

Chemical gradients across martensite/austenite phase boundaries in precipitation hardened TRIP steel studied by atom probe tomography

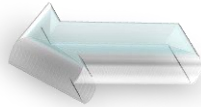
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- **Motivation and alloy design**
- **Atom probe tomography**
- **Mechanical properties and microstructure**
- **Mn partitioning and simulations**
- **Conclusions**



- **TRIP:** deformation-stimulated transformation of unstable austenite into martensite and accommodation plasticity (e.g. Mn, Ni, low C)
- **Maraging effect:** hardening of heavily strained martensite via nano-sized (intermetallic) precipitates (Ni, Al, Ti, Mo)

(see also conventional Maraging steels)

What is **maraging-TRIP** ?

Quenched austenite: ductile low carbon martensite

Retained austenite (TRIP or TWIP)

Aging: austenite reversion (increase TRIP / TWIP volume)

Aging: controlled precipitation hardening

* *TRIP: transformation-induced plasticity*

* *Maraging: martensite aging*

Steel	C	Ni	Co	Mo	Ti	Al	Mn	Fe
Maraging	0.01	18	12	4	1.6	0.15	0.05	Balance
09MnPH	0.01	2	-	1	1.0	0.15	9	Balance
12MnPH	0.01	2	-	1	1.0	0.15	12	Balance
15MnPH	0.01	2	-	1	1.0	0.15	15	Balance



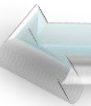
Low carbon: ductile martensite

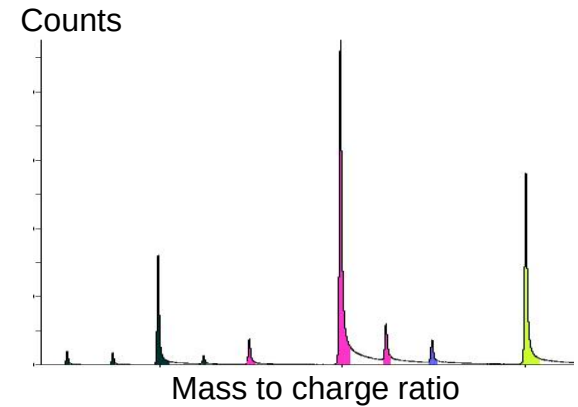
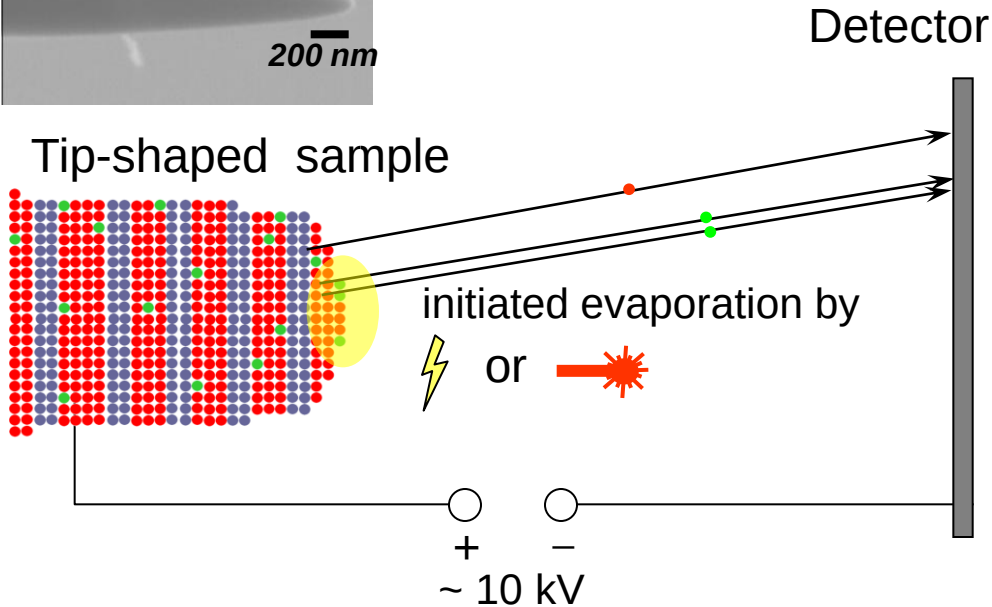
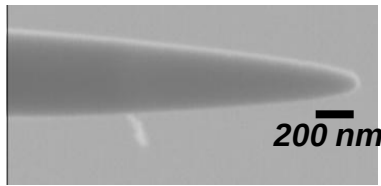
Precipitation Hardenable

