresponding statistics analyzing the ruptured area or the equivalent magnitude with four different scale-realtions (see calcuMagniSpaceTimeSinglets.jl script) by using the file "EventID-MagnitudeStatisticalResults.dat"

Examples

This section includes two examples using different input values in order to show how the user can configure the code. To execute the examples the path directory **"YOURPATH"** must be modify only in two scripts

YOURPATH/TREMOL_singlets/examples/ExampleMw7_0-1982/TREMOL_Singlets.jl and

YOURPATH/TREMOL_singlets/examples/ExampleMw8_2-2017/TREMOL_Singlets.jl

Example 1: Modelling the magnitude and ruptured area of the earthquake Mw=7.0, 07/06/1982

After define YOURPATH, to run this example is required to executed the following command in the bash

\$ julia YOURPATH/TREMOL_singlets/examples/ExampleMw7_0-1982/preprocessing.jl

The value described in the `YOURPATH/TREMOL_singlets/examples/ExampleMw7_0-1982/preprocessing.jl` are:

```
VecID = ["Mw7-1982"]

fhi_bkg = 0.67

fhi_asp = 0.90

strength_asp = 4

nbox = 100

nk = 10

DurTeo = [40]

VelAsp = [3.2]

Leff = [34.47]

Weff = [17.81]

VectorSeff = [Leff[1].*Weff[1]]

SaOri = [0.23]

YsizeOri = [Leff[1].*Weff[1]]
```

Values of: *DurTeo*, *VelAsp*, *SaOri*, *Leff* and *Weff* are computed in Rodríguez-Pérez & Ottemöller, (2013). While *fhi_bkg*, *fhi_asp*, *strength_asp*, *nbox*, are defined by the user in order to explore the behavior of the algorithm TREMOL_Singlets.

The output plots and file are saved in the folder "/YOURPATH/TREMOL_singlets/examples/ExampleMw7_0-1982/Results_PlotsFilesMw7_1982/"

To compare the user results with the expected you can use the files saved in the folder "~/YOURPATH/TREMOL_singlets/examples/ ExampleMw7_0-1982 /Results_PlotsFilesMw7_1982/Expected_Output_Mw7_0_1982"

Example 2: Modelling the magnitude and ruptured area of the earthquake Mw=8.2, 08/09/2017

To run this example is requiered to executed the following command in the bash

```
$ julia YOURPATH/TREMOL_singlets/examples/ExampleMw8_2-2017/preprocessing.jl
```

The input value described in the `YOURPATH/TREMOL_singlets/examples/ExampleMw8_2-2017/preprocessing.jl` file are:

```
VecID = ["EventMw8_2"]

fhi_bkg = 0.67

fhi_asp = 0.90

strength_asp = 4

nbox = 100

nk = 1

DurTeo = [40]

VelAsp = [2.0]

Leff = [125.95]

Weff = [71.13]

VectorSeff = [Leff[1].*Weff[1]]

SaOri = [0.34]

YsizeOri = [Leff[1].*Weff[1]]
```

Code documentation

Following sub-sections are dedicated to described the **parameters** and **return** parameters of scripts of the code documentation TREMOL_Singlets.

preprocessing/

Define the input parameters defined arguments

preprocessing.jl: Define the input by the user and call the function TREMOL_Singlets to pass the input values as arguments

```
Import modules
using PyPlot
using PyCall
@pyimport matplotlib.colors as matcolors
@pyimport matplotlib as mpl
@pyimport matplotlib.patches as patches
```

Import functions

- include(joinpath("/YOURPATH/TREMOL singlets/main/","TREMOL Singlets.jl"))

Parameters that assigned the initial conditions to the simulated domain

- `VecID ::String: Name to identify the simulated event.
- `fhi_asp::Float`: percentage of load to be transferred for any ruptured cell in the asperity domain
- `fhi_bkg::Float or Vector`: percentage of load to be transferred for any ruptured cell in the background
- `strength_asp::Integer`: strength value to define the asperity cells.
- `nbox::Integer`: Number of cells assigned to each Seismic Source.
- `nk::Integer`: Number of times that will execute a same experiment to obtain statistics

Parameters coming from the finite fault source method

- `DurTeo::Float Vector`: rupture duration given by finite fault computation [seconds]
- `VelAsp :: Float Vector`: rupture velocity given by finite fault computation [km/s]
- `Leff::Float Vector`: Effective length size of the effective rupture area [km]

- `Weff::Float Vector`: Effective wide size of the effective rupture area [km]
- `VectorSeff::Float Vector`: Effective area [km²]
- `SaOri::Float Vector`: Ratio of the asperity size
- `YsizeOri::Float Vector`: Asperity area [km²]

main/

TREMOL_Singlets.jl

Main program that define the size and shape to the effective domain and asperity domain. Also this script gives the input values as arguments to the FBM algorithm. Lastly the output values goes to the postprocessing function.

