

Vademecum: list of functions of the software StrainModeler developed in the paper:

StrainModeler: A MATHEMATICA™ –based program for 3D analysis of finite and progressive strain

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Basis Change (BC__)

BLOCK DESCRIPTION

This is a straightforward method to make a basis change.

The new basis is defined from the old one using 2 + 2 angles

{ α_1 , θ_1 } { α_2 , θ_2 }

(see figure)

BLOCK FUNCTIONS

Code: BC01

Description: It computes the matrix of the basis change.

Syntax: **generalBasisChangeMatrix**[*axisIpAngles*, *directionUpAngles*]

Inputs:

axisIpAngles: they are the angles corresponding to the direction of the axis i' of the new basis.

directionUpAngles: they are the angles of a direction u' that is also contained in the plane i - k .

Output: 3x3 base change matrix.

Code: BC02

Description: It computes the matrix of the deformation gradient in the old basis.

Syntax: **computeGradientOldBasis**[*changeMatrix*, *gradientNewBasis*]

Inputs:

changeMatrix: matrix of basis change.

gradientNewBasis: matrix of the deformation gradient in the new basis

Output: 3x3 matrix of the deformation gradient in the old basis

Code: BC03

Description: It computes the matrix of a vector in the old basis.

Syntax: **computeVectorOldBasis**[*changeMatrix*, *vectorNewBasis*]

Inputs:

changeMatrix: matrix of basis change.

vectorNewBasis: column matrix of a vector in the new basis.

Output: column matrix of a vector in the old basis.

Code: BC04

Description: It computes the matrix of the deformation gradient in the new basis.

Syntax: **computeGradientNewBasis**[*changeMatrix*, *gradientOldBasis*]

Inputs:

changeMatrix: matrix of basis change.

gradientOldBasis: matrix of the deformation gradient in the old basis.

Output: 3x3 matrix of the deformation gradient in the new basis.

Code: BC05

Description: It computes the matrix of a vector in the new basis.

Syntax: **computeVectorNewBasis**[*changeMatrix*, *vectorOldBasis*]

Inputs:

changeMatrix: matrix of basis change.

vectorOldBasis: column matrix of a vector in the old basis.

Output: column matrix of a vector in the new basis.

Matrix n - roots (MR)

BLOCK DESCRIPTION

Given the accumulated shear/flattening/exponential matrix, it computes its n-root.

BLOCK FUNCTIONS

Code: MR01

Description: It computes the n-root of a 3x3 simple shear matrix.

Syntax: **shearMatrixRoot**[*Ac*, *n*]

Inputs:

Ac: Matrix of the total simple shear.

n: Number of steps in which the process of simple shear is divided (order of the root).

Output: Matrix of the incremental deformation gradient (the n-root of the input matrix)

Code: MR02

Description: It computes the n-root of a 3x3 flattening matrix.

Syntax: **flatteningMatrixRoot**[*Ac*, *n*]

Inputs:

Ac: Matrix of the total flattening.

n: Number of steps in which the process of flattening is divided (order of the root).

Número de pasos en que quiero descomponer el proceso de flattening (orden de la raíz).

Output: Matrix of the incremental deformation gradient (the n-root of the input matrix)

Code: MR03

Description: It computes the n -root of a 3×3 exponential matrix.

Syntax: **exponentialMatrixRoot**[A_c, n]

Inputs:

A_c : Matrix of the total exponential process.

n : Number of steps (order of the root)

Output: Incremental gradient (the n -root of the input matrix)

Gradient Sequence Computations (GSC)

BLOCK DESCRIPTION

Computations associated to an Incremental Gradient Sequence (IGS).

BLOCK FUNCTIONS

Code: GSC01

Description: Builds an IGS composed by n identical matrices M :

$M, n \rightarrow \{M, M, M, \dots, M\}$

Syntax: **buildIGS**[M, n]

Inputs:

M : Repeated matrix.

n : Number of steps.