Mn (+Ni): austenite (TRIP)

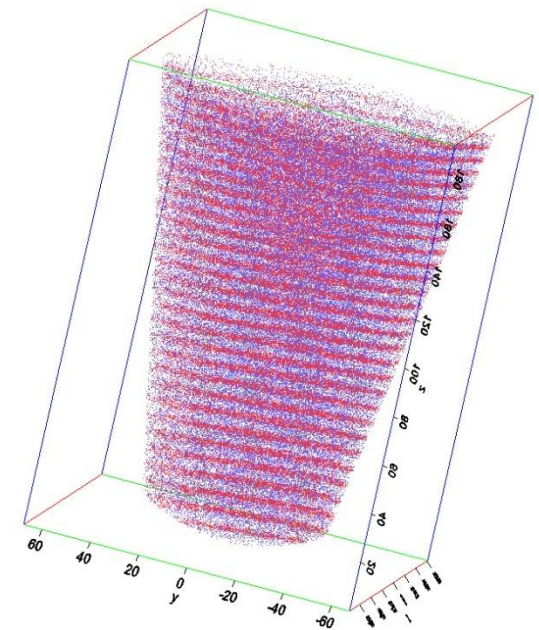
Martensite aging after quenching at 450°C

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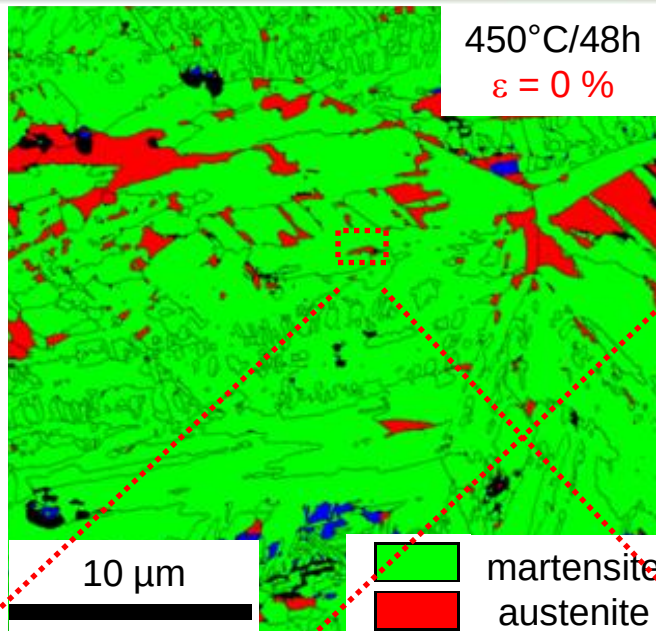
3D atom map



APT: Time of flight measurement
(chemical identification)
+
Ion projection microscopy
(determination of position)

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Hierarchy of maraging TRIP / TWIP steels