Import functions

- include(joinpath("/YOURPATH/TREMOL_singlets/main/","FBM_Singlets.jl"))
- include(joinpath("/YOURPATH/TREMOL_singlets/postprocessing/","postprocessing.jl"))

Parameters

- `VecID::String`: Name to identify the simulated event.
- `fhi_asp::Float`: percentage of load to be transferred for any ruptured cell in the asperity domain
- `fhi_bkg::Float or Vector`: percentage of load to be transferred for any ruptured cell in the background
- `strength_asp::Integer`: strength value to define the asperity cells.
- `DurTeo::Float Vector`: rupture duration given by finite fault computation [seconds]
- `VelAsp :: Float Vector`: rupture velocity given by finite fault computation [km/s]
- `Leff::Float Vector`: Effective length size of the effective rupture area [km]
- `Weff::Float Vector`: Effective wide size of the effective rupture area [km]
- `VectorSeff::Float Vector`: Effective area [km²]
- `SaOri::Float Vector`: Ratio of the asperity size
- `YsizeOri::Float Vector`: Asperity area [km²]
- `nbox::Integer`: Number of cells assigned to each Seismic Source.
- `nk::Integer`: Number of times that will execute a same experiment to obtain statistics

return

- `VecMagni::Vector`: Array that contains the results of the magnitude analysis coming from the function postprocessing.jl

FBM Singlets.il

This function carry out the FBM asperity algorithm

Import function

- $-include (joinpath ("/YOURPATH/TREMOL_singlets/main/", "contaravalan_Asp.jl")) \\$
- $-include (joinpath ("/YOURPATH/TREMOL_singlets/main/", "distmayor LLS_Asp.jl")) \\$
- include(joinpath("/YOURPATH/TREMOL_singlets/main/","distmenorLLS_Asp.il"))

Parameters

- `nbox_x::Integer`: number of cells in X-axis of the domain \Omega
- `nbox_y::Integer`: number of cells in Y-axis of the domain \Omega
- `smin::Integer`: number of steps that realize the algorithm

- `VecPosi::Array`: size(nbox_x,nbox_y), Matrix of the load values. Dynamical matrix because changes the value of some cells at each step.
- `MatrizStrengtInitial::Array`: size(nbox_x,nbox_y), Initial Matrix of the strength value.
- `VecAsperi::Array`: size(nbox_x,nbox_y), Matrix of the strength value. This matrix evolves at each step
- `fhiFuera::Float`: load-transfer value assigned to the cells located in the domain
- `fhiDentro::Float`: load-transfer value assigned to the cells located in the asperity domain
- `VecPhi::Array`: size(nbox_x,nbox_y), Matrix of the load-tranfer values.

return

- `vectk1::Array`: size(smin+10,12). Raw data that contains the results of the transfer, accumulation and rupture process following the FBM rules. At each step each row contains:
 - 1. k:number of step
 - 2. acumt: cumulative time $(T_k = sum(tiempo[1:k]))$
 - 3. tiempo[k]: inter-event time [dimensionless]
 - 4. (0 or 1): counter that indicates if is normal or avalanche event
 - 5. suma: sum of the load in all the cells
 - 6. sumarho= 1/tiempo[k] (rupture rate)
 - 7. parametrosigma: load value of the cell chosen to fail
 - 8. a: coordinate in the X-axis, of the cell chosen to fail
 - 9. b: coordinate in the Y-axis, of the cell chosen to fail
 - 10. fhi: load-transfer value
- 11-12. (iorigi,jorigi): coordinates of the cell chosen to fail at the first stage of the searching algorithm defined in the function contaravalan_Asp.
- `vectparamestad::Array`: size(smin+10,2). Mean and Standard deviation of the load in the system, computed at each time step considering only the cells active
- `VecPosiFinal::Array`: size(nbox_x,nbox_y). Final configuration of VecPosi.

contaravalan Asp.jl

Compute the number of cells that overpass the threshold load value (STAGE 1). But also in this version the cells that fails is choosen considering not only its load but also the strength criterion defined in TREMOL algorithm (STAGE 2)

