



# Precise prediction of NNS HIP components through DEM and FEM modelling

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## Introduction

- Purpose of the study
- Process description

## Simulation of HIP Process with “Image Analyzed” Densities Distribution

- “Image Analyze” for investigation of initial density distribution
- FEM Simulation

## Simulation of HIP Process with DEM simulated Densities Distribution

- Introduction of DEM and Modelling Approach
- Validation, Simulation and Comparison

## Summary and Outlook



**Welded Capsule (SS304) filled with powder (SS316L) before HIP Process**



$T_{HIP} = 1125^{\circ} \text{ C}$   
 $p_{HIP} = 110 \text{ MPa}$   
 $t_{Holding} = 2 \text{ h}$

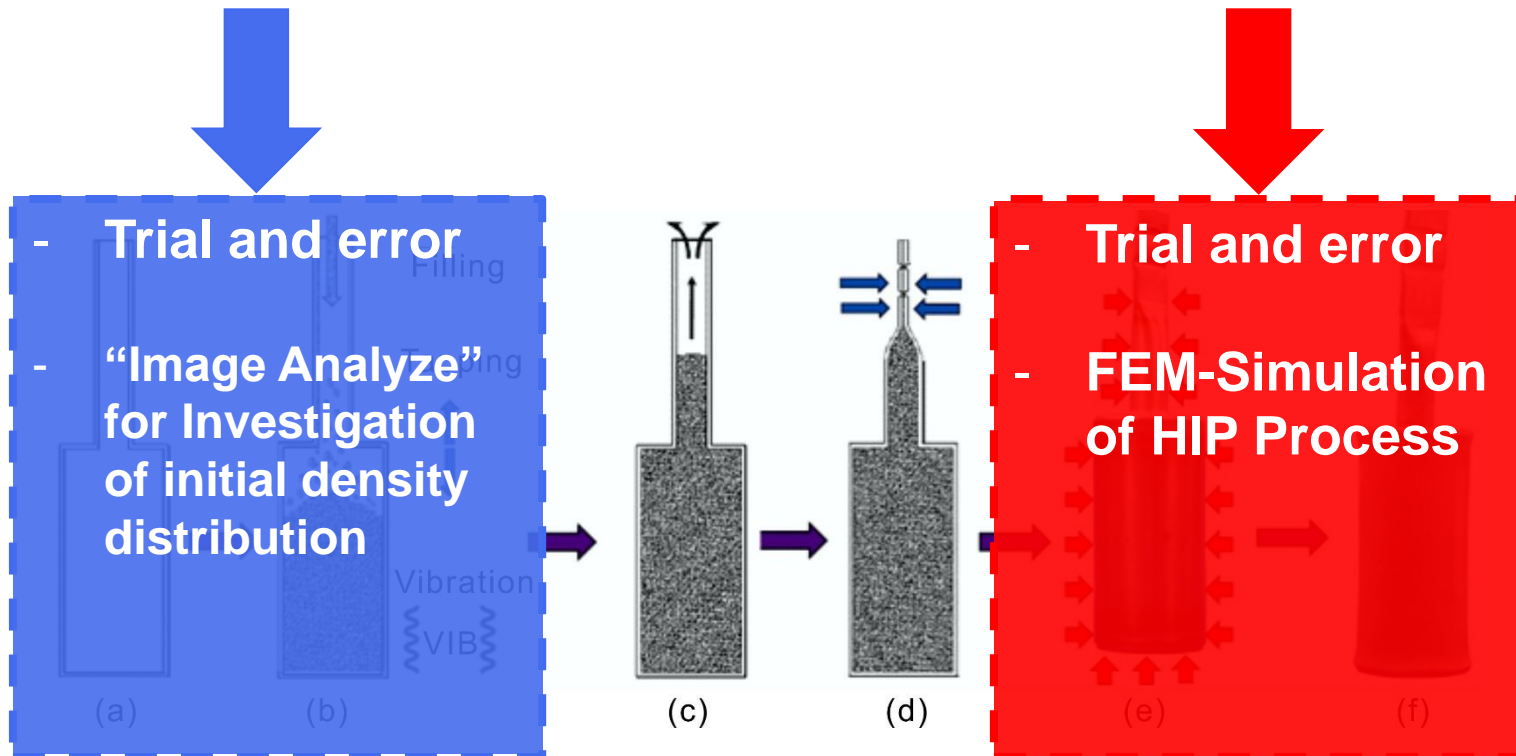


**Capsule (SS304) filled with powder (SS316L) consolidated to full dense after HIP Process**

- **Anisotropic shrinkage of capsule during PM HIP:**
  - Inhomogeneous powder densities distribution in the capsule
  - Temperature gradients and inhomogeneity in the HIP Unit
  - Imperfection of material

- Capsule design
- Particle size
- Initial relative density and distribution

- Material properties
- Process parameters  
(Time, Pressure, Temperature, etc.)



**HIP process in general:** (a) Capsule production; (b) Filling powder, tapping and vibration; (c) Evacuation; (d) Sealing capsule; (e) Applying high temperature and high pressure; (f) Full density HIPed component

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## **Simulation of HIP Process with “Image Analyzed” Densities Distribution**

- **“Image Analyze” for investigation of initial density distribution**
- **FEM Simulation**

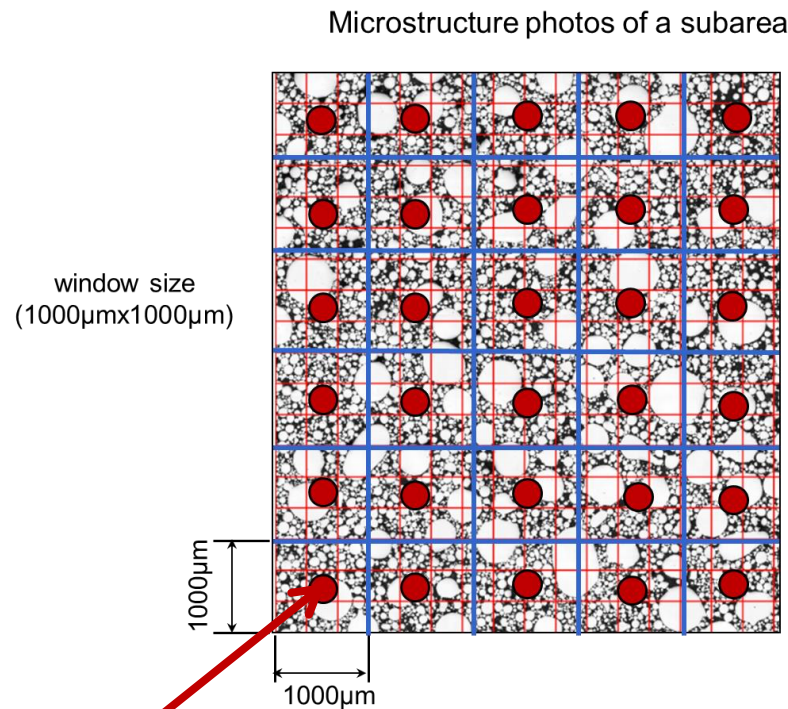
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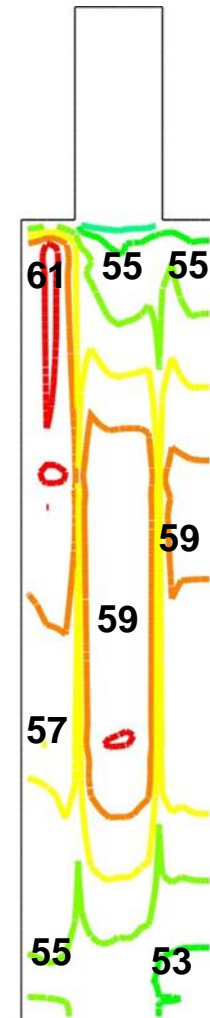
# “Image Analyze” for investigation of initial density distribution