Output: An IGS list of matrices, $\{M, M, M, \dots, M\}$.

Code: GSC02

Description: It computes an Accumulated Gradient Sequence (AGS) from an IGS:

$F = \{M, M, M, \dots, M\} \rightarrow \{M^1, M^2, \dots, M^n\}$

Note: The elements of F may be different.

Syntax: **computeAGS**[F]

Inputs:

F : The IGS list of matrices.

Output: The AGS list $\{M^1, M^2, \dots, M^n\}$.

Code: GSC03

Description: It computes a sequence of items composed of the axis ratios of the incremental strain ellipsoid $\{R_{13}, R_{12}, R_{23}\}$ from an incremental gradient sequence (IGS).

Syntax: **computeSeqR13R12R23**[F]

Inputs:

F : The IGS list of matrices (the elements of F may be different).

Output: The list $\{ \{R_{13}, R_{12}, R_{23}\}_1, \{R_{13}, R_{12}, R_{23}\}_2, \dots, \{R_{13}, R_{12}, R_{23}\}_n \}$

Code: GSC04

Description: It computes a sequence of items composed of the axis ratios of the incremental strain ellipsoid $\{R_{13}, R_{12}, R_{23}, A\}$ from an incremental gradient sequence (IGS).

A : The product eigenvalue1 \times eigenvalue3.

Syntax: **computeSeqR13R12R23A**[*F*]

Inputs:

F: IGS list of matrices (the elements of *F* may be different).

Output: The list { {*R*₁₃, *R*₁₂, *R*₂₃,*A*}₁, {*R*₁₃, *R*₁₂, *R*₂₃,*A*}₂,..., {*R*₁₃, *R*₁₂, *R*₂₃,*A*}_{*n*}}

Code: GSC05

Description: It obtains the sequence of the three principal directions of the strain corresponding to the accumulated gradients.

Syntax: **computeSeqPrincVectors**[*F*]

Inputs:

F: IGS list of matrices (the elements of *F* may be different).

Output: The list { {*V*₁,*V*₂,*V*₃}₁, {*V*₁,*V*₂,*V*₃}₂, ..., {*V*₁,*V*₂,*V*₃}_{*n*} }.

Progressive Transformations (PT)

BLOCK DESCRIPTION

A progressive transformation is defined by an IGS *F*. We provide functions to transform directions and planes.

BLOCK FUNCTIONS

Code: PT01

Description: Starting from the initial vector *v*₀, we compute the successive transformed vectors;

F, *v*₀ -> {*v*₀, *v*₁, ..., *v*_{*n*}}

F: IGS

*v*₀: Initial vector

Caveat: The resulting sequence has *n*+1 elements.

Syntax: **progressiveVectorTransform**[*F*, *v*₀]

Inputs:

F: IGS *F*={*F*₁, *F*₂, ..., *F*_{*n*}}

*v*₀: Initial direction vector

Output: The list {*v*₀, *v*₁, ..., *v*_{*n*}}, where *v*_{*i*}=*F*_{*i*}**v*_(*i*-1).

Code: PT02

Description: It calculates two vectors {*V*₁,*V*₂} that define the plane from the dip and the dip direction, see fig. 4.

*V*₁ is the trace of the plane with the horizontal plane; *V*₂ is the vector corresponding to the dip.

Syntax: **calculatePlaneBasis**[*alpha*, *theta*]

Inputs:

alpha, *theta* are the angles that characterize the dip direction.

Output: The list {*V*₁,*V*₂}.

Code: PT03

Description: The function transforms the plane by a deformation defined by its gradient.

Syntax: **transformsPlaneParameters**[*basis*, *grad*]

Inputs:

basis: Vectors that define the original plane

grad: Gradient of the deformation.

