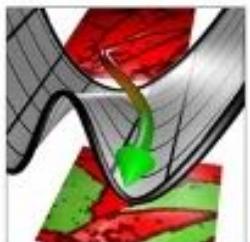


Overview presentation-Internship in MPIE

C. Zhu, D. Yan, C.C. Tasan, D. Raabe



Department Microstructure Physics and Alloy Design



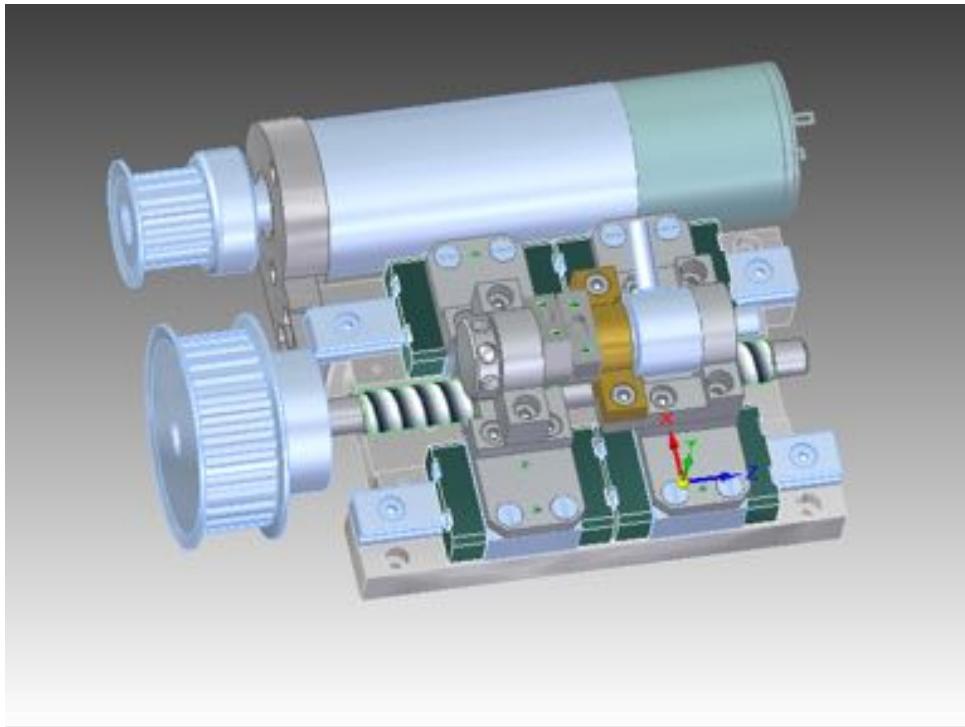
15-09-2014

Outline

- In-situ Tensile Stage (ITS)
 1. Displacemnet sensor
 2. Load sensor
 3. ITS setup
 4. Tensile Stage Control
 5. The Control Software
 6. Experiment I-BSE/SE movie
- DAMASK in (Experiment II)
 1. Introduction of DAMASK (CPFEM) / how it is used to solve the boundary value problems
 2. Four files used in DAMASK
 3. Process of DAMASK
- Results
 1. Experiment II results obtained through DIC
 2. Simulation Results obtained through DAMASK
 3. Comparison of experimental results and simulation results
- Previous Work

In-situ Tensile Stage-'ITS'

Solid Edge ST6 animation (1000X)



Specifications:

Teeth Ratio: 15:30

Mass: 500g

| load | \leq 500N

Max Specimen Length \leq 20mm

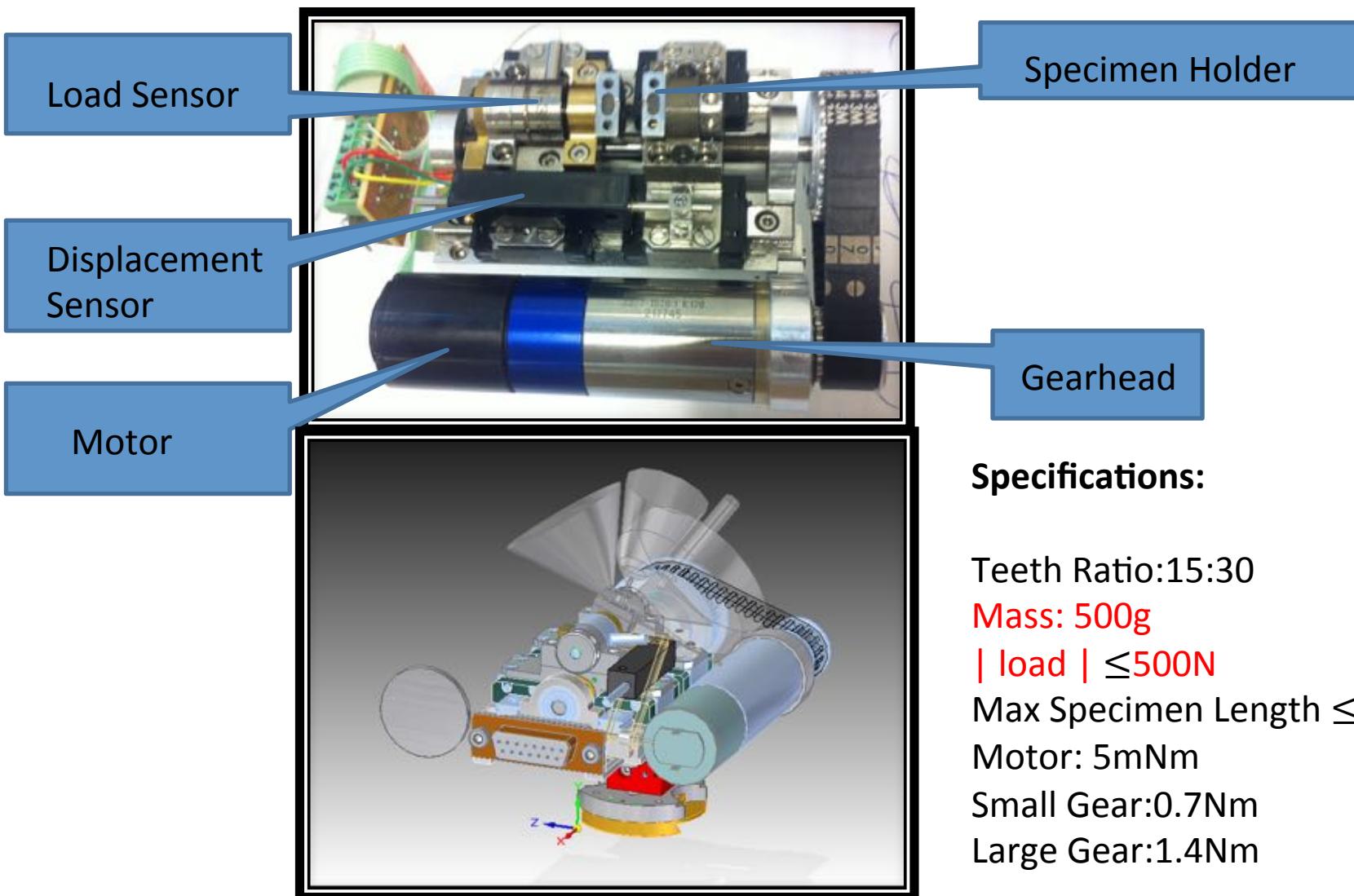
Motor: 5mNm

Small Gear: 0.7Nm

Large Gear: 1.4Nm

- 2 revolution of small gear
- =1 revolution of large gear
- =0.5mm linear movement of carriages
- =1mm elongation of specimen

In-situ Tensile Stage-'ITS'



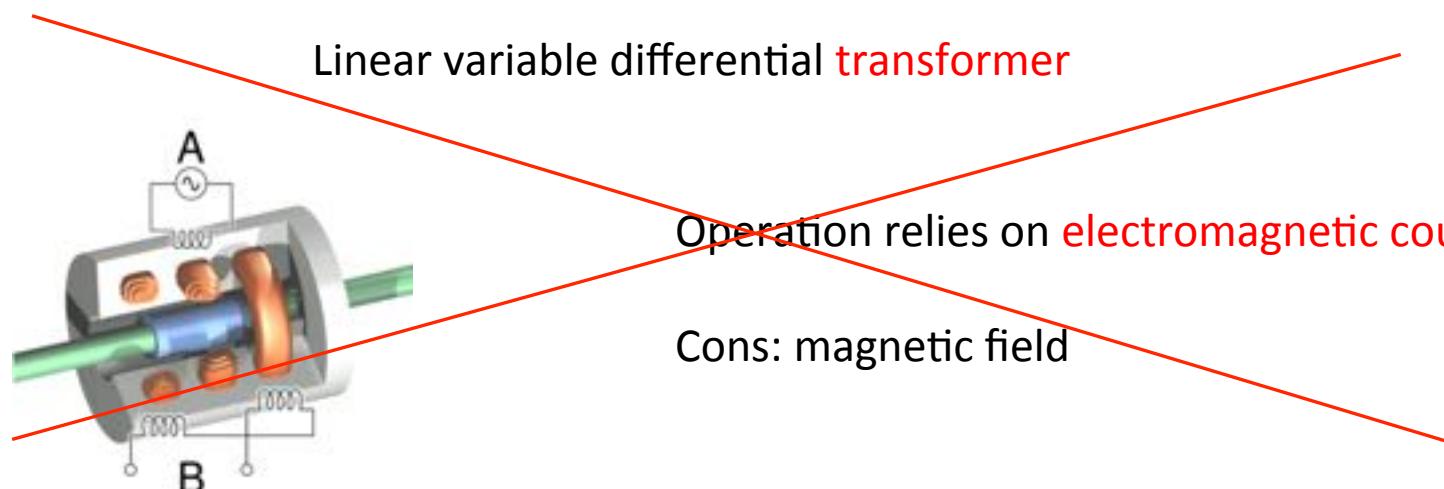
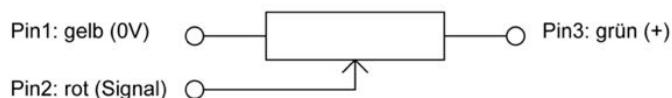
Displacement Sensor

Weg Sensoren: Serie MM10 - Potentiometric Linear Transducer

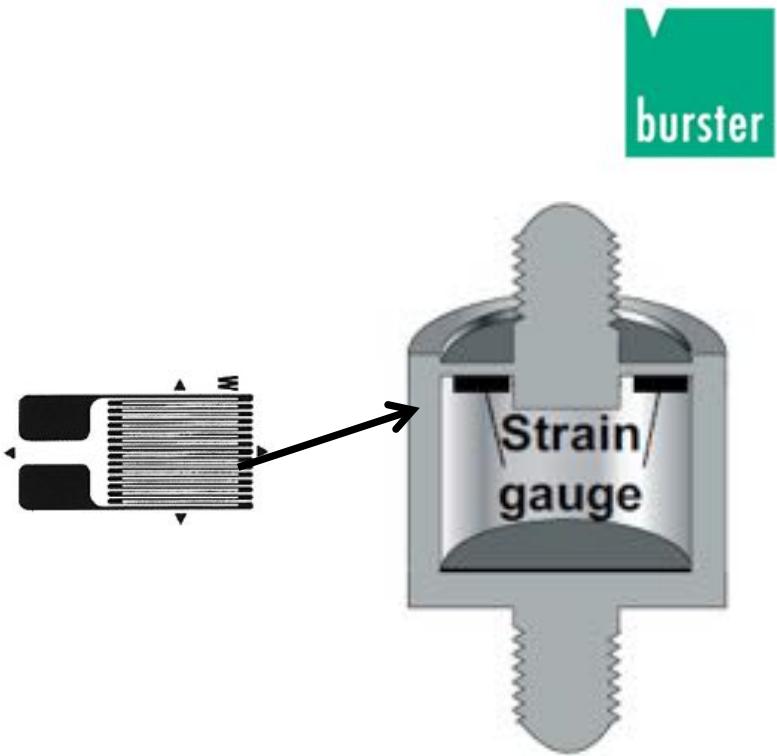


Electrical contact travels from 8 mm to 15 mm
Conductive plastic element

Cons: friction



Load Sensor: Model 8417



Force

- > deflection of elastic membrane
- > change of sensor's overall height
- > tension or compression in the spring element
- > resistance change in strain gauge