20			
67	66	59	← (L1)
1a	1b	1c	
68	67	59	← (L2)
2a	2b	2c	
69	68	59	← (L3)
3a	3b	3c	
70	70	59	← (L4)
4a	4b	4c	
71	70	59	← (L5)
5a	5b	5c	



$$\text{Relative Density (RD) of each window} = \frac{\text{Amount of powder}}{\text{Total area of the window}}$$

RD [%]



## ABAQUS

Get initial Values:

$\sigma_{ij}, \Delta \varepsilon_{ij}, T, t, \Delta t$   
 $\rho, \rho_0, \sigma_0, r_1, f_1, \varepsilon_{ij}^{inel}, \varepsilon$

$r_1$  plastic hardening parameter

Calculate:

$E, \nu, \mathbf{C}^{el}, \sigma_{eq1}, \sigma_{eq2}, f_1$

$f_1 = \sigma_{eq1} - r_1 - \sigma_y$

If  $f_1 \geq 0$

No

Yes

Plastic Deformation:  $\Delta \varepsilon_{ij}^{pl}$

Viscoplastic Deformation:  $\Delta \varepsilon_{ij}^{cr}$

Update:  $\Delta \varepsilon_{ij}^{el}, \Delta \sigma_{ij}$

$\Delta \varepsilon_{ij}^{el} = \Delta \varepsilon_{ij} - \Delta \varepsilon_{ij}^{pl} - \Delta \varepsilon_{ij}^{cr}$

$\Delta \sigma_{ij} = \mathbf{C}_{ijkl}^{el} \cdot \Delta \varepsilon_{kl}^{el}$

## User-defined Material Model

Calculate:

$\Delta \varepsilon_{kk}^{inel}$

$\Delta \varepsilon_{kk}^{inel} = \Delta \varepsilon_{kk} - \Delta \varepsilon_{kk}^{el}$

Calculate:

$\rho$

RD

$\rho^{(t+\Delta t)} = \rho^t \cdot \Delta \varepsilon_{kk}^{inel}$

Update Jacobi-Matrix:  $\mathbf{J}$

End of Increment

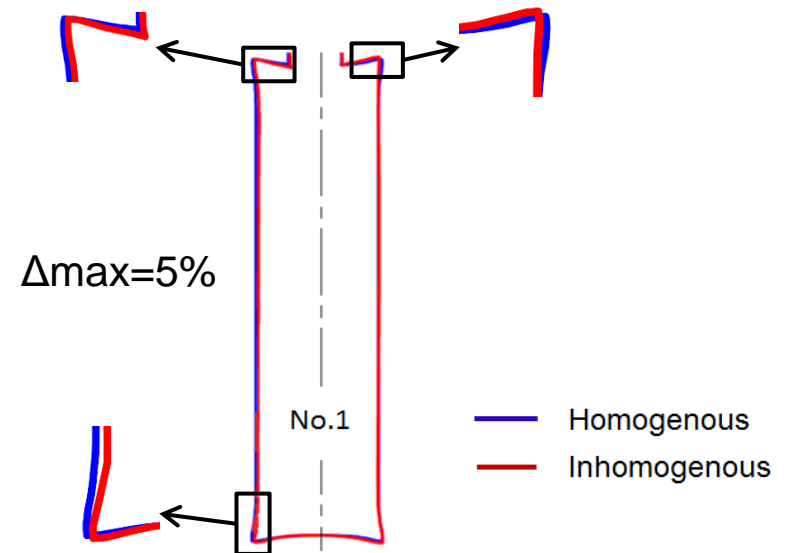
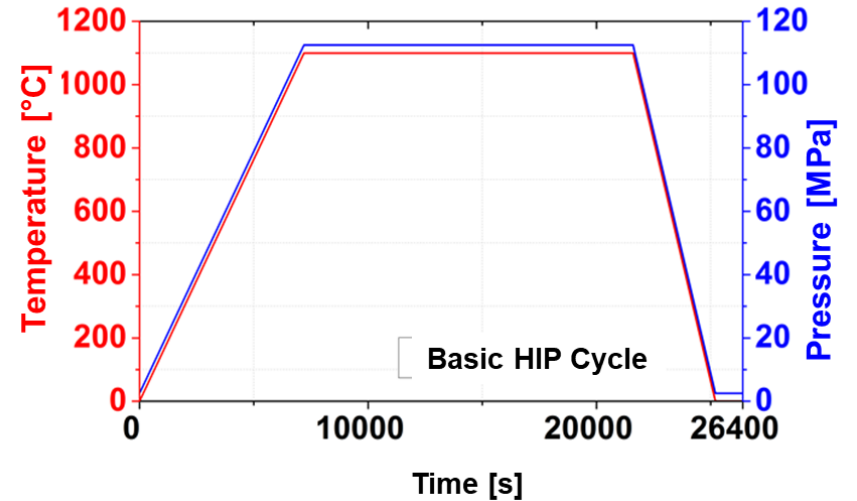
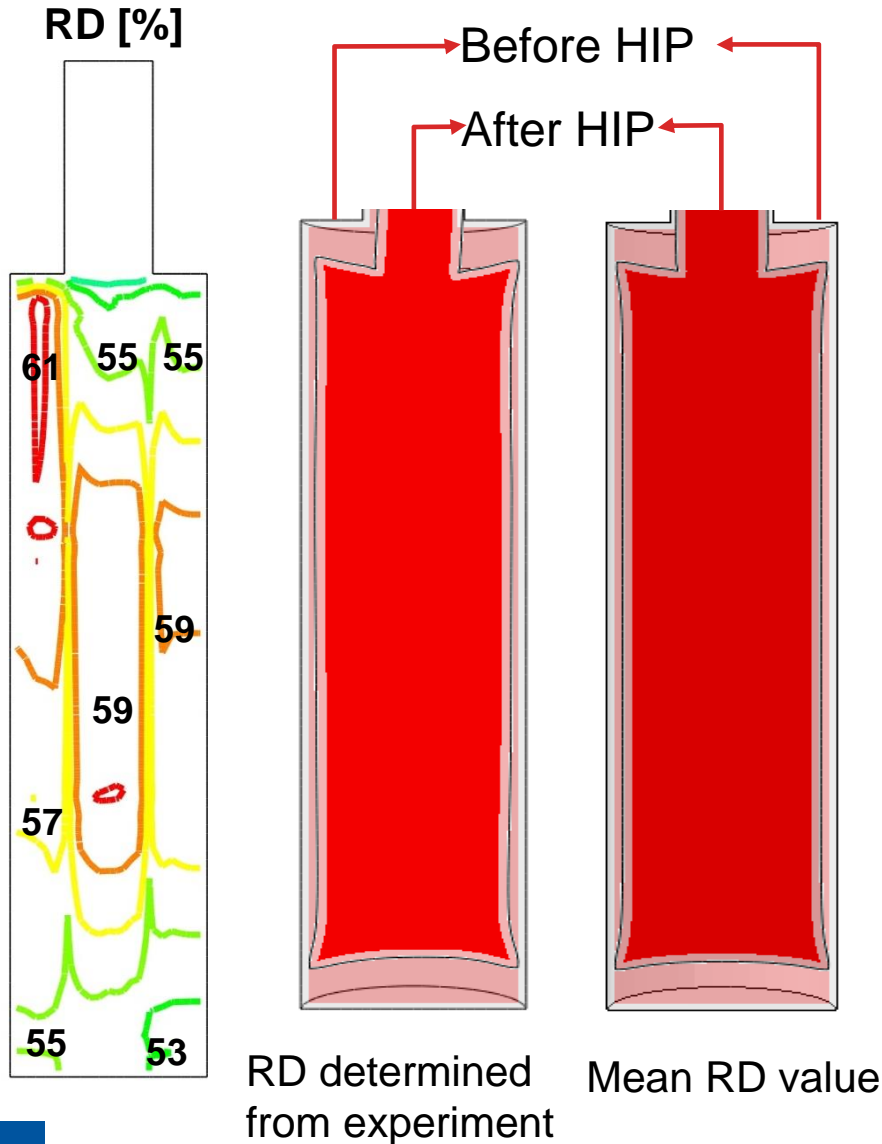
No, to the next increment

