

# Prediction of microstructure for AISI316L steel from numerical simulation of laser powder bed fusion

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## Introduction

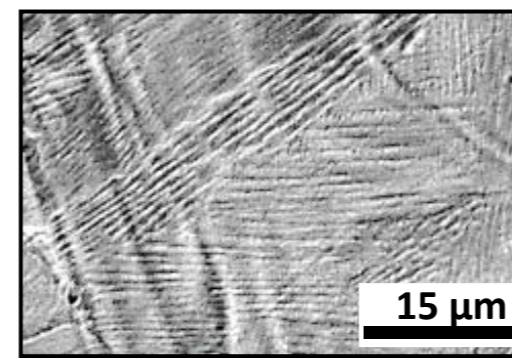
**Laser powder bed fusion (L-PBF)** is the most promising additive manufacturing technology for metals.

Many **numerical simulation** software provide solidification data useful for **microstructure prediction**, therefore they can represent a powerful tool for L-PBF improvement.

The **cooling rate** ( $\dot{T}$ ) of the process is the key parameter determining the **microstructure** of the final component, directly responsible for the mechanical properties.

**Austenitic stainless steel 316L** commonly exhibits a **cellular microstructure** when produced by L-PBF.

**Primary cell arm spacing (PCAS)** is the characteristic feature of this microstructure, on which the mechanical properties such as **microhardness (H)** depend.



✓ **Objective:**

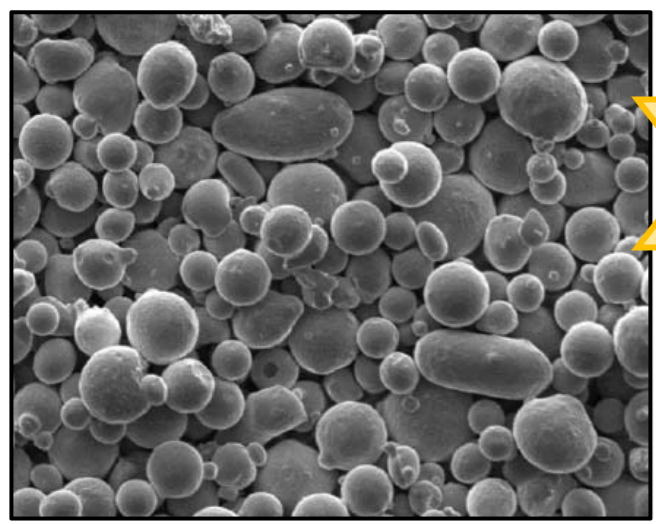
**Estimate the PCAS and the microhardness** of 4 single scan tracks of 316L stainless steel by using **FLOW-3D AM**, a commercial CFD software, and **validate** the model