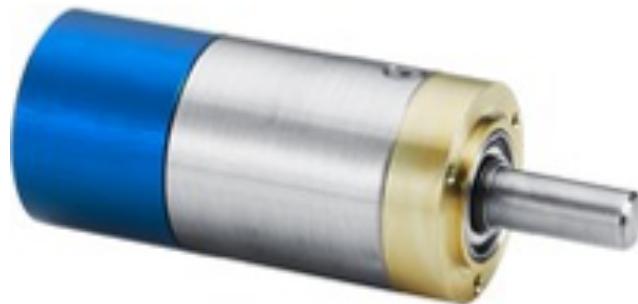


Motor+Gearhead



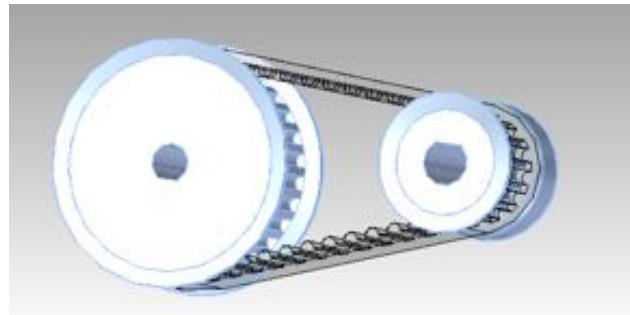
FAULHABER

DC-Micromotors (5mNm, steel, black coated) + Planetary Gearhead (0.7Nm, Metal housing)



