

The model of austenite decomposition kinetics based on CCT curves and its application in 22MnB5 steel

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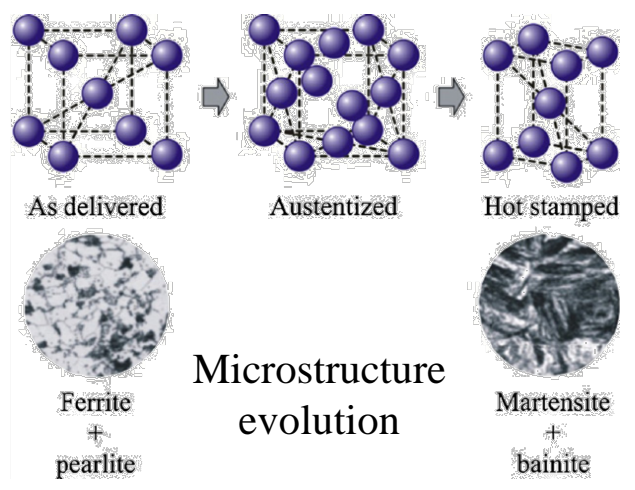
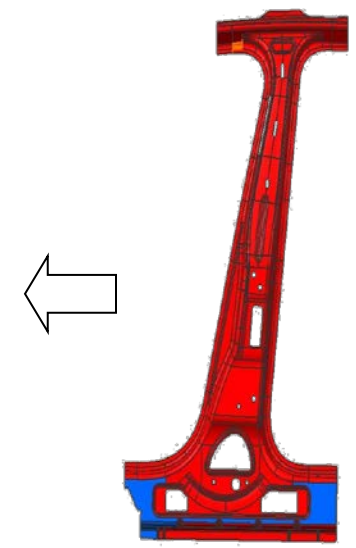
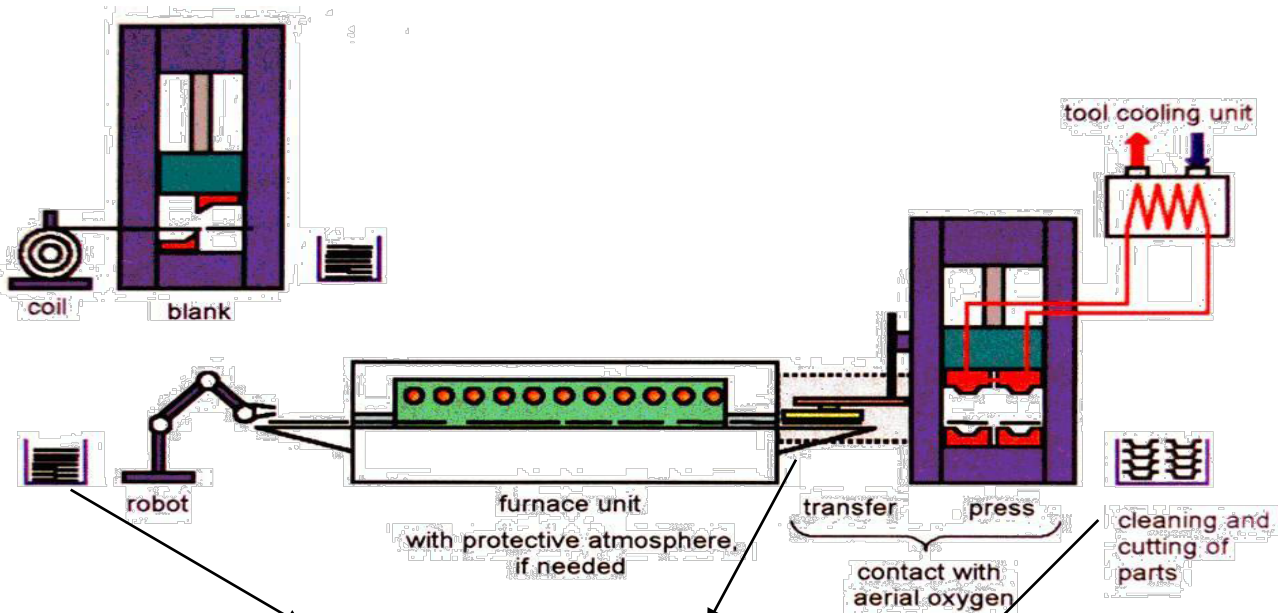
- Introduction to hot stamping**
- Thermal-mechanical-metallurgical coupled FEM model**
- Theoretical analysis of diffusional austenite decomposition**

The mathematical relationship between isothermal and non-isothermal transformation