Yes

End



# HIP Simulation with determined densities field in the capsule





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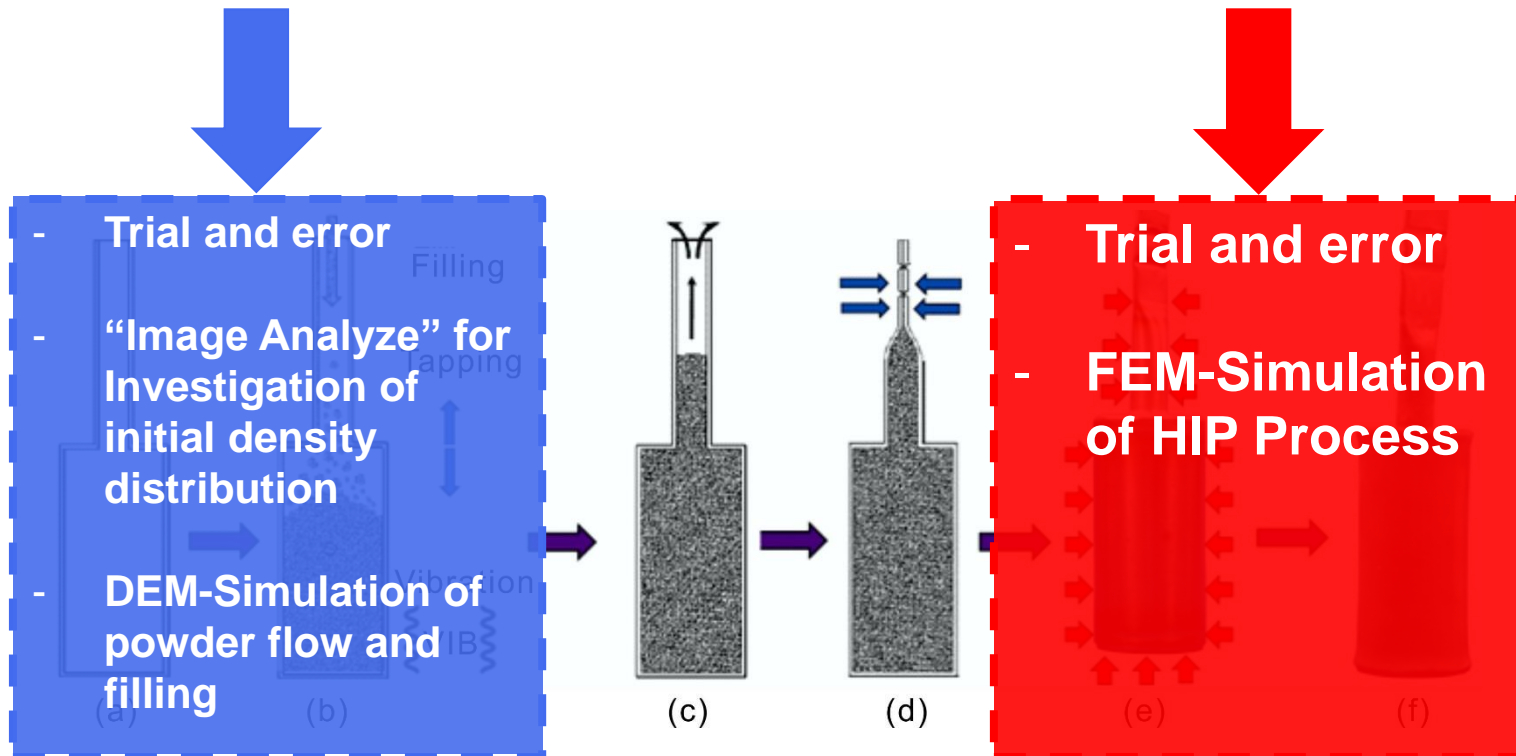
## **Simulation of HIP Process with DEM simulated Densities Distribution**

- **Introduction of DEM and Modelling Approach**
- **Validation, Simulation and Comparison**

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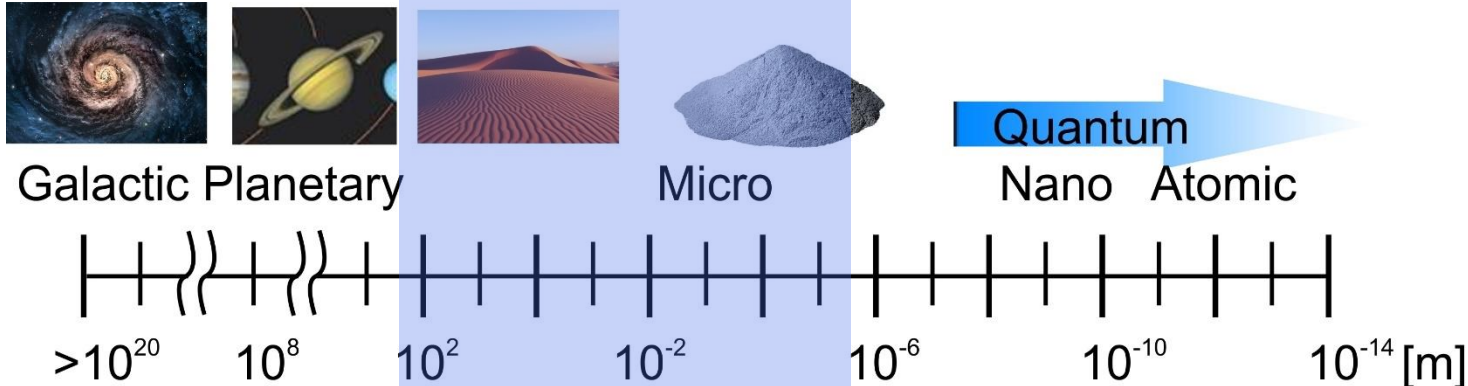
- Capsule design
- Particle size
- Initial relative density and distribution

- Material properties
- Process parameters  
(Time, Pressure, Temperature, etc.)



**HIP process in general:** (a) Capsule production; (b) Filling powder, tapping and vibration; (c) Evacuation; (d) Sealing capsule; (e) Applying high temperature and high pressure; (f) Full density HIPed component