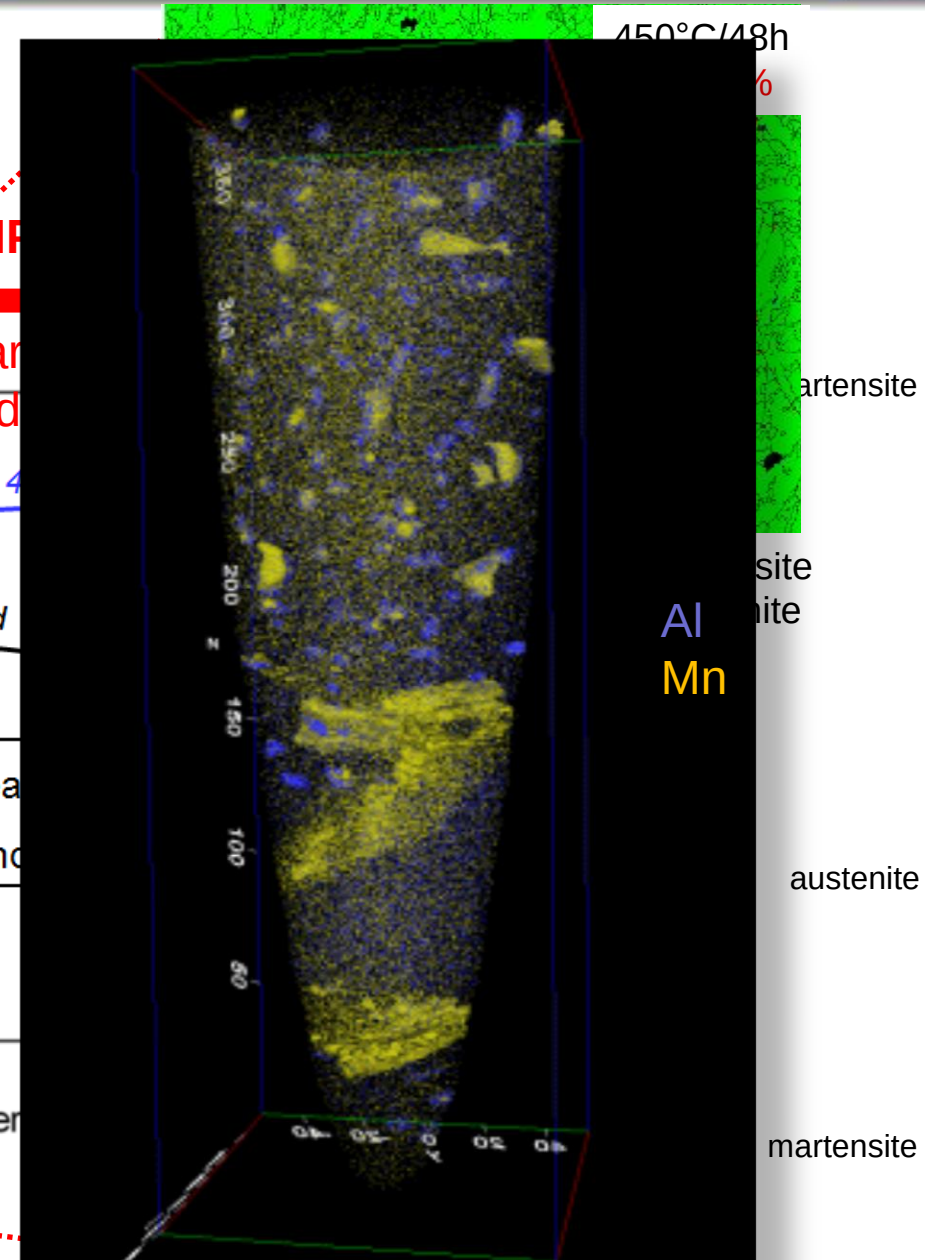
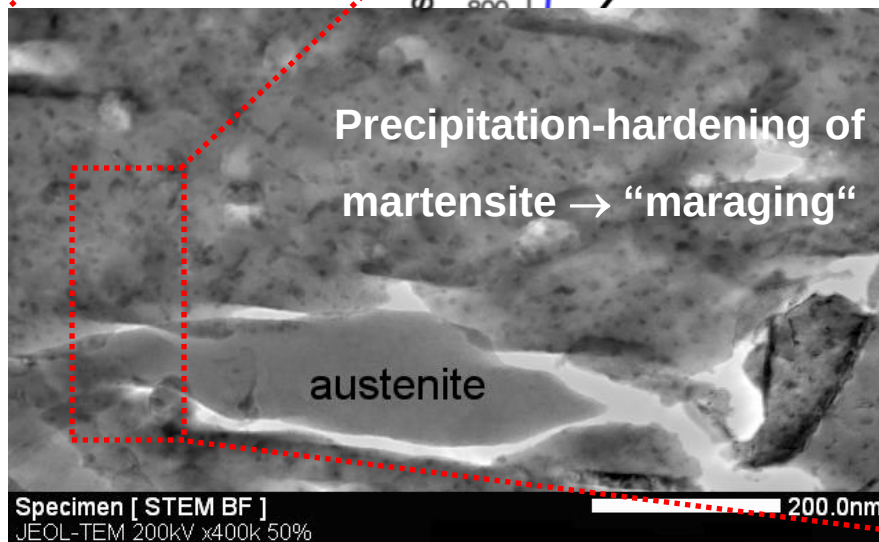
TRIP

