

DYALOG

Elsinore 2023

APL and Metallurgy

Jesús Galán López (Spain)

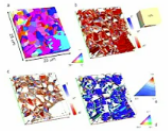
Dyalog / Ghent University



Computational problems
Materials Science and Engine

Microstructural Analysis

- Large amounts of SEM/EBSD and simulation data (2D/3D)
- Study properties, relationships, topology
- Calculate distributions
- Crystallographic texture analysis (boundaries, ODFs)
- Bridge between simulations and experiments



20:22

Dyalog and Academia // Jesús Galán López // Dyalog '22

232 views • 10 months ago



Dyalog Usermeeting

Jesús Galán López Jesús and Gitte introduce the plan for bringing Dyalog back into academia. The final objective is to position ...



Introduction and background | STEM in academia and where APL fits | Project goals | Action plan |... 9 chapters ▾



Dyalog and Academia

- Get visibility in academic and research environments
- Learn about them and from them
- Study and improve the use of APL as a tool for education and research of technical topics
- Introduce a new generation to APL
- Establish fruitful relationships with the academic world



APL and Metallurgy

- Use APL to teach / learn about complex topics
 - Mathematical methods
 - Crystallography and crystal plasticity
 - Modelling of thermo-mechanical processes
- Evaluate capacities of APL as a teaching tool when compared with mainstream languages
- More presence in the classroom and the student community



APL and Metallurgy

- 1. Initial plan**
- 2. Results**
- 3. Conclusions**
- 4. Next**

What happened

- ✦ Kept goals
- ✦ Merged some tasks, split some others
- ✦ Amount of work and time sometimes different to expectations
- ✦ Results



Evaluation of state of the art

- ◆ Programming languages and software at the university
 - ◆ Review current literature
 - ◆ Personal point of view and real cases
 - ◆ Programming languages strengths and weaknesses
 - ◆ Report



Evaluation of state of the art

- ◆ APL at the faculty
 - ◆ Unheard of
 - ◆ Looks hard (those symbols!)
 - ◆ How do I ... in APL?
 - ◆ How does APL compare to ...?
 - ◆ Can APL solve *my* problem?



Development

- Introduction tutorial
 - No previous programming experience
 - Cover: syntax, primitives, dfns, system
 - Tradfns, namespaces, classes only mentioned
 - Simple examples
 - Jupyter notebook and slides



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```
++-x÷*@[]!?!|_|⊥T-H=≠≤<>≥≠vλλ̃v†‡c>ε[]⊥ψi⊥εuη~/\ƒλ,τρφθϕ''""*.öë@[]@[]@[]I±∓⊕⊖+ωα∇&~θΔΔ{}[]()
```

↑

Arithmetic

	3+2
5	
	3-2
1	
	3×2
6	
	3÷2
1.5	

jgl@dyaLog.com



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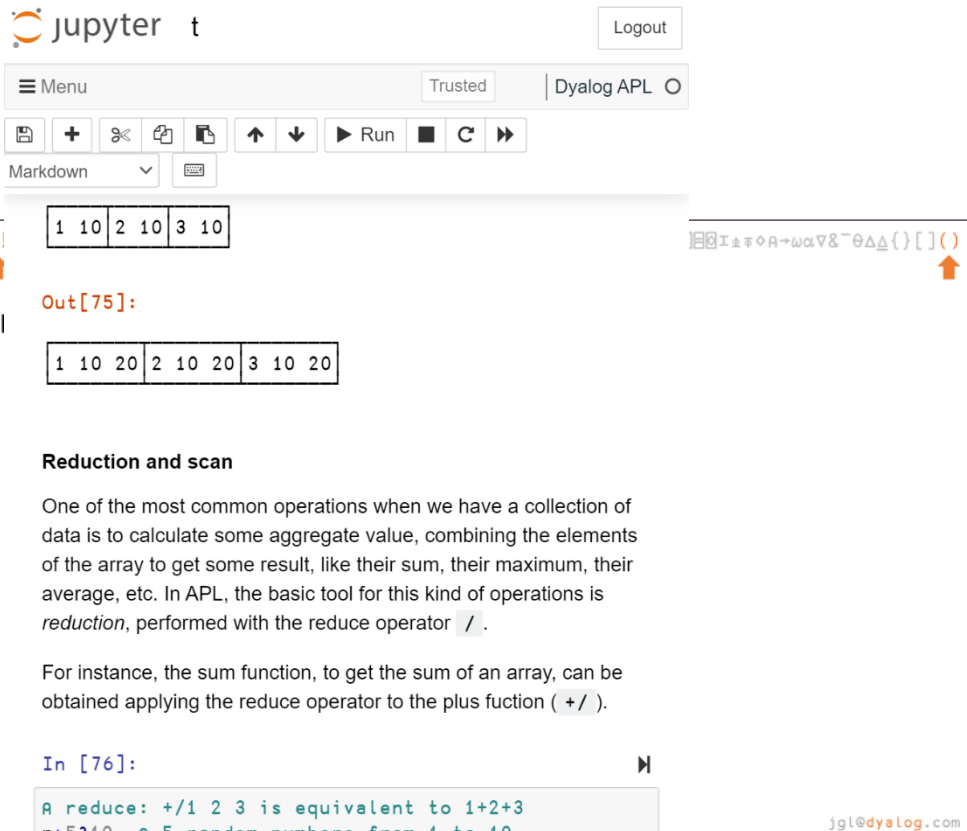
```
++-x÷*⊞o!?!|_|⊥T→H=≠≤<>≥≡≠vλã∨†‡c>ε[]⊞ψι⊥εεun~/\fλ,;ρφθϕ''''".°öë@□□⊞⊞⊞I±∓⊕⊖+ωα∇&^-θΔΔ{}[]()
```

Trains

(i j f h g)

jgl@dyaLog.com





The screenshot shows a Jupyter notebook interface. At the top, there is a 'Logout' button and a 'Trusted' status indicator. Below that is a 'Menu' button and a 'Dyalog APL' logo. The main toolbar contains icons for file operations, navigation, and execution. The code cell contains the following APL code:

```
1 10 2 10 3 10
```

The output of the code is:

```
Out[75]:
```

1 10 20	2 10 20	3 10 20
---------	---------	---------

The code cell also includes a title 'Fu' and a toolbar with mathematical symbols. Below the code cell, there is a section titled 'Reduction and scan' with the following text:

One of the most common operations when we have a collection of data is to calculate some aggregate value, combining the elements of the array to get some result, like their sum, their maximum, their average, etc. In APL, the basic tool for this kind of operations is *reduction*, performed with the reduce operator `/`.

For instance, the sum function, to get the sum of an array, can be obtained applying the reduce operator to the plus function (`+/`).

In [76]:

```
A reduce: +/1 2 3 is equivalent to 1+2+3
r+5210 A 5 random numbers from 1 to 10
(c,+,x/,[,[/,-/,+/,+/#) A array, summation, pr
A dyadic: reduce on windows (overlapping subarrays o
2,/r 2-~/r A concatenate and subtract elements o
3,/r 3+/r A concatenate and sum elements on win
```

The output of the code is:

```
Out[76]:
```

1 2 5 7 4	19	280	1	7	1	4	3.8
-----------	----	-----	---	---	---	---	-----

The screenshot also shows a footer with the email address 'jgl@dyaLog.com' and a logo of a traditional APL hat.

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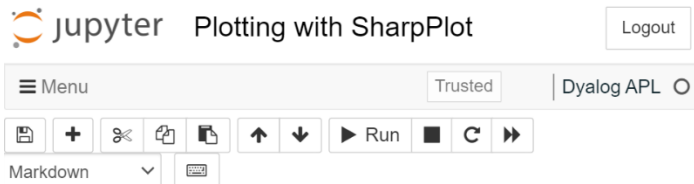
Development

- ◆ Specialized tutorials: engineering (jupyter)
 - ◆ Plotting
 - ◆ Linear fitting
 - ◆ Simple calculus
 - ◆ Geometric algebra
 - ◆ Formulae (WIP)



- Specialized tutorials: engineering (jupyter)

- Plotting
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- Formulae (WIP)



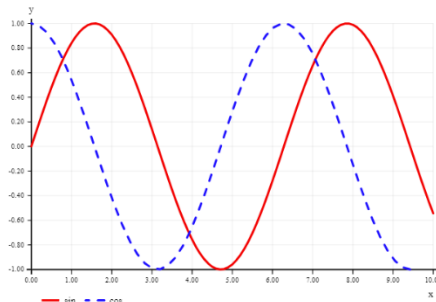
```
A Plot and display
:For i :In 1#data
  sp.SetColors i[]colors
  sp.DrawLineGraph i>data
:EndFor
svg←sp.RenderSvg Causeway.SvgMode.FixedAspect
```

The return of `Plot` is the plot as SVG, so we use `]html` to display it.

In [33]:

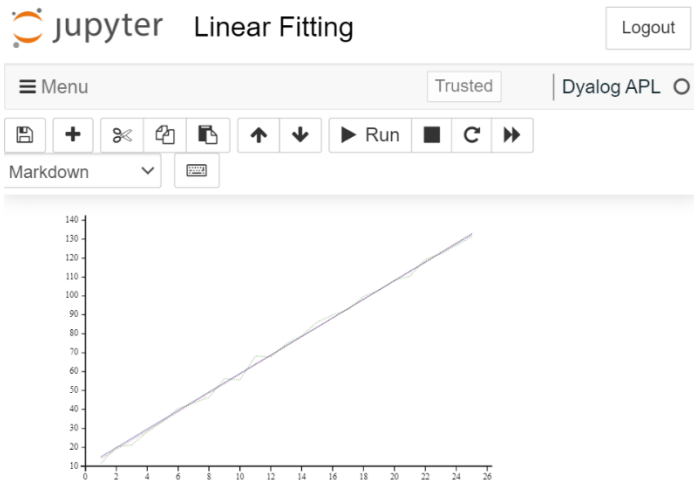
```
A eg
]html 'red' 'blue' ('x'Plot'y') ('sin' sin) ('cos'
]html 'purple' ('x'Plot'y') 'log' ((⊖1+x) x)
]html ('Time [s]'Plot'Distance [m]') ('a = 20' ((x
```

Out[33]:



- Specialized tutorials: engineering (jupyter)

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R-squared value

We still need to determine the [R-squared](#) value, which can be calculated as one minus the ratio of the sums of squares ($\sum (y - \hat{y})^2$) of differences between the real and estimated values and between the real values and their mean:

In [8]:

```
1 - (sum((y - a + b * x) ** 2) / sum((y - mean(y)) ** 2))
```

Out[8]:

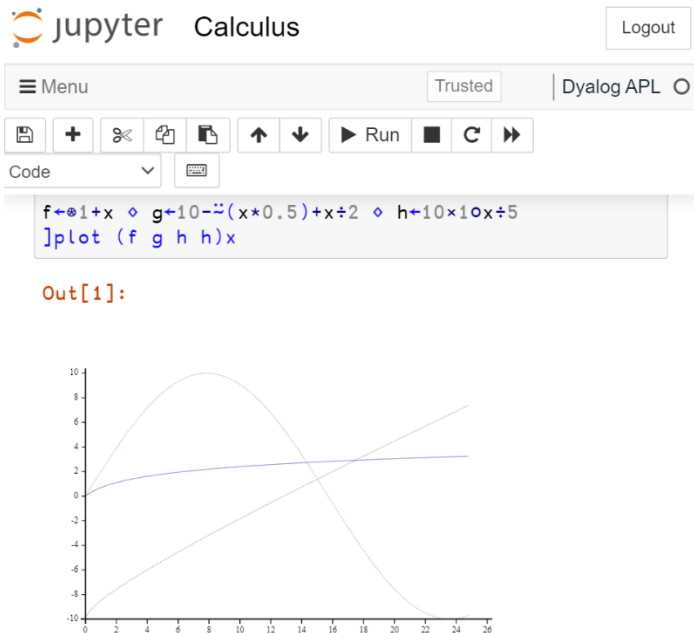
```
0.9971435494
```

The R-squared value is defined not only for linear fittings, but also for any other fitting we perform on some data. Therefore, we can define an operator that takes the fitted function as operand (in our



Specialized tutorials: engineering (jupyter)

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Derivative

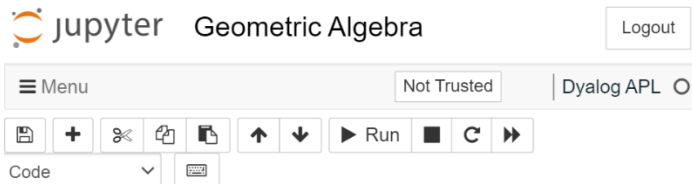
The derivative will be calculated as a simple ratio of the difference between two consecutive values. The differences will be calculated using the n-wise reduction operator ([dyadic /](#)).

In [2]:

```
dy←2-/f ◊ dx←2-/x ◊ d←dy÷dx  
]plot d x
```

Out[2]:





The result of squaring a vector is the square of its magnitude. To multiply two parallel vectors v and w defined, respectively, as $a \times u$ and $b \times u$:

$$\begin{aligned}(v \Delta v w) &\equiv (a \times u) \Delta v b \times u \\(v \Delta v w) &\equiv a \times b \times u \Delta v u \\(v \Delta v w) &\equiv a \times b\end{aligned}$$

So, the product of two parallel vectors is the scalar that results from multiplying their magnitudes.

Geometric product of perpendicular vectors

Every vector can be decomposed in two perpendicular components in some base. For example:

```
In [24]:  
  
v1 v2 ← 3 2   A eg  
Assert (⇒ v1 v2 +.x (1 0) (0 1)) ≡ (M×U) v1 v2   A  
Assert 0 ≡ v1 0 +.x 0 v2   A  
'v' 'v1' 'u1' 'v2' 'u2' Table (3 2) 3 (1 0) 2 (0 1)  
'v1 0 +.x 0 v2' 'm ← M v' 'u ← U v' 'm × u' Table (
```

Out[24]:

v	v1	u1	v2	u2
3 2	3	1 0	2	0 1

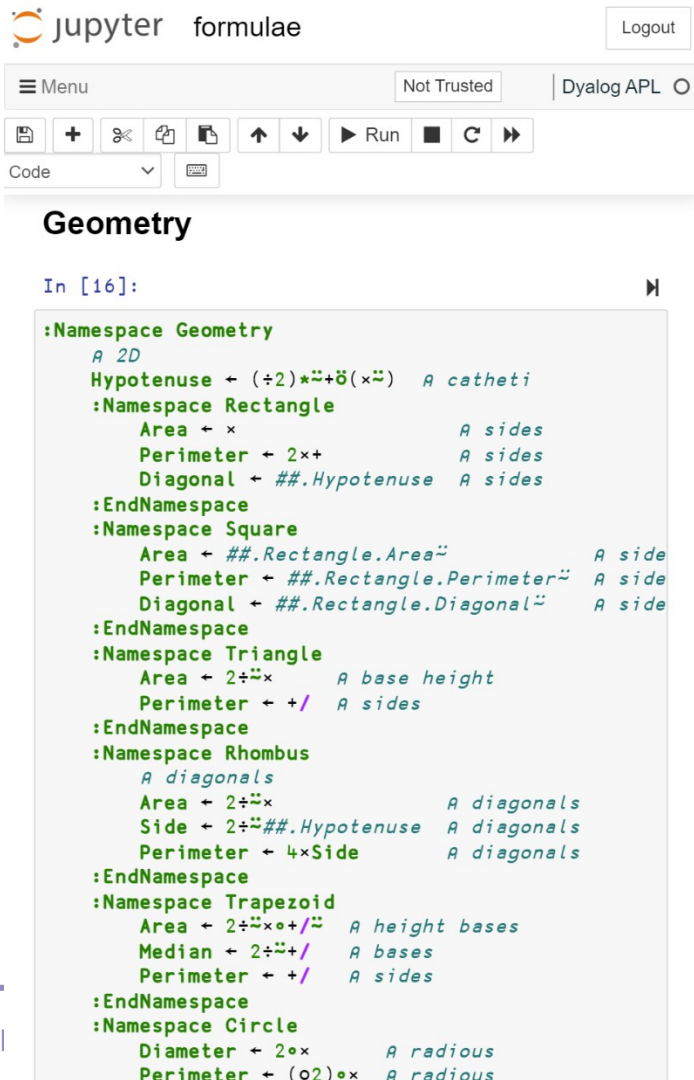


Development

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Rectangle.Area $\leftarrow x$
Square.Area $\leftarrow x^2$

A uniform acceleration
 $t \perp (a \div 2) v_0 x_0$



The screenshot shows a Jupyter Notebook interface with the following elements:

- Top bar: Jupyter logo, "formulae", and a "Logout" button.
- Navigation: "Menu", "Not Trusted", and "Dialog APL" options.
- Toolbar: Icons for file operations, navigation, and execution (Run, Stop, Refresh, Next).
- Code cell: A code cell titled "Geometry" containing the following code:

```
In [16]:  
  
:Namespace Geometry  
  A 2D  
  Hypotenuse  $\leftarrow (\div 2) * \sqrt{x^2 + y^2}$  A catheti  
:Namespace Rectangle  
  Area  $\leftarrow x * y$  A sides  
  Perimeter  $\leftarrow 2 * (x + y)$  A sides  
  Diagonal  $\leftarrow \#.Hypotenuse$  A sides  
:EndNamespace  
:Namespace Square  
  Area  $\leftarrow \#.Rectangle.Area$  A side  
  Perimeter  $\leftarrow \#.Rectangle.Perimeter$  A side  
  Diagonal  $\leftarrow \#.Rectangle.Diagonal$  A side  
:EndNamespace  
:Namespace Triangle  
  Area  $\leftarrow 2 \div x * y$  A base height  
  Perimeter  $\leftarrow + /$  A sides  
:EndNamespace  
:Namespace Rhombus  
  A diagonals  
  Area  $\leftarrow 2 \div x * y$  A diagonals  
  Side  $\leftarrow 2 \div \#.Hypotenuse$  A diagonals  
  Perimeter  $\leftarrow 4 * Side$  A diagonals  
:EndNamespace  
:Namespace Trapezoid  
  Area  $\leftarrow 2 \div (x + y) * h$  A height bases  
  Median  $\leftarrow 2 \div (x + y)$  A bases  
  Perimeter  $\leftarrow + /$  A sides  
:EndNamespace  
:Namespace Circle  
  Diameter  $\leftarrow 2 * r$  A radius  
  Perimeter  $\leftarrow (Q2) * r$  A radius
```



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Development

- ◆ Specialized tutorials: materials (jupyter)
 - ◆ Analysis of tensile experiments
 - ◆ Crystallographic orientations and misorientations
 - ◆ Grain growth



- Specialized tutorials: materials (jupyter)
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Introduction

This notebook presents a short tutorial on how to use Dyalog APL for the analysis of tensile experimental data. The tutorial is directed towards scientists, engineers and students who have already learnt the basics of the APL language and intend to use Dyalog APL for their data analysis work.

The topic of tensile data analysis is chosen because it represents a well known task for many engineers of different disciplines, and it is a perfect example of the more general procedure of reading data from a file, doing some basic processing, and plotting the obtained results, which is familiar to many engineering researchers and students.

Tensile analysis

Tensile testing is one of the most common experimental methods for the determination of the mechanical properties of materials. Materials scientists employ tensile tests to find different properties that define the mechanical behaviour of materials, and mechanical, aeronautical, and civil engineers, use these properties for the design of structures.

In a uniaxial tensile test (the most common kind of tensile test), a specimen with a predefined geometry is subjected to a uniaxial load that deforms the material under **controlled conditions**. The force during the experiment is measured with a load cell, while the deformation on the sample is measured either directly on the specimen using an extensometer or a strain gauge, or indirectly based on the crosshead displacement of the testing machine. From this data, strain and stress are calculated, respectively, as displacement with respect to the original length and force divided by cross sectional area (width multiplied by thickness):

$$e_{eng} = \frac{\Delta l}{l_0} = \frac{l - l_0}{l_0} \quad S_{eng} = \frac{F}{A} = \frac{F}{w \cdot t}$$

It is customary to present this data as a **tensile diagram** or strain-stress curve.

During a tensile experiment, both the length and the area of the specimen change. When strain and stress are calculated with respect to the initial values, as they are in the formulas presented above, we will call them *engineering strain* and *engineering stress*. When they are calculated with respect to the instantaneous values, we will call them *true strain* and *true stress*. Since true values are not dependent on initial conditions, true values are not dependent on a specific specimen geometry, at difference of the corresponding engineering values. Assuming that the volume is conserved, true and engineering strain and stress will be correlated by the equations:

$$e_{true} = \ln(1 + e_{eng}) \tag{1}$$

$$S_{true} = S_{eng} (1 + e_{eng}) \tag{2}$$



Development

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```
[-0.000397524 4.758206844]
```

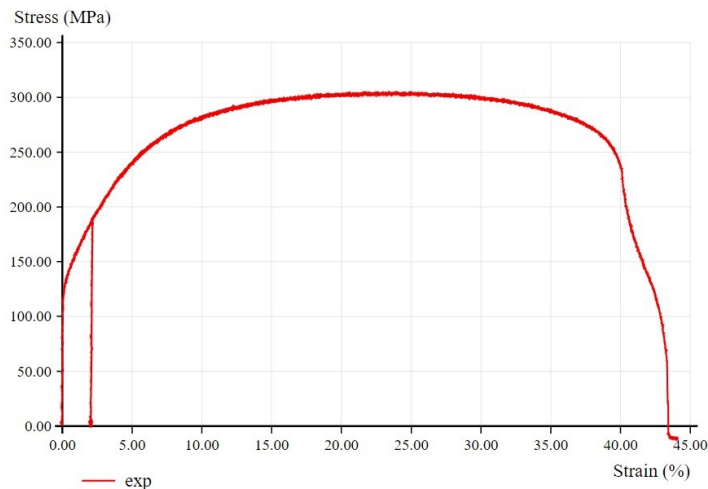
We will store the strain-stress data as a vector with two columns. Since it is a trivial operation, we will not define specific functions to extract the strain and stress data, and will instead use `>` for strain and `>φ` for stress.

Data files might be quite large, and there could be a significant amount of experimental noise:

```
In [10]: 'Number of rows:', #=>e  
        ]html PlotTD 'exp' e
```

```
Out[10]: Number of rows: 20040
```

```
Out[10]:
```



We observe that the data does not only contain a very large number of points and experimental scatter. Other issues are that it does not exactly start at zero, there are some suspicious points around 2% strain which are assumed to be erroneous and need to be discarded, and the measurement continues after fracture of the sample, with data that is not relevant.