## *Discrete Element Method (DEM)*



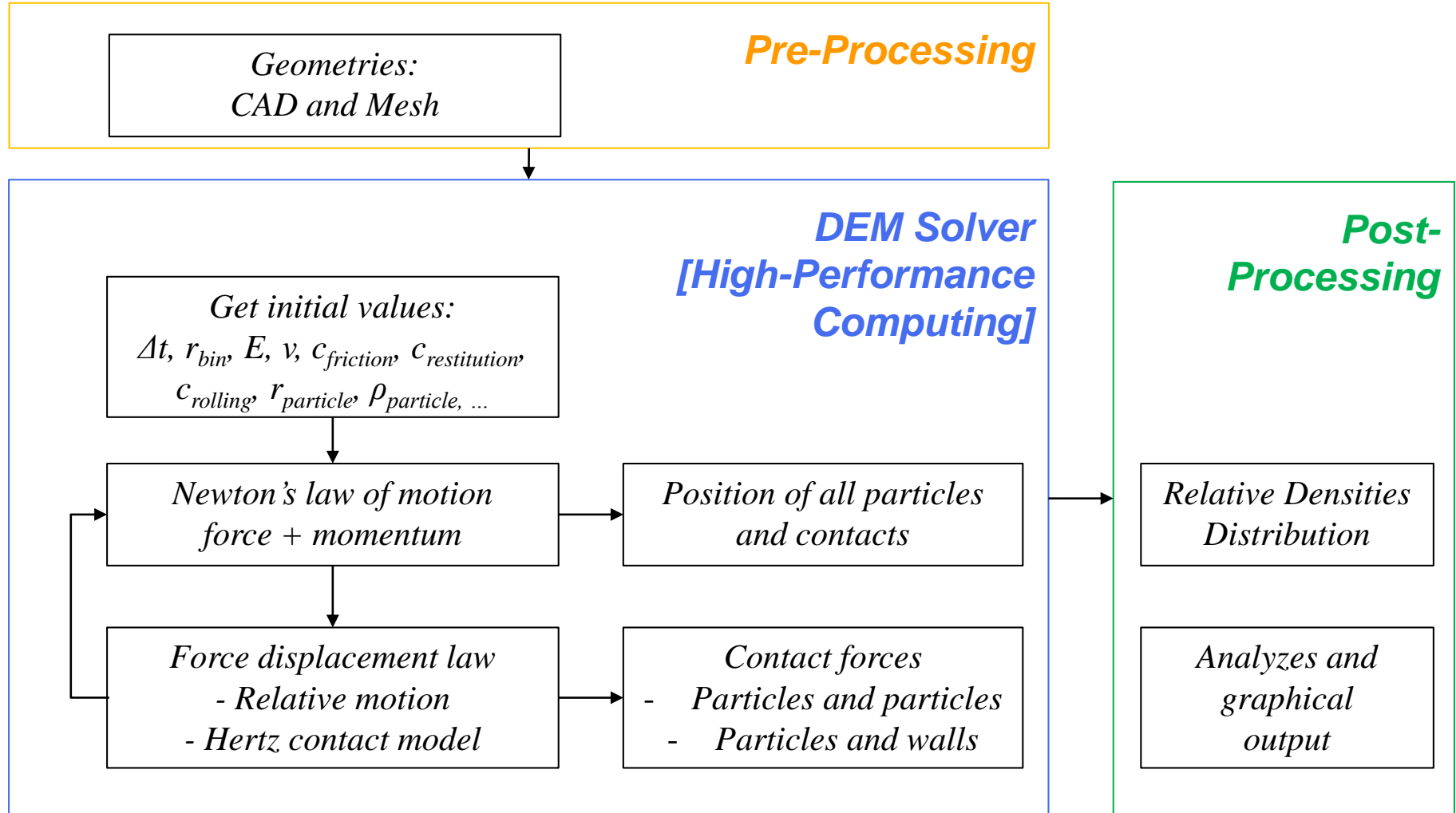
- mathematical description of single objects / particles / granules
- meshless method compared to FEM
- interactions between objects only at contact points

### **Advantages:**

- Few equations which have to be solved for each object
- Forces and motions can be investigated for each single object, which are usually not measureable

### **Limits in usage:**

- Computing time-consuming for real particle systems consists of billions of elements



Materials:

- Stainless Steel 316L (Powder)
- SS304 (Underlayment)

Particles: radius 250  $\mu\text{m}$

Density  $\rho$  7800 kg/m<sup>3</sup>

Particle radius  $r$  250  $\mu\text{m}$

Young's modulus  $E$

Poisson's ratio  $\nu$

Coefficient of restitution  $c_{restitution}$

Coefficient of static friction  $c_{friction}$

Coefficient of rolling friction  $c_{rolling}$

Capacity: 10<sup>9</sup> Particles

- Iron powder 500 kg
- Ti-Alloys powder 300 kg

Rayleigh time increment:

$$\Delta t_R = \frac{\pi r \sqrt{\rho / G}}{(0.1631 \nu + 0.8766)}$$

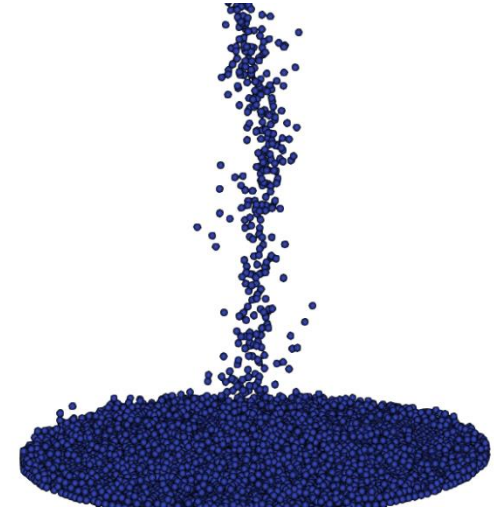
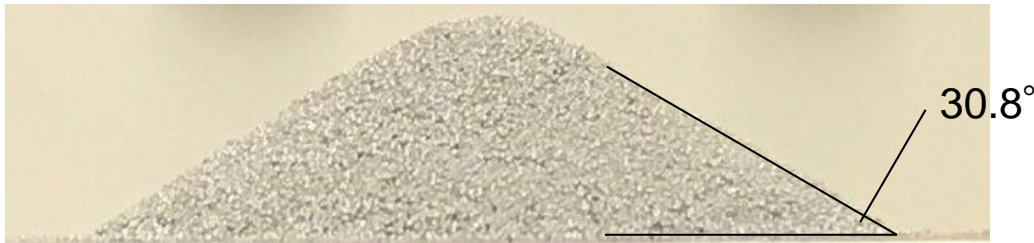
Hertz time increment:

$$\Delta t_H = 2.87 \left[ \frac{(\rho (4/3) \pi r^3)^2}{r E^2 v_{max}} \right]^{0,2}$$

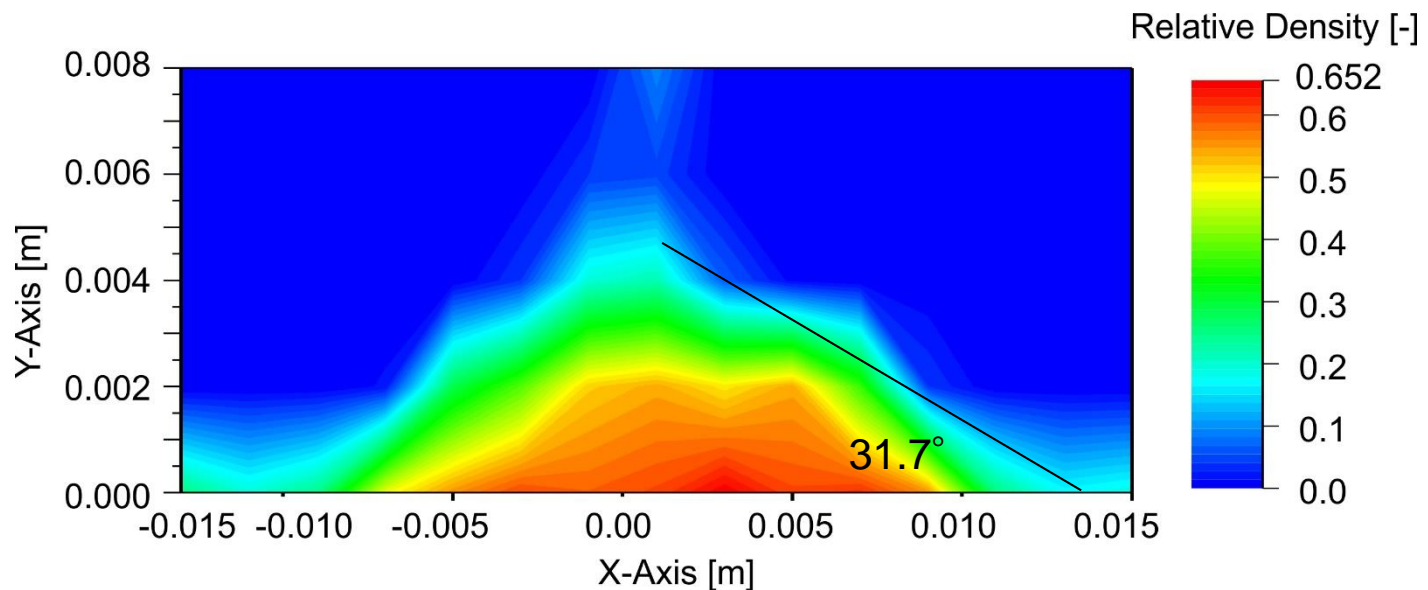
$$\Delta t = 0,20 \min(\Delta t_R, \Delta t_H)$$