hard

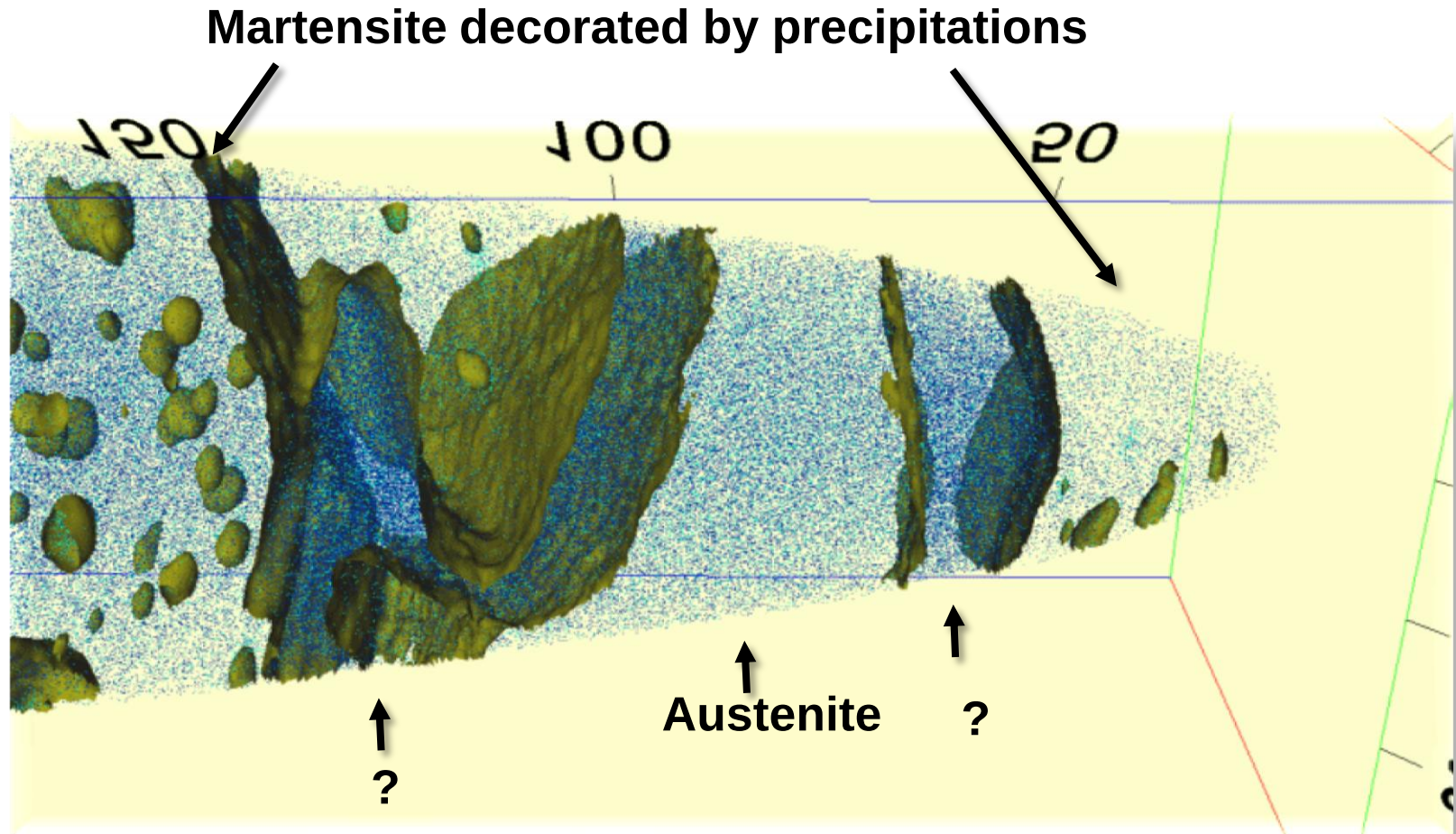
+ d

aged 4

enched



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Mn atoms

Ni atoms

Mn iso-concentration: 18 at.%

C	Ni	Mo	Ti	Al	Mn	Fe
0.01	2.0	1.0	1.0	0.15	12	bal.