Parameters

- `nx::Integer`: number of cells in X-axis of the domain \Omega
- `ny::Integer`: number of cells in Y-axis of the domain \Omega
- `VectorP::Array`: size(nbox_x,nbox_y), Matrix of the load values. Dynamical matrix because changes the value of some cells at each step.
- `VecAsperi::Array`: size(nbox_x,nbox_y), Matrix of the strength value. This matrix evolves at each step
- `VecPhi:Array`: size(nbox_x,nbox_y), Matrix of the load-tranfer values.
- `rho::Integer`: Weibull exponent, we take it as a constant 30
- `k::Integer`: step of number

return

- `contador::Integer`: number of cells that overpass the threshold load value
- `Nflag1::Integer` : code that indicates if the rupture is normal or avalanche
- `maxtempo1::Float` : maximum load value
- `iout1::Integer`: x coordinate in the array of the chosen cell to fail
- `jout1::Integer`: y coordinate in the array of the chosen cell to fail
- `VecAsperi::Array`: updated matrix of the strength
- `iout::Integer`: x coordinate in the array of the chosen cell to fail at STAGE 1

- `jout::Integer`: Y coordinate in the array of the chosen cell to fail at STAGE 1

distmayorLLS_Asp.jl

Function that distribute the load following the avalanche events algorithm (avalanche event is the cell that overpass their threshold values)

Import function

- include(joinpath("/YOURPATH/TREMOL_singlets/main/","vecinosLLS.jl"))

Parameters

- `vector::Array`:vector of nine positions that contains the neighbors and failed cell, being: vector[9], vector[3], vector[7] and vector[1] diagonal neighbors; vector[8], vector[2], vector[4] and vector[6] perpendicular neighbors. Finally vector[5] is the failed cell.
- `fhi::Float`: load-transfer value assigned to the failed cell.

return

- `vector::Array`: updated vector of nine positions after the load transfer of the failed cell.

distmenorLLS_Asp.jl

Function that distributes the load following the normal-events algorithm (when any cell ooverpass their threshold values)

Import function

- include(joinpath("/YOURPATH/TREMOL_singlets/main/","vecinosLLS.jl"))

Parameters

- `vector1::Array`:vector of nine positions that contains the neighbors and failed cell, being: vector1[9], vector1[3], vector1[7] and vector1[1] diagonal neighbors; vector1[8], vector1[2], vector1[4] and vector1[6] perpendicular neighbors. Finally vector1[5] is the failed cell.
- `fhi::Float`: load-transfer value assigned to the failed cell

return

- `vecreload::Array`: updated vector of nine positions after the load transfer of the failed cell.

vecinosLLS.jl

This function distributed the load for the orthogonal neighbors (N,S,E,W) that are allowed to received load.

Import function

- include(joinpath("/YOURPATH/TREMOL_singlets/main/","veciProhiDiag.jl"))

Parameters

- `vector::Array`:vector of nine positions that contains the neighbors and failed cell, being: vector[9], vector[3], vector[7] and vector[1] diagonal neighbors; vector[8], vector[2], vector[4] and vector[6] perpendicular neighbors.
- `vecEsfTot::Float`: Load value of the failed cell.

return

- `vectclon::Array`: Updated vector of nine positions after the load transfer of the failed cell.

veciProhiDiag.jl

Compute the number of cells that overpass the threshold load value. But also in this version it choose the cells that is ruptured because the strength criterion defined in TREMOL algorithm

Parameters

- `vector::Array`:vector of nine positions that contains the neighbors and failed cell, being: vector[9], vector[3], vector[7] and vector[1] diagonal neighbors; vector[8], vector[2], vector[4] and vector[6] perpendicular neighbors.`
- `vecEsfDiag::Float`: Load value will be transfer to the Diagonal neighbors. `
- `A::Array`:vector of four positions that contains the diagonal neighbors

return

- `B::Array`: updated vector of four positions that contains the new amount of load given to the diagonal neighbors

postprocessing/

postprocessing.jl

Main program that assigns the size and shape to the effective domain and asperity domain. Also this script gives the input values to the FBM algorithm. Lastly the output values goes to the postprocess function.