Time increment 0.000001s

## *Angle of repose*

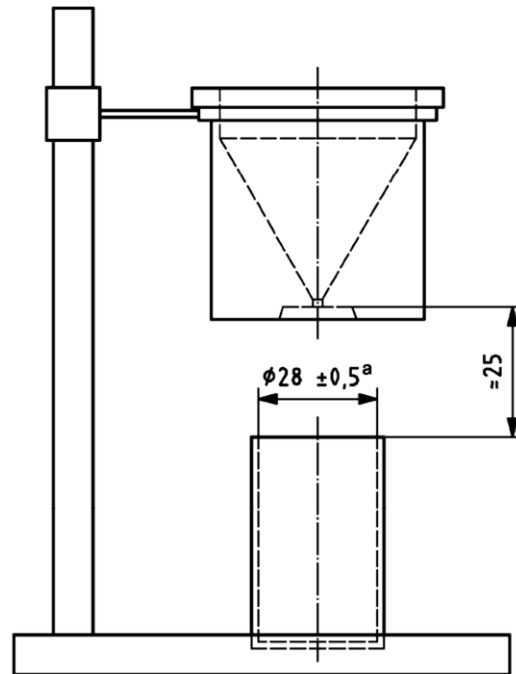


Deviation: 2.9%



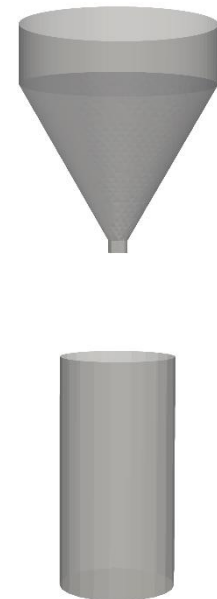
**Flowmeter**

**Experiment**



**Simulation Model**

Simplified



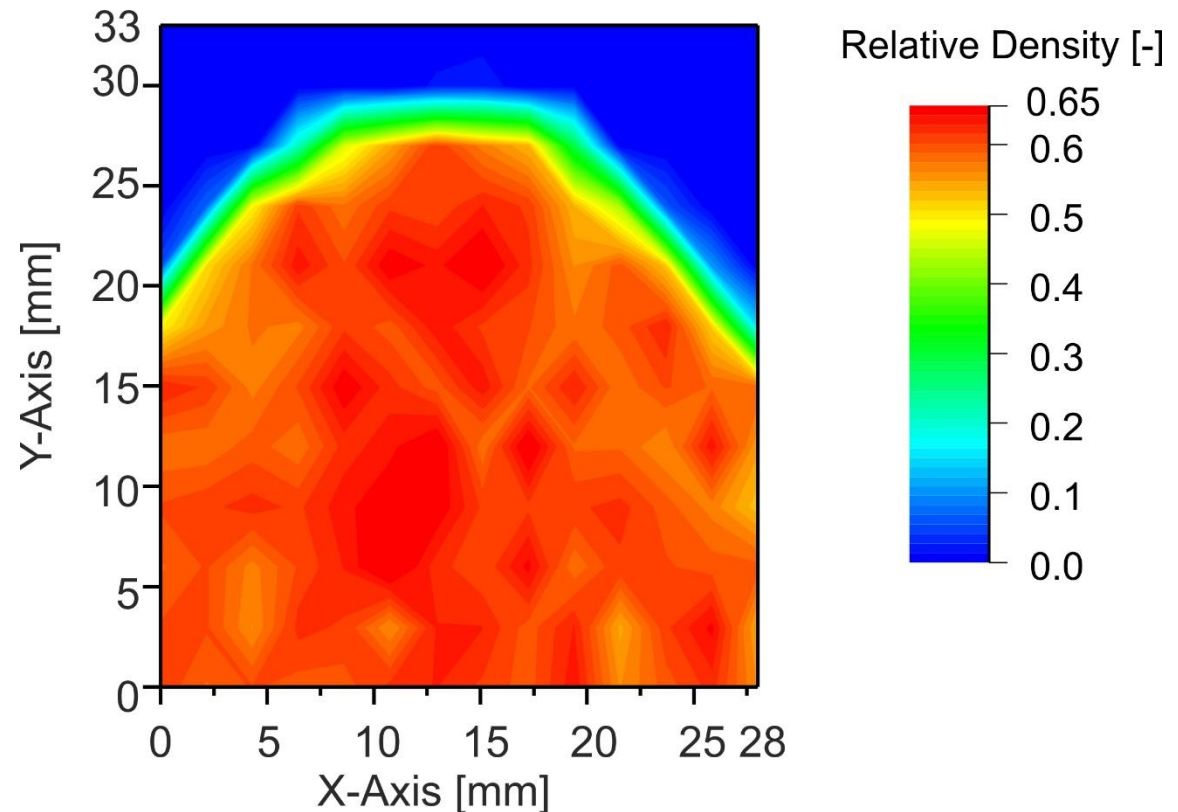




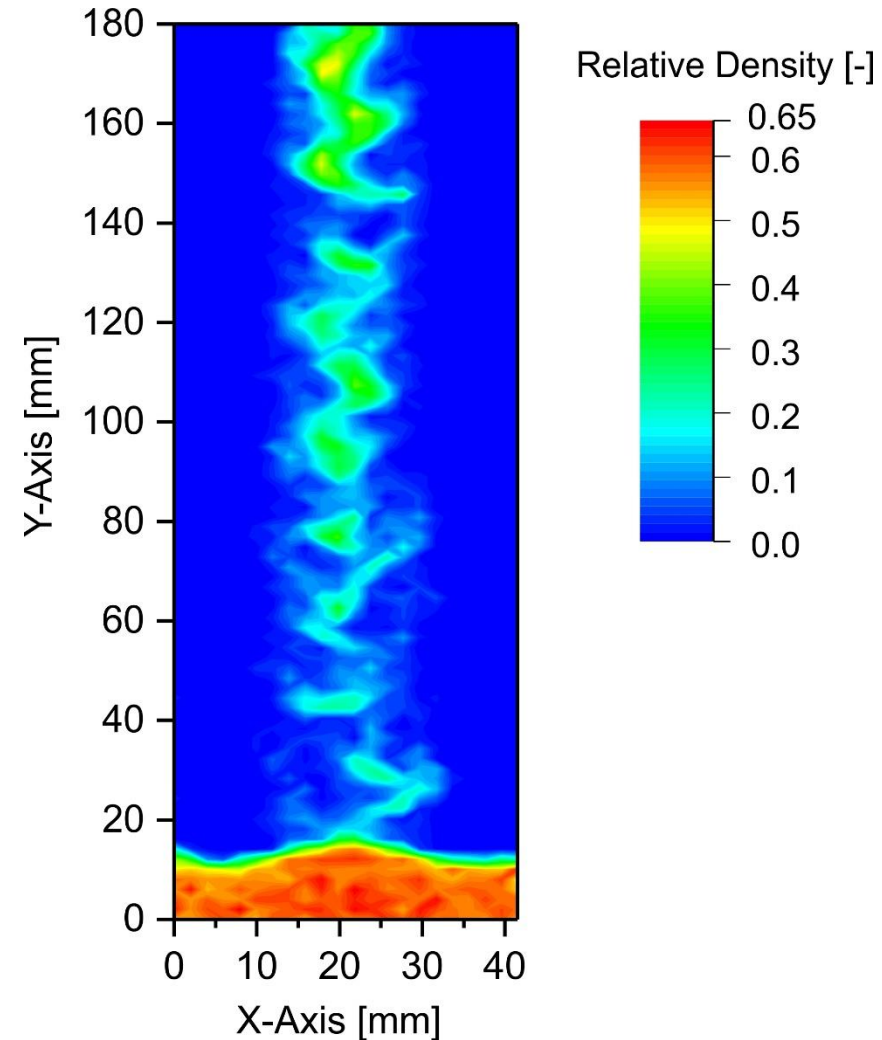
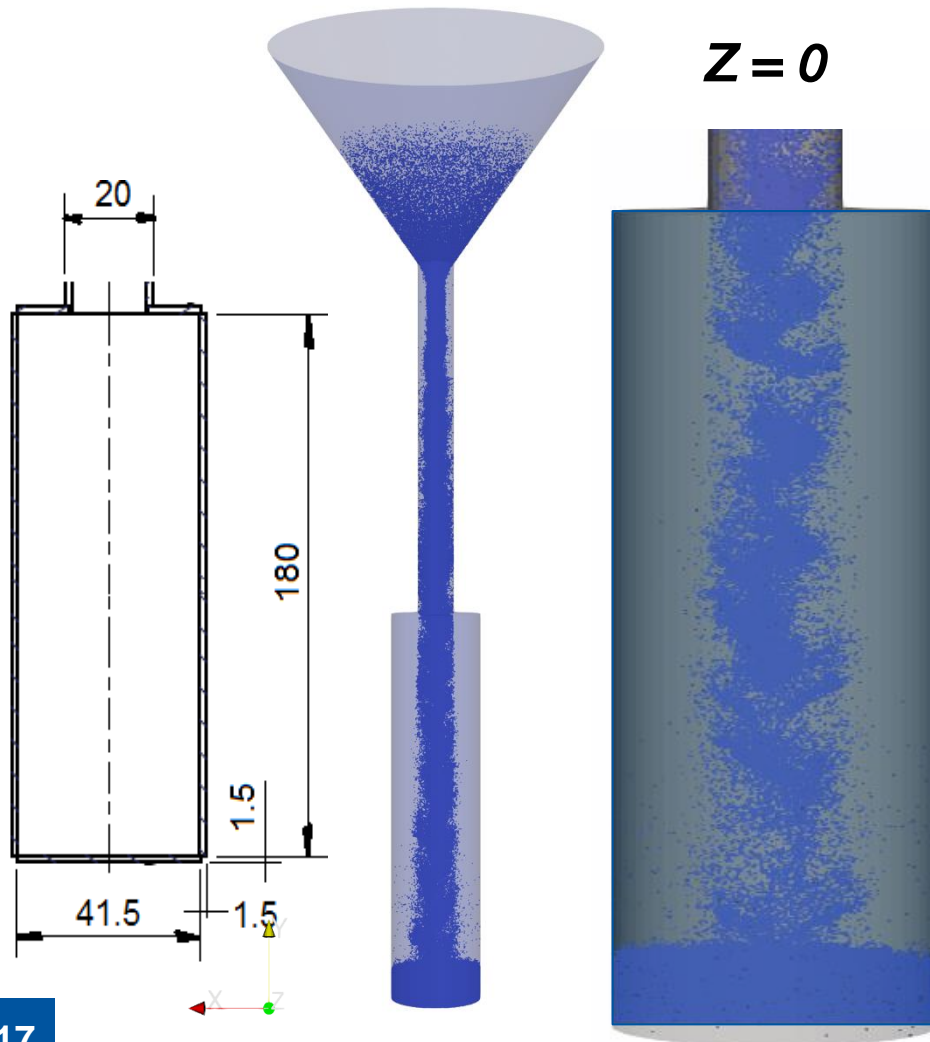
**SS316L Powder**