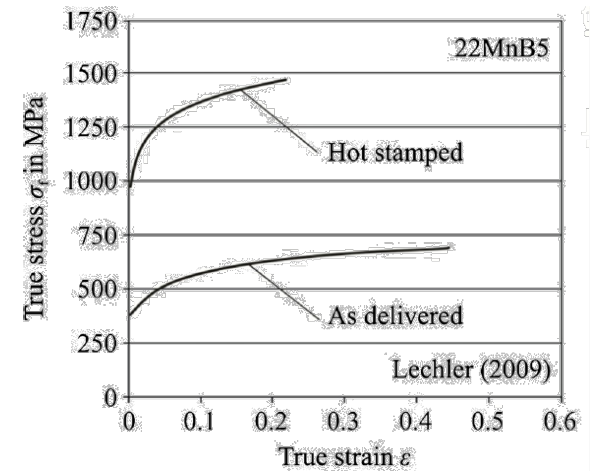
Generalized additive rule

- Summary**

Introduction to Hot Stamping



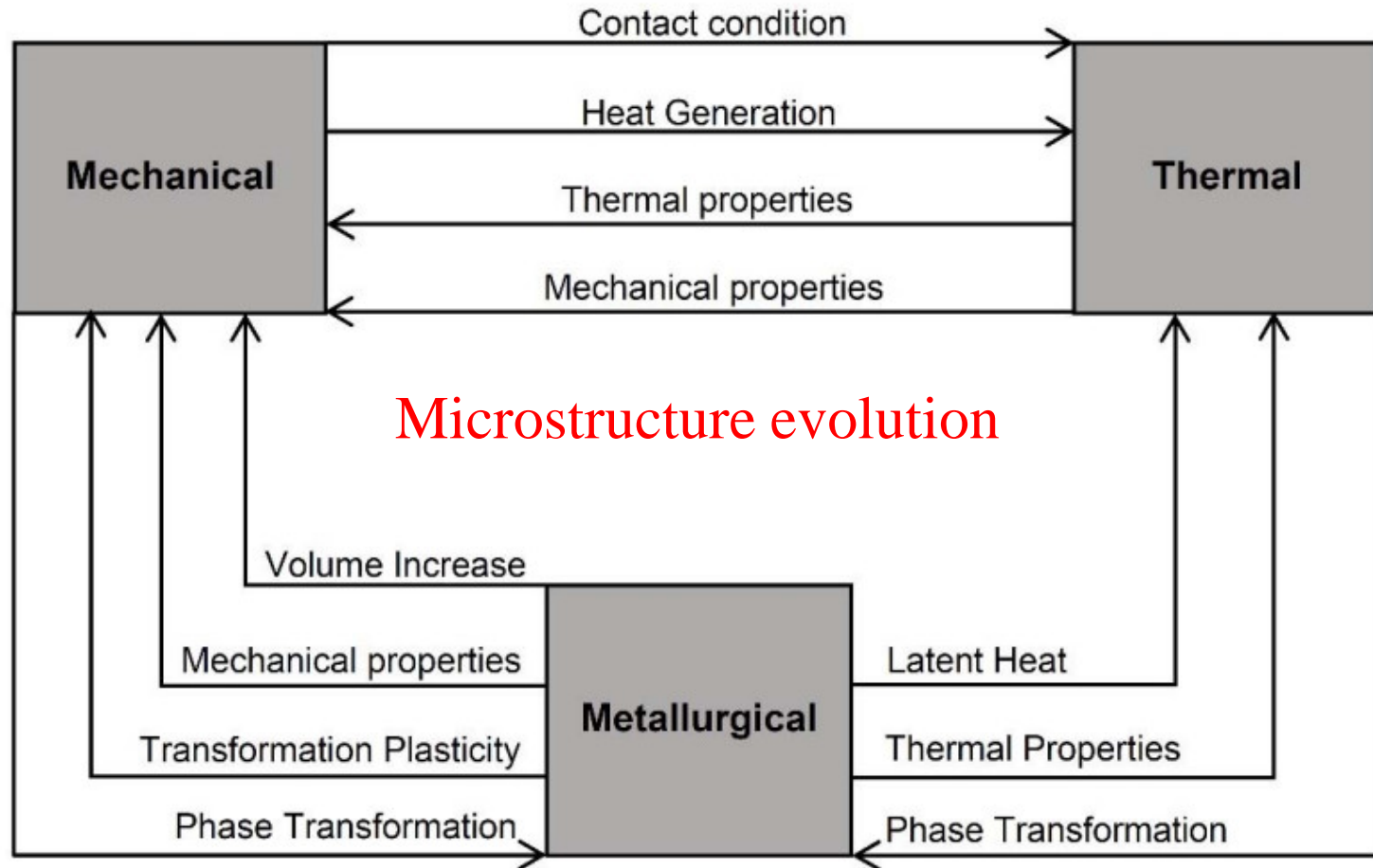
Mechanical properties



Thermal-mechanical-metallurgical coupled FEM model



Thermal-mechanical-metallurgical coupled FEM model



Thermal-mechanical-metallurgical coupled model during hot stamping

Austenite decomposition kinetics

1, Diffusional phase transformation

JMAK type model (1) $F = 1 - \exp(-b \times t^n)$

K-V type model (2)
$$\frac{\partial F_i}{\partial t} = \frac{f(T)f(F_i)}{f(G)f(C)}$$

{

- T*---Temperature
- F*---Current fraction formed
- G*---Austenite grain size
- C*---Alloy composition

Scheil additivity hypothesis (**Calculating non-isothermal transformation from isothermal kinetics**)

$$\sum_i^m \frac{\Delta t_i}{\tau_i} = 1 \qquad t_j = \Delta t_j + \left[-\frac{\ln(1 - F^{j-1})}{b} \right]^{\frac{1}{n}}$$

2, Diffusionless phase transformation

Koistinen-Marburger equation

(1) M. Avrami., et al. The Journal of Chemical Physics.1939 (7),1103.

(2) J.S. Kirkaldy, et al. International Conference on Phase Transformations in Ferrous Alloys.1983, 125-148.