Import function

- include(joinpath("/YOURPATH/TREMOL_singlets/main/","calcuMagniSpaceTimeMultiSinglets.jl"))
- include(joinpath("/YOURPATH/TREMOL_singlets/postprocessing/","plotcoordenadasSingletesbis.jl"))

Parameters

- `datos::Array`: size(smin,12) raw data coming from the FBM algorithm. This data base contains the rupture information of model

Each row contains:

- 1. k:number of step
- 2. acumt: tiempo acumulado $(T_k = sum(tiempo[1:k]))$
- 3. tiempo[k]: inter-event time [dimensionless]
- 4. (0 or 1): identifier to normal or avalanche event
- 5. suma: sum of the load in all the cells
- 6. sumarho= 1/tiempo[k] (rupture rate)
- 7. parametrosigma: load value of the cell chosen to fail
- 8. a: coordinate in the X-axis, of the cell chosen to fail
- 9. b: coordinate in the Y-axis, of the cell chosen to fail
- 10. fhi: load-transfer value
- 11-12. (iorigi,jorigi): coordinates of the cell chosen to fail at the first stage of the searching algorithm defined in the function contaravalan_Asp.
- $\ensuremath{^{\cdot}}$ VectorCenterAperities::Vector Integer $\ensuremath{^{\cdot}}$: size(1), Central coordinates of the aspertity (x,y)
- `VectorSaOriLateralSize_y::Vector`: Asperity lateral size in the Y-axis
- `VectorSaOriLateralSize_x::Vector`: Asperity lateral size in the X-axis

- `VectorCoordsAperities_x::Vector Integer`: coordinates of the asperity vertex in the X-axis
- `VectorCoordsAperities y::Vector Integer`: coordinates of the asperity vertex in the Y-axis
- `nbox_x::Integer`: number of cells in X-axis of the domain \Omega
- `nbox_y::Integer`: number of cells in Y-axis of the domain \Omega
- `smin::Integer`: number of steps that realize the algorithm
- `vecEstadistico::Array`: size(smin+10,2). Mean and Standard deviation of the load in the system, computed at each time step considering only the cells active
- `VectorSaOri::Float Vector`: Random size of the asperity
- `SaOri::Float Vector`: Original Ratio of the asperity size
- `TotalNumberCells::Integer`: number of cells in the domain \Omega (nbox_x * nbox_y)
- `a::Float`: random number
- `CellSize::Float`: Size of a cell in km²
- `velAsp::Float`: rupture velocity
- `Weff::Float`: width of the effective area in km
- `Leff::Float`: length of the effective area in km
- `DurTeo::Float`: rupture duration

return

- `vecMagniLoop::Array`: results of tha magnitude analysis
- Each row contains:
- 1. vecMagniLoop[1,1]: b-value computed trough the function bmemag.jl using Somerville relation
- 2. vecMagniLoop[1,2]: maximum magnitude using Somerville relation
- 3. vecMagniLoop[1,3]: b-value computed trough the function bmemag.jl using Mai relation
- 4. vecMagniLoop[1,4]: maximum magnitude using Mai relation
- 5. vecMagniLoop[1,5]: b-value computed trough the function bmemag.jl using Mai-VL relation
- 6. vecMagniLoop[1,6]: maximum magnitude using Mai-VL relation
- 7. vecMagniLoop[1,7]: b-value computed trough the function bmemag.jl using Ramirez relation
- 8. vecMagniLoop[1,8]: maximum magnitude using Ramirez relation
- 9. vecMagniLoop[1,9]: ratio of the largest simulated earthquake [cells] and the total number of cells
- 10. vecMagniLoop[1,10]: largest simulated earthquake in [cells]
- 11. vecMagniLoop[1,11]: VectorSaOri
- 12. vecMagniLoop[1,12]: a
- 13. vecMagniLoop[1,13]: SaOri
- 14. vecMagniLoop[1,14]: smin
- 15. vecMagniLoop[1,15]: maximum magnitude using Somerville relation
- 16. vecMagniLoop[1,16]: length of the area earhquake (in km²) divided by the rupture velocity (km/s)
- 17. vecMagniLoop[1,17]: equivalent rupture time in seconds, considering the longest length divided by the rupture velocity
- 18. vecMagniLoop[1,18]: equivalent rupture time in seconds considering the root square of the area divided by the rupture velocity
 - 19. vecMagniLoop[1,19]: rupture velocity, velAsp
 - 20. vecMagniLoop[1,20]: area of the largest simulated earthquake
 - 21. vecMagniLoop[1,21]: sqrt(area of the largest simulated earthquake)/velAsp
 - 22. vecMagniLoop[1,22]: sqrt(area of the largest simulated earthquake)/DurTeo
 - 23. vecMagniLoop[1,23]: velAsp*DurTeo
 - 24. vecMagniLoop[1,24]: sqrt(area of the largest simulated earthquake)/velAsp

calcuMagniSpaceTimeSinglets.jl

Cluster the avalanches considering the time and space criterion. Also is computed the equivalent magnitude at each new group.