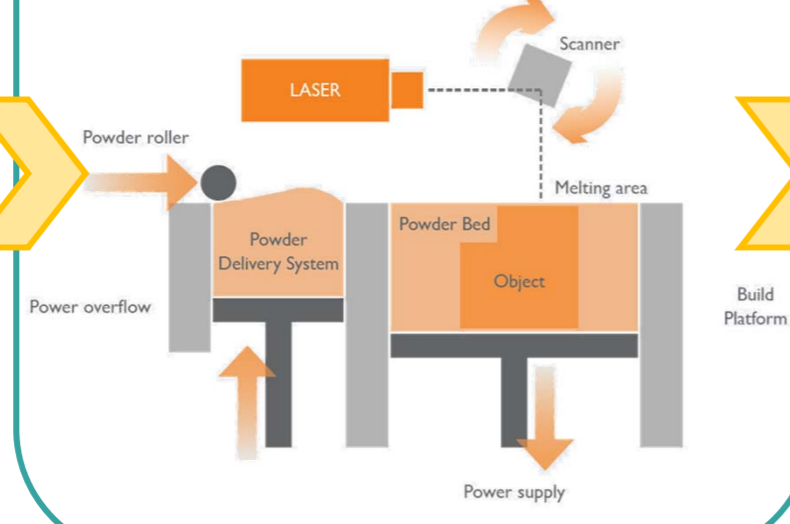
## Materials and Methods

### Powders

Gas atomized 316L stainless steel



### L-PBF process



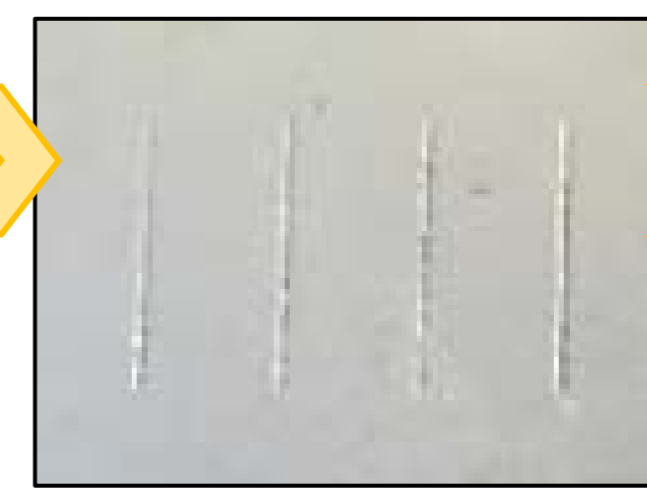
### Process parameters

Laser Power = 140 W Laser beam spot size = 36  $\mu$ m

Track	Scanning speed	Layer thickness
A	50 cm/s	60 $\mu$ m
B	100 cm/s	
C	50 cm/s	90 $\mu$ m
D	100 cm/s	