Thermal-mechanical-metallurgical coupled FEM model

Decomposition of austenite

isothermal ferrite, pearlite and bainitic transformation

$$X = 1 - \exp\left[-b \left(\frac{D\gamma_{TTT}}{D\gamma}\right)^m (t - \tau_o)^n\right]$$

Martensitic transformation

$$F_m = F_a \times \{1 - \exp[-0.011 \times (M_s - T)]\}$$

Heat transfer model

$$\rho c_p \frac{\partial T}{\partial t} = \lambda \nabla T + Q$$

$$Q^{th} = \Delta H_i \frac{\Delta F_i}{\Delta t}$$

Mechanical response

$$d\varepsilon_{ij} = d\varepsilon_{ij}^e + d\varepsilon_{ij}^p + d\varepsilon_{ij}^{th} + d\varepsilon_{ij}^{tr} + d\varepsilon_{ij}^{tp}$$

$$\begin{cases} d\varepsilon_{ij}^{th} \rightarrow \text{Thermal strain increment} \\ d\varepsilon_{ij}^{tr} \rightarrow \text{Transformation strain increment} \\ d\varepsilon_{ij}^{tp} \rightarrow \text{Transformation induced plasticity} \end{cases}$$

$$d\varepsilon_{ij}^{th} = \alpha_{ij} dT + \frac{\partial \alpha_{ij}}{\partial T} \delta_{ij} T dT + \sum_{l=0}^N \frac{\partial \alpha_{ij}}{\partial \xi_l} \delta_{ij} T d\xi_l \quad \text{Volume change due to the thermal process}$$

$$d\varepsilon_{ij}^{tr} = - \sum_{l=0}^N \left(\frac{\partial \beta_{ijkl}^{-1}}{\partial \xi_l} \sigma_{kl} + \beta_l \delta_{ij} + \frac{\partial \beta_{ij}}{\partial \xi_l} \delta_{ij} T \right) d\xi_l \quad \text{Volume change due to phase transformation}$$

$$d\varepsilon_{ij}^{tp} = - \sum_{l=0}^N 3K_l s_{ij} (1 - \xi_l) d\xi_l \quad \text{Local plastic flow below the yield stress, Greenwood-Johnson model}$$

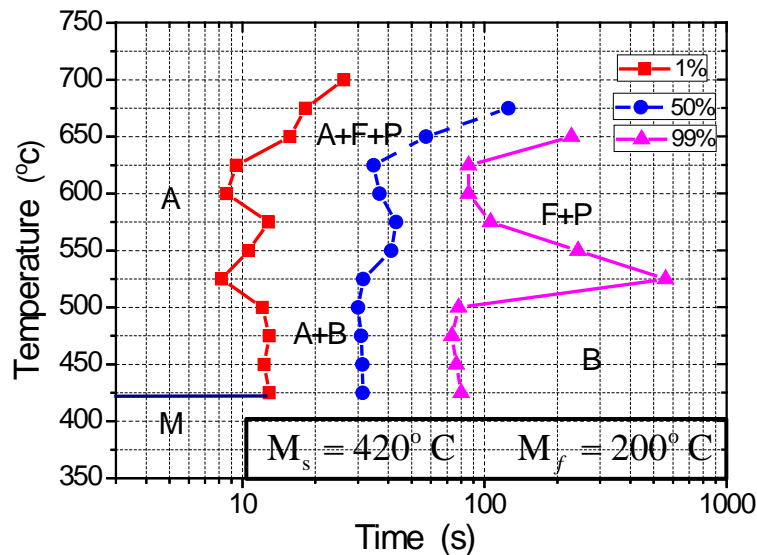
Thermal-mechanical-metallurgical coupled FEM model



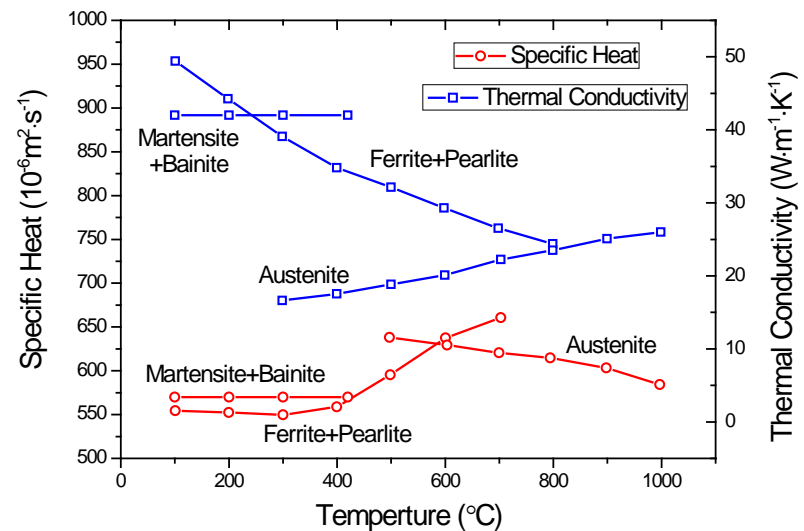
Calculation parameters

Chemical composition for 22MnB5 in weight percent

	C (%)	Si (%)	Mn (%)	P (%)	S (%)	Cr (%)	Ni(%)	B(ppm)
Measured	0.27	0.29	1.25	0.007	<0.005	0.22	0.013	39