Output: The list

{singular,norm1,norm2,normsRatio,dp1,dp2,normalPlano,normalPlanoAxial,norm1*norm2}

- a) “singular” is an indicator if the ellipse is actually a circle.
singular=False means that it is not a circle.
If singular=True, the remaining entries are meaningless.
- b) “norm1”: magnitude of the major semi-axis of the strain ellipse on the deformed plane.
- c) “norm2”: magnitude of the minor semi-axis of the strain ellipse on the deformed plane.
- d) “normsRatio” : semi-axes ratio of the strain ellipse on the deformed plane.
- e) “dp1”: orientation (α and θ values) of the major axis of the strain ellipse on the deformed plane.
- f) “dp2” :orientation (α and θ values) of the minor axis of the strain ellipse on the deformed plane.
- g) “PlaneNormal” : α and θ values for the transformed plane.
- h) “AxialPlaneNormal”: orientations (α and θ values) of the axial planes of the folds formed on the deformed plane. It is assumed that these axial planes contain the major axis of the finite strain ellipse on the plane and are perpendicular to the deformed plane.
- i) “norm1 \times norm2”: area change on the deformed plane.

Code: PT04

Description: Starting from the initial plane defined by basis0, it computes the data of the successive transformed planes. The successive transformations are defined by the IGS F.

Syntax: **progressivePlaneTransform**[*F*, *basis0*]

Inputs:

F: IGS $F=\{F_1, F_2, \dots, F_n\}$

basis0: Initial plane defined by two vectors (see the function PT02).

Output: The list { D_1, D_2, \dots, D_n }, where D_i are the data calculated by PT03.

Code: PT05

Description: Given a sequence of vectors $V = \{v_0, v_1, \dots, v_n\}$, the function gives a sequence constituted by the corresponding lengths of the vectors of V.

Syntax: **lengthVectorSeq**[*V*]

Inputs:

V: a sequence of vectors, $\{v_0, v_1, \dots, v_n\}$.

Output: The list { L_0, L_1, \dots, L_n }, where $L_i = \text{Norm}[L_i]$.

Interface: Angles/Cartesian Coordinates (ACC)

BLOCK DESCRIPTION

Functions that compute the angular parameters (α , θ) of a vector from its Cartesian components and viceversa (see Fig. 2).

BLOCK FUNCTIONS

Code: ACC01

Description: Given a direction, defined by angles (α , θ), it computes the corresponding vector in Cartesian components.

Syntax: **fromAnglesToCartesian**[*alpha*, *theta*]

Inputs:

alpha, *theta*: The two angles that define a direction.

Output: The list $\{v_1, v_2, v_3\}$: Cartesian components of the unit vector associated to the direction.

Code: ACC02

Description: Given a direction, defined by a vector in Cartesian components, it computes the corresponding angles (vertical vectors have not well defined angles). This function is the inverse of ACC01.

Syntax: **evalDirAngles**[*v*]

Inputs:

v: The list $\{v_1, v_2, v_3\}$: Cartesian components of the unit vector associated to the direction.

Output : The list $\{alpha, theta\}$: the two angles that define a direction.

Code: ACC03

Description: A sequence of vectors in Cartesian coordinates is transformed in a sequence of directions in angle form (using ACC02).

Syntax: **fromVectorSeqToAngleSeq**[*seqv*]

Inputs:

seqv: The list $\{v_1, v_2, \dots, v_n\}$ of vectors in Cartesian components.

Output: The list $\{d_1, d_2, \dots, d_n\}$ of the corresponding directions in angle form, where $d_i = \{alpha_i, theta_i\}$ are the two angles that define the direction d_i .

Code: ACC04

Description: Given a sequence of vectors normal to planes, it computes another sequence corresponding to the same planes defined by the two angles that define their dip direction.

Syntax: **fromNormalVectorSeqToAngleSeq**[*seqv*]

Inputs:

seqv: The list $\{v_1, v_2, \dots, v_n\}$ of normal vectors in Cartesian components.