**Time: Exp. 18.5 s / 50g, Sim. 20.1 s / 50g**

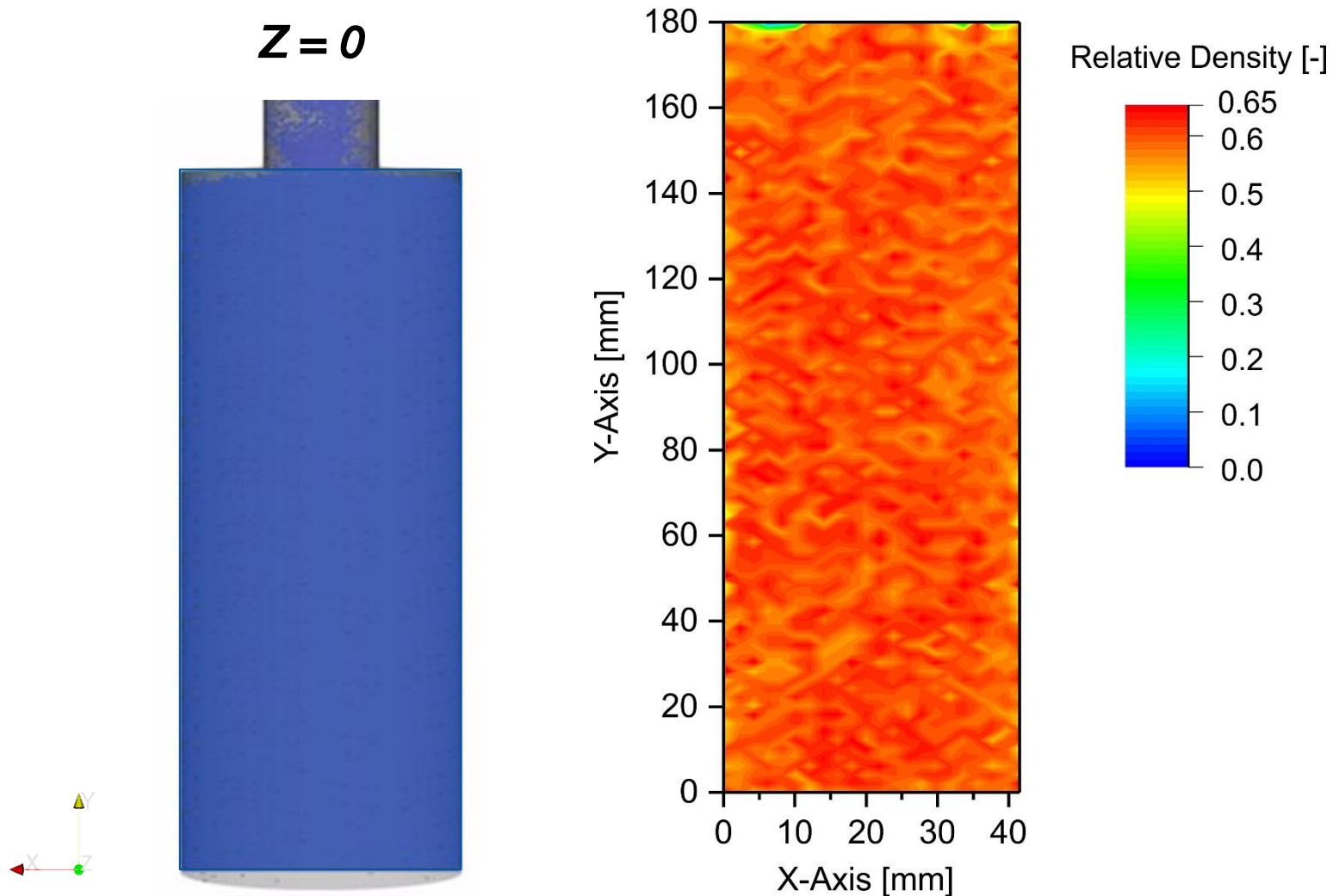
**Deviation: 1.6 s , 10.2 %**



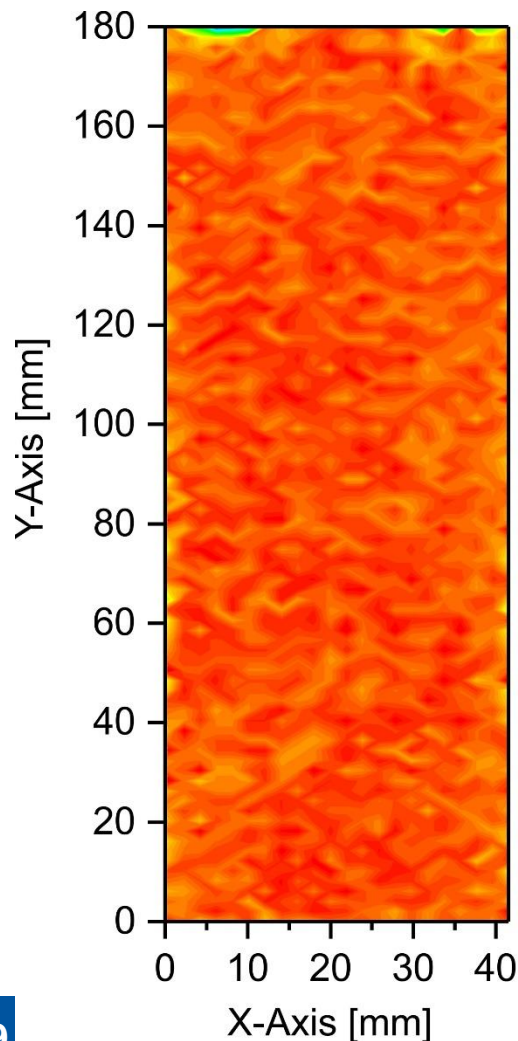
## *Filling process using DEM Simulation*



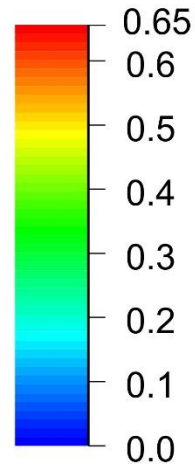
## *Filling process using DEM Simulation*



# HIP Simulation with determined densities field in the capsule



Relative Density [-]

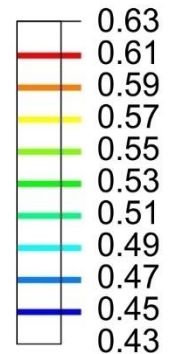