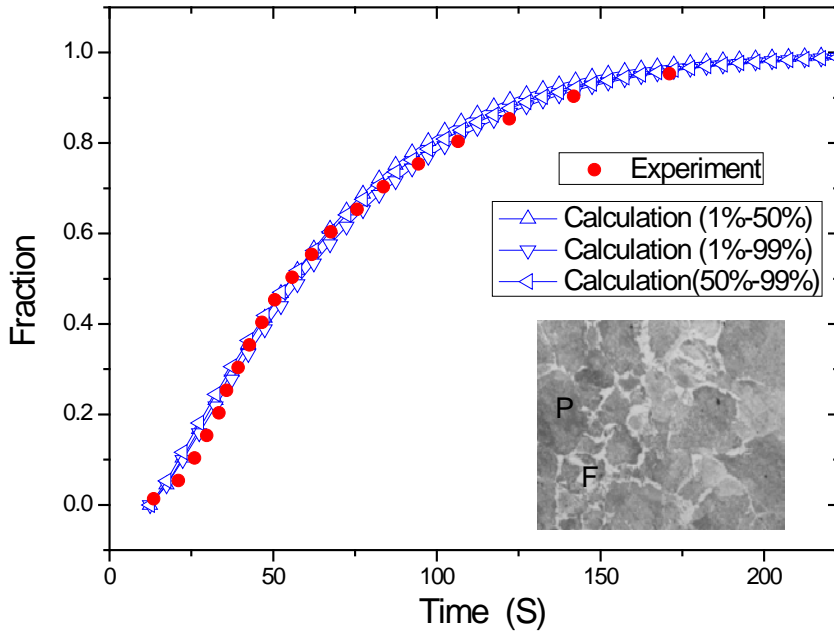
The TTT diagram for the 22MnB5



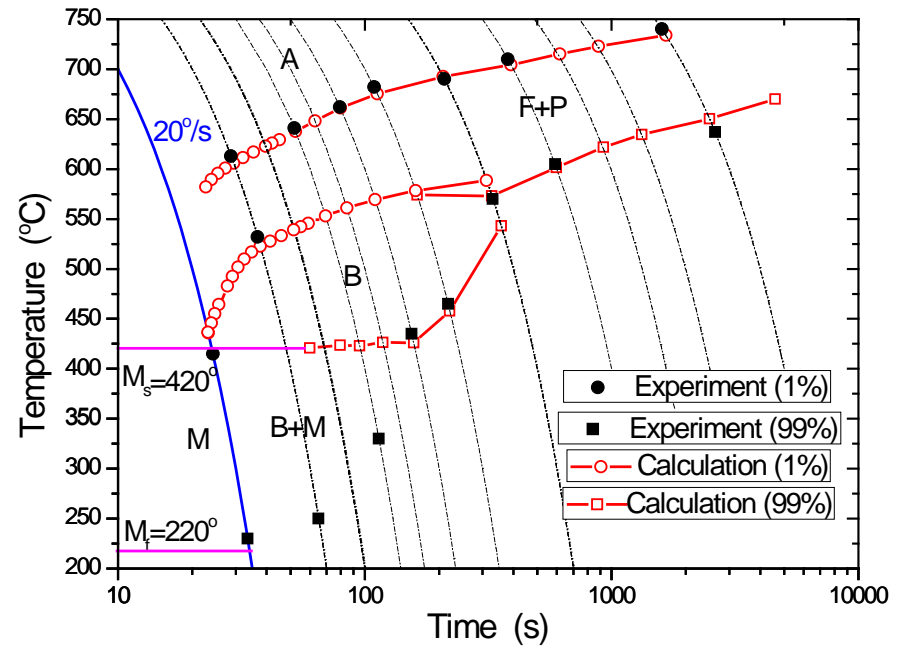
The Thermal conductivity and specific heat for the 22MnB5

Thermal conductivity, specific heat and heat transfer coefficient are experimentally measured as the function of temperature.

Thermal-mechanical-metallurgical coupled FEM model



The phase transformation kinetics considering incubation time in JMAK equation and the corresponding microstructure metallograph (F: ferrite, P: pearlite)