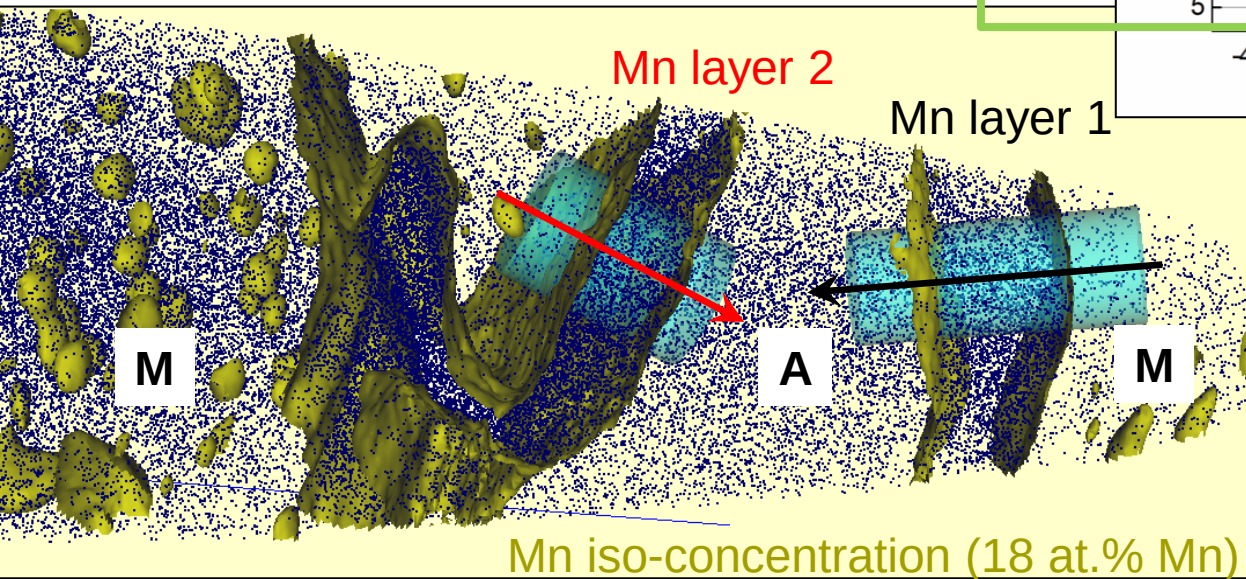
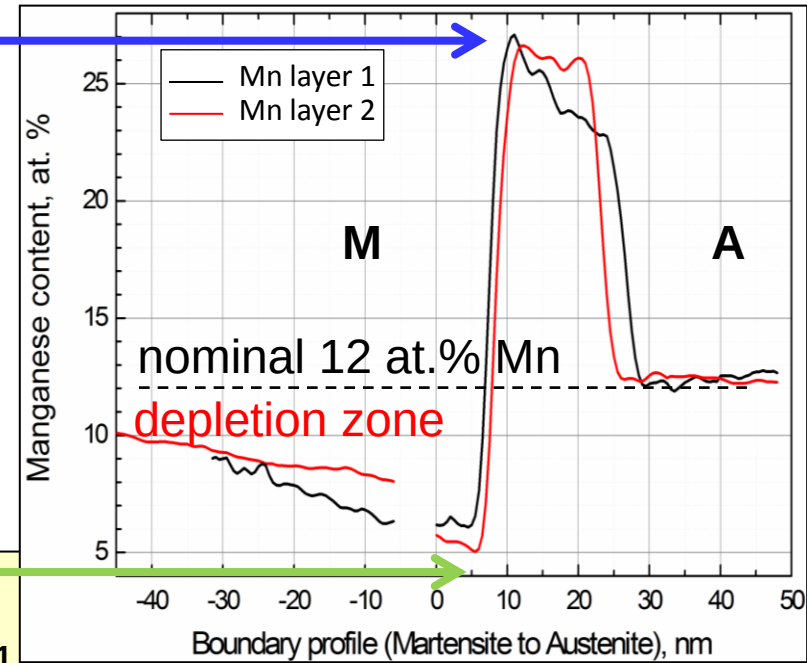
70 million ions
Laser mode
(0.4nJ, 54K)