**DEM -> FEM**

**three dimensional  
densities field**

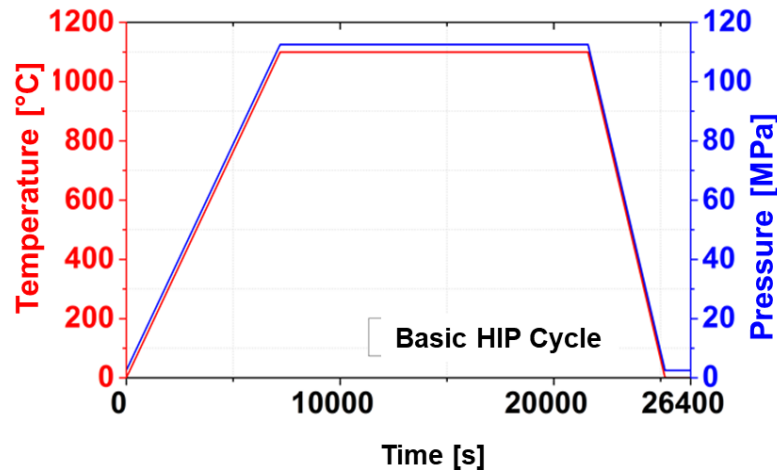
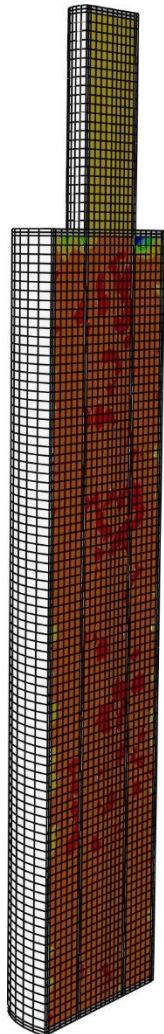


RD [-]

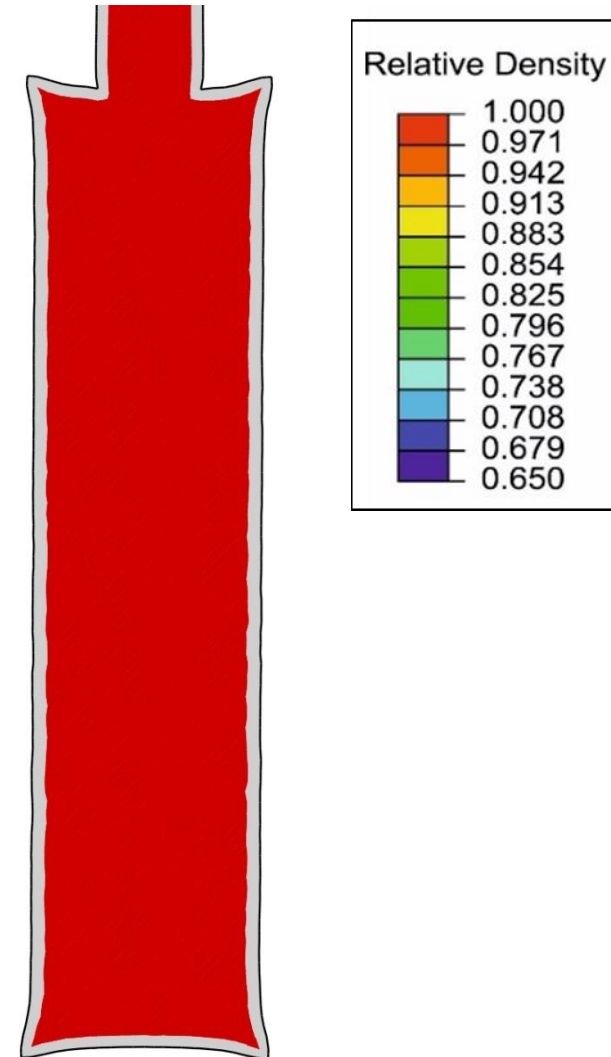




# HIP Simulation with determined densities field in the capsule



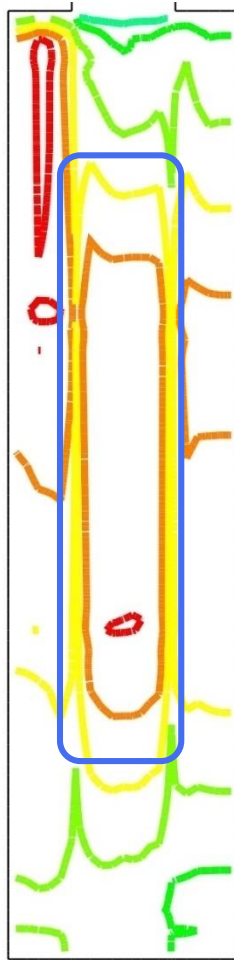
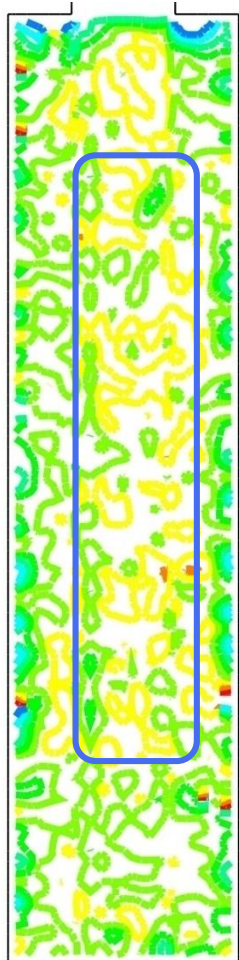
Cylinder Capsule  
Thickness: 1.5 mm  
Element size~2mm  
Capsule material: SS304  
Powder material: SS316L