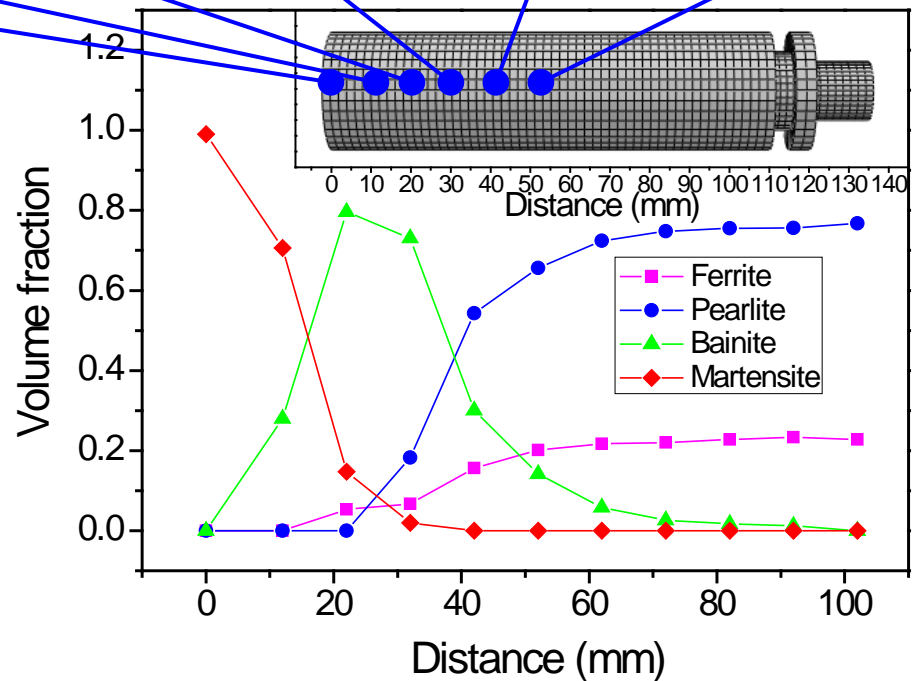
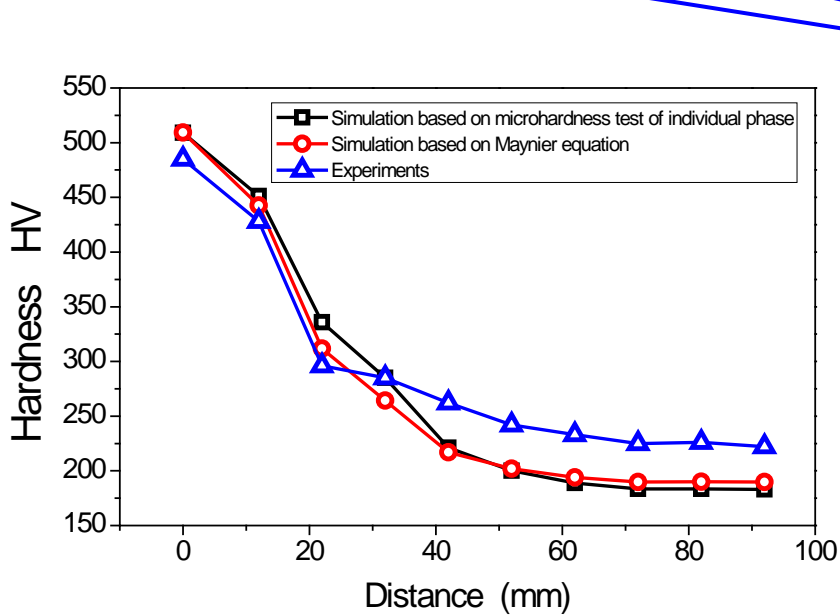
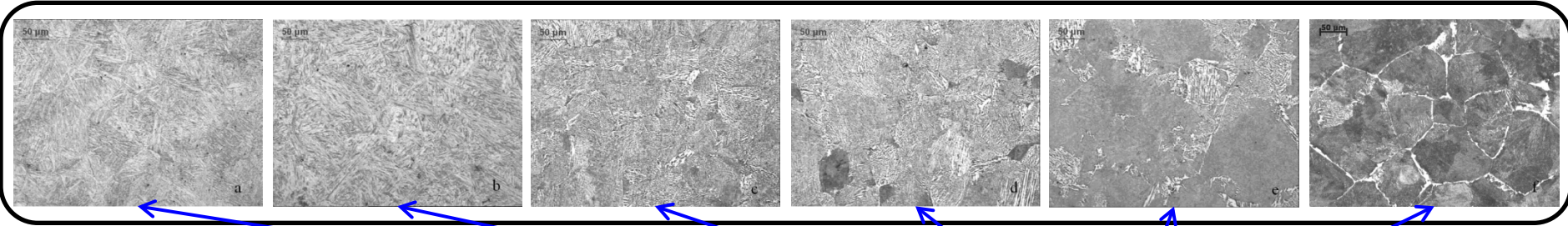


The CCT diagram prediction and comparison with dilatational experiments

Thermal-mechanical-metallurgical coupled FEM model



The Jominy end-quenching test and the microstructure prediction



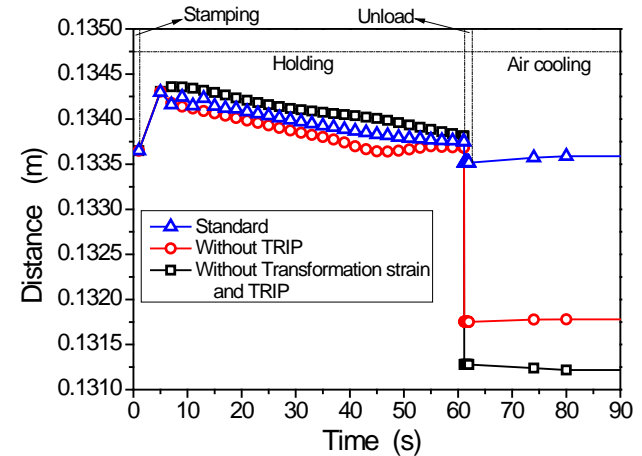
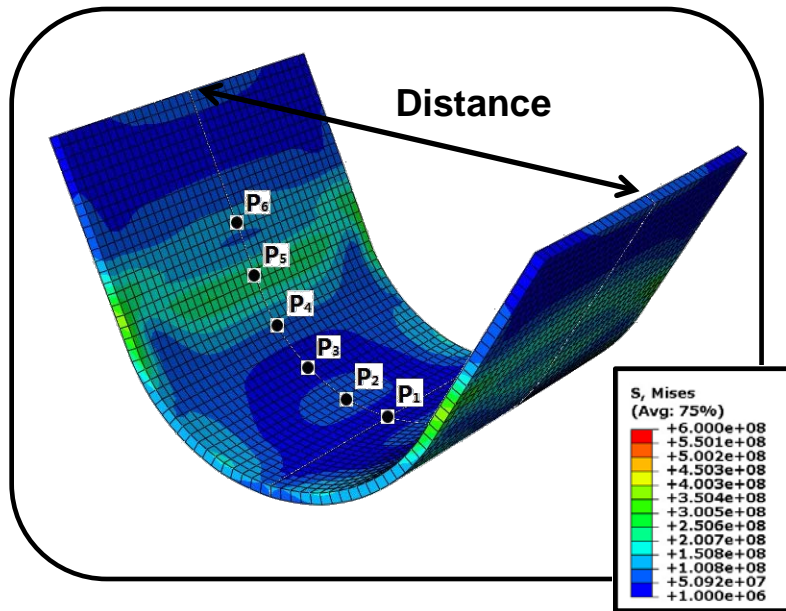
Measured and calculated hardness as function of distance from the quenched end

The simulated final microstructure components and the corresponding fraction along the Jominy bar