Young modulus and initial strain offset

In order to find the Young modulus, we need to perform a linear fitting. The curve should start at strain zero,



Development

Specialized tutorials: materials (jupyter)

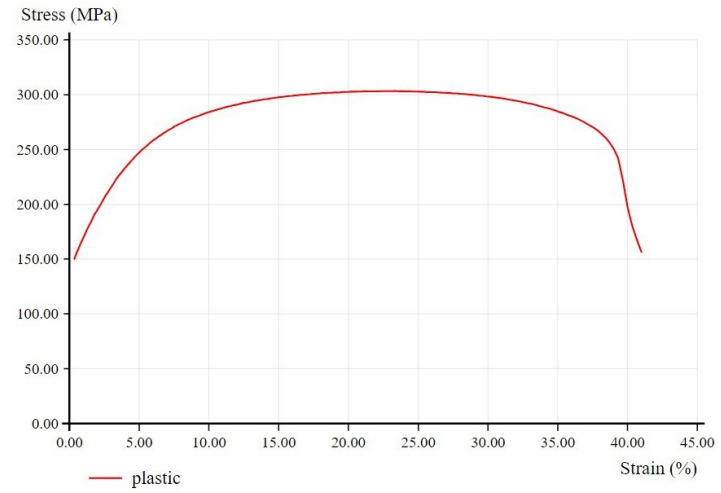
- Analysis of tensile experiments
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(x+y).

The `Plastic` operator allows to specify what is the yielding offset as `αα` (usually defined as the point of 0.2% or 0.02% of plastic strain), then uses `Clean` to remove the points with plastic strain higher than specified.

```
In [18]: Plastic ← {x y ← α α α ≤ 0 → Clean ω - (y + x → φ ω) 0} A get plastic curve from total tensile curve  
A plastic curve with 0.2% yield offset  
p ← x y (0.002 Plastic) c  
]html PlotTD 'plastic' p
```

Out[18]:



True tensile curve, hardening rate, and UTS point

After calculating the plastic strain, our next goal is to calculate the true curve and the strain hardening rate. The true strain can be calculated from the engineering strain as $\ln(1 + \epsilon)$. To calculate the true stress, we need the engineering stress and the engineering strain. If the engineering strain is the left argument and the engineering stress the right one, it is calculated as $\frac{\sigma}{1 + \epsilon}$. The strain hardening rate is calculated as the derivative of the true plastic curve using `D`, from the [Calculus notebook](#). The UTS point is the point where the true curve and the strain hardening rate intercept.

The `TrueRate` function is used to calculate both the true curve until the UTS point (the uniform deformation point) and the strain hardening rate. The `TrueRate` function is used to calculate both the true curve until the UTS point (the uniform deformation point) and the strain hardening rate.



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(x+y).
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```
In [18]: Plastic ← {x y ← α α α ≤ 0 ⇒ Clean ω - (y - x + φ ω) 0} A get plastic curve from total tensile curve
A plastic curve with 0.2% yield offset
p ← x y (0.002 Plastic) c
]html PlotTD 'plastic' p
```

Out[18]:



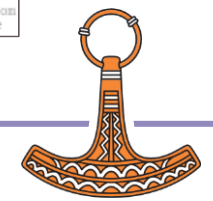
Resampling and cleaning

```
Resample ← {(α-1) {(σ ω), ÷ ÷ n + / ω p ~ α, n + [ α ÷ ~ ≠ ω ] ~ ω} A average rows
Smooth ← {α+3 ÷ ω Resample ~ [ α ÷ ~ ≠ ω ] A average rows
Clean ← {α+1 ÷ (α α α ω) ÷ / ~ ~ ~ (c * 2 Δ ÷ ÷) ~ ω / ~ ~ ~ c ≠ ÷ ω} A select points
c ← ((0.02 > ÷) v 150 < ÷ ÷ φ) Clean 250 Resample e A strain < 0.02 or stress > 150
]html PlotTD 'clean' c
```

jgl@dyalog.com
jesus.galanlopez@ugent.be

true strain can be calculated from the engineering strain as $\epsilon \ln(1 + \epsilon)$. To calculate the true stress, we need the engineering stress and the engineering strain. If the engineering strain is the left argument and the engineering stress the right one, it is calculated as $\sigma(1 + \epsilon)$. The strain hardening rate is calculated as the derivative of the true plastic curve using `D`, from the [Calculus notebook](#). The UTS point is the point where the true curve and the strain hardening rate intercept.

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```

:Namespace Euler
  A quaternion product
  c ← t(0 1 2 3)(1 0 3 2)(2 3 0 1)(3 2 1 0)
  u ← t(1 -1 -1 -1)(1 1 1 -1)(1 -1 1 1)(1 1 -1 1)
  QP ← u{+/αα[-2t+r+1]α×[(i(≠ρα)-1),r+IO](cIO+ωω)[(r+≠ρω)-1-IO]ω}c
  QC ← ×i1o(1,-3p1)
  QD ← +.×i1oQC

  A quaternion from Euler angles
  RD ← 180÷ω
  UV ← {ω×[(i(≠ρω)-1)÷(+/-ω)]*÷2}
  QA ← {(2ω), (1ω)◦.×UVα}◦(÷2) ◊ QAD ← QA◊RD
  QE ← {α+(0 0 1)(1 0 0)(0 0 1) ◊ αQP.QA=[-1+i≠ρω]ω} ◊ QED ← QE◊RD

  A cubic symmetry
  cs ← c1 0 0 0
  cs,+, (1 0 0)(0 1 0)(0 0 1) ◊ QAD 90 180 270
  cs,+, (1 1 1)(-1 1 1)(1 -1 1)(1 1 -1) ◊ QAD 120 240
  cs,+, (1 1 0)(1 0 1)(0 1 1)(1 -1 0)(-1 0 1)(0 1 -1)◊.QAD 180

  A misorientations
  ML ← {α+0.5 ◊ [(180×ω)÷α]}
  MC ← cs{2×-2o1[≠f|αα◦.QD=αQP=QCω]}
  IM ← {IO+1-≠(α[ω]+/-≠i(α[ω]-1)}
  CM ← {c+INSθ ◊ c.m+1p≠(-IM-1)≠, tω ◊ c}
  _M_ ← {α=ω:0 ◊ 0sm+(i+αIMω)÷ωω.m:m ◊ t(i=ωω.m)+MLα(MC)δ(=αα)ω}

  A namespace with function M to calculate misorientations
  M ← {α+0.5
    m+INS'QD' 'QP' 'QC' 'IM' 'MC' ◊ m.ML+α◊ML
    m.M+(α÷2)+α×(+ω)_M_(CMω) ◊ 2≠ρω: m
    m.M+m.Mδ((-1+ρω)◊{IO+α±ω-IO}) ◊ m
  }

  Space ← {t>.,/(ω÷2)+ω×(i"360 90 90÷ω)-1}
  Random ← {F+{(1 2o2<c◊?αp0)×cω÷2} ◊ UVt[1]ω(F◊(1◊-),F)?ωp0}
:EndNamespace
  
```

A product components $\omega=0$
 A product unit factors
 A quaternion product
 A quaternion conjugate
 A quaternion dot product

A radians from degrees
 A unitary vector
 A quaternion from axis-angle
 A quaternion from Euler angles (zxz)

A identity
 A 4-fold around <001>
 A 3-fold around <111>
 A 2-fold around <110>

A misorientation level from radians
 A misorientation with cubic symmetry
 A index of misorientation
 A cache of misorientations
 A memoization

A misorientation step
 A namespace with curried ML
 A misorientation function
 A over list index for higher rank

A Euler space of given step
 A random orientations of given step



Development

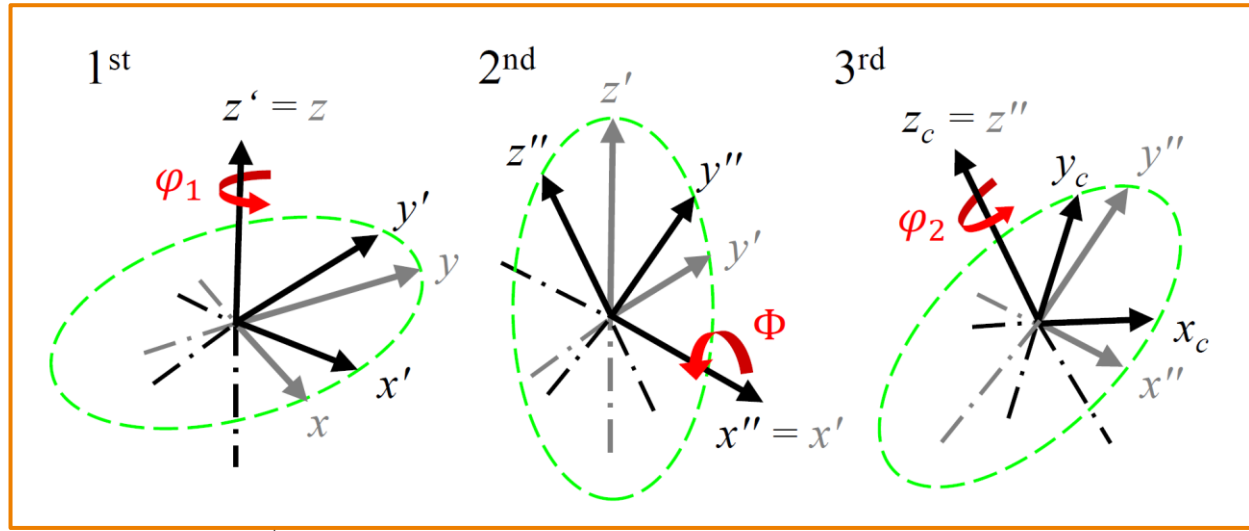
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  u ← t(1 -1 -1 -1)(1 1 1 -1)(1 -1 1 1)(1 1 -1 1)
  QP ← u{+/αα×[-2t+r+1]α×[(i(≠ρα)-1),r+ΠIO](c-ΠIO+ωω)[(r+≠ρω)-1-ΠIO]ω}c
  QC ← ×∂1o(1,-3p1)
  QD ← +.×∂1oQC
  
```

A product components $\square IO=0$
 A product unit factors
 A quaternion product
 A quaternion conjugate
 A quaternion dot product



angle
angles (zcx)

from radians
cubic symmetry
ns

ML
n
gher rank

```

Space ← {t>o.,/(ω≠2)+ω×(i''360 90 90÷ω)-1}
Random ← {F+{(1 2o2×c0?ωρ0)×cω*≠2} ∘ UVt[1]ω(Fo(1o-),F)?ωρ0}
:EndNamespace
  
```

A Euler space of given step
 A random orientations of given shape



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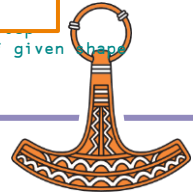
1st

```

:Namespace Euler
  A quaternion product
  c + t(0 1 2 3)(1 0 3 2)(2 3 0 1)(3 2 1 0)
  u
  QF
  QQ
  QQ
      
```

```

      A product components []IO=0
      s (z x z)
      adians
      ymmetry
      ank
      A random orientations of given haph
      
```



Development

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1st

```

:Namespace Euler
  A quaternion product
  c + t(0 1 2 3)(1 0 3 2)(2 3 0 1)(3 2 1 0)
  u
  QF
  QQ
  QQ
        
```

A product components $\square I=0$

(z x z)

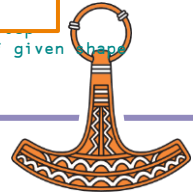
adians
ymmetry

rank