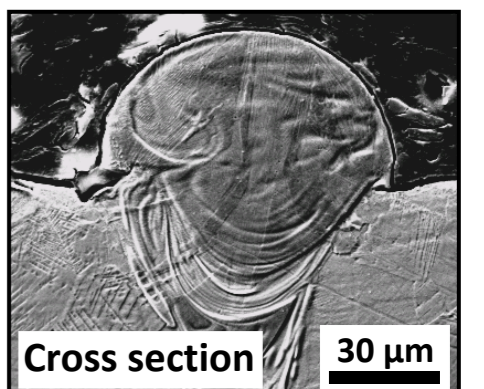
### Real samples

4 single scan tracks of 10 mm



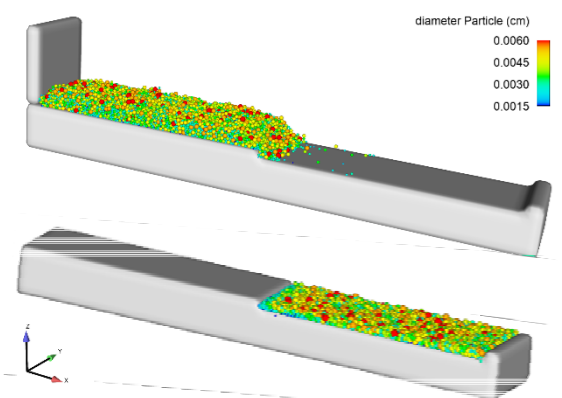
### Characterization

**PCAS and microhardness** measurements on the cross section

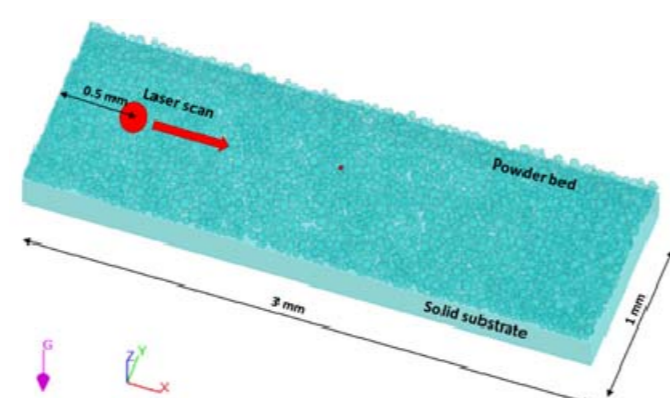


### Model

**Discrete Element Method (DEM)** of **FLOW-3D AM** was used to **create the powder bed**

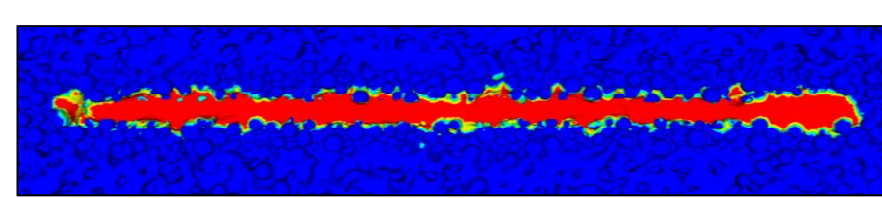


**Finite Volume Method (FVM)** of **FLOW-3D AM** was used to reproduce powder melting and subsequent solidification



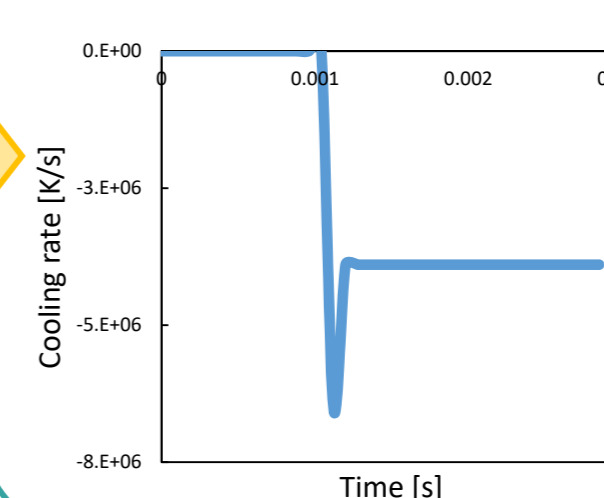
### Simulated samples

A preliminary phase of calibration for the **absorption coefficient** allowed the obtaining of the **first 2 mm** of the track



### Characterization

**Cooling rate data**  $\dot{T}$  evaluation



**Estimation of PCAS and H** through the **empirical formula**:

$$PCAS = 80 \dot{T}^{-0.33}$$

$$H = 152 + 498d^{-0.5}$$

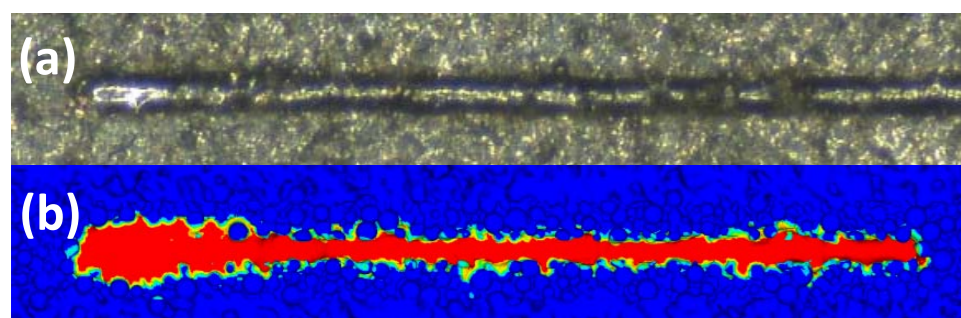
where d is the grain width, calculated in turn from  $\dot{T}$

## Results and conclusions

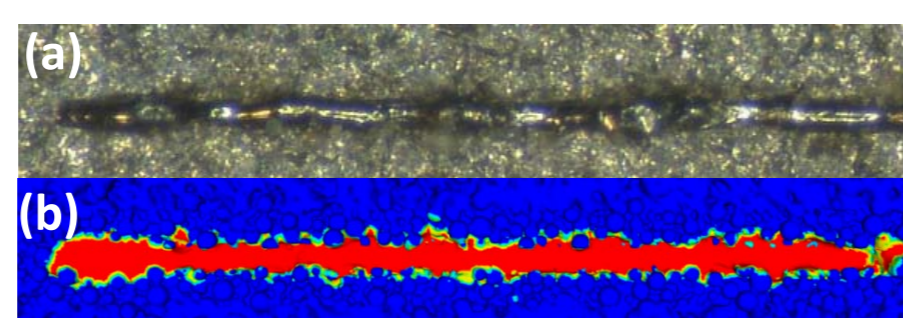
### Model calibration

The confrontation between the top view of the track shows **good agreement** between results, as the morphology of the real track (a) is well predicted by simulations (b). This denotes a correct calibration of the absorption coefficient, set to **0.6** for these combinations of process parameters.