Thermo-Calc $\Rightarrow$

equilibrium Mn-conc.:

27 at. % Mn in austenite (A)

3 at. % Mn in ferrite (martensite) (M)



precipitates in α'

$$x_{Diff} \cong 2\sqrt{Dt} \cong 30nm$$

no precipitates in
austenite

$$x_{Diff} \cong 2nm$$

men [STEM BF]
STEM 200kV x400k 50%

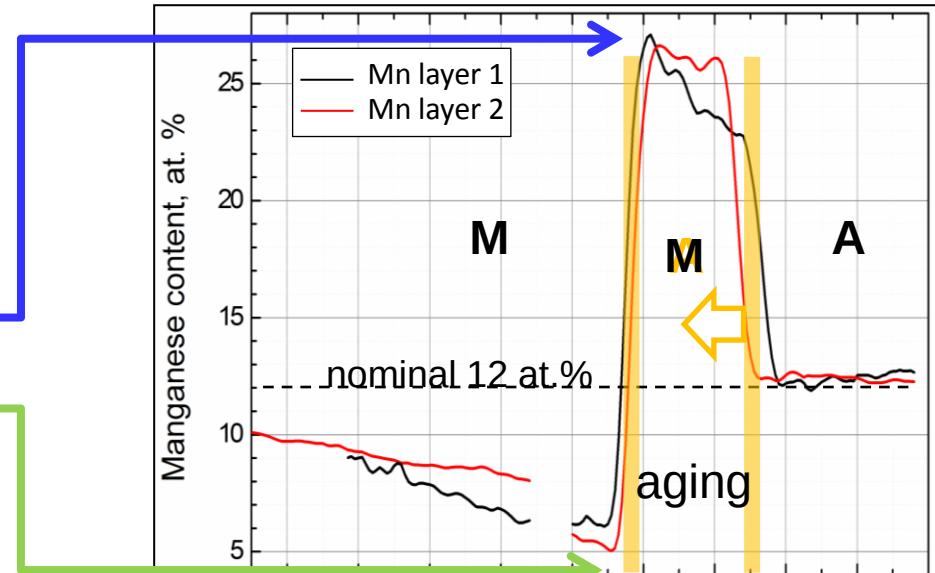
200.0nm

Thermo-Calc $\Rightarrow$

equilibrium Mn-conc.:

27 at. % Mn in austenite (A)

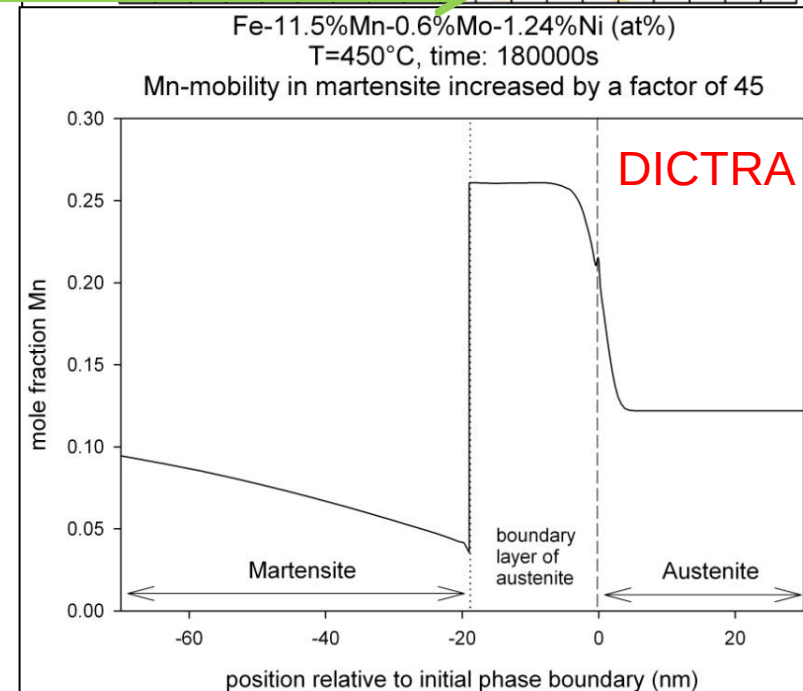
3 at. % Mn in ferrite (martensite) (M)