```

Sp
Random + {F+((1 2o2<co?ap0)<w*+2) o UVt[1]w(Fo(1o-),F)?wp0}
:EndNamespace
        
```

A random orientations of given shape



Development

- Specialized tutorials: materials (jupyter)

- Analysis of tensile experiments
- Crystallographic orientations and misorientations
- Grain growth

1st

```

:Namespace Euler
  A quaternion product
  c + t(0 1 2 3)(1 0 3 2)(2 3 0 1)(3 2 1 0)
  u
  QF
  QQ
  QQ
        
```

A product components $\square I=0$

(z x z)

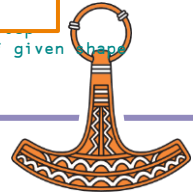
adians
ymmetry

rank

```

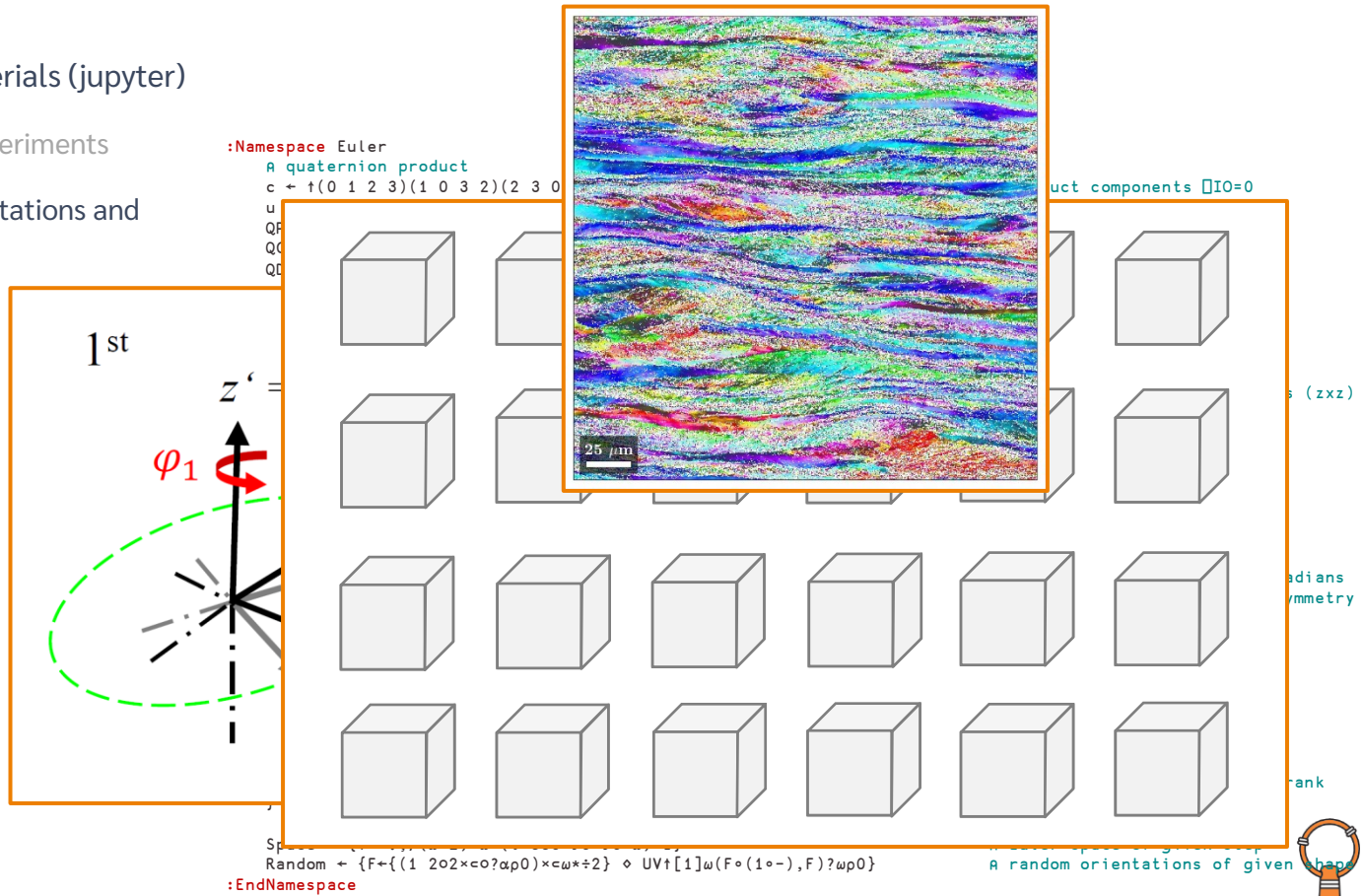
Sp
Random + {F+((1 2o2xco?ap0)x<w*+2) o UVt[1]w(Fo(1o-),F)?wp0}
:EndNamespace
        
```

A random orientations of given shape



Development

- Specialized tutorials: materials (jupyter)
 - Analysis of tensile experiments
 - Crystallographic orientations and misorientations
 - Grain growth



Development

- Specialized tutorials: materials (jupyter)
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 - Crystallographic orientations and misorientations
- Grain growth



Grain Growth Modelling in APL



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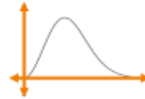
Development

- Specialized tutorials: materials (jupyter)
 - Analysis of tensile experiments
 - Crystallographic orientations and misorientations
- Grain growth

Grain Growth (Geiger, 2001)

[https://doi.org/10.1016/S1359-6454\(00\)00352-9](https://doi.org/10.1016/S1359-6454(00)00352-9)

- Thermal energy



Maxwell-Boltzmann
distribution

$$G_T(T) = -R T \log(x)$$

- Boundary energy



Read-Shockley
equation

$$\Delta\theta_{ij} = \frac{\pi(q_i - q_j)}{2 q_{max}}$$
$$G_{Bij}(\Delta\theta_{ij}) = G_0 \sin(\Delta\theta_{ij}) (1 - \log(\sin(\Delta\theta_{ij})))$$

- Activation energy $G_A = 10000 \text{ J/mol}$

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Development

- Specialized tutorials: materials (jupyter)

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- Grain growth

Grain Growth (Geiger, 2001)

https://

Grain growth operator

```
gg={
  A α: parameters ω: repetitions
  A α: temperature ω: initial orientations
  (GA GO SF RP)+αα      A simulation parameters
  qmax+[/,ω             A maximum orientation
  A solve:
  (q a)+α(next*ωω)ω(area ω) A get final orientations and list of areas
  q(2*(a+01)*+2)        A return final orientations and list of diameters
}
```

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Development

- Specialized tutorials: materials (jupyter)
 - Analysis of tensile experiments
 - Crystallographic orientations and misorientations
 - Grain growth

Grain Growth (Geiger, 2001)

https://github.com/jgalanlopez/grain-growth

Grain growth operator

```
gg←{  
  A←(q←q)⊖(q←q)  
  (q←q)⊖(q←q)  
}
```

Results

The first plot, 'Average Diameter', shows a linear increase from 0 to approximately 17.5 over 175 time units. The second plot, 'Circle Diameter', shows a non-linear decrease from approximately 24 to 0 over 100 time units.

The top row shows five heatmaps at time steps 0 (0.0), 10 (0.396), 20 (0.63), 30 (0.86), and 40 (1.03). The bottom row shows five heatmaps at time steps 50 (1.41), 60 (1.74), 70 (2.06), 80 (2.38), and 90 (2.61). The heatmaps show a transition from a noisy, multi-colored state to a smooth, single-colored state.

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Development

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 - Grain growth

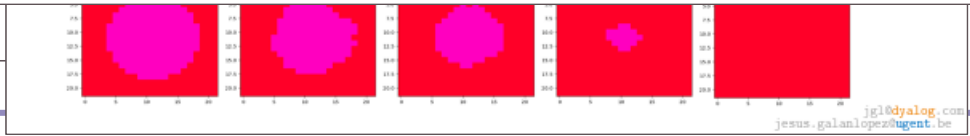
DIALOG
Elsinore 2023

Grain Growth and Array Programming

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jesus.galanlopez@ugent.be



Grain
https://
●
●
●



Development

- ◆ Specialized tutorials: materials (jupyter)
 - ◆ Analysis of tensile experiments
 - ◆ Crystallographic orientations and misorientations
 - ◆ Grain growth



Research

- Geometric algebra
- Grain growth
- Crystal plasticity



- Geometric algebra
- Grain growth
- Crystal plasticity

```

:Namespace ga A
  @IO = 0
  @CT = 1e-10
  Assert = {a~assertion failure' o Oe:ia [signal 8 o shy=1]

  A vectors
  M = 0,5+2i, u = U = +M; o I = +(a+u); o BV = {1+2i+1i}
  AV = c^2i[-+2] o VA = c^2i[-+2]

  A vector extension
  C = (a~@CT o @CT+u o 1#u; u o a~.|e:u:0 o u)
  S = (8#u:0 o (1#u) o (0#u) o 1#u; u o u) o US = +(u/v;u);
  I = (a~@CT o @CT+u o S u; +1-(1+u+1+e) u) o E = #B2BC
  X = (a~2+Du o a~u,0p^0[a~#u] o _X = (a~+ o a(aaB((aI@#u)*X)u)
  _V = (a~+ o 2+c(a aa_X)u) o _I = (0#u:0 o 0#u:0 o a(aa)u)

  A multivectors
  B = (a~2+1#u o r=a|6p u o v=0p^1+|f, a~rta o v[a]++rtu o v)
  B8 = +i+k2+1i; C = ((f/r)/(O#C)@B8#u) o D = [2#B8
  GB = (b~2((ccu)B~/B8#u))C, u o a=i#b o (aa)]](1+/,a)X b)

  A product tables
  PB = 2+u o SB = 1+(a~1+1)u, a1+ o MS = +1 1 0, p^2+3#
  FM = ((aSBu)*(aa)u, a~u) o TS = ((a, PB, B8, ((MSu)_FM))B8 2+u/v)
  TX = TS 0, +/ o TI = 0 0#u+o#TS o TR = SB^2B8# o BT = +V/

  A geometric operators
  _d = (p s+(a~u^2)u o (rtp)V(rts)+r(r#^au)[pp]ta+,{aa}2)B,u)
  _dI = (O V X (aa~u,aa^2)u) o _dX = [(aa,aa,u)u)
  _d = _d (TS^D) o _dX = _d (TX^D) o _dI = _dI (TIFD)
  _d = (u LC a aaB((a(u^2)u+RC)u) o _d = _d D
  _I = (a~1 o 2+Ca aa c+h^2)B(1+1#B^2[1-+2]2aa^BV#)#)USuX^2+Du)
  _C = (a~Du o (a~Vv/aa~+B^2BV#)#u X^2+u)

  A geometric functions
  d = _d o dX = _dX o dI = _dI
  d = d _d o dX = dX _d o dI = dI _d
  LC = d _C o RC = d^2_C o DC = LC^2
  aa = d _I o dd = d _I o R = +*TR; o PS = 1+2+2+2#