Diameter (mm)	Length (mm)
22	24

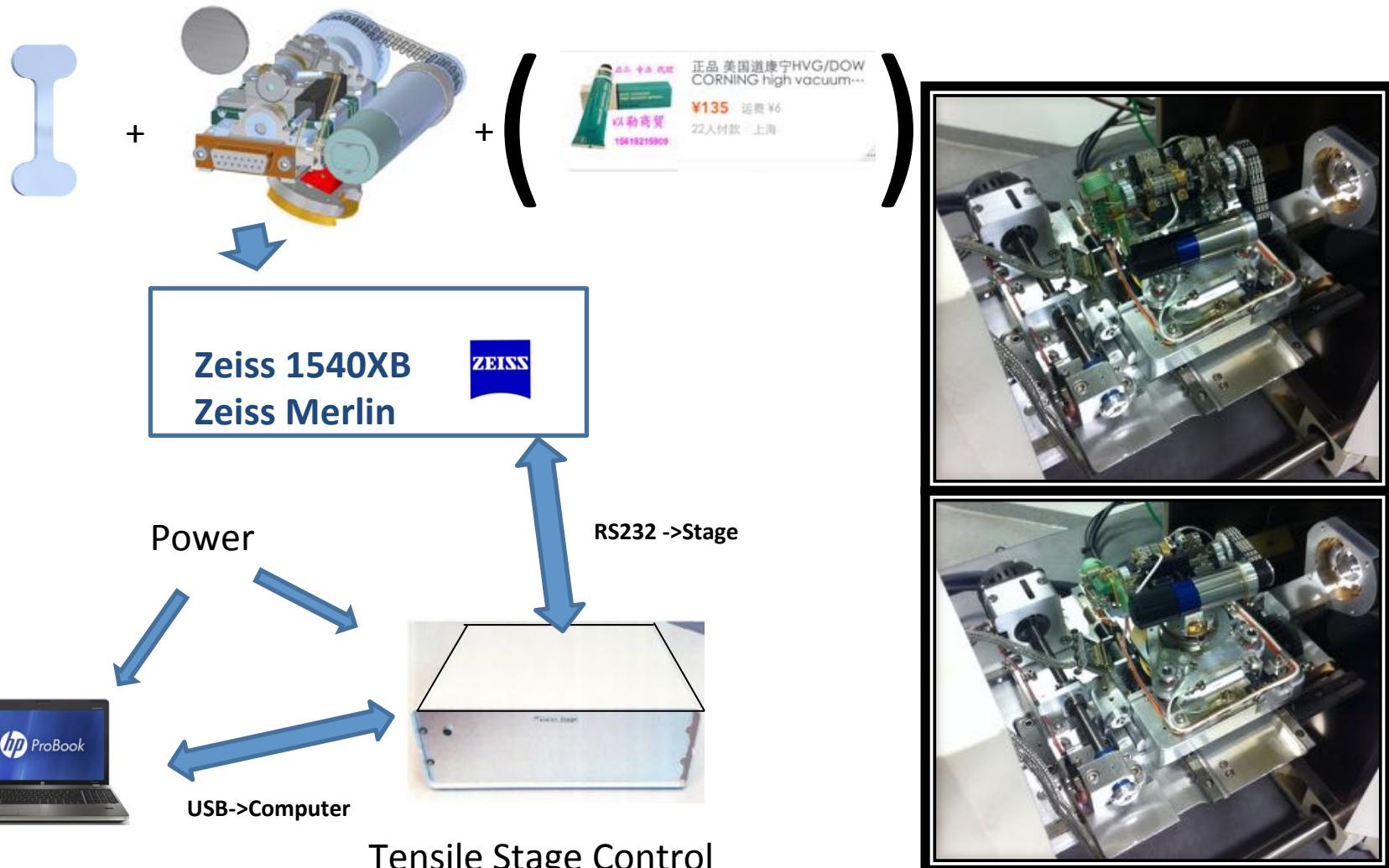
For high torque



1.4Nm

0.7Nm

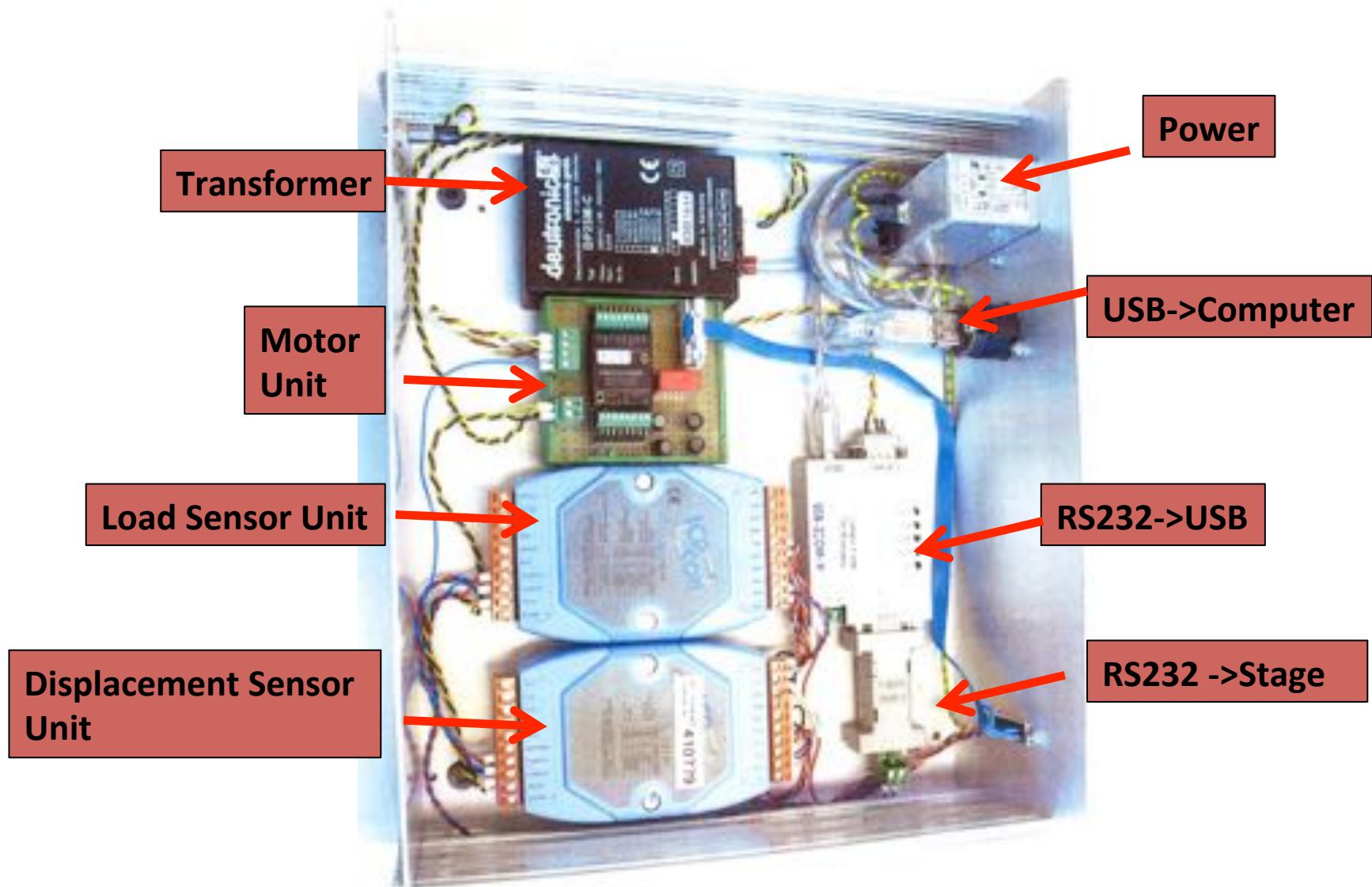
ITS Setup



E
B
S
D

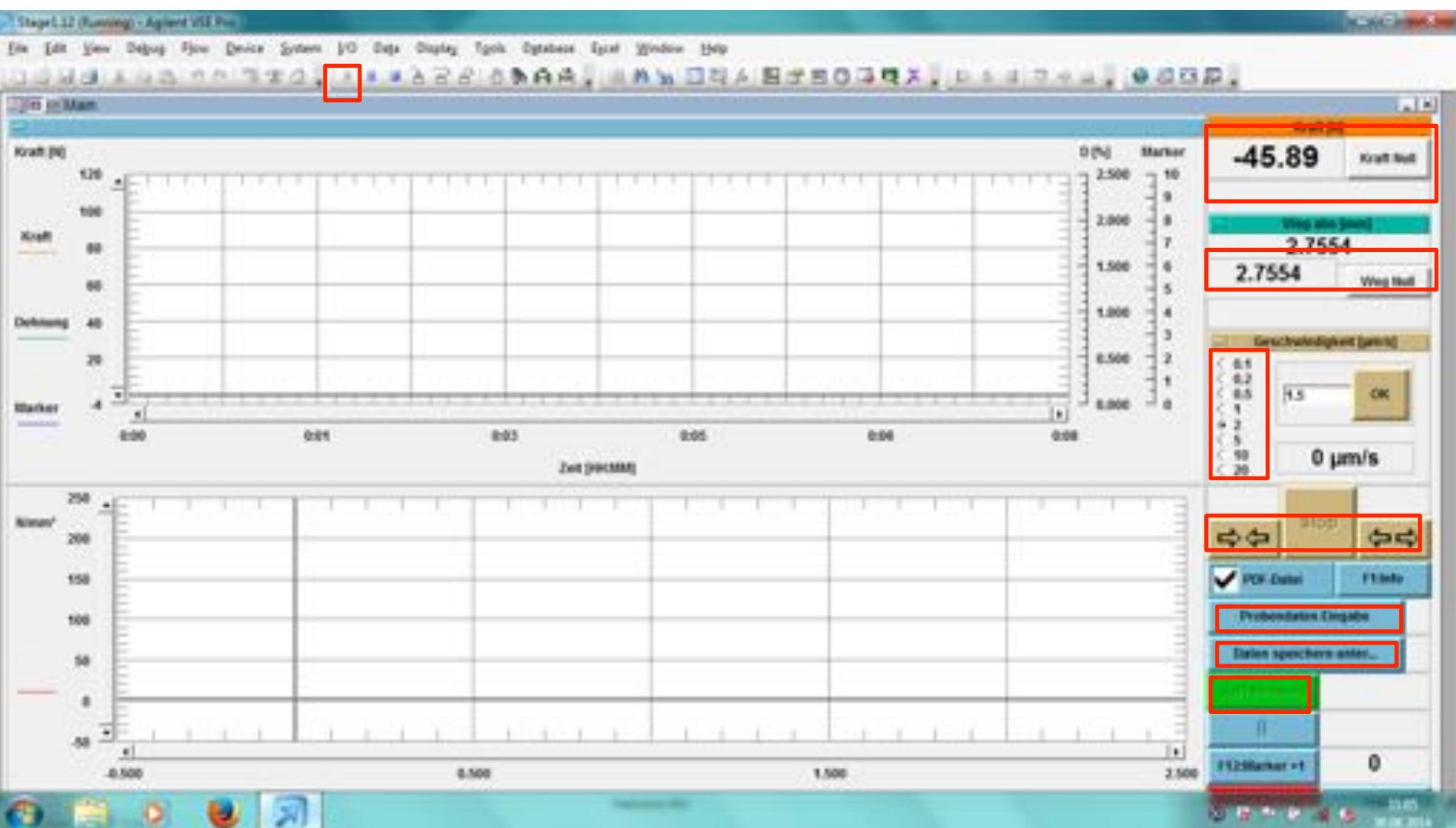
S
E

Tensile Stage Control

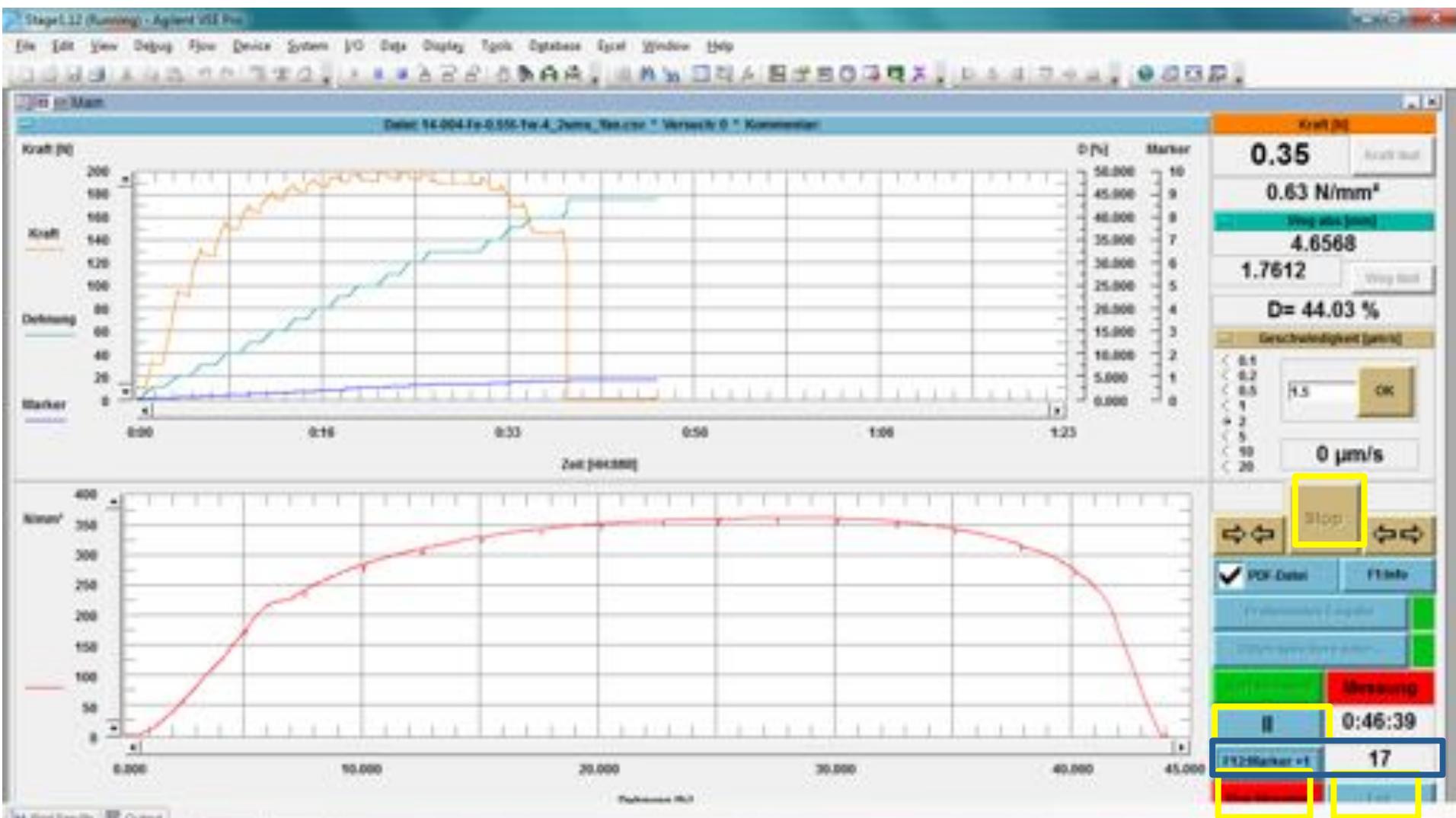




The Control Software



The Control Software





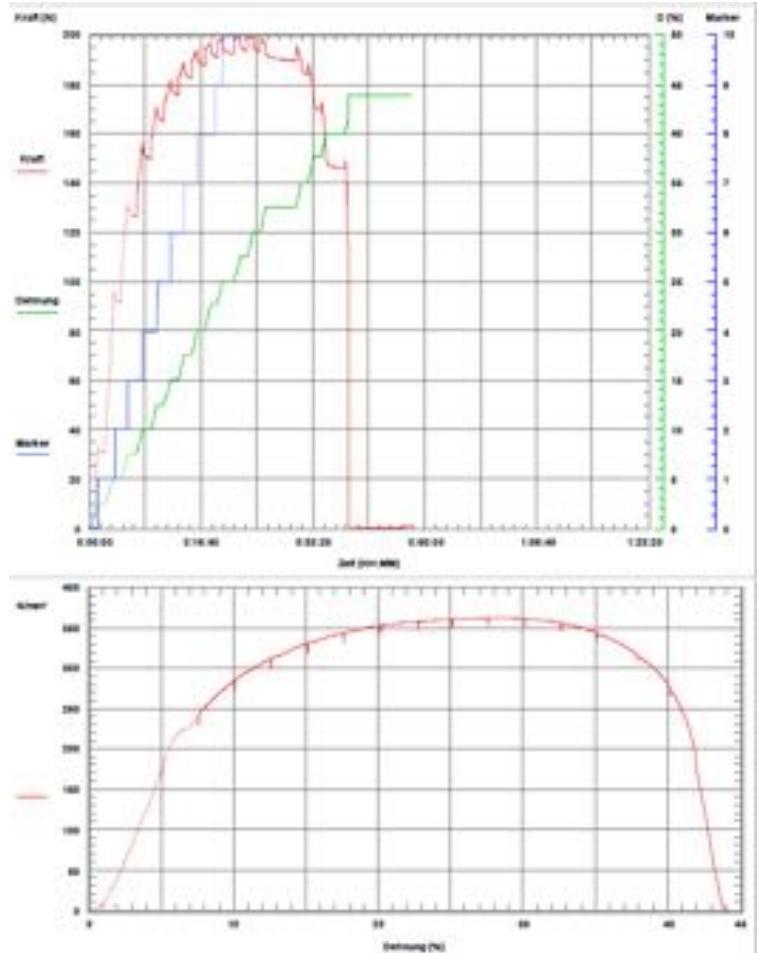
The Control Software



14-005-Fe-0.55t-1w-4_20ums_Yan.pdf
Adobe Acrobat Document



14-005-Fe-0.55t-1w-4_20ums_Yan.csv
Microsoft Excel Comma Separate...



A screenshot of Microsoft Excel showing a spreadsheet with experimental data. The columns are labeled A through G. Row 1 contains the header "Versuchsnummer". Rows 2 through 7 contain experimental parameters: Datum (12.12.2012), Zeit (21:48), Probenbr. (1), Probendic (0,55), akt. Probe (4), and Kommentar (empty). Rows 9 through 26 contain the measured data, with columns for Messzeit (Marker), Kraft (N), Weg (mm), Spannung (MPa), and Dehnung (%). The data shows a cyclic loading and unloading process.

