

Thermo-Mechanical Fatigue of Cast 319 Aluminum Alloys

Huseyin Sehitoglu,¹ Carlos C. Engler-Pinto Jr.²,
Hans J. Maier³, Tracy J. Foglesong⁴

¹ University of Illinois at Urbana- Champaign, USA,
² Ford Motor Company, Dearborn, ³ Universität-
GH Paderborn, Germany, ⁴ Exxon , Houston

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Outline

- Use of Aluminum Alloys in engine blocks and cylinder heads
- Thermo-Mechanical Fatigue Results
- Summary
- Modeling Studies (Precipitation hardened aluminum alloys)

Percentage of Vehicles with Aluminum Engine Blocks and Heads (*)

	1994	2000	2005
Heads			
Passenger cars	78%	85%	95%
Light trucks	20%	40%	60%
Blocks			
Passenger cars	13%	30%	50%
Light trucks	5%	10%	20%

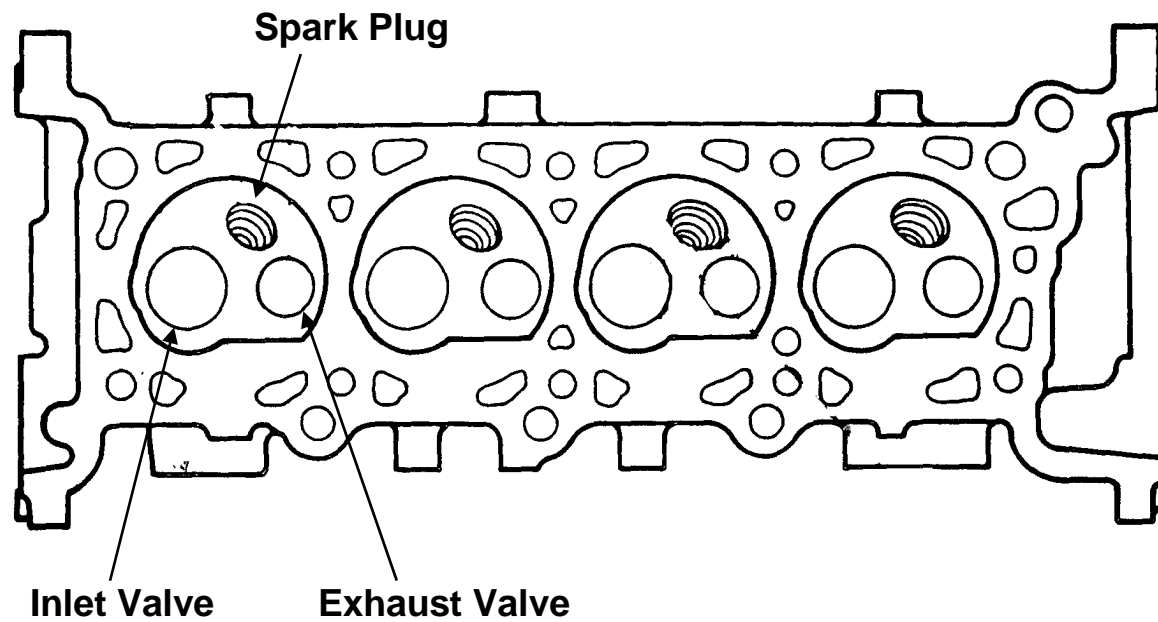
(*) Delphi VIII Study, 1996

Advantages of cast aluminum

- Lightweight
 - V-8 Engine Block: 150 lbs Cast Iron vs. 68 lbs Aluminum
- Cast into complex shapes
- Increased thermal conductivity

Practical Application

- Cylinder Heads



Al319-T7B

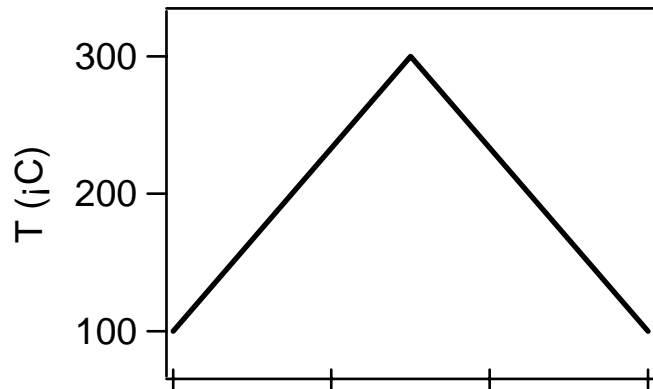
- Nominal Composition in weight percentage

Al	Si	Cu	Mg	Mn	Fe	Zn	Ti	Cr
Bal.	7.2- 7.7	3.3- 3.7	0.25- 0.35	0.20- 0.30	*	0.25 max	0.25 max	0.05 max

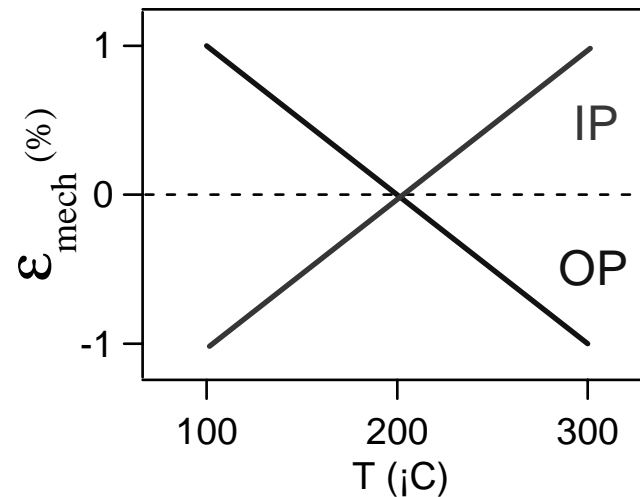
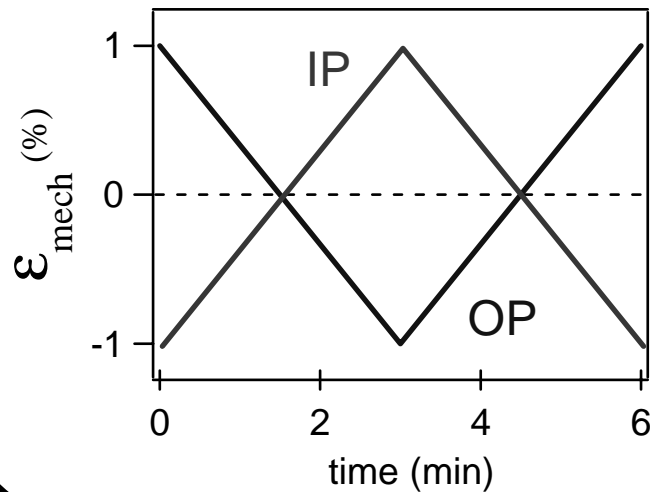
– (*) WAP319: max 0.4% Fe - EAP319: max 0.8% Fe

- Thermal treatment
 - solutionizing at 495°C for 8 hours followed by precipitating at 260°C for 4 hours)

Thermo-Mechanical Fatigue Cycles



- Simultaneously changing strain and temperature (T)
- **In-Phase:** max-strain at max- T
- **Out-of-Phase:** max-strain at min- T