  GA = {
    g = D#B8 o g_s g_b = u (S BV 2+d+u)
    g_I = 2 o g_u X o g_I = 2 o g_S o g_US = US o g_C = C
    g_V = V o g_V = V o g_D = D o g_BV = BV o g_d = d
    g_d = _d (TSu) o g_dX = _d (TXu) o g_dI = _dI (TIu)
    g_d = g_d _d o g_dX = g_dX o g_dI = g_dI
    g_dV = g_dBV o g_dXV = g_dXV o g_dIV = g_dIV
    g_dV = g_dV o g_dXV = g_dXV o g_dIV = g_dIV
    g_d = g_d _d o g_dX = g_dX o g_dI = g_dI
    g_dV = g_dV o g_dXV = g_dXV o g_dIV = g_dIV
    g_LC = g_d _C o g_RC = g_d^2_C o g_DC = g_LC^2
    g_R = *(TRu) o g_X = X(a~B o a(aa)u) o g
  }

  g0 g1 g2 g3 c q pgs = GA^(14),(0 1) (0 2) (3 0 1)

  VS = {
    - Assert (e0) A, # p^" a b = a
    - Assert (c,1) A, z p^p^" u v w = w
    - Assert v E v v o
    - Assert v E I x v
    - Assert 0 E v +_V =v
    - Assert (v +_V u) E v +_V v
    - Assert (a x v) E v x a
    - Assert (u +_V v +_V u) E u +_V u +_V v
    - Assert (v x a + b) E (v x a) + b
    - Assert (a + v +_V u) E v +_V (a+v) u
    - Assert (v x a + b) E a +_V (v x) b
  }

  _GA = {
    - Assert (a b = a) VS (u v v + u)
    - Assert (u aa v aa w) E w aa^2 u aa v
    - Assert (u aa v +_V v) E (v +_V u aa v)
    - Assert (aa v +_V v) E (v +_V aa v)
    - Assert (a aa v aa b) E a x b v
    - Assert (v +_V v) E v +_V v
  }
}
:EndNamespace

```

A important
 A avoid rounding issues
 A Assert by Roger Hui
 A magnitude, unitary, inverse
 A array from vector, vector from array, base
 A collapse equal values
 A scalar if possible and unscalar
 A remove trailing zeros and equivalent
 A extend and apply extending arguments
 A apply vector extension or zero function
 A multivector
 A binary unitary base, grade and dimensions
 A group in blades
 A position, sign and metric
 A factor from metric, tables from signature
 A exterior, inner, reversion and base
 A geometric product
 A inner and exterior product
 A geometric, inner and exterior product
 A antigeometric product
 A inverse
 A complement
 A geometric, exterior and inner product
 A antigeometric, exterior and inner product
 A left, right and double complement
 A inverse, antiinverse, reverse, pseudoscalar
 A GEOMETRIC ALGEBRA
 A namespace, signature, base
 A geometric, exterior and inner operator
 A products
 A products of vectors
 A products by vector
 A antiproducts
 A inverse and division
 A complements: left, right, double
 A reverse and extend
 A 0-3D, complex, quaternions, pgs
 A Vector Space
 A a and b are scalars
 A u, v and w are vectors (rank 1)
 A identity element for addition
 A identity element for multiplication
 A inverse element for addition
 A commutative addition
 A commutative product with scalars
 A associative addition
 A associative product with scalars
 A distributive over addition of vectors
 A distributive over addition of scalars
 A Geometric Algebra (geometric product aa)
 A vector space: a b scalars, u v w vectors
 A associative product
 A distributive by left
 A distributive by right
 A extension of scalar product
 A scalar product of parallel vectors
 A jgl@dyaalog.com 2022



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

```

GA = {
  g = [NSB ◊ g.s g.b - u ($ BV 2+d+1/a)
  g.I=I ◊ g._K_X ◊ g.I=I ◊ g.S+S ◊ g.US+US ◊ g.C+C
  g.V+V ◊ g._V-V ◊ g.D+D ◊ g.BV+BV ◊ g._d-d
  g._d = _d(TSu) ◊ g._dX = _d(TXu) ◊ g._dI = _dI(TIu)
  g.d = "g._d ◊ g.dX = "g._dX ◊ g.dI = "g._dI
  g.dV = g.dBV ◊ g.dXV = g.dXBV ◊ g.dIV = g.dIBV
  g.dV = g.dV ◊ g.dXv = g.dXv ◊ g.dIv = g.dIv
  g.d = g.d g._d ◊ g.dX = g.dX g._d ◊ g.dI = g.dI g._d
  g.dA = g.d I ◊ g.dAV = g.dABV ◊ g.dAV = g.dAV
  g.LC = g.d _C ◊ g.RC = g.d_C ◊ g.DC = g.LC^2
  g.R = *(TRu)◊ ◊ g.X = X(a*b ◊ a(aa)u) ◊ g
}

```

```

R GEOMETRIC ALGEBRA
R namespace, signature, base

R geometric, exterior and inner operator
R products
R products of vectors
R products by vector
R antiproducts
R inverse and division
R complements: left, right, double
R reverse and extend

```



- Geometric algebra
- Grain growth
- Crystal plasticity

ω)
 $.US \leftarrow US \quad \diamond \quad g.C \leftarrow C$
 $. \underline{\Delta} \leftarrow \underline{\Delta} \quad d$
 $\diamond \quad g. \underline{\Delta} I \leftarrow \underline{\Delta} I (T I \omega)$
 $\diamond \quad g. \Delta I \leftarrow \times g. \underline{\Delta} I$
 $\diamond \quad g. \Delta I V \leftarrow g. \Delta I \circ V$
 $\diamond \quad g. \Delta I v \leftarrow g. \Delta I \circ V$
 $\diamond \quad g. \underline{\Delta} I \leftarrow g. \Delta I \quad g. \underline{\Delta}$
 $\diamond \quad g. \Delta \Delta v \leftarrow g. \Delta \Delta \circ V$
 $\diamond \quad g. DC \leftarrow g. LC * 2$
 $\alpha(\alpha\alpha)\omega \} \quad \diamond \quad g$

A GEOMETRIC ALGEBRA

A namespace, signature, base

A geometric, exterior and inner operator products

A products of vectors

A products by vector

A antiproducts

A inverse and division

A complements: left, right, double

A reverse and extend



- Geometric algebra
- Grain growth
- Crystal plasticity



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- Recent changes
- Random page
- Help about MediaWiki
- Tools
- What links here
- Related changes
- Special pages
- Printable version
- Permanent link
- Page information

ω

$\cdot US \leftarrow US \quad \diamond \quad g \cdot C \leftarrow C$

$\cdot \underline{\Delta} \leftarrow \underline{\Delta} \quad d$

$\diamond \quad g \cdot \underline{\Delta} I \leftarrow \underline{\Delta} I$

$\diamond \quad g \cdot \Delta I \leftarrow \times g \cdot$

$\diamond \quad g \cdot \Delta IV \leftarrow g \cdot \Delta$

$\diamond \quad g \cdot \Delta I v \leftarrow g \cdot \Delta$

$\diamond \quad g \cdot \underline{\Delta} I \leftarrow g \cdot \Delta$

$\diamond \quad g \cdot \Delta \Delta v \leftarrow g \cdot \Delta$

$\diamond \quad g \cdot DC \leftarrow g \cdot l$

$\alpha(\alpha\alpha)\omega \} \quad \diamond \quad g$

Main Page

Rigid Geometric Algebra

This wiki is a repository of information about Rigid Geometric Algebra (RGA), and specifically the four-dimensional Clifford algebra $\mathcal{G}_{3,0,1}$. This wiki is associated with the following websites:

- [Projective Geometric Algebra overview site](#)
- [Conformal Geometric Algebra companion site](#)

Rigid geometric algebra is a mathematical model that naturally incorporates representations for Euclidean points, lines, and planes in 3D space as well as operations for performing rotations, reflections, and translations in a single algebraic structure. It completely subsumes conventional models that include homogeneous coordinates, Plücker coordinates, quaternions, and screw theory (which makes use of dual quaternions). This makes rigid geometric algebra a natural fit for areas of computer science that routinely use these mathematical concepts, especially computer graphics and robotics. [Conformal Geometric Algebra](#) (CGA) is a larger algebra that contains the complete RGA and also includes round objects like circles and spheres.

Rigid geometric algebra is an area of active research, and new information is frequently being added to this wiki.

If you are experiencing problems with the LaTeX on this site, please clear the cookies for rigidgeometricalgebra.org and reload.

Introduction

In the four-dimensional rigid geometric algebra, there are 16 graded basis elements. These are listed in Table 1.

There is a single *scalar* basis element that we denote by **1**, in bold, and its multiples correspond to the real numbers, which are values that have no dimensions.

There are four *vector* basis elements named $e_1, e_2, e_3,$ and e_4 that have one-dimensional extents. A general vector $\mathbf{v} = (v_x, v_y, v_z, v_w)$ has the form

Type	Values	Grade / Antigrade	
Scalar	1	0 / 4	□□□□
Vectors	e_1 e_2 e_3 e_4	1 / 3	■□□□ □■□□ □□■□ □□□■
Bivectors	$e_{23} = e_2 \wedge e_3$ $e_{31} = e_3 \wedge e_1$ $e_{12} = e_1 \wedge e_2$ $e_{43} = e_4 \wedge e_3$ $e_{42} = e_4 \wedge e_2$ $e_{41} = e_4 \wedge e_1$	2 / 2	■□■□ ■□□■ □■□■ □■□■ □■□■ □■□■
Trivectors / Antivectors	$e_{321} = e_3 \wedge e_2 \wedge e_1$ $e_{412} = e_4 \wedge e_1 \wedge e_2$ $e_{431} = e_4 \wedge e_3 \wedge e_1$ $e_{423} = e_4 \wedge e_2 \wedge e_3$	3 / 1	■□■□ ■□■□ ■□■□ ■□■□
Antiscalar	1 = $e_1 \wedge e_2 \wedge e_3 \wedge e_4$	4 / 0	■□■□

Table 1. The 16 basis elements of the 4D rigid geometric algebra.

<https://rigidgeometricalgebra.org/wiki/>

nature, base

erior and inner operator

ctors
ctor

vision
eft, right, double
tend



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

A geometric product operator

$$_GP_ \leftarrow \{t \leftarrow \alpha(\omega\omega\ddot{\omega})\omega \diamond +f\uparrow, \alpha(t\times\circ c''(\rho t)\uparrow\circ.(\alpha\alpha_Z))\omega\}$$

$$_GP_ \leftarrow _GP_ (BN\uparrow\ddot{\neq})$$

v1		v2		x_GP			
1	2	3	4	11	0	0	2



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

A geometric product operator