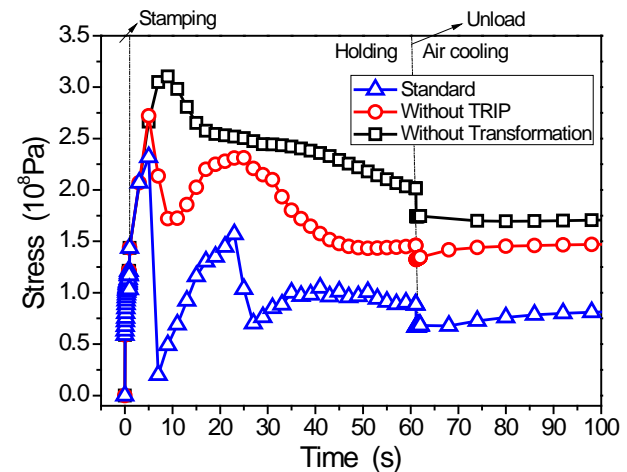
Thermal-mechanical-metallurgical coupled FEM model



Unloading spring-back and Residual stress



The effect of transformation strain and transformation plasticity on the opening distance



The effect of transformation strain and transformation plasticity on the residual stress

The expansion due to martensitic transformation is benefit for controlling the unloading spring-back.

The expansion due to martensitic transformation results in the smaller distribution of residual stress.

Theoretical analysis of diffusional austenite decomposition

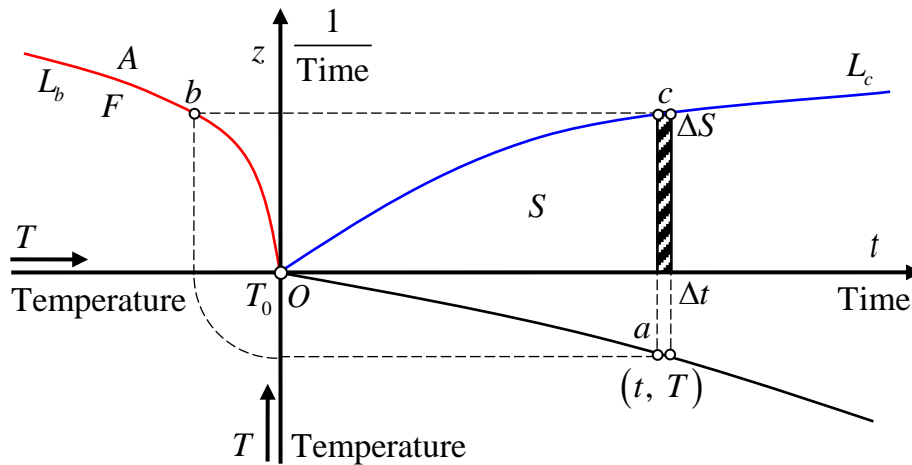
- The CCT curves are easily to be obtained in most industrial design
- The isothermal kinetics curves are hard to be obtained When considering the influence of the deformation as well as the heating process

Is it possible to calculate the transformation kinetics undergoing arbitrary thermal path from CCT curves?

- The mathematical relationship between isothermal and non-isothermal transformation
 - Generalized additivity rule of transformation kinetics based on the CCT curves
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Theoretical analysis of diffusional austenite decomposition

The mathematical relationship between isothermal and non-isothermal transformation – incubation time



- L_a : The non-isothermal path curve
- L_b : The isothermal transformation curve ($f(x) = 0$)
- L_c : The incubation contribution curve

$$\tau_{TTT}(T) = \left. \frac{\partial T_{CCT}(v)}{\partial v} \right|_T$$

$$\sum_i^m \frac{\Delta t_i}{\tau_i} = 1$$

$$\Delta S = \frac{1}{\tau_{TTT}(T)} \Delta t$$

$$S = \int_0^t \frac{1}{\tau_{TTT}(T)} dt \quad \leftarrow \quad T = T_0 + v(t) \cdot t$$

Incubation contribution

$$S = \int_{T_0}^T \frac{1}{\tau_{TTT}(T) \times [v'(t) \cdot t + v(t)]} dT$$

For CCT, $v = \text{constant}$

$$\int_{T_0}^{T_{CCT}} \frac{1}{\tau_{TTT}(T) \cdot v} dT = 1$$

The calculation of TTT incubation time from non-isothermal curves

Theoretical analysis of diffusional austenite decomposition

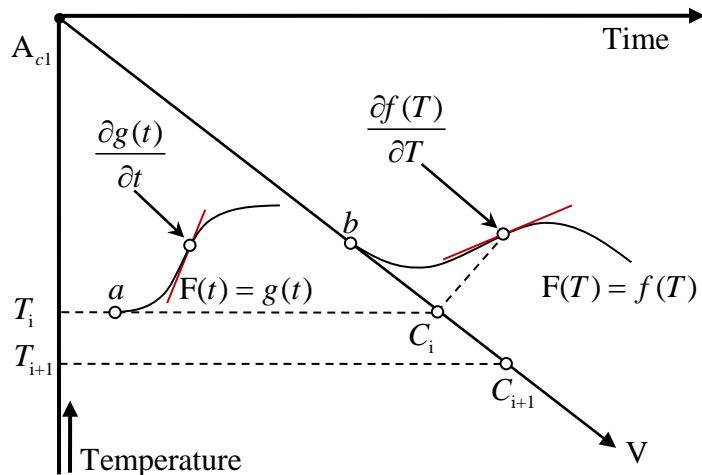


The mathematical relationship between isothermal and non-isothermal transformation – growth kinetics