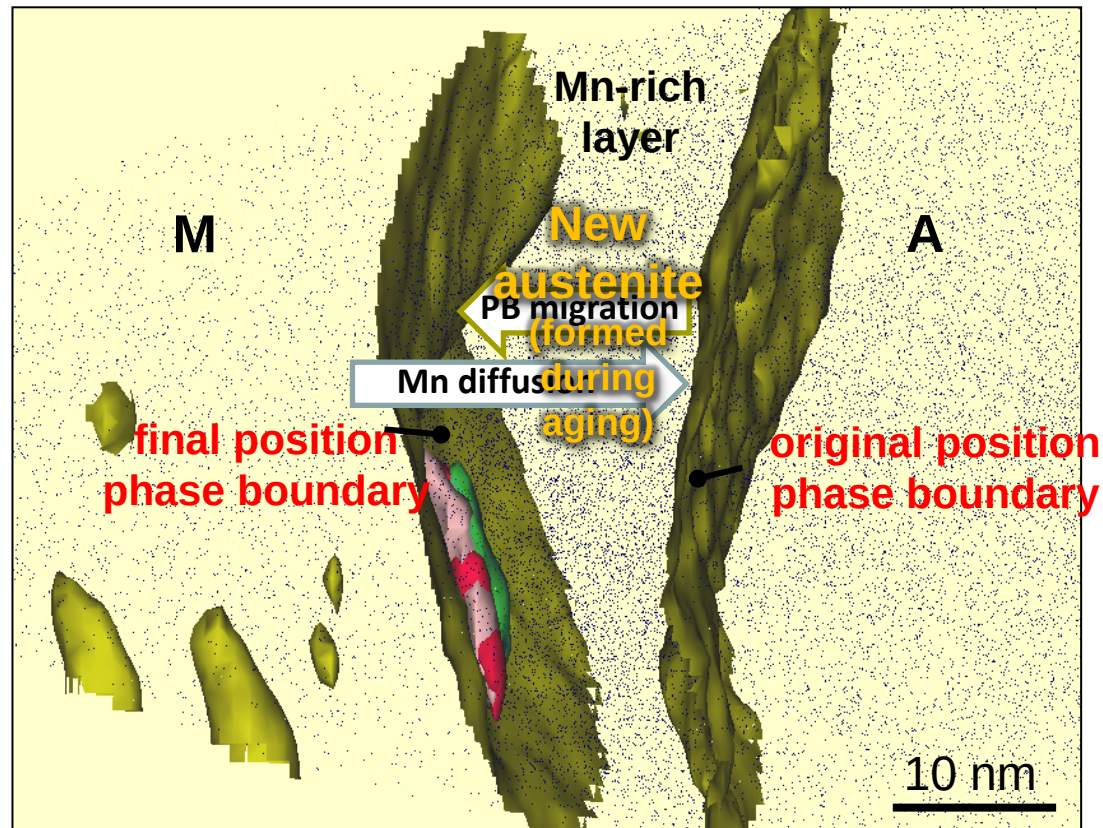


**Excellent agreement between
experiment & simulation !**



**Kinetic freezing and
associated austenite reversion !**





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- Maraging-TRIP as a new GPa steel design approach
- Unexpected simultaneous increase in strength and elongation
- Mn partitioning, predicted by ThermoCalc/DICTRA
- Austenite stability predicted using ab initio methods (see talk of T. Hickel)
- B2 and L21 nano-precipitates
- Next steps: lean composition, interstitials, alloy variants, equilibrium partitioning, nano-precipitates