```

_GP_ ← {t ← α(ωω̃)ω ⋄ +f↑, α(t×c̃(ρt)↑∘.(αα_Z))ω}
_GP ← _GP_(BN[ö≠])

```

v1	v2	x_GP																
<table border="1"> <tr> <td>1</td> <td>2</td> <td>3</td> <td>4</td> </tr> </table>	1	2	3	4	<table border="1"> <tr> <td>5</td> <td>6</td> <td>7</td> <td>8</td> </tr> </table>	5	6	7	8	<table border="1"> <tr> <td>26</td> <td>44</td> <td>0</td> <td>0</td> <td>0</td> <td>0</td> <td>8</td> <td>8</td> </tr> </table>	26	44	0	0	0	0	8	8
1	2	3	4															
5	6	7	8															
26	44	0	0	0	0	8	8											



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

A geometric product operator

$$_GP_ \leftarrow \{t \leftarrow \alpha(\omega\omega\ddot{\sim})\omega \diamond +/ \uparrow, \alpha(t \times \circ c''(pt) \uparrow \circ .(\alpha\alpha_Z))\omega\}$$

$$_GP_ \leftarrow _GP_ (BN \uparrow \ddot{\neq})$$

v1	v2	x_GP	o . x_GP	+ . x_GP																																
<table border="1"> <tr> <td>1</td><td>2</td><td>3</td><td>4</td> </tr> </table>	1	2	3	4	<table border="1"> <tr> <td>5</td><td>6</td><td>7</td><td>8</td> </tr> </table>	5	6	7	8	<table border="1"> <tr> <td>26</td><td>44</td><td>0</td><td>0</td><td>0</td><td>0</td><td>8</td><td>8</td> </tr> </table>	26	44	0	0	0	0	8	8	<table border="1"> <tr> <td>26</td><td>30</td><td>0</td><td>0</td><td>0</td><td>0</td><td>8</td><td>10</td> </tr> <tr> <td>38</td><td>44</td><td>0</td><td>0</td><td>0</td><td>0</td><td>6</td><td>8</td> </tr> </table>	26	30	0	0	0	0	8	10	38	44	0	0	0	0	6	8	70 0 0 16
1	2	3	4																																	
5	6	7	8																																	
26	44	0	0	0	0	8	8																													
26	30	0	0	0	0	8	10																													
38	44	0	0	0	0	6	8																													



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

A geometric product operator

$$_GP_ \leftarrow \{t \leftarrow \alpha(\omega\omega\ddot{\omega})\omega \diamond +f\uparrow, \alpha(t\times\circ c''(\rho t)\uparrow\circ.(\alpha\alpha_Z))\omega\}$$

$$_GP_ \leftarrow _GP_ (BN\lceil\ddot{\neq})$$

z1	z2	x_GP	o.x_GP	+ .x_GP																				
<table border="1"> <tr><td>1</td><td>2</td><td>3</td><td>4</td></tr> </table>	1	2	3	4	<table border="1"> <tr><td>5</td><td>6</td><td>7</td><td>8</td></tr> </table>	5	6	7	8	<table border="1"> <tr><td>-16</td><td>-20</td><td>22</td><td>40</td></tr> </table>	-16	-20	22	40	<table border="1"> <tr><td>-16</td><td>-18</td><td>22</td><td>26</td></tr> <tr><td>-18</td><td>-20</td><td>34</td><td>40</td></tr> </table>	-16	-18	22	26	-18	-20	34	40	-36 62
1	2	3	4																					
5	6	7	8																					
-16	-20	22	40																					
-16	-18	22	26																					
-18	-20	34	40																					



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

z1	z2	×	∘.×	+ .×
1J3 2J4	5J7 6J8	-16J22 -20J40	-16J22 -18J26 -18J34 -20J40	-36J62

z1	z2	×_GP	∘.×_GP	+ .×_GP																				
<table border="1"> <tr> <td>1</td><td>2</td><td>3</td><td>4</td> </tr> </table>	1	2	3	4	<table border="1"> <tr> <td>5</td><td>6</td><td>7</td><td>8</td> </tr> </table>	5	6	7	8	<table border="1"> <tr> <td>-16</td><td>-20</td><td>22</td><td>40</td> </tr> </table>	-16	-20	22	40	<table border="1"> <tr> <td>-16</td><td>-18</td><td>22</td><td>26</td> </tr> <tr> <td>-18</td><td>-20</td><td>34</td><td>40</td> </tr> </table>	-16	-18	22	26	-18	-20	34	40	-36 62
1	2	3	4																					
5	6	7	8																					
-16	-20	22	40																					
-16	-18	22	26																					
-18	-20	34	40																					



Research

- Geometric algebra
- Grain growth
- Crystal plasticity

GrainGrowth



- Geometric algebra
- Grain growth
- Crystal plasticity

GrainGrowth

```
:Namespace GrainGrowth
R←8.314 ⋄ Area←#.EBSD.Area
_B←{>+/G0×s×1-⊖@(0<)}''s←10180÷∞(t×GB<t)▷α,ω
B1←{α(∓1⊖α)ω(∓1ϕω)}_B
B2←{α(∓1⊖1ϕα)ω(∓1⊖1ϕω)}_B
Step←{
  gt←-R×α×⊖?∞ω ⋄ q1←1 ∓1(⊖'',ϕ'')cω ⋄ gb←ω
  v1←1 ∓1(⊖'',ϕ'')cν ⋄ t←(∓v/v1)^(∓(cω)ν.∞q)
  gb1←(1+RP×?∞q1)×{ω+SF×q1 B2''{(1ϕα)(∓1⊖
  (v1 gb1 q1)←{t/ö,ϕ†ϕ†ω}}''v1 gb1 q1 ⋄ m←∞
}
Next←{q Step∞←ω ⋄ a,←Area q ⋄ q}

FromEBSD←{
  α←10000 3000 5 0.1 0.1 1000 0.25 0.01 ⋄
  gg.g gg.q←gg.DA #.EBSD.Orientations ω ⋄
  gg.v←(ρgg.q)ρ(gg.IQ<#.EBSD.IQ ω)∧gg.CI<∞
}
ToEBSD←{ebsd←ω ⋄ ebsd[ι3]←†ϕ†g[,q] ⋄ ebsd}

Solve←{a←Next∞≡ω}
_Solve_←{
  α←1 ⋄ f←α ⋄ n←0 ⋄ ebsd←αα ⋄ name←ωω
  a←Next∞{n←+1 ⋄ (α≡ω)←(ToEBSD_ebsd)#.EBSD
}
:EndNamespace
```



Research

- Geometric algebra
- Grain growth
- Crystal plasticity



```

:Namespace Euler
  A quaternion product
  c ← t(0 1 2 3)(1 0 3 2)(2 3 0 1)(3 2 1 0)
  u ← t(1 -1 -1 -1)(1 1 1 -1)(1 -1 1 1)(1 1 -1 1)
  QP ← u{+/αα×[-2t+r+1]α×[(i(≠ρα)-1),r+IO]}(c[IO+ωω][(+≠ρω)-1])
  QC ← ×ö1°(1,-3p1)
  QD ← +.×ö1°QC

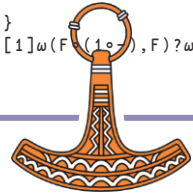
  A quaternion from Euler angles
  RD ← 180÷ω
  UV ← {ω×[i(≠ρω)-1]÷(+×ω)*÷2}
  QA ← {(2ω), (1ω)°×UVα}°(÷2) ◊ QAD ← QA°RD
  QE ← {α+(0 0 1)(1 0 0)(0 0 1) ◊ αQP.QA=[-1+i≠ρω]ω} ◊ QED ← Q

  A cubic symmetry
  cs ← c1 0 0 0
  cs,+, (1 0 0)(0 1 0)(0 0 1) °.QAD 90 18
  cs,+, (1 1 1)(-1 1 1)(1 -1 1)(1 1 -1) °.QAD 120 2
  cs,+, (1 1 0)(1 0 1)(0 1 1)(1 -1 0)(-1 0 1)(0 1 -1)°QAD 180

  A misorientations
  ML ← {α+0.5 ◊ [(180×ω)÷α]}
  MC ← cs{2×-2o1[>[≠|αα°QD=αQP°QCω]}
  IM ← {IO+1-ω(α|ω)+/1-ωi(α|ω)-1}
  CM ← {c+NSθ ◊ c.m+1p(-IM-1)≠, tω ◊ c}
  _M_ ← {α=ω:0 ◊ 0≤m+(i+αIMω)ωω.m:m ◊ t(iωωω.m)+MLα(MC)ö(÷αα)}

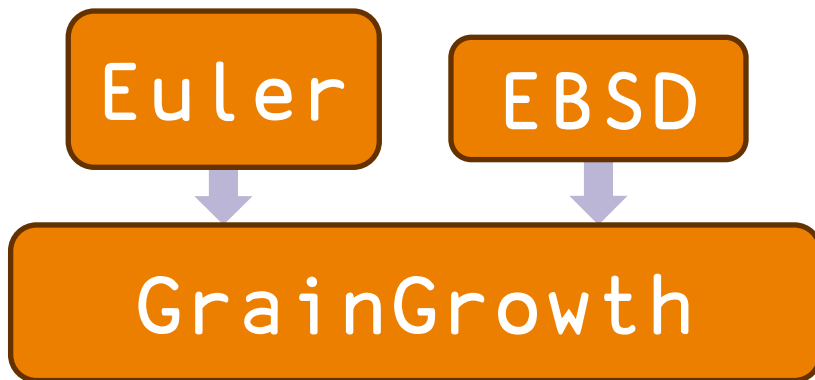
  A namespace with function M to calculate misorientations
  M ← {α+0.5
    m+NS'QD' 'QP' 'QC' 'IM' 'MC' ◊ m.ML+α°ML
    m.M+(α÷2)+α×(,tω)_M_(CMω) ◊ 2≥≠ρω: m
    m.M+m.Mö((-1↓ρω)°{IO+α.1ω-IO}) ◊ m
  }

  Space ← {t°.,/(ω÷2)+ω×(i°360 90 90÷ω)-1}
  Random ← {F+(1 2o2×c°?αρ0)×cω*÷2} ◊ UVt[1]ω(F(1°-),F)?ωρ0}
:EndNamespace
  
```



Research

- Geometric algebra
- Grain growth
- Crystal plasticity



:Namespace EBSD

```
Read←{ A read whitespace separated data f
      f←80 ~1 MAP ω A map
      lf cr←UCS 10 13 A line
      clean←,/c~(≠'#'^1,-1↓=o{f})~ A spli
      s←'\s+'R', 't'(^\\s+)|\\s+$'R' 't→clea
      CSV'Invert' 2t-s'S' 2 0
}
```

```
Write←({'#','α')CSV('Invert' 2)('Separato
```

```
Crop←{x y←α ◇ s←((x>5o)∧y>4o)ω ◇ s/'w}
```

```
Orientations←{
  x←5oω ◇ nx←[0.5+1+(f/x)÷f/|2-/x ◇ ny←[
  α←0 ◇ ea←↓φ↑3↑ω ◇ ea←α([0.5+~)×(α>0)↑
}
```