Output: The list $\{dip_1, dip_2, \dots, dip_n\}$ of the corresponding directions in angle form, where $dip_i = \{alpha_i, theta_i\}$ are the two angles that define the dip direction dip_i .

Code: ACC05

Description: Given a plane defined by a normal vector, it computes the angles of

the corresponding dip direction.

Syntax: **evalPlaneAngles**[*r*]

Inputs:

r: The list {*r*₁, *r*₂, *r*₃} of the normal vector in Cartesian components.

Output: The list {*alpha*, *theta*} with the dip direction angles.

Code: ACC06

Description: A sequence of angles in radians is transformed to degrees.

Syntax: **radianSeqToDegreeSeq**[*rs*]

Inputs:

rs: A list of angles in radians.

Output: The input list converted to degrees.

Plotting (P)

BLOCK DESCRIPTION

Graphical utilities.

BLOCK FUNCTIONS

Code: P01

Description: A sequence *seq* is paired with evenly spaced real numbers of a specified interval. The resulting sequence is adequate for the *Mathematica* ListPlot command.

Example:

{a,b,c}, {3,5} -> {{3,a}, {4,b}, {5,c}}

Syntax: **putOneSequence**[*seq*, *interval*]

Inputs:

seq: Sequence.

interval: {start, end} interval of real numbers.

Output : A sequence of pairs in which *seq* is paired with evenly spaced real numbers of the specified interval.

Code: P02

Description: This function generates a linear sequence of values.

Example:

the inputs

{3,5}, 5

generates the list of values

{3, 3.5, 4, 4.5, 5}

Example:

{a,b,c}, {3,5} -> {{3,a}, {4,b}, {5,c}}

Syntax: **generateLinParameterSeq** [*interval*, *n*]

Inputs:

interval: {start, end} interval of real numbers.

n: number of points.

Output : A sequence with evenly spaced real numbers of the specified interval.

Code: P03

Description: In order to plot seq2 vs seq1, this function constructs a seq of pairs; for example,

$\{2,7,8\}, \{6,3,1\} \rightarrow \{\{2,6\}, \{7,3\}, \{8,1\}\}$

If one of the sequences has an initial extra term, it is deleted:

$\{0,2,7,4\}, \{6,3,1\} \rightarrow \{\{2,6\}, \{7,3\}, \{4,1\}\}$.

The resulting sequence is adequate for the *Mathematica* ListPlot command.

Syntax: **putTwoSequences**[sx, sy]

Inputs:

sx: Sequence.

sy: Sequence.

Output : A sequence of pairs.

Code: P04

Description: From the three lists

paramSeq, alphaSeqRad, thetaSeqRad

It generates a new list with the values of the parameter and the coordinates (x,y) in the net for each point corresponding to the specified angles.

Syntax: **genParametrizedCurveSeq**[paramSeq,alphaSeqRad,thetaSeqRad]

Inputs:

paramSeq: The list $\{p_1, p_2, \dots, p_n\}$

alphaSeqRad: The list $\{\alpha_1, \alpha_2, \dots, \alpha_n\}$

thetaSeqRad: The list $\{\theta_1, \theta_2, \dots, \theta_n\}$

Output: The list $\{t_1, t_2, \dots, t_n\}$, where $t_i = \{p_i, x_i, y_i\}$.

Code: P05

Description: A list is sectioned into a sequence of lists overlapping one element.

Syntax: **listSectioning**[longList, p]

Inputs:

longList: a list

p: The length of the sublists.

Output: The list of the sublists.

Code: P06

Description: In order to plot a curve with 'tics' over the EPN(*), a list with graphic elements of subarcs is generated.

Syntax: **putCurve**[parCurveSeq, p]

Inputs:

parCurveSeq: A parameterized curve (generated by P04)

p: The length of the subarcs.

Output: A list with graphic elements of subarcs.

The successive values of the parameter at the 'tics' are printed.

(*) EPN is for Equiareal Projection Net.

Code: P07

Description: It generates a list with graphic elements of subarcs, including parameter values at the 'tics'.