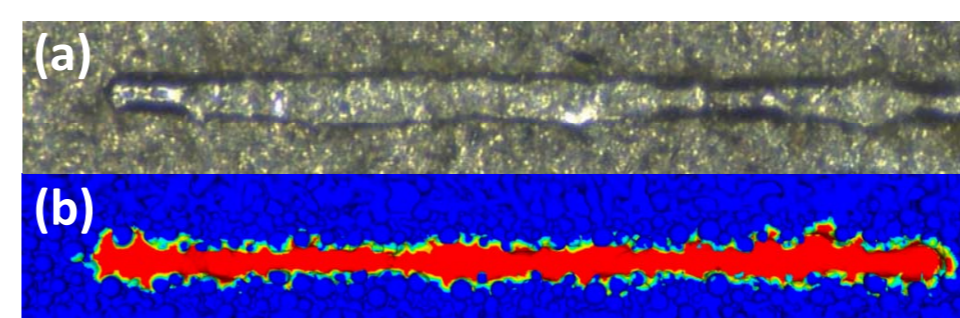
Track A



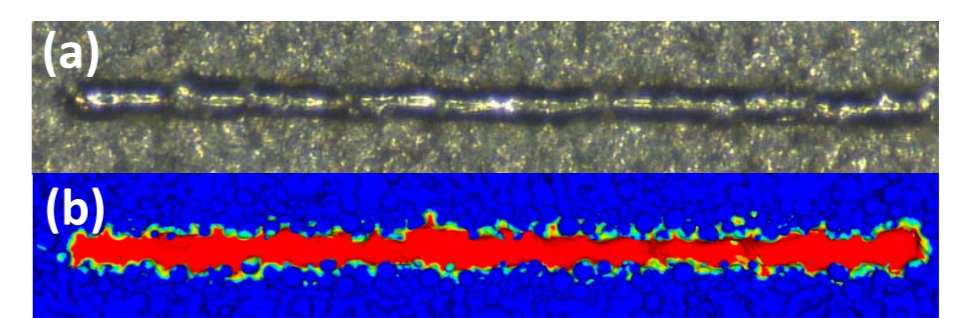
Track B



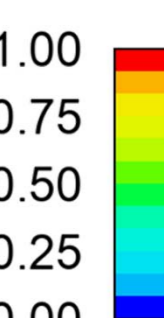
Track C



Track D



Melt region



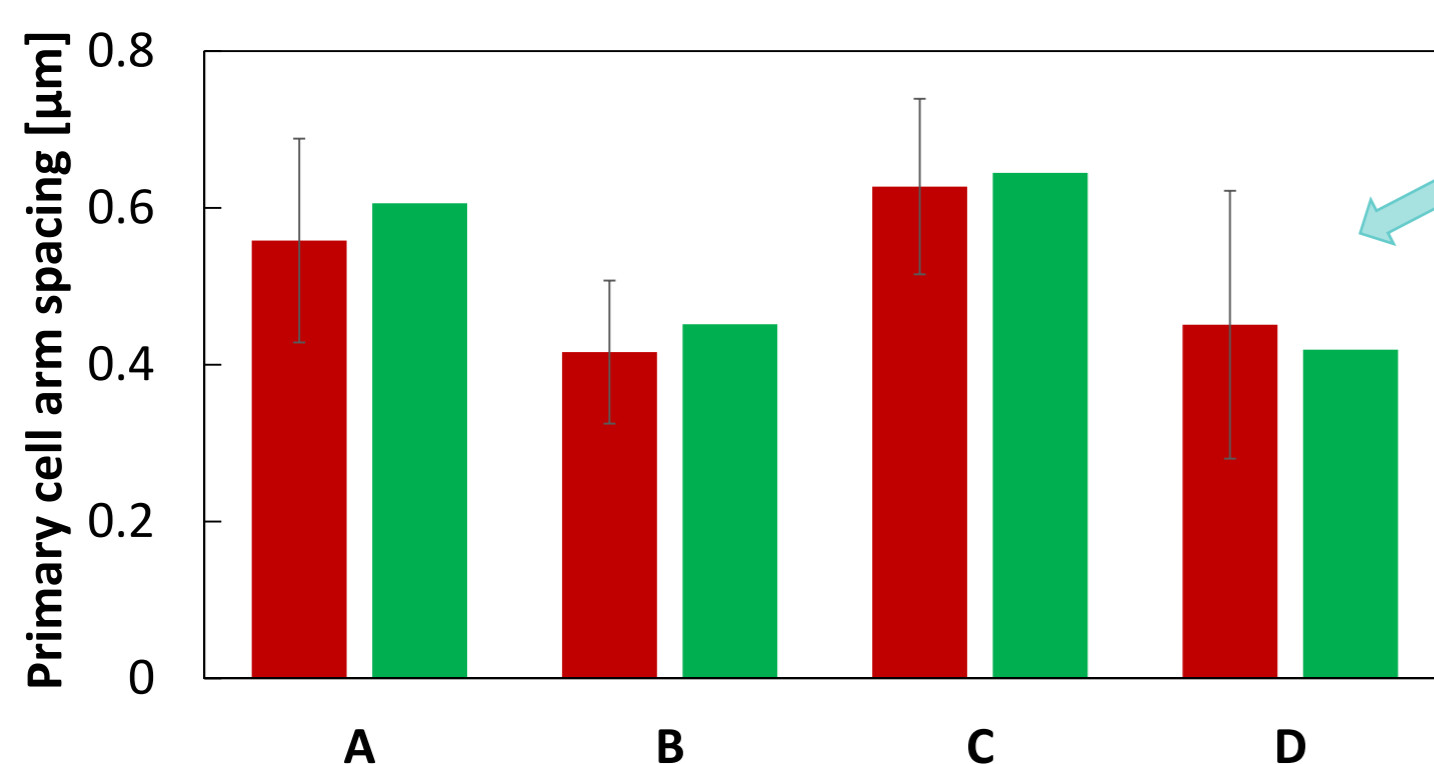
0.5 mm

Good morphological agreement