Import function

- include(joinpath("/YOURPATH/TREMOL_singlets/postprocessing/","gutenberRichAspDef.il"))
- include(joinpath("/YOURPATH/TREMOL_singlets/postprocessing/","bmemag.jl"))

Parameters

- `Y::Array`:raw data coming from the FBM algorithm. This data base contains the rupture information of model Each row contains:
 - 1. k:number of step
 - 2. acumt: tiempo acumulado (T_k = sum(tiempo[1:k]))
 - 3. tiempo[k]: inter-event time [dimensionless]
 - 4. (0 or 1): identifier to normal or avalanche event
 - 5. suma: sum of the load in all the cells
 - 6. sumarho= 1/tiempo[k] (rupture rate)
 - 7. parametrosigma: load value of the cell chosen to fail
 - 8. a: coordinate in the X-axis, of the cell chosen to fail
 - 9. b: coordinate in the Y-axis, of the cell chosen to fail
 - 10. fhi: load-transfer value
- 11-12. (iorigi,jorigi): coordinates of the cell chosen to fail at the first stage of the searching algorithm defined in the function contaravalan_Asp.
- `Nbox_x::Integer`: number of cells in X-axis of the domain \Omega
- `Nbox_y::Integer`: number of cells in Y-axis of the domain \Omega`
- `smin::Integer`: number of steps that realize the algorithm
- `VectorCenterAperities::Vector Integer`: size(1), Central coordinates of the aspertity (x,y)
- `VectorSaOriLateralSize_y::Vector`: Asperity lateral size in the Y-axis
- `VectorSaOriLateralSize x::Vector`: Asperity lateral size in the X-axis
- `VectorCoordsAperities_x::Vector Integer`: coordinates of the asperity vertex in the X-axis
- `VectorCoordsAperities_y::Vector Integer`: coordinates of the asperity vertex in the Y-axis
- `CellSize::Float`: Size of a cell in km2`
- `velAsp::Float`: rupture velocity
- `Weff::Float`: width of the effective area in km
- `Leff::Float`: length of the effective area in km

return

- `vecMagni::Array`: matrix that contains the results of the magnitude analysis
- VecNewAvalSpaceTime::Array`: matrix that contains the data of the regrouping algorithm coming from the Y matrix

Each row contains:

- 1. VecNewAvalSpaceTime[contNumAvalSpaceTime,1] = contadorAval: number of elements considering in the regrouping algorithm
- 2. VecNewAvalSpaceTime[contNumAvalSpaceTime,2] = original number in the raw catalog Y
- 3. $VecNewAvalSpaceTime[contNumAvalSpaceTime,3] = cumulative time (T_k = sum(tiempo[1:k])) computed in$

the raw catalog Y

- 4. VecNewAvalSpaceTime[contNumAvalSpaceTime, 4] = inter-event time [dimensionless] computed in the raw catalog Y
- 5. VecNewAvalSpaceTime[contNumAvalSpaceTime,5] = normal-avalanche code considered in the new regrouping
- 6. VecNewAvalSpaceTime[contNumAvalSpaceTime,6] = x coordinate
- 7. VecNewAvalSpaceTime[contNumAvalSpaceTime,7] = y coordinate
- 8. VecNewAvalSpaceTime[contNumAvalSpaceTime,8] = rupture rate

bmemag.jl

Function calculates the mean magnitude, the b-value based on the mean and the standard deviation

Parameters

- `VecMag ::Array`: Vector that contains the equivalent magnitudes

return

- `meanm1::Float`: mean value of VecMag vector

- `b1::Float`: b-value

- `sig1::Float`: standard deviation

- `av2::Float`: a-value

gutenberRichAspDef.jl

Function that computes the Gutenberg-Richter fit relation using the method of least squares

Parameters

- `VecMagHB1::Array`: Vector that contains the equivalent magnitudes

- `nflag::Integer`: Identifier

- `AreaSUB::Float`: Cell area ion km2

return

- `paramGRAreaWY::Array`: results of the Gutenber-Richter fitting using the least square method
- $1.\ paramGRA reaWY [1] = magMax$
- 2. paramGRAreaWY[2]=magMin
- 3. paramGRAreaWY[3]=p1 (a-value)
- 4. paramGRAreaWY[4]=p2 (b-value)
- 5. paramGRAreaWY[5]=rho (correlation coeficient)
- `cuenrep::Integer`: number of times that a same frequency of magnitudes is repeated for different magnitudes
- `vecWrite::Array`: vector that contains the frequency-magnitude data.