```
IQ←6o A image quality
```

```
CI←7o A confidence index
```

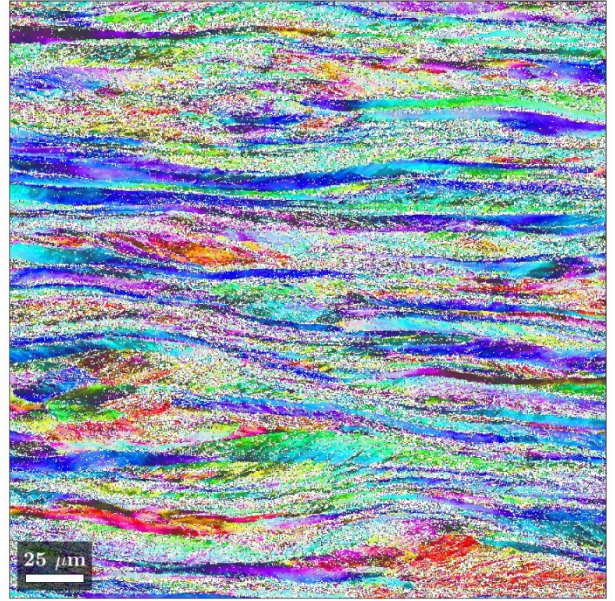
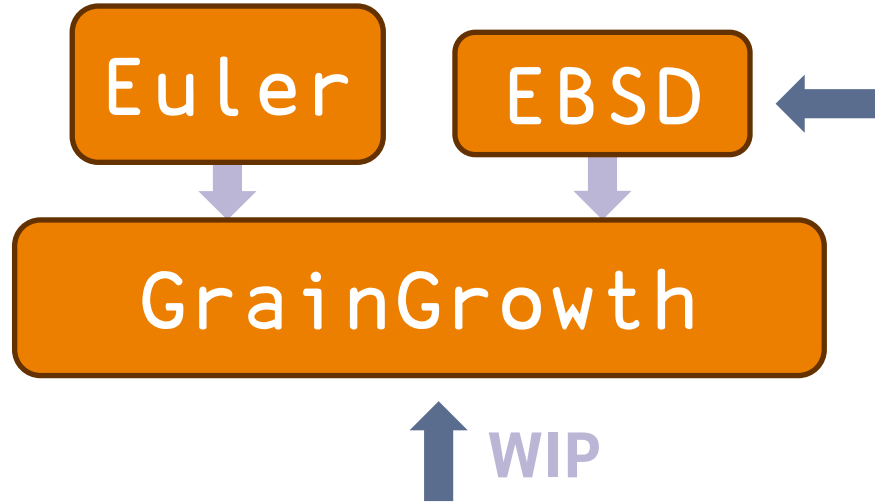
```
Area←{(i2)×.((≠,ω){αα÷1[+/,ω≠1φ[α]ω})cω}
```

:EndNamespace



Research

- Geometric algebra
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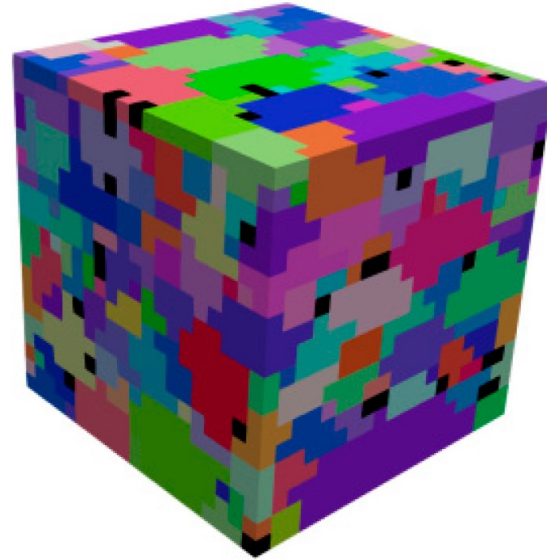
Sample preparation, microscopy: Estefanía Sepulveda
MTEX processing: JGL

$$r_{0i} = \frac{M_{0i}}{V_0^2} \sqrt{\sum_d \left(\sum_j \delta_{ij} A_{0j} |\bar{u}_j \cdot \bar{u}_d| \right)^2} \sum_j (\gamma_{0j} - (1 - \delta_{ij}) c_{0ij} \gamma_{ij}) A_{0j}$$



Research

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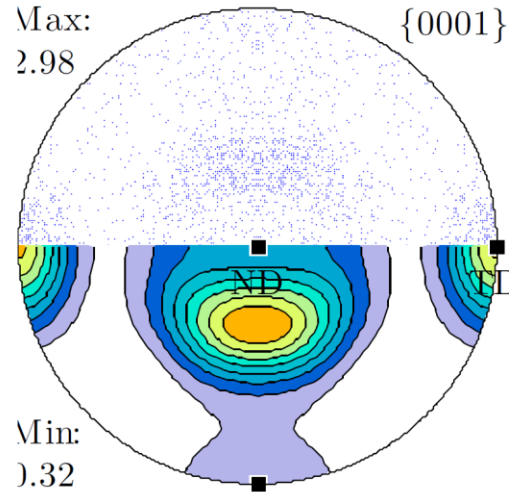
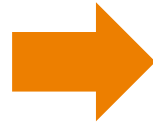
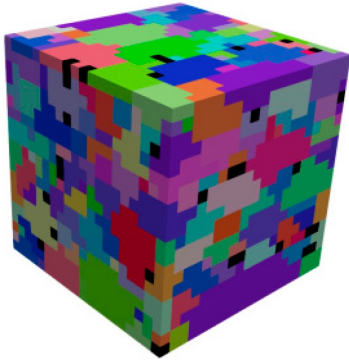


<https://doi.org/10.3390/cryst10090819>



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- Grain growth
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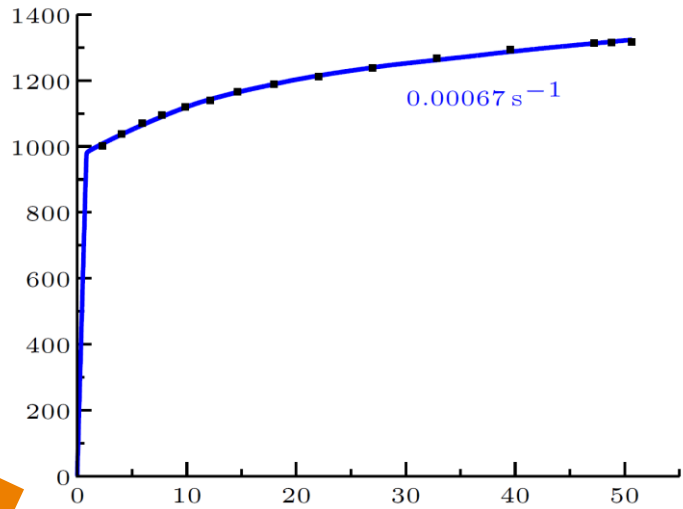
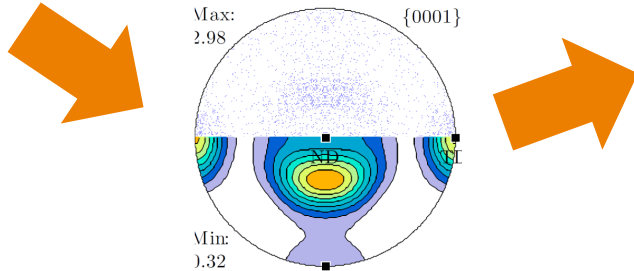
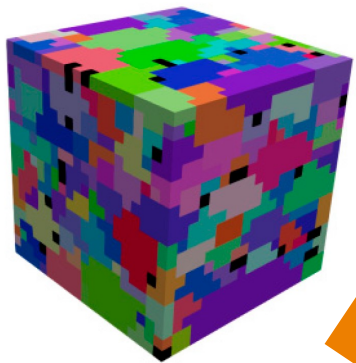


<http://hdl.handle.net/1854/LU-4388117>



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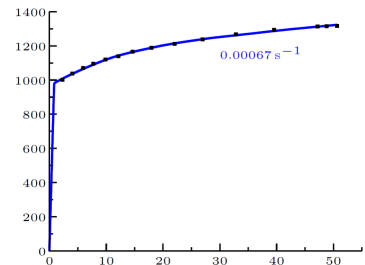
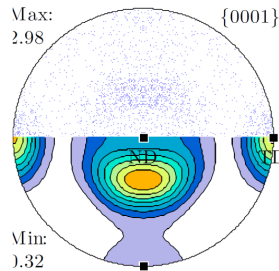
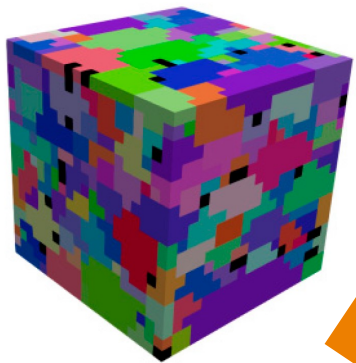


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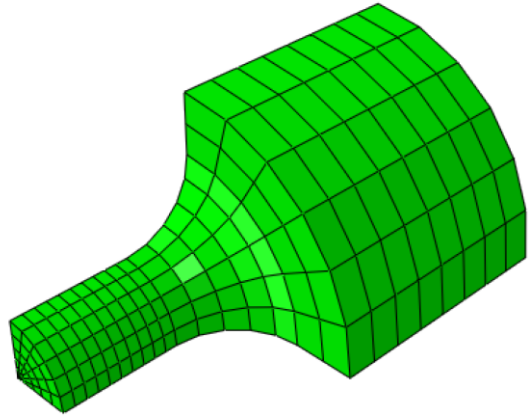


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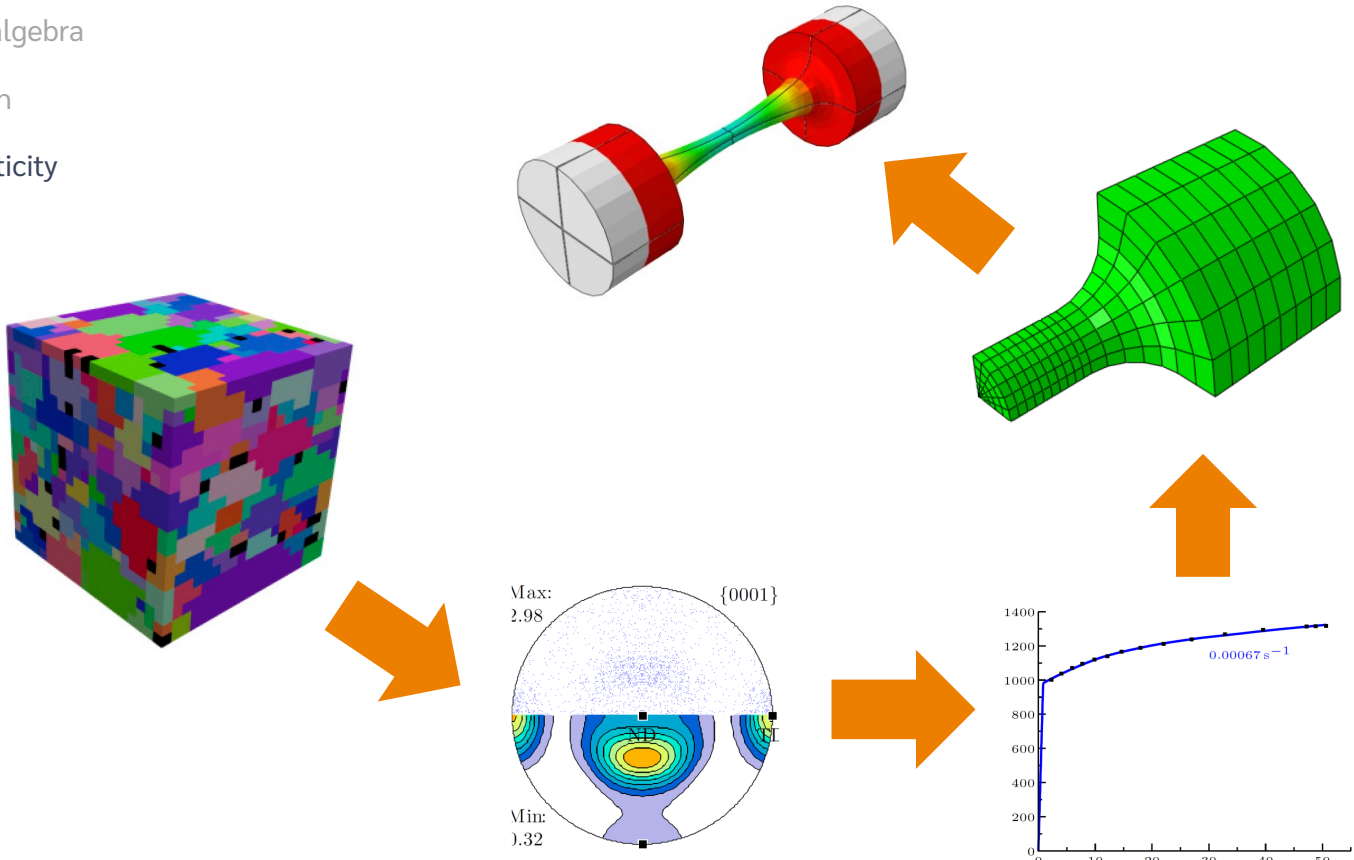


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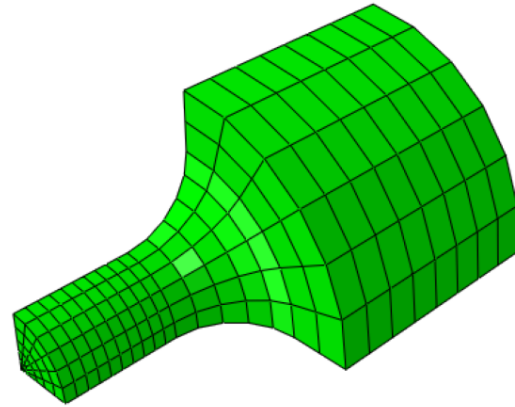
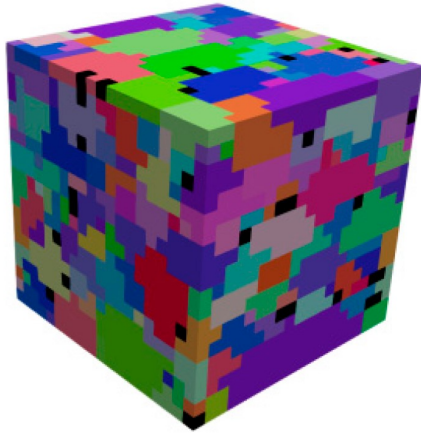


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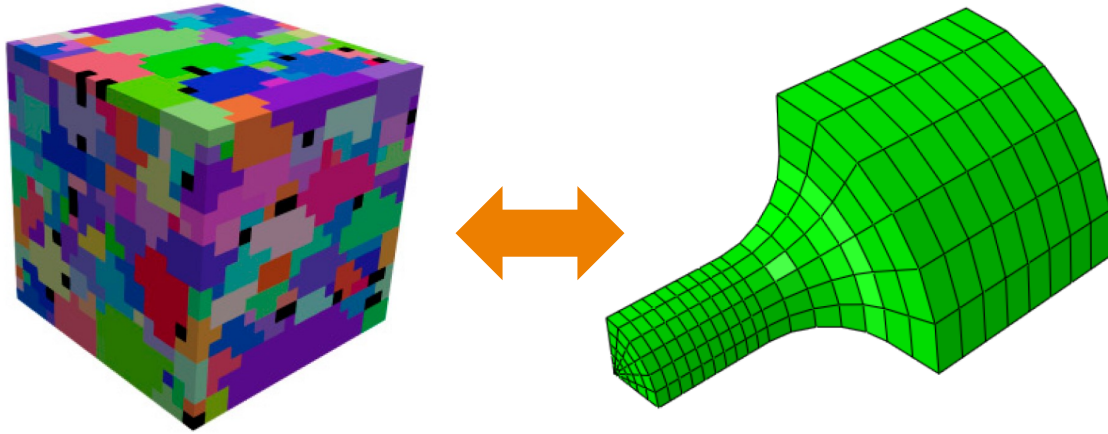
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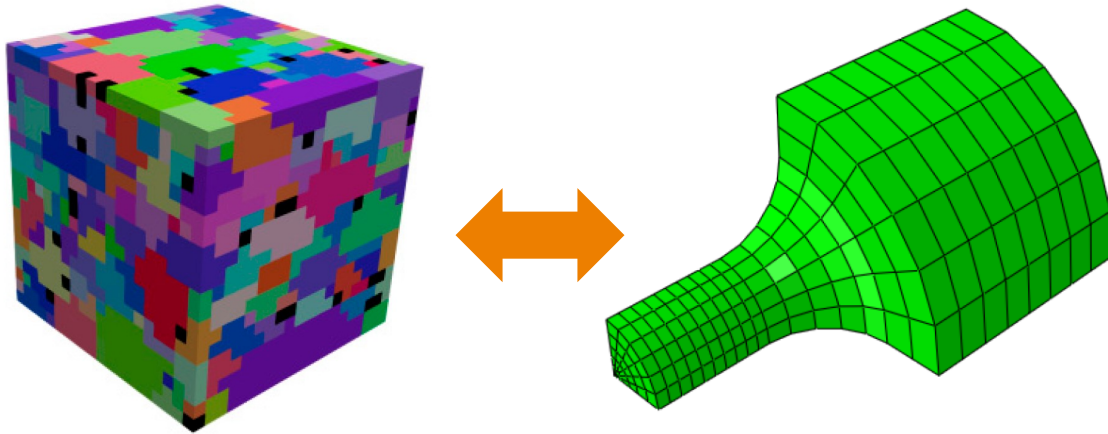
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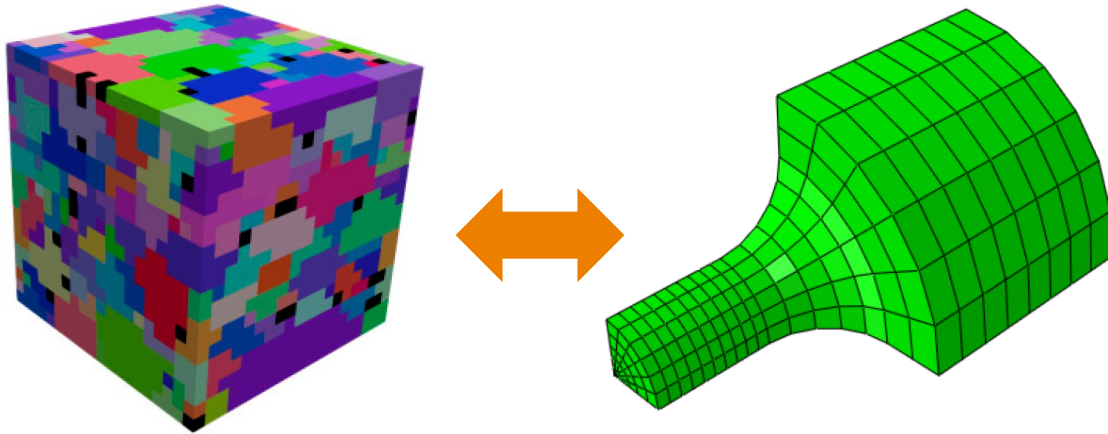
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CPFEH



- Geometric algebra
- Grain growth
- Crystal plasticity

C_{rystal} P_{lasticity} F_{inite} E_{lement} H_{omogenisation}



- Geometric algebra
- Grain growth
- Crystal plasticity

Crystal Plasticity Finite Element Homogenisation