DO NOT consider incubation time in JMAK equation: only for constant cooling rate (Rios P R. *Acta Materialia* 2005)

$$F(t, T) = 1 - \exp[k(T) \times t^{n(T)}] \longrightarrow t(F, T) = \left. \frac{\partial T_{CCT}(F, v)}{\partial v} \right|_{F, T} = \left\{ \frac{\ln\left[\frac{1}{1-F(t, T)}\right]}{k(T)} \right\}^{\frac{1}{n(T)}} \longrightarrow k(T) = \frac{\ln\left[\frac{1}{1-F(t, T)}\right]}{\left[\left. \frac{\partial T_{CCT}(F, v)}{\partial v} \right|_{F, T} \right]^{n(T)}}$$

CONSIDER incubation time in JMAK equation: for arbitrary cooling path



$$g(t)|_T = 1 - \exp[k(T) \times (t - \tau)^{n(T)}]$$

$$\left. \frac{\partial g(t)}{\partial t} \right|_{t=g^{-1}[f(T_i)]} = \left. \frac{\partial f(T)}{\partial T} \right|_{T=T_i} \times v$$

$K(T)$, $n(T)$ and τ can be obtained

$f(T)$ and $g(T)$ are the non-isothermal and isothermal transformation functions respectively

Schematic mathematical method of transition from CCT to TTT

Theoretical analysis of diffusional austenite decomposition

The generalized additivity rule of CCT curve

$$T = T_0 + v(t) \cdot t$$

$$T_{\text{CCT}}(v) = T_0 + v \cdot \tau_{\text{CCT}}(v)$$

$$v = \frac{\tau_{\text{CCT}}(v)}{T_{\text{CCT}}(v) - T_0}$$

$$S = \int_{T_0}^T \frac{1}{\tau_{\text{TTT}}(T) \times v} dT$$

$$\Delta S = \int_T^{T+\Delta T} \frac{\tau_{\text{CCT}}(v)}{\tau_{\text{TTT}}(T)} \times \frac{dT}{T_{\text{CCT}}(v) - T_0} \leftarrow \tau_{\text{TTT}}(T) = \left. \frac{\partial T_{\text{CCT}}(v)}{\partial v} \right|_T$$

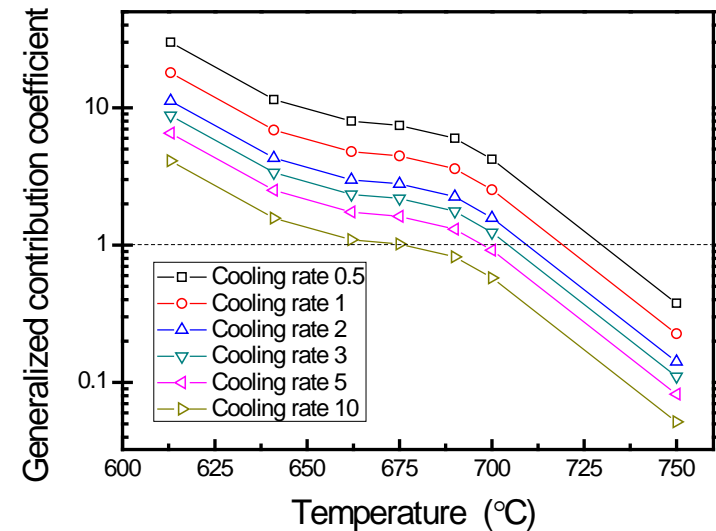
$$\int_{T_0}^{T_f} \left[K(T, v) \times \frac{dT}{T_{\text{CCT}}(v) - T_0} \right] = 1$$

$$K(T, v) = \frac{\tau_{\text{CCT}}(v)}{\left(\left. \frac{\partial T_{\text{CCT}}(v)}{\partial v} \right) \right|_T}$$

Kinetics contribution coefficient

$$\int_{T_0}^{T_f} \left[K(T, v) \times \frac{dT}{T_{\text{CCT}}(v) - T_0} + \frac{\Delta t}{\left(\left. \frac{\partial T(v)}{\partial v} \right) \right|_T} \right] = 1$$

Generalized additivity rule



Kinetics contribution coefficient as the function of temperature

Kinetics contribution coefficient increases with the decreasing transformation temperature, implying that higher cooling degree results in the shorter incubation time.

Theoretical analysis of diffusional austenite decomposition



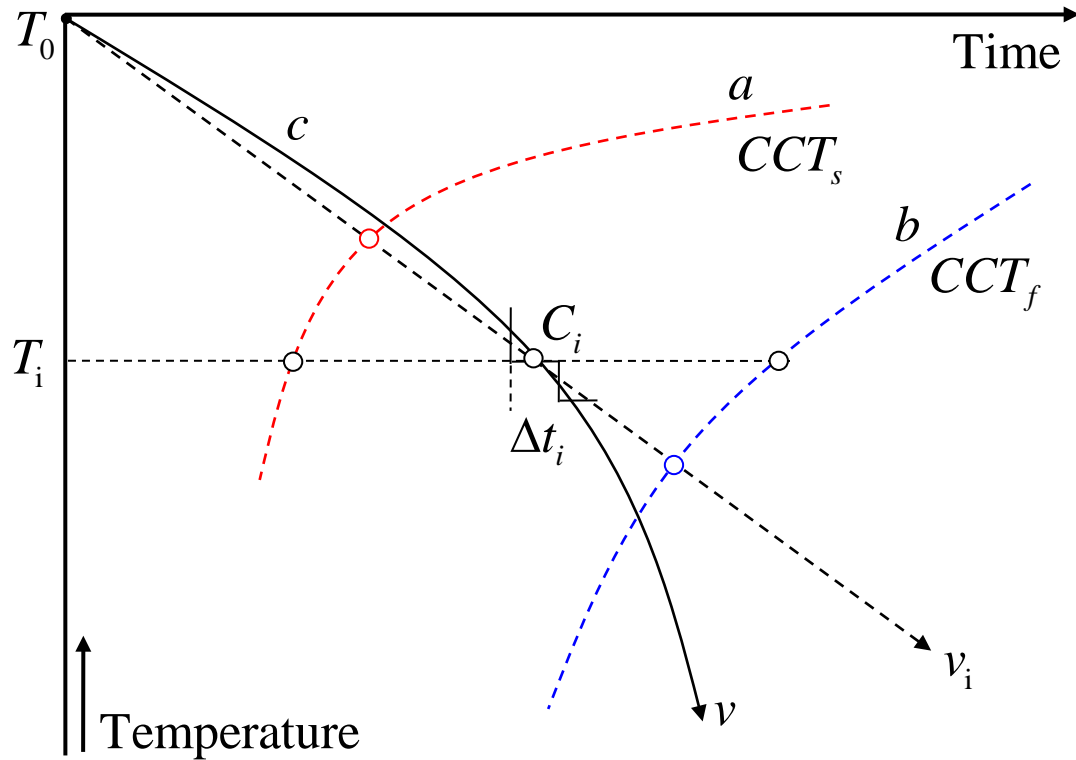
The generalized additivity rule of CCT curve