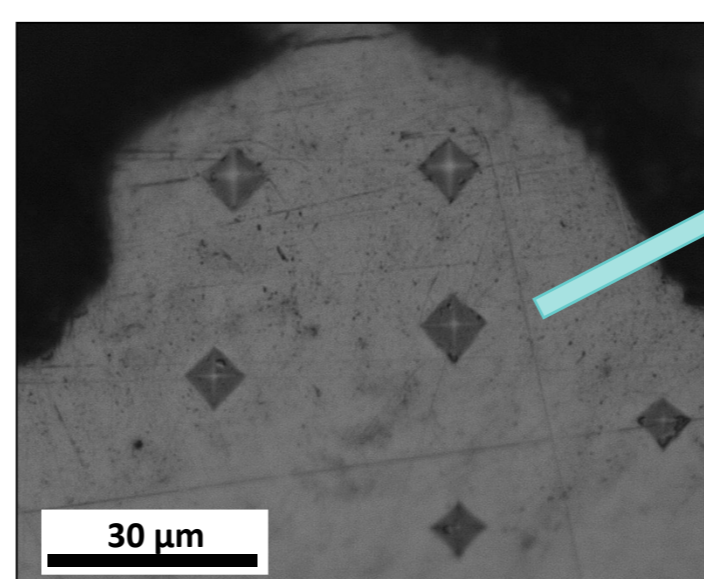
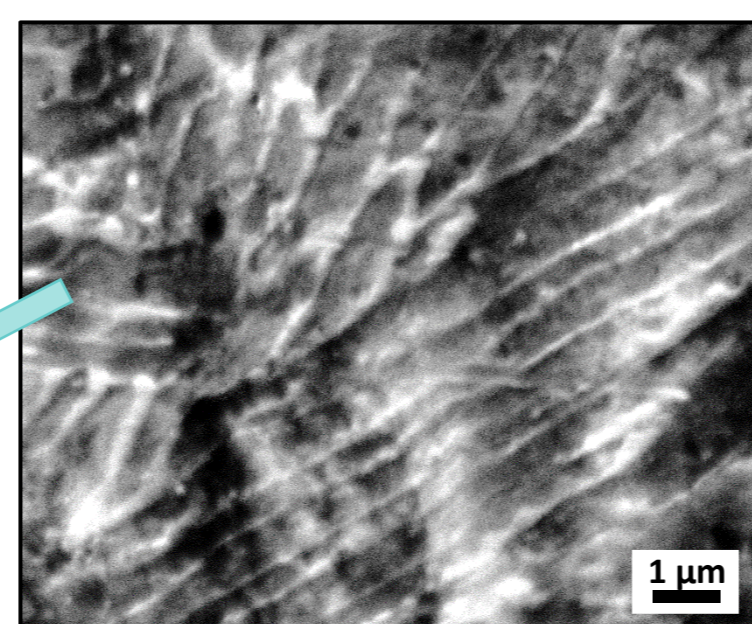
### Prediction of Primary Cell Arm Spacing

The simulated results show **high agreement** with the experimental data. The deviation between real and simulated results varies **from 2% to 8%**.