Syntax: **putCurveWithLabels**[*parCurveSeq*, *p*]

Inputs:

parCurveSeq: A parameterized curve (generated by P04)

p: The length of the subarcs.

Output: A list with graphic elements of subarcs, including parameter values at the 'tics'.

The successive values of the parameter at the 'tics' are printed.

Code: P08

Description: It generates a graphic element for a labeled point.

Syntax: **myPointWithLabel**[*t*, *x*, *y*]

Inputs:

t: Numeric label.

x: Point abscissa.

y: Point ordinate.

Output: Graphic element for the point including label.

Code: P09

Description: It generates a graphic element for a point.

Syntax: **myPoint**[*x*, *y*]

Inputs:

x: Point abscissa.

y: Point ordinate.

Output: Graphic element for the point.

Code: P10

Description: It generates a list with graphic primitives for plotting an arc without 'tics'.

Syntax: **putSimpleArc**[*parCurveSeq*]

Inputs:

parCurveSeq: A parameterized curve (generated by P04), a list {{*t*,*x*,*y*}, ...}.

Output: A list with graphic elements of the arc (without 'tics').

Code: P11

Description: It generates a graphic primitive list to draw a EPN(*).

Syntax: **generatesTemplate**[*deltaTheta*, *smallDeltaTheta*, *deltaAlpha*, *smallDeltaAlpha*, *thin*, *thick*, *pale*, *dark*]

Inputs: Parameters defining the style of the template.

deltaTheta, *smallDeltaTheta*, *deltaAlpha*, *smallDeltaAlpha*, *thin*, *thick*, *pale*, *dark*.

Example:

degrees=Pi/180.;

deltaTheta=15 degrees;

smallDeltaTheta=5 degrees;

deltaAlpha=15 degrees;

smallDeltaAlpha=5 degrees;

(* Absolute thickness *)

thin=0.25;

thick=0.35;
(* Gray levels *)
pale=0.75;
dark=0.35;
Output: The list of graphic elements of the EPN.
(*) EPN is the Equiareal Projection Net.

Code: P12

Description: It generates a list with graphic primitives for plotting an arc without 'tics'.

Syntax: **plotCurveAndTemplate**[*curve*, *template*]

Inputs:

curve: List of graphic primitives of the curve.

template: List of graphic primitives of the EPN (generated by P11).

Output: Combined plot of curve and template.

Sequence Processing (SP)

BLOCK DESCRIPTION

Sequence manipulations.

BLOCK FUNCTIONS

Code: SP01

Description: From a list *s*, whose elements are also lists, this function constructs another list with the *na*_th elements of the elements of *s*.

Example:

With *s*={{3,2,9},{6,7,-4},{1,8,9},{5,3,4}} and *na*=2 the function gives the list {2,7,8,3}.

Syntax: **selectArgument**[*s*, *na*]

Inputs:

s: A list whose elements are also lists.

na: Ordinal number of the selected element.

Output: The list with the *na*_th elements of the elements of *s*.

Code: SP02

Description: It composes two sequences.

Syntax: **composeSequences**[*G1*, *G2*]

Inputs:

G1: A sequence.

G2: A sequence.

Output: A sequence made by the elements of *G1* first and *G2* last.

Change of reference (CHR)

BLOCK DESCRIPTION

IJK are the kinematic axes for the description of the deformation and NET the

geographical axes. The function defined in this block can be used to obtain IJK coordinates from NET coordinates.

BLOCK FUNCTIONS

Code: CHR01

Description: This function computes the change matrix from NET axis to IJK axis.

Syntax: **computeBasisChangeMatrix**[*axis1Angles*, *alternateDirection13Angles*]

Inputs:

axis1Angles: Angles corresponding to the direction of the kinematic axis 1.

alternateDirection13Angles: Angles corresponding to a direction u contained in the plane 13.

Output: 3x3 change matrix.