A	B	C	D	E	F	G
1	Versuchsnummer	0				
2	Datum	12.12.2012				
3	Zeit	21:48				
4	Probenbr.	1				
5	Probendic	0,55				
6	akt. Probe	4				
7	Kommentar					
9	Messzeit (Marker)	Kraft (N)	Weg (mm)	Spannung (MPa)	Dehnung (%)	
10	0,0	0 -0,39	-0,0003	-0,72	0,000	
11	0,4	0 -0,35	0,0000	-0,63	0,000	
12	0,8	0 -0,30	0,0000	-0,54	0,000	
13	1,2	0 -0,30	0,0000	-0,54	0,000	
14	1,6	0 -0,30	0,0000	-0,54	0,000	
15	2,0	0 -0,35	0,0000	-0,63	0,000	
16	2,4	0 -0,39	0,0000	-0,72	0,000	
17	2,8	0 -0,44	0,0000	-0,81	0,000	
18	3,2	0 -0,44	0,0000	-0,81	0,000	
19	3,6	0 -0,39	0,0000	-0,72	0,000	
20	4,0	0 -0,35	0,0000	-0,63	0,000	
21	4,4	0 -0,30	0,0000	-0,54	0,000	
22	4,8	0 -0,35	0,0000	-0,63	0,000	
23	5,2	0 -0,35	0,0000	-0,63	0,000	
24	5,6	0 -0,10	0,0000	-0,18	0,000	
25	6,0	0 1,43	0,0003	2,60	0,008	
26	6,4	0 3,90	0,0007	7,09	0,017	

Experiment I-BSE movie

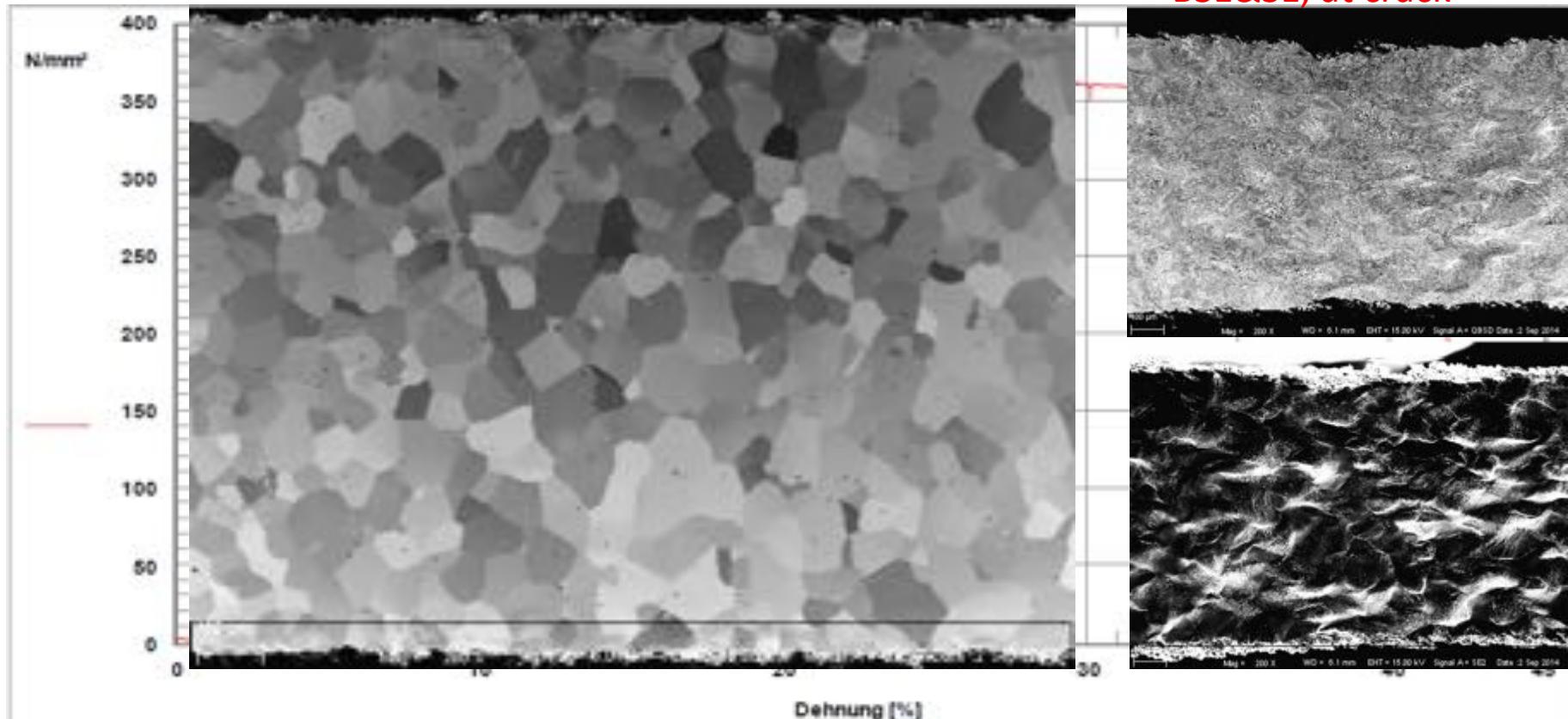
RD Tensile

$\varepsilon \downarrow global = 0\%-32.5\%$, BSE

$\varepsilon \downarrow global$

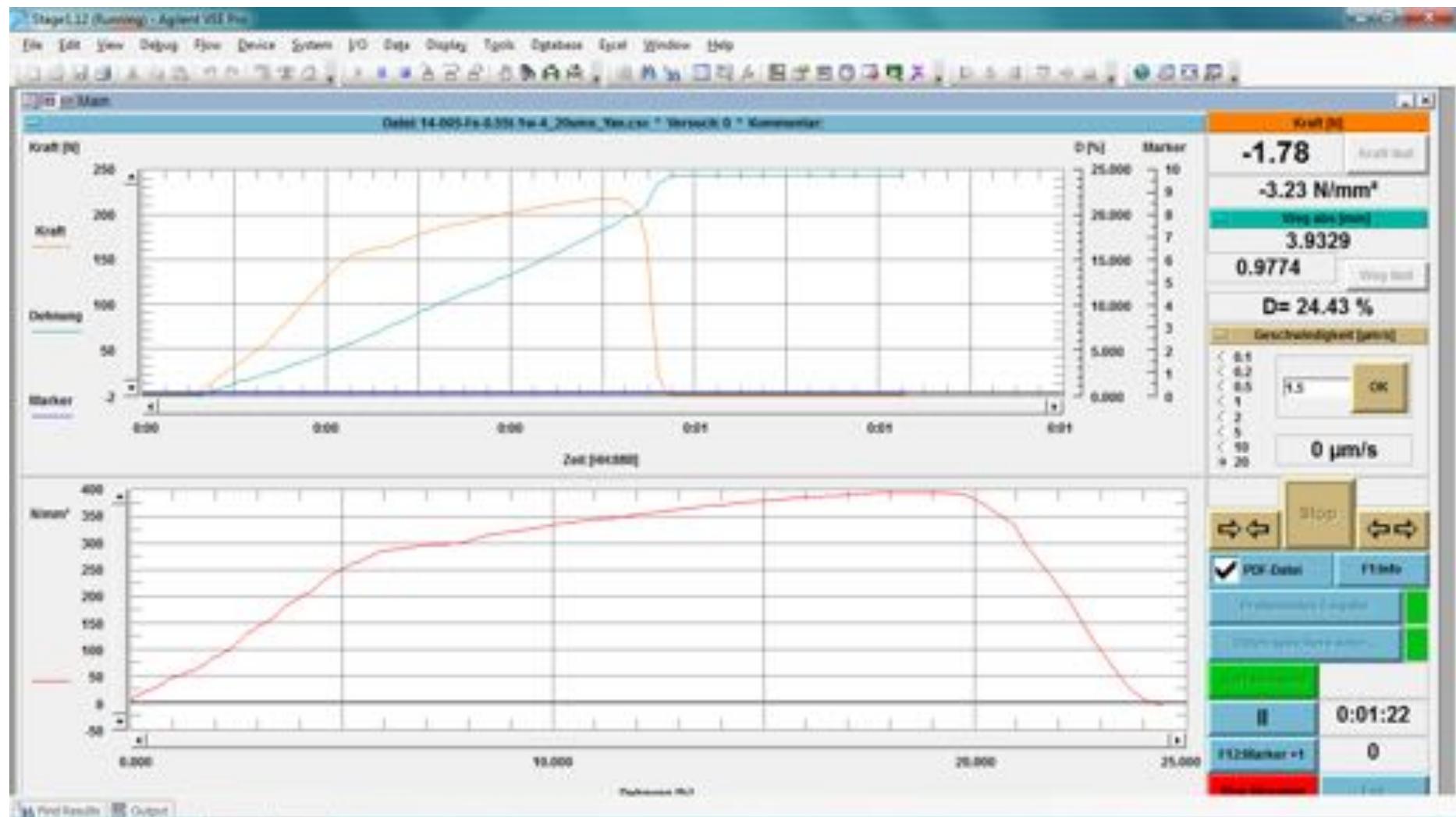
$= 32.5\%-44.3\%$,

BSE&SE, at crack



Versuchsnummer:	0
Datum:	02/Sep/2014
Uhrzeit:	15:50
Probenbreite [mm]:	1
Probendicke [mm]:	0,55
akt. Probenlänge [mm]:	4
Kommentar:	

Experiment II- $\varepsilon_{global} \downarrow$ global =0% to 19%





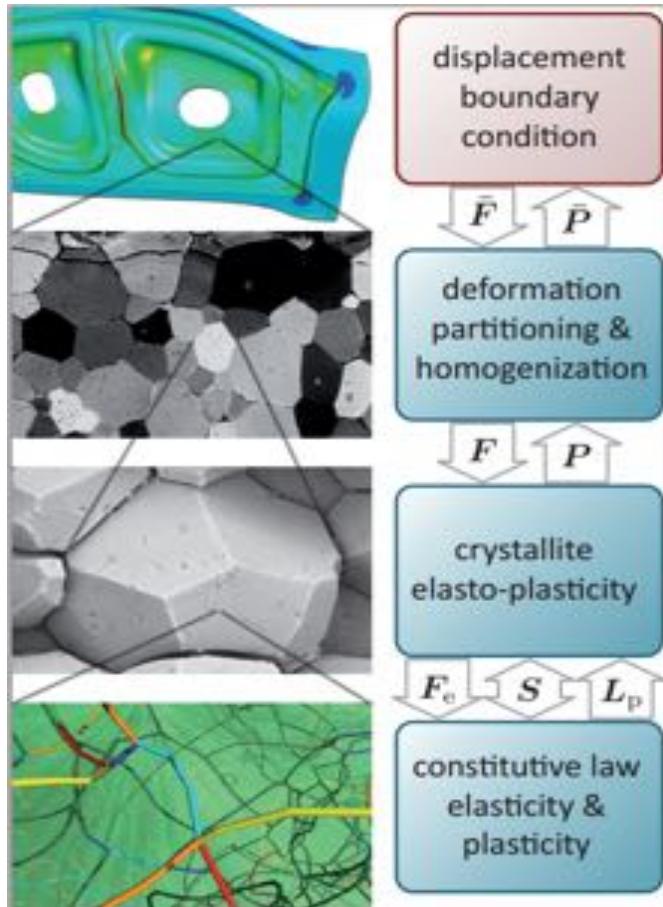
The screenshot shows the DAMASK software interface. At the top, there is a navigation bar with tabs like "Home", "About", "Contact", and "Help". Below the navigation bar, there is a large blue header area with some graphical elements. The main content area contains several sections of text and tables. One section is titled "Details of the simulation showing flow conditions". This section includes a table with columns "Flow direction", "Flow velocity", and "Flow angle". Another section is titled "Material properties". This section includes a table with columns "Material", "Type", "Value", and "Unit". There are also other sections and tables below, though they are less clearly legible due to the blurriness of the screenshot.