■ Experimental result ■ Simulation result



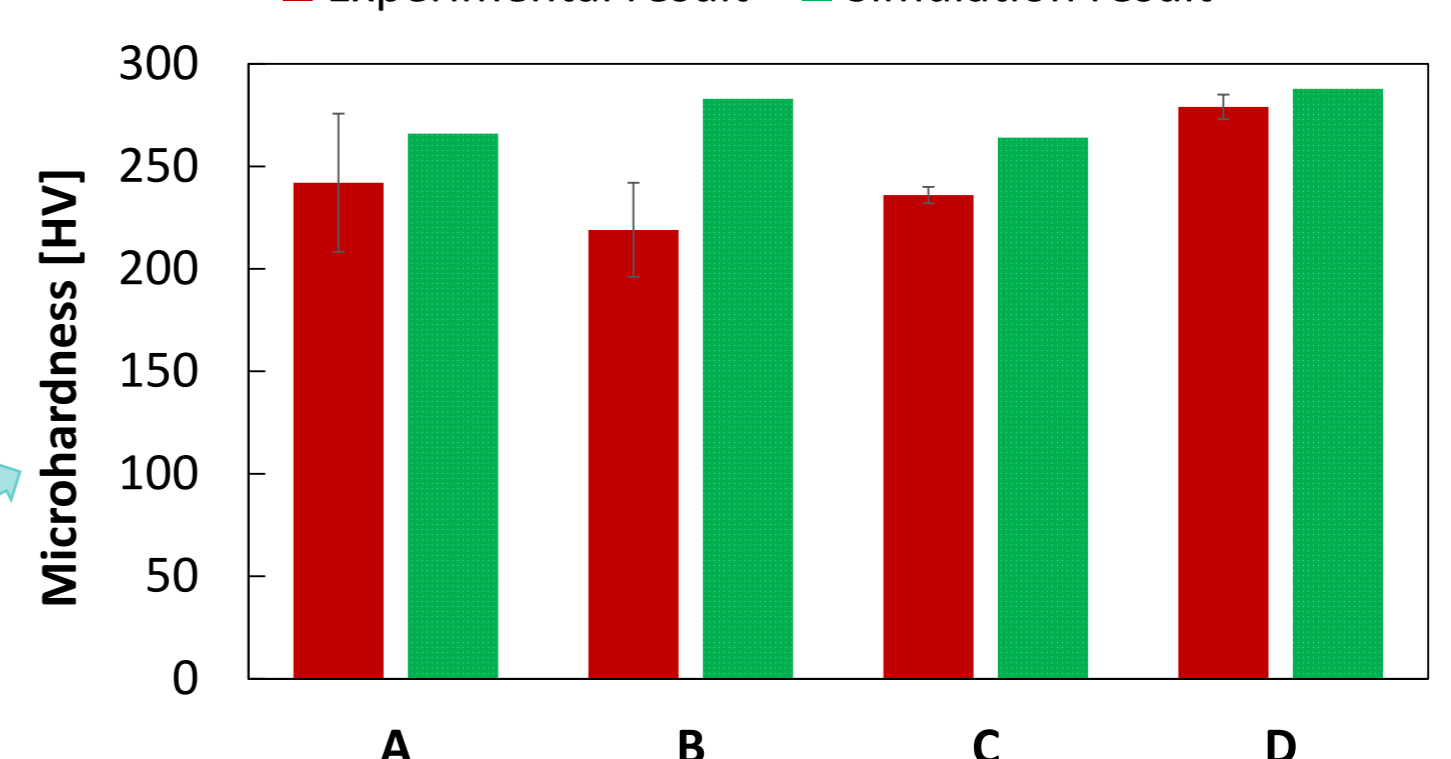
Confirmation of the model validity and its potential in obtaining reliable results for microstructure prediction



### Prediction of Microhardness

The predicted values **overestimate** the microhardness of the samples. The deviation between real and simulated results varies **from 3% to 22%**.

■ Experimental result ■ Simulation result



Overestimation of the model with respect to the real microhardness of the tracks

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