Code: CHR03

Description: A plane can be defined by its dip direction or by its normal vector. This function calculates the dip direction angles from the normal vector angles. The same function is used in the opposite conversion.

Syntax: **LMPvsNanglesChange**[*inputAngles*]

Inputs:

inputAngles: Angles of the normal vector (respectively dip direction).

Output: Dip direction angles (respectively normal vector).

Code: CHR04

Description: Using the NET coordinate system, this function obtains the Cartesian coordinates of a direction vector specified by its angles {alpha, theta}.

Syntax: **computeNETcoordinates**[*angles*]

Inputs:

angles: The list {alpha, theta} of the angles that define a direction u.

Output: The list {u₁, u₂, u₃} of the Cartesian components of the direction vector u.

Code: CHR05

Description: This function obtains the IJK coordinates of a direction vector from the NET coordinates.

Syntax: **fromNETcoordinatesToIJKcoordinates**[*NETcoordinates*, *changeMatrix*]

Inputs:

NETcoordinates: The list of the NET coordinates of a direction u.

changeMatrix: The change matrix (obtained by CHR01).

Output: The list of IJK coordinates of the same direction u.

Auxiliary functions (AF)

BLOCK DESCRIPTION

Several auxiliary functions.

BLOCK FUNCTIONS

Code: AF01

Description: Given two directions by the corresponding angles, it calculates the acute angle between them.

Syntax: **angleBetweenDirections**[*par1*, *par2*]

Inputs:

par1: The list {alpha1, theta1} of angles of the first direction.

par2: The list {alpha2, theta2} of angles of the second direction.

Output: The angle between the directions (degrees).

Code: AF02

Description: This function computes the principal directions of the strain corresponding to a given gradient matrix.

Syntax: **principalDirections**[*mC*]

Inputs:

mC: The 3x3 gradient matrix.

Output: It prints several geometrical data of the principal directions.

Code: AF03

Description: A principal direction is represented by the angles {*alpha*, *theta*}. This function calculates the *alpha* angle of the principal direction number *pdN*.

Syntax: **alphaPrincipalDirection**[*mC*, *pdN*]

Inputs:

mC: The 3x3 gradient matrix.

pdN: The number of the principal direction for which the angle is calculated.

Output: The alpha angle of the principal direction number *pdN*.

Code: AF04

Description: A principal direction is represented by the angles {*alpha*, *theta*}. This function calculates the *theta* angle of the principal direction number *pdN*.

Syntax: **thetaPrincipalDirection**[*mC*, *pdN*]

Inputs:

mC: The 3x3 gradient matrix.

pdN: The number of the principal direction for which the angle is calculated.

Output: The *theta* angle of the principal direction number *pdN*.

Code: AF05

Description: Given a gradient matrix *mC*, this function calculates the ratio $R_{13} = \sqrt{\lambda_1} / \sqrt{\lambda_3}$ between eigenvalues.

Syntax: **ratioR13**[*mC*]

Inputs:

mC: The 3x3 gradient matrix.

Output: The ratio $R_{13} = \sqrt{\lambda_1} / \sqrt{\lambda_3}$.

Code: AF06

Description: Given a gradient matrix *mC*, this function calculates the ratio $R_{12} = \sqrt{\lambda_1} / \sqrt{\lambda_2}$ between eigenvalues.

Syntax: **ratioR12**[*mC*]

Inputs:

mC: The 3x3 gradient matrix.

Output: The ratio $R_{12} = \sqrt{\lambda_1} / \sqrt{\lambda_2}$.

Code: AF07

Description: Given a gradient matrix *mC*, this function calculates the ratio

$R_{23} = \sqrt{\lambda_2} / \sqrt{\lambda_3}$ between eigenvalues.

Syntax: **ratioR23**[*mC*]

Inputs:

mC: The 3x3 gradient matrix.

Output: The ratio $R_{23} = \sqrt{\lambda_2} / \sqrt{\lambda_3}$.