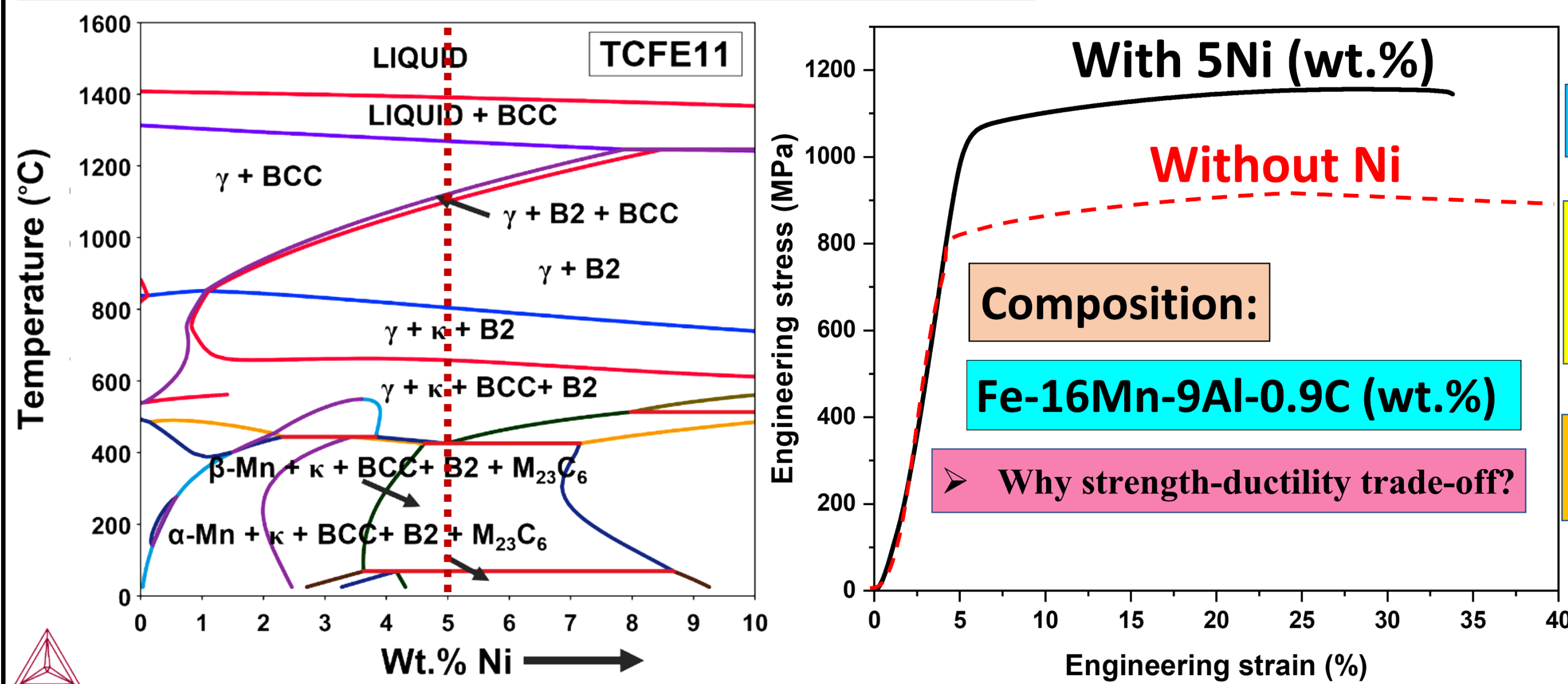


## Introduction and objectives

### Design of Ni-alloyed Fe-Mn-Al-C low-density steel



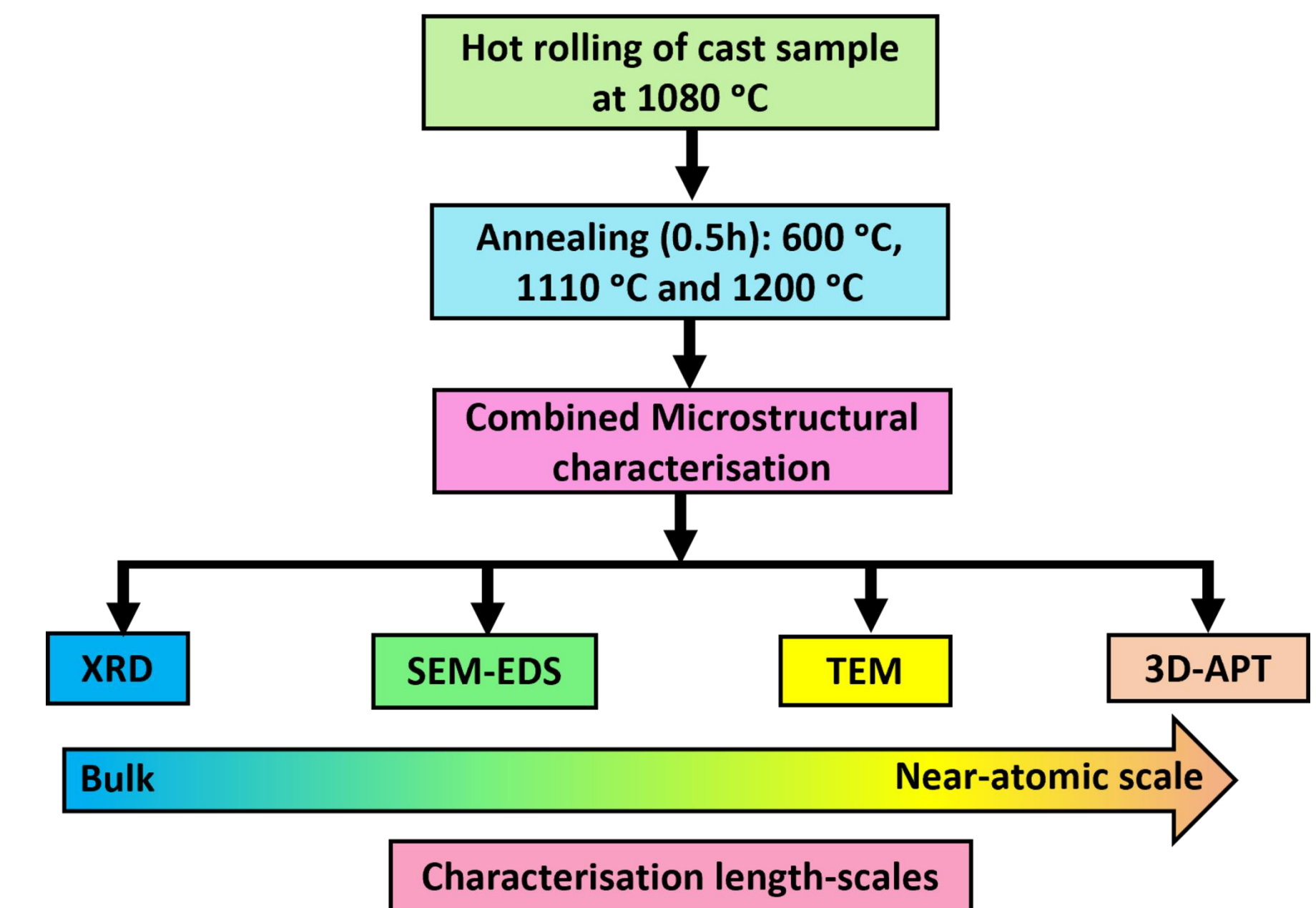
To investigate:

- Ni-induced B2 evolution sequence vs annealing temperature

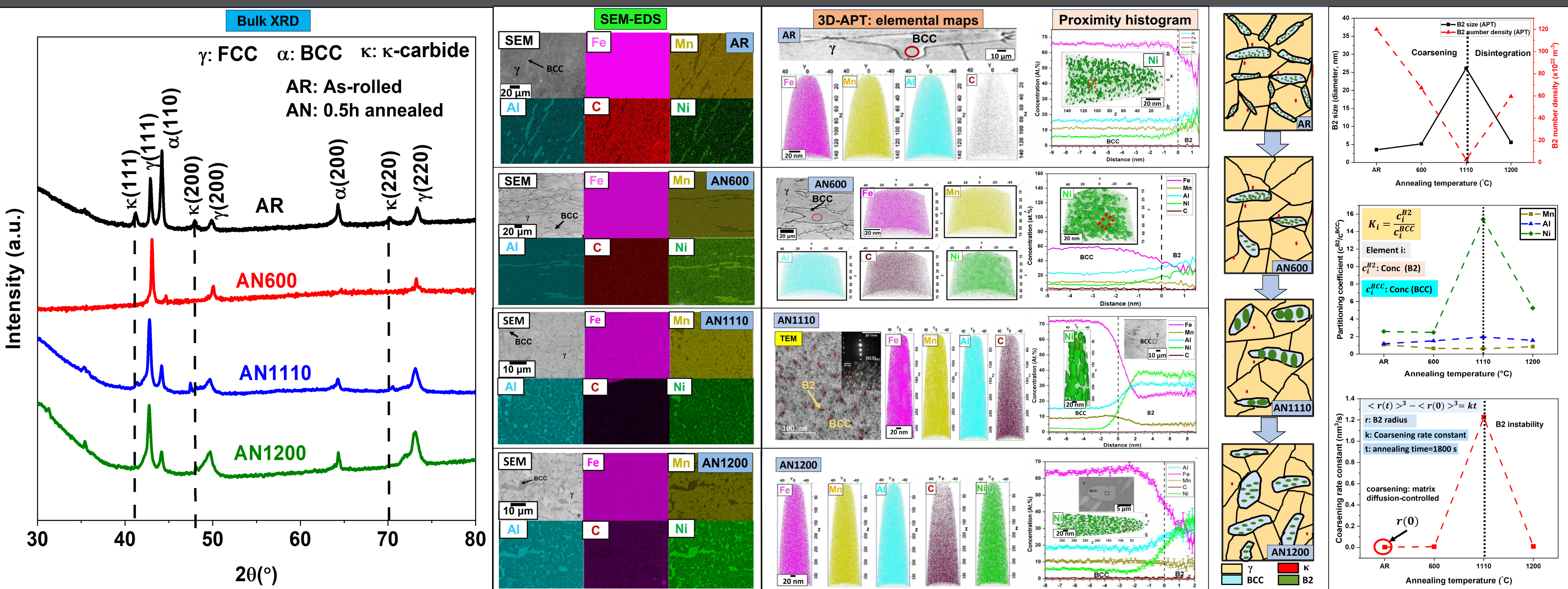
- B2 stability and composition

## Methodology – Combined microscopy

### Fe-16Mn-9Al-0.9C-5Ni (wt.%)



## Results



## Conclusions

1. Near-atomic scale analysis using 3D-APT reveals the localized precipitation of nano-scale B2 precipitates within BCC phase.
2. B2 coarsening kinetics follows modified LSW model and hence is bulk diffusion controlled.
3. Partitioning of Ni between BCC and B2 phases controls the rate of B2 coarsening.
4. B2 disintegration at 1200 °C indicates its limited stability at elevated temperatures, as per thermodynamic estimates.
5. B2 localization within BCC phase in  $\gamma$ -based Ni-alloyed Fe-Mn-Al-C steel leads to enhanced strength-ductility trade-off.

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