DAMASK simulation requirements (Franz Roters):

- arbitrary mechanical boundary value problems
- continuum mechanics
- accounting for crystal plasticity



Crystal Plasticity
Finite Element Method
(CPFEM)



DAMASK/boundary value problems solved by spectral solver

Solving with iteration through a convolution in Fourier space

Loop is finished when a given tolerance is reached (error stress BC, error $F \downarrow BC$, error $P \downarrow BC$)

Phenomenological law: critical resolved shear stress

Physics law: dislocation density

tensionX-variant3.load defines the loading conditions with deformation gradient rate $F \downarrow BC$ specified.

tensionX-variant3.load :

```
Fdot 5e-3 0 0 0 * 0 0 0 * stress *** * 0 * * * 0 time 50 incs 1000 freq 20
```

Strain tensor:

X: $F \downarrow BC$ (strain rate)

Y: arbitrary

Z: arbitrary

Stress Tensor:

Unknown

numerics.config defines the maximum and minimum iteration steps for the spectral solver.

numerics.config:

```
itmax 60
itmin 4
myspectralsolver basicPETSc
#myfilter cosine
```

material.config

contains information about the material separated by five `<part>` (`<output>`, `<phase>`, `<homogenization>`, `,` `<microstructure>`, `<texture>`) and phenomenological law used as the constitutive model for ferrite phase.

material.config:

```
#-----#
<crystallite>
#-----#
Output
#-----#
<phase>
#-----#
plasticity :phenopowerlaw/elasticity: hooke/slip systems/twin
#---
<homogenization>
#---
Ngrains 1
#---
<microstructure>
#---
[Grain00000001]
crystallite 1
(constituent) phase 1 texture 00000001 fraction 1.0
1 to 61401

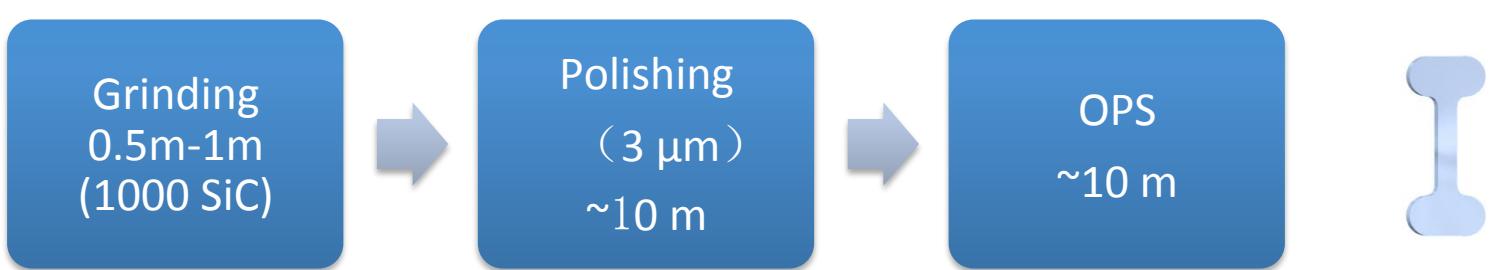
#---
<texture>
#---
[Grain00000001](gauss) phi1 115.7272 Phi 154.8848 phi2 75.8287 scatter 0.0 fraction 1.0
1 to 61401
```

fe.geom maps the microstructure information of the model to the location in the volume element (VE). It also contains the resolution, dimension, origin coordinate, applied homogenization scheme and number of grains in the model.

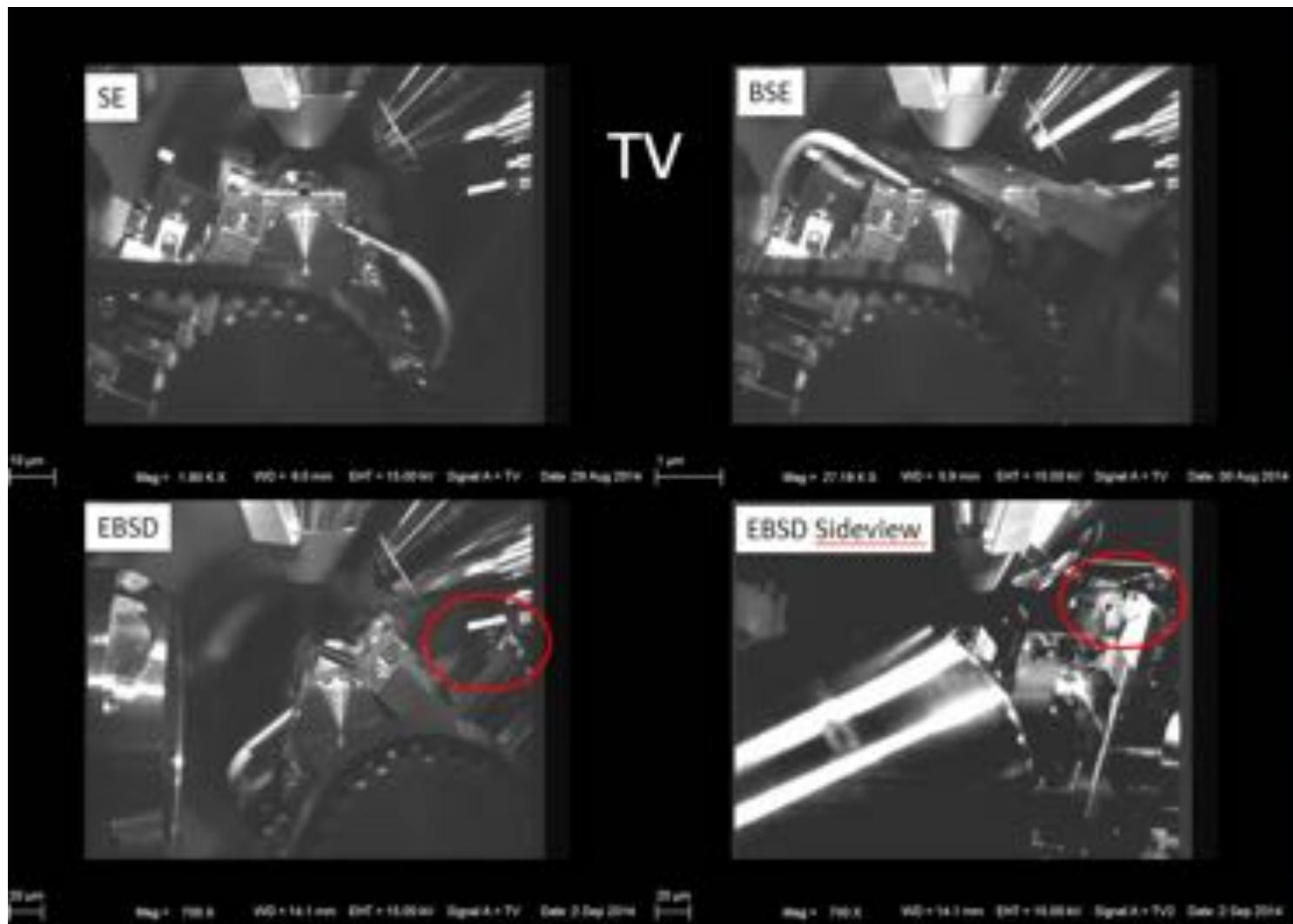
fe.geom:

```
5 Header
resolution          a 291      b 211      c1
dimension          x 291      y 211      z1
origin            x 0        y 0        z0
homogenization 1
maxGrainCount 0
1 to 61401
```

Pure Iron Sample Preparation

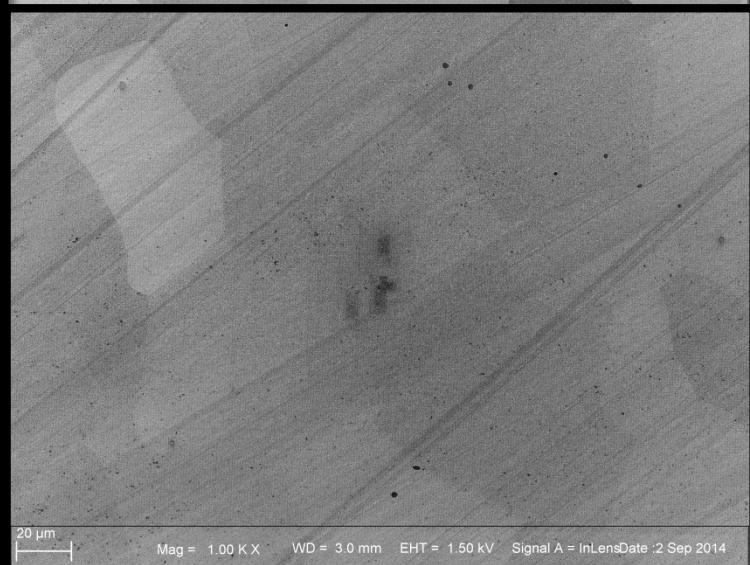
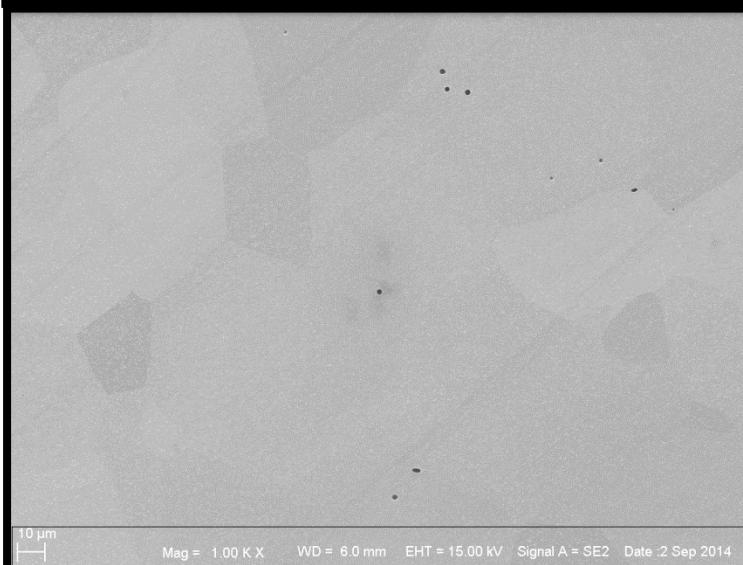
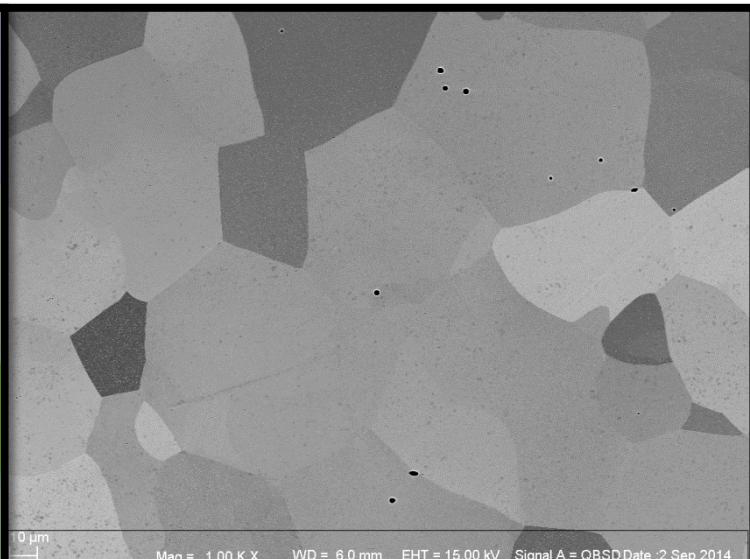
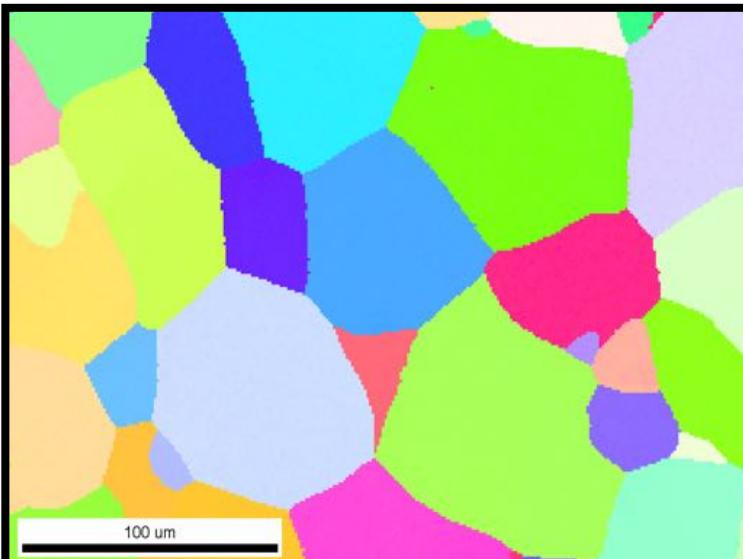