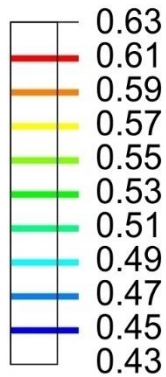
# Comparison of initial and final shapes

DEM  
Simulated

„Image  
Analyse“



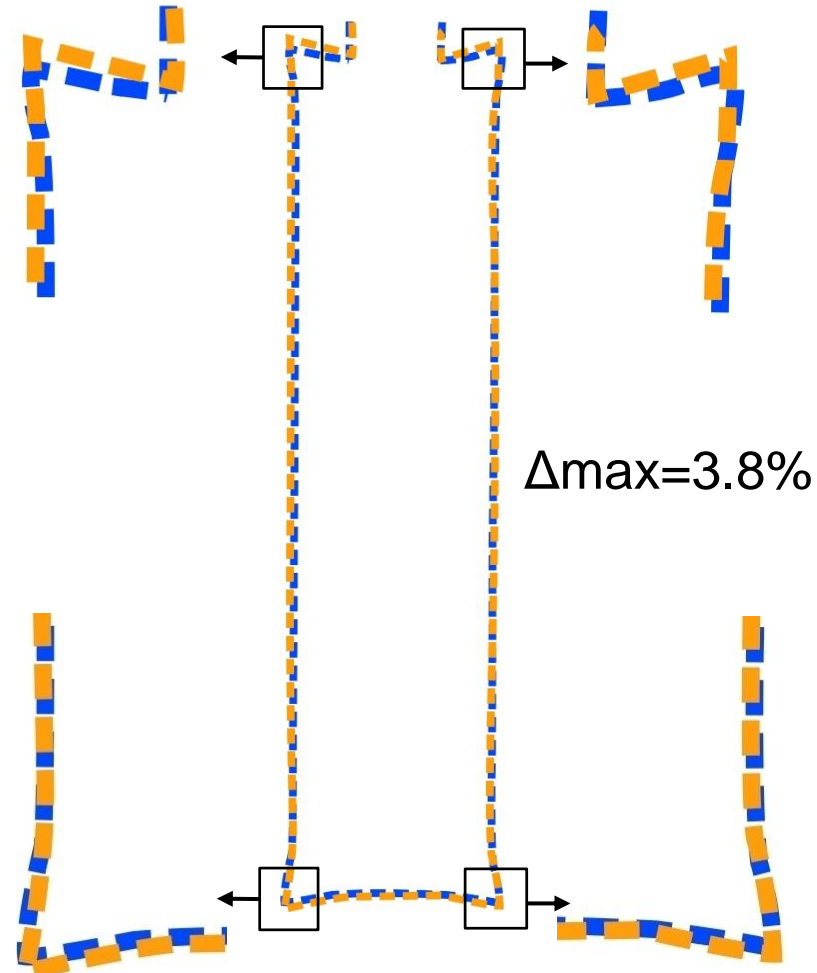
RD [ - ]



Difference:  
~2%

Contours of  
final shape

Initial RD input from  
DEM Simulation  
Image Analyse



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- **Summary**

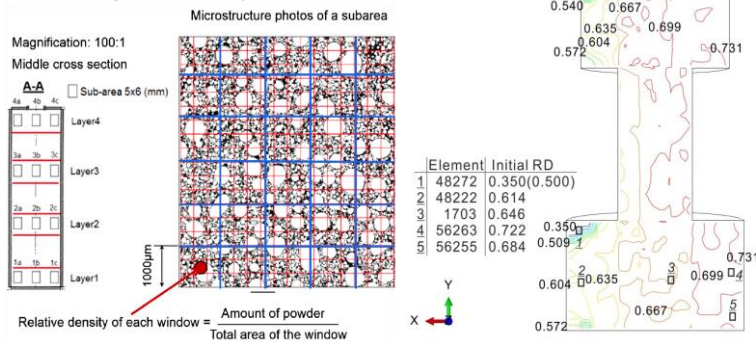
- Simulation approach with coupled DEM and FEM Modelling has been used to calculate the capsule filling densities and predict the densification behavior of HIP Process.
- The simulation results correspond well with the experimental measurement.

- **Outlook**

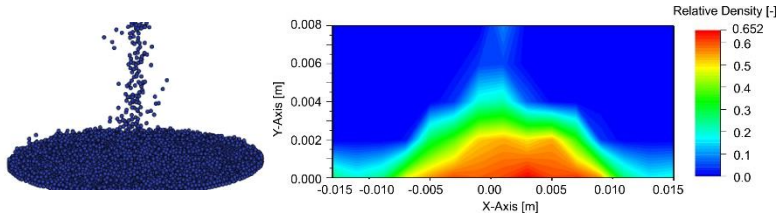
- Influences of particle size distribution
- Influences of pre-desification process, vibration and tapping
- Influences of more complex shape

## Inhomogeneous initial density distributions

### "Image Analyze"



### DEM Simulation



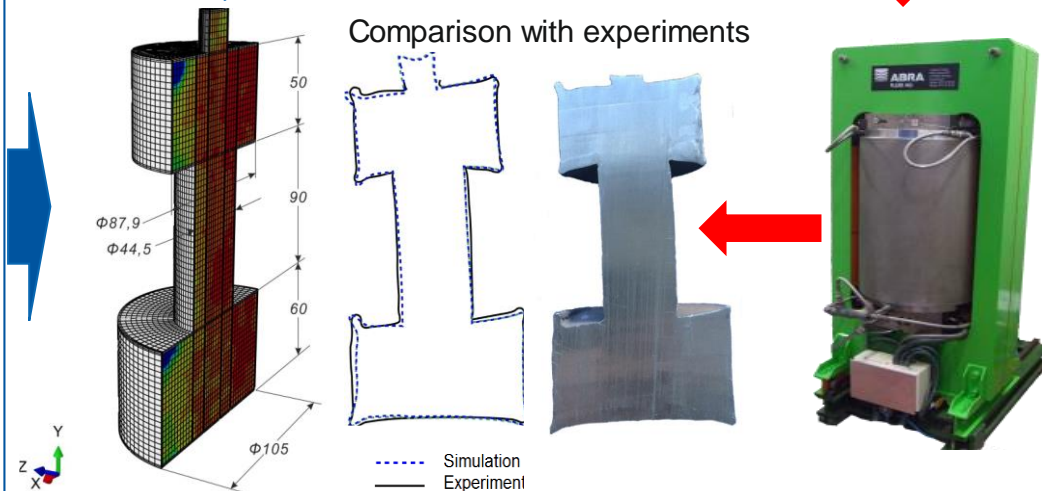
## Temperature gradient on the capsule surfaces

## Capsule thickness

Properties of powder and capsule materials

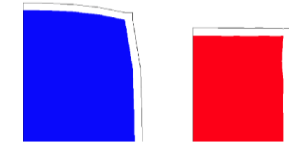
Kuhn-Donwey and Abouaf-Chenot Model

Temperature and pressure profile



Capsule shape optimization

Optimized capsule



Before and after HIP



## **Acknowledgement**

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Thank you very much for your attention!

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