$$\left. \frac{\partial g(t)}{\partial t} \right|_{t=g^{-1}[f(T_i)]} = \left. \frac{\partial f(T)}{\partial T} \right|_{T=T_i} \times v$$

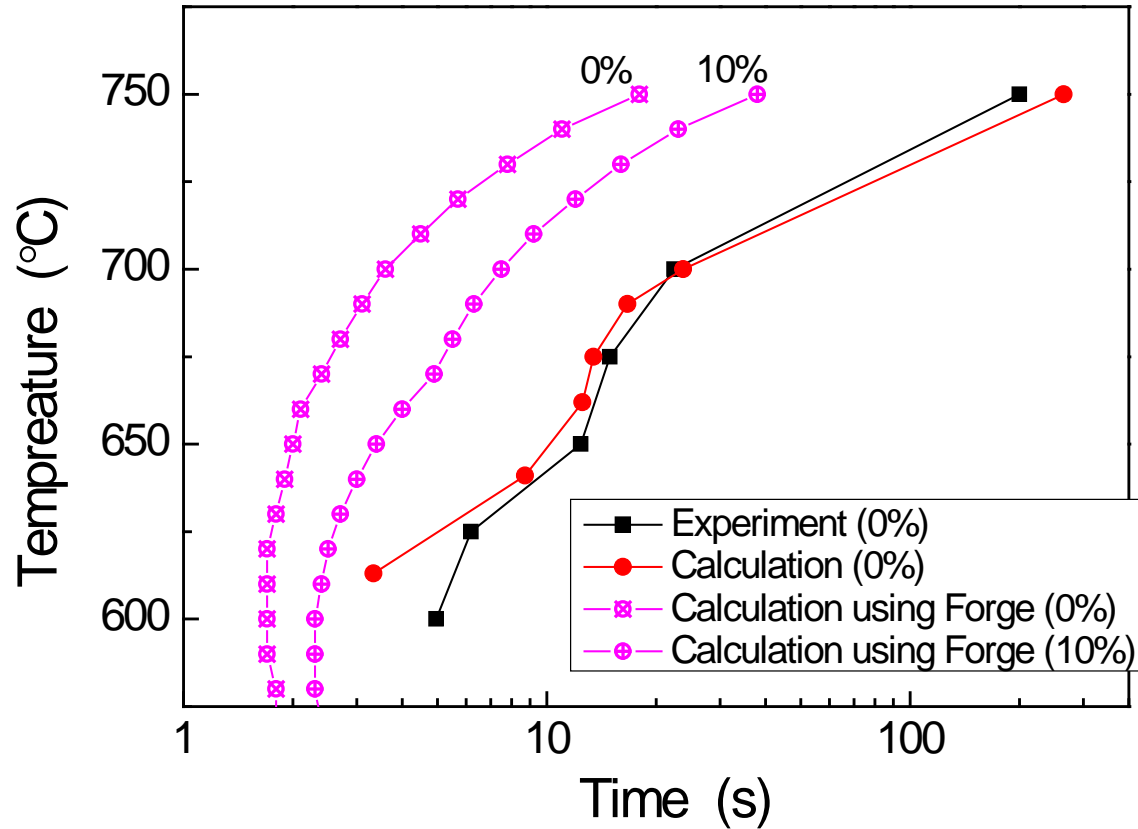
$$\Delta F_i = \left. \frac{\partial F}{\partial t} \right|_{F_i, T_i} \times \Delta t_i$$

$$\Delta F_i = \left. \frac{\partial f(T)}{\partial T} \right|_{T_i} \times v_i \cdot \Delta t_i$$

$$F_{i+1} = F_i + \Delta F_i$$



Schematic generalized additivity rule using CCT curves



The calculated TTT incubation time using CCT data and comparison with experiments

- A computational model which describes the interactive influence of thermal history, phase transformation and its mechanical responses is developed for simulating hot stamping high strength boron steel.
- The introduction of incubation time in JMAK type equations provide the more reasonable simulation results of phase transformation kinetics. The hardness prediction agrees well with the Jominy test, especially at the range of middle and high cooling rate.
- A theoretical relationship between the isothermal and non-isothermal transformation kinetics and the generalized additivity rule for CCT curve are built.
- This generalized model can be used to calculate the transformation kinetics undergoing arbitrary thermal path (including the cooling and heating) using the experimental CCT or CHT curves.

Thank you for your attention!