Experiment I&II



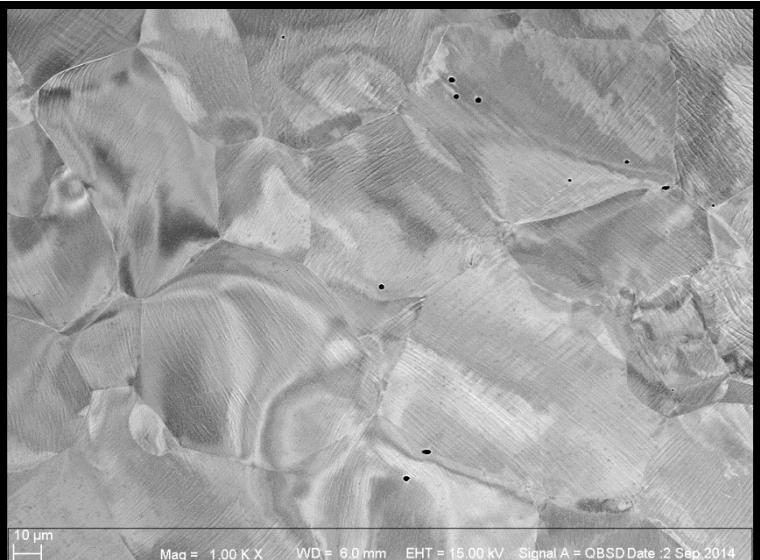
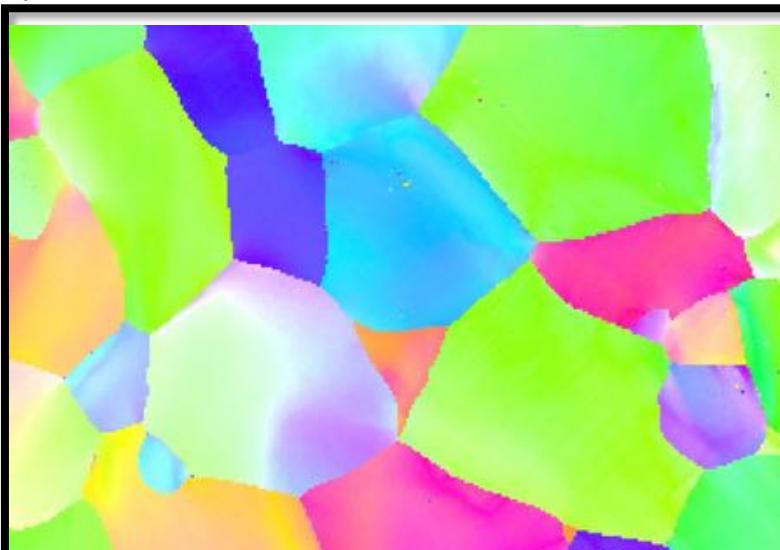
Experiment II- $\varepsilon \downarrow global = 0\%$

RD Tensile



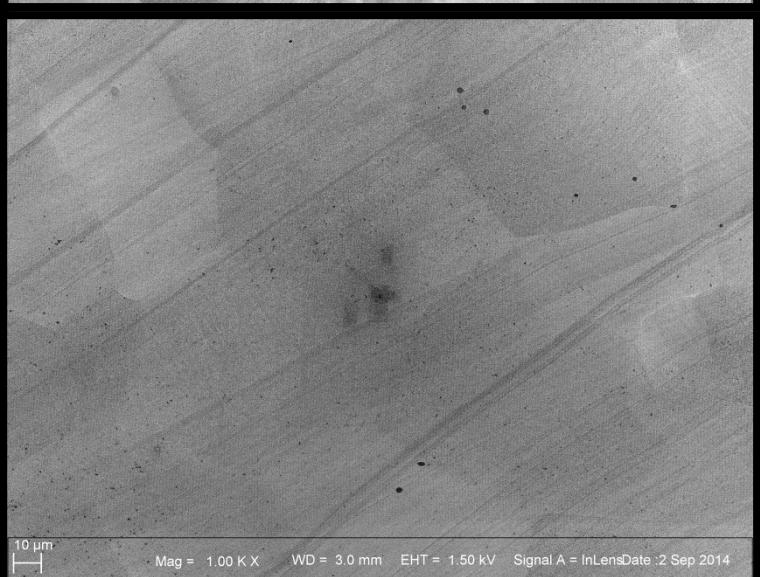
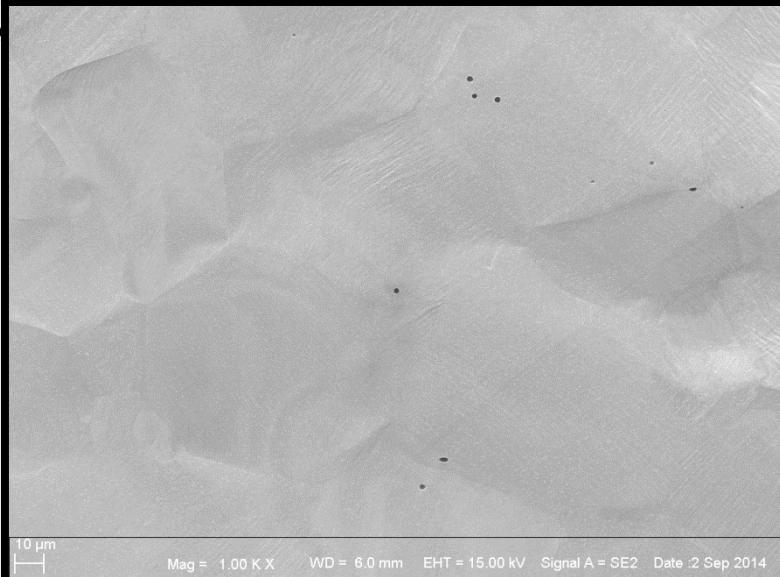
Experiment II- $\varepsilon \downarrow$ global =19%

RD Tensile



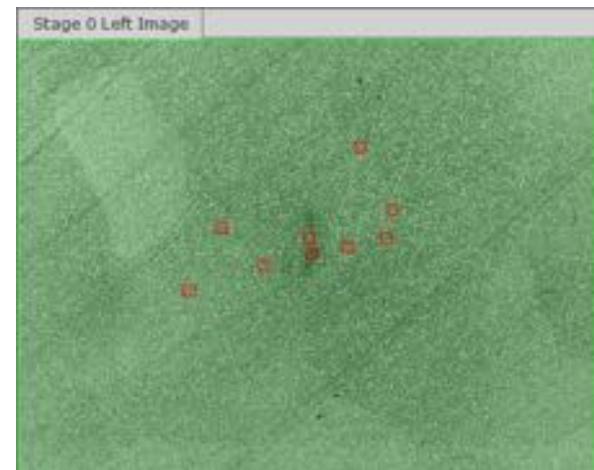
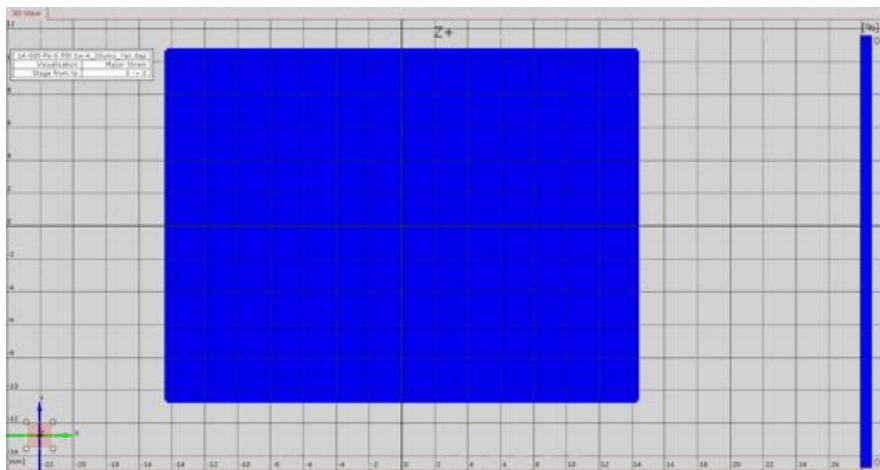
EBSD

Binning:4*4
Gain:21.27
Black:-0.28
Exposure:5.18

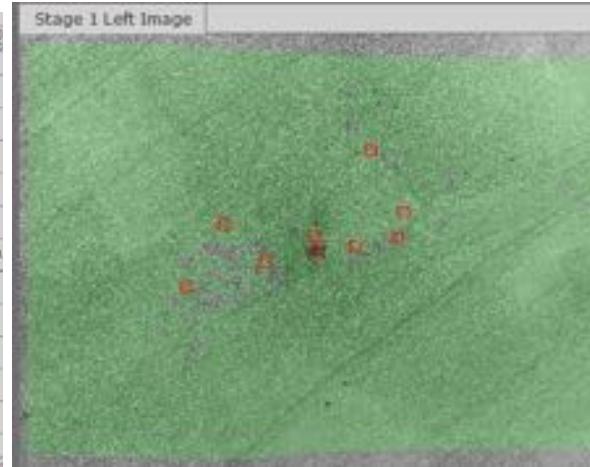
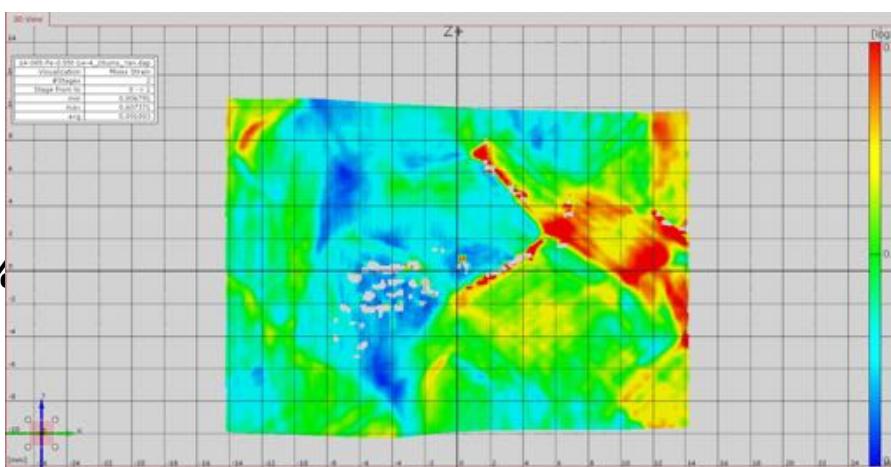