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Superposition principle

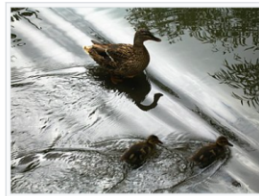
39 languages

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From Wikipedia, the free encyclopedia

This article is about the superposition principle in linear systems. For other uses, see Superposition (disambiguation).

The **superposition principle**,^[1] also known as **superposition property**, states that, for all **linear systems**, the net response caused by two or more stimuli is the sum of the responses that would have been caused by each stimulus individually. So that if input *A* produces response *X* and input *B* produces response *Y* then input (*A* + *B*) produces response (*X* + *Y*).



Superposition of almost plane waves (diagonal lines) from a distant source and waves from the wake of the ducks. Linearity holds only approximately in water and only for waves with small amplitudes relative to their wavelengths.

A function $F(x)$ that satisfies the superposition principle is called a **linear function**. Superposition can be defined by two simpler properties: **additivity**

$$F(x_1 + x_2) = F(x_1) + F(x_2)$$

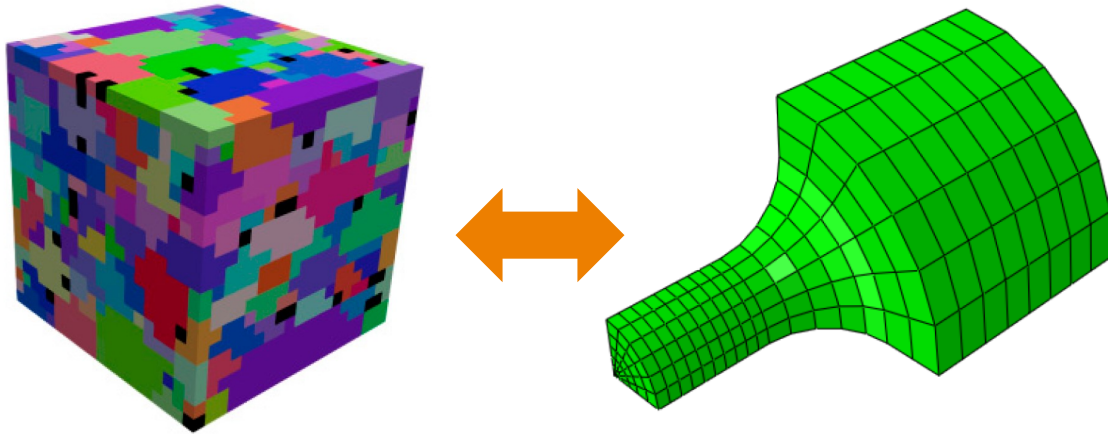
and **homogeneity**



Research

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Research

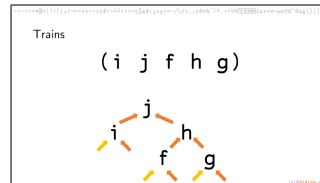
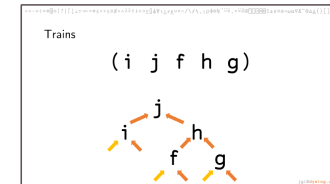
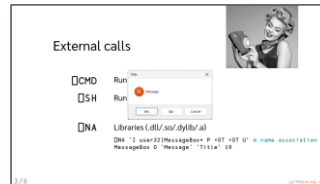
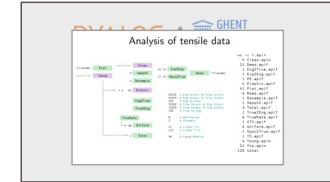
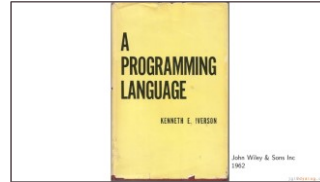
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Dissemination and promotion



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Conclusions

- ◆ APL is an interesting choice for the solution of materials science and engineering problems
- ◆ Convincing other researchers is not easy
- ◆ But there is potential
- ◆ We will keep trying



Plans

- ◆ Apply feedback
- ◆ More tutorials and publish them
- ◆ Continue research
 - ◆ Geometric algebra, grain growth, crystal plasticity
 - ◆ Articles and conferences
- ◆ Networking (more universities)



Thank you



Elsinore 2023

APL and Metallurgy

- TensileAnalysis ←
- Crystallography ←
- GrainGrowth ←
- CrystalPlasticity ←
- ...

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