Experiment II-DIC calculation (Aramis)

Stage 0
 $\varepsilon_{\text{global}} = 0\%$

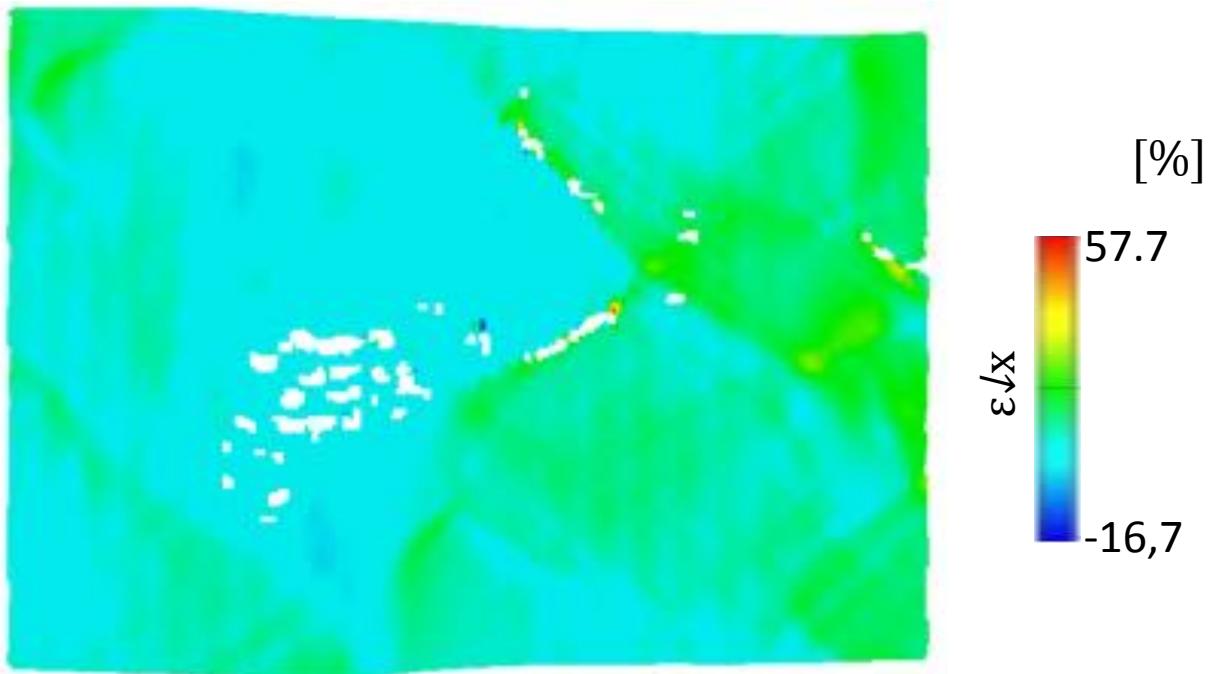


Stage 1
 $\varepsilon_{\text{global}} = 19\%$



Experiment II Results-DIC

Visualisation: $\varepsilon \downarrow x$

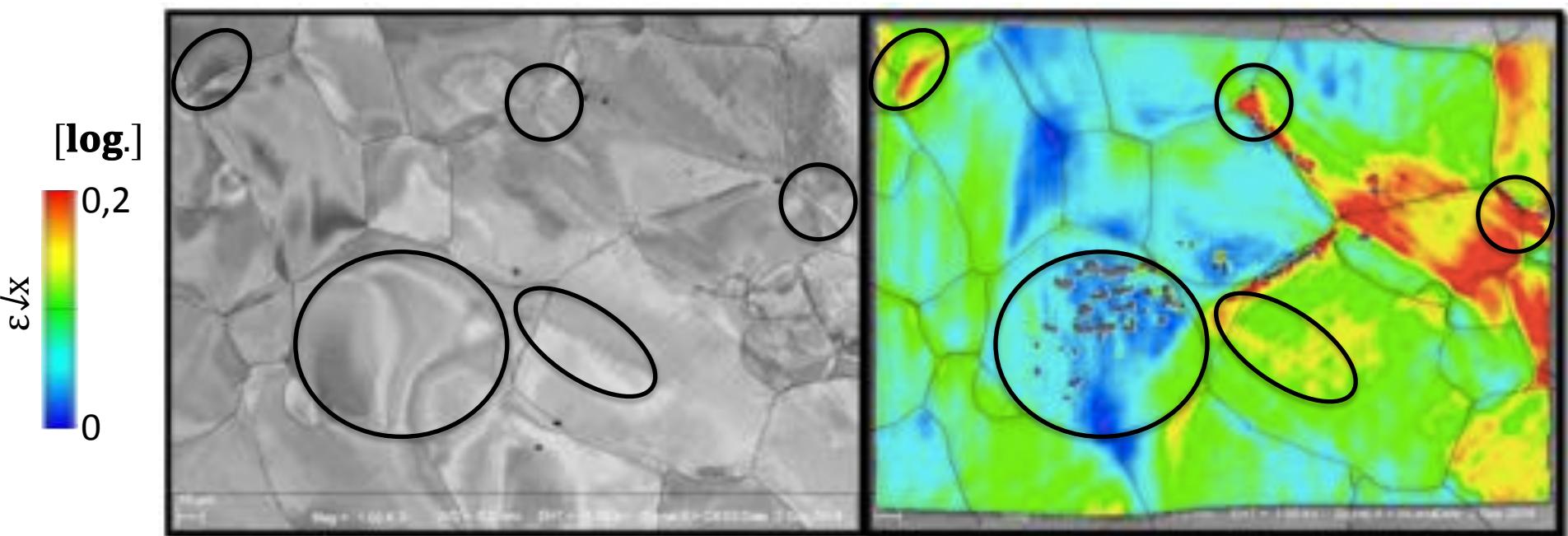


$\varepsilon \downarrow local$

min	-16.775
max	57.756
avg	8.568

Experiment II Results-DIC

Visualisation: $\varepsilon_{\downarrow x}$



Simulation Process- CPFEM



Start Simulation

- cd into /DAMASK, make spectral processing install
- wsLoad
- ssh maws02.mpie.de
- screen
- cd into working folder
- **DAMASK_spectral -l xxx.load -g xxx.geom > monitor**
- ctrl+A+D
- close window

```
MPIE\d.yan@maws01:~$ wsLoad
maws01      14:05:47 up 113 days, 15:41,  3 users, load average: 2.08, 2.08, 2.06  6%  51% (32 @ 3.0 GHz 251 GB)
maws02      14:06:05 up 72 days, 22:51,  2 users, load average: 0.06, 0.04, 0.05  0%  1% (32 @ 3.0 GHz 251 GB)
maws03      14:06:07 up 65 days, 1:49,  3 users, load average: 0.14, 0.07, 0.06  0%  1% (32 @ 3.0 GHz 251 GB)
```

Prepare Output

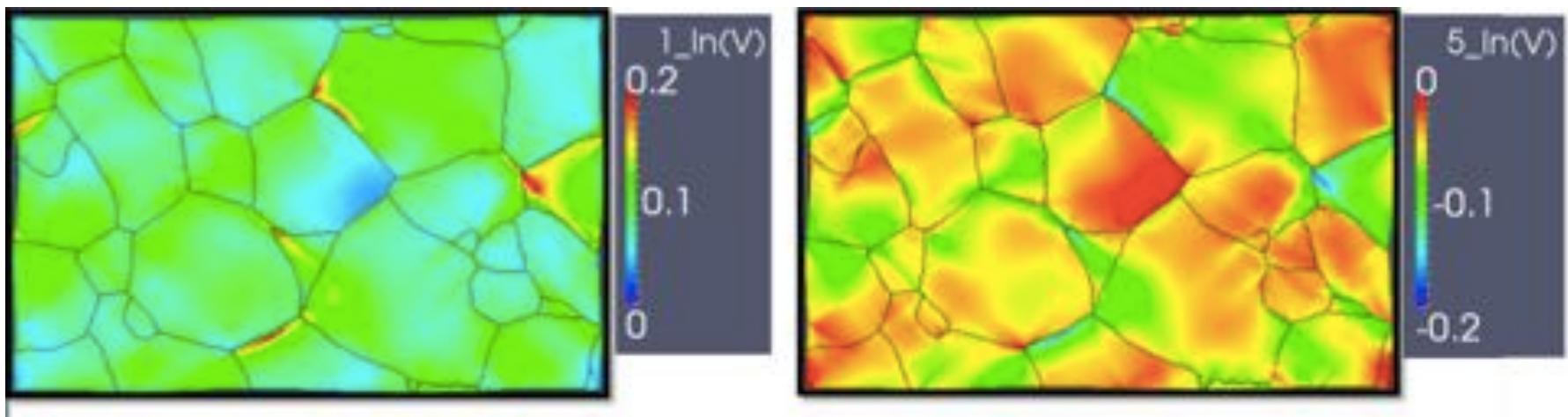
- postResults **xxx.spectralOut** --separation x,y,z --split -r 0 100 1 --increments --cr f,p,eulerangles
- cd into /postProc
- addStrainTensors -v -l **xxx_inc*.txt** [True strain]
- addCauchy **xxx_inc*.txt** [True stress]
- addMises -s Cauchy -e "ln(V)" **xxx_inc*.txt** [Equivalent stress/strain]
- addDeterminant -t f **xxx_inc*.txt** [Dilatation strain]
- addDeviator -t Cauchy -s Cauchy **xxx_inc*.txt** [Hydrostatic stress]
- For IPF, first Manually change degree to radius by excel and save as txt
- addIPFcolor -p 1 0 0 -e eulerangles **xxx_inc*.txt** [Inverse pole figure]

For Visualize in Paraview

- 3Dvisualize -s [1-9]_f,[1-9]_p,[1-9]_"ln(V)",
[1-9]_Cauchy,"Mises(ln(V))","Mises(Cauchy)","det(f)","sph(Cauchy)",[1-3]_IPF_100 **xxx_inc*.txt**
- vtk_addVoxelgridData --vtk **mesh_xxx_inc*.vtk** --color **xxx_inc*.txt**

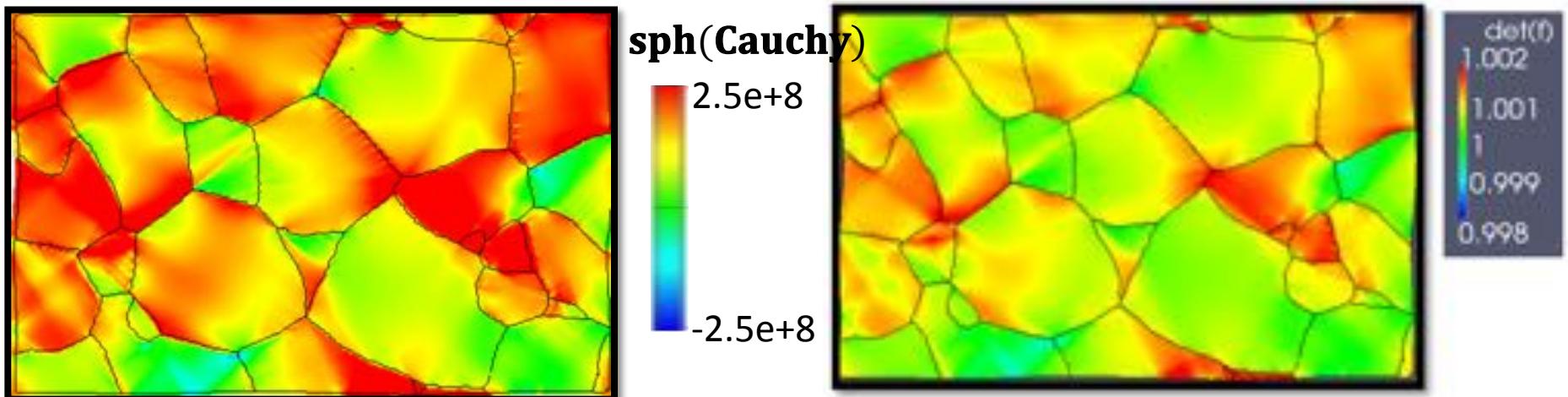
Simulation Results- CPFEM

DAMASK: True Local Strain Partitioning of Pure Fe sample in x and y directions



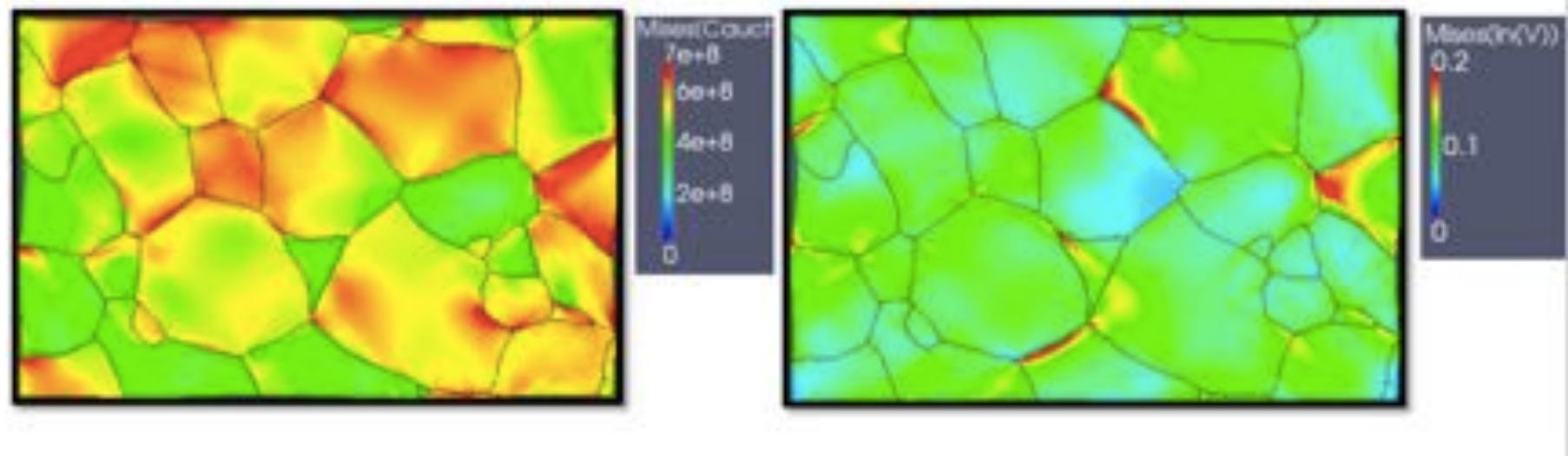
Simulation Results- CPFEM

DAMASK: Hydrostatic Stress and Determinant of the Deformation Gradient in Pure Fe



Simulation Results- CPFEM

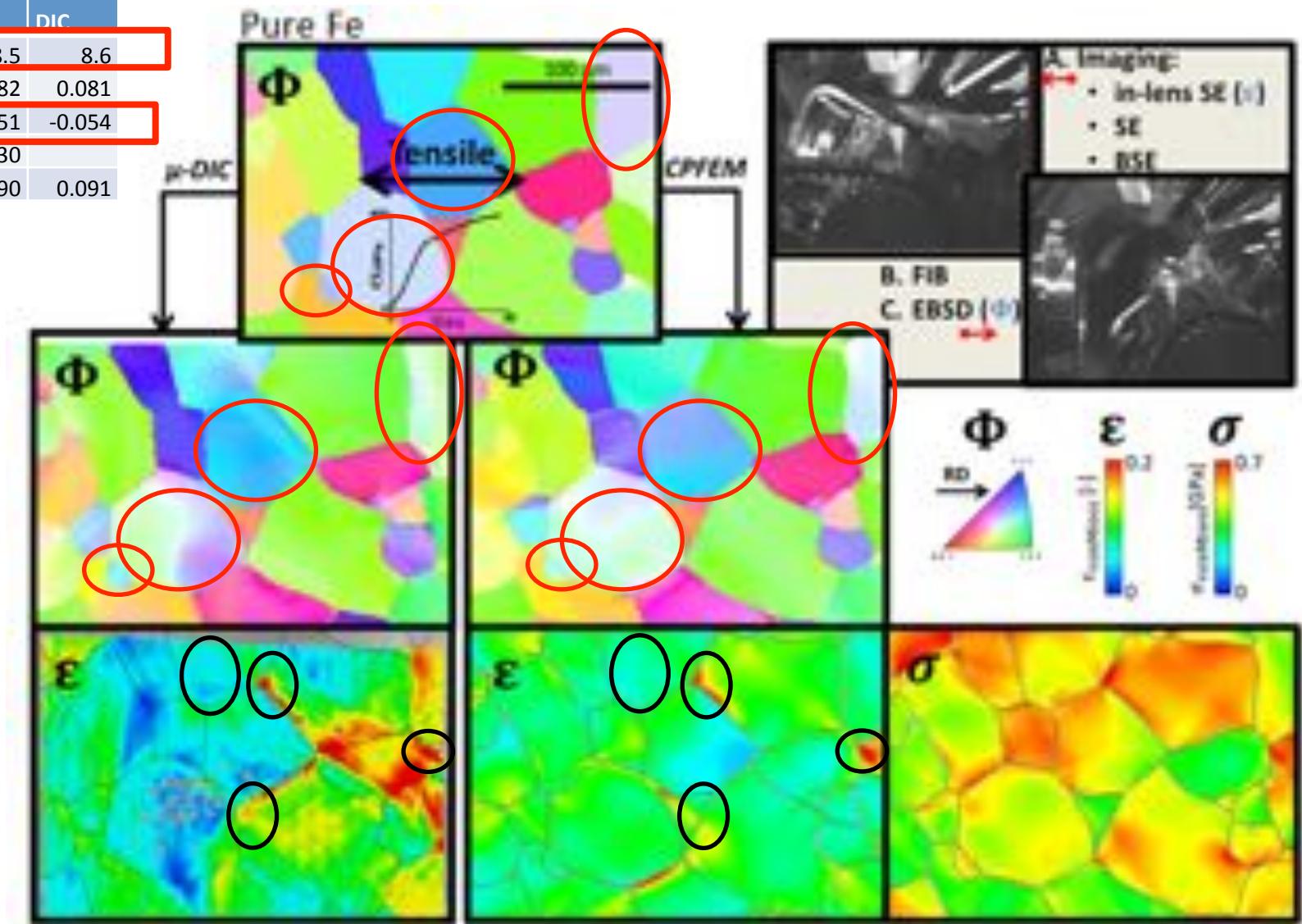
DAMASK: Local Equivalent Stress and Strain Partitioning in Pure Fe



DIC and CPFEM Comparison

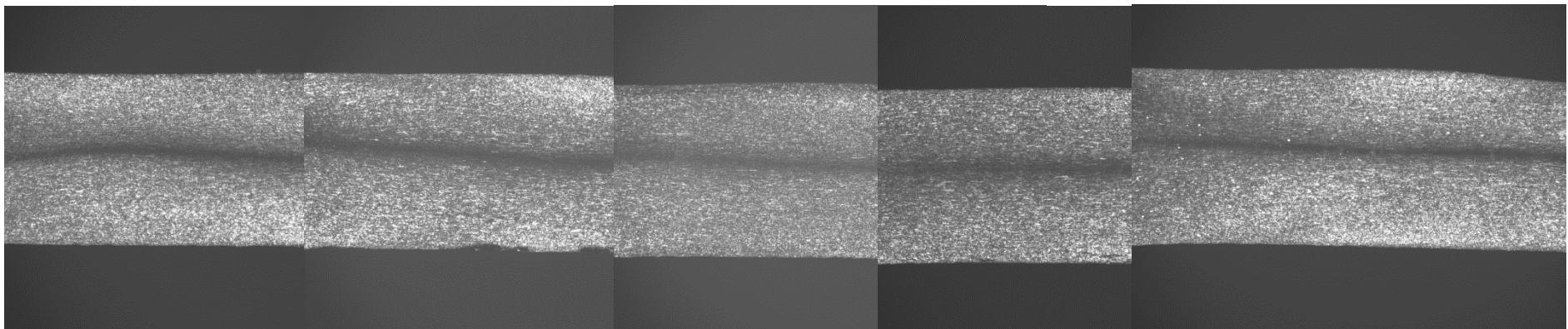
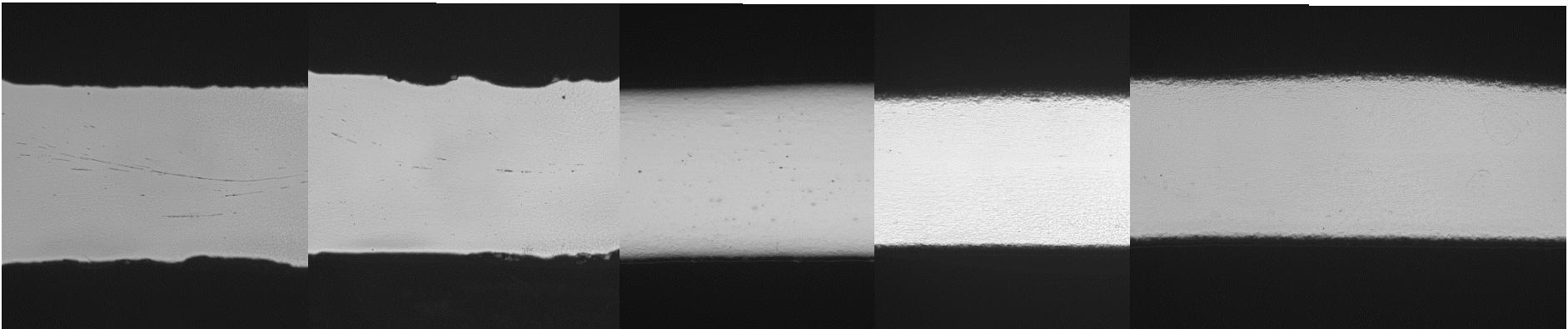


Ave.	Damas k	DIC
x	8.5	8.6
lnx	0.082	0.081
lny	-0.051	-0.054
lnz	-0.030	
mises	0.090	0.091



Previous Work

Optimisation of the UFG steel sample



Summary

- Introduction to DAMASK and ITS
- Experimental Results from DIC
- Simulation Results from DAMASK
- Experiment and Simulation Comparison
 1. Strain with a grain
 2. Strain at triple points
 3. slip transition across grain boundary
- Previous Work



End

- Thank you all!
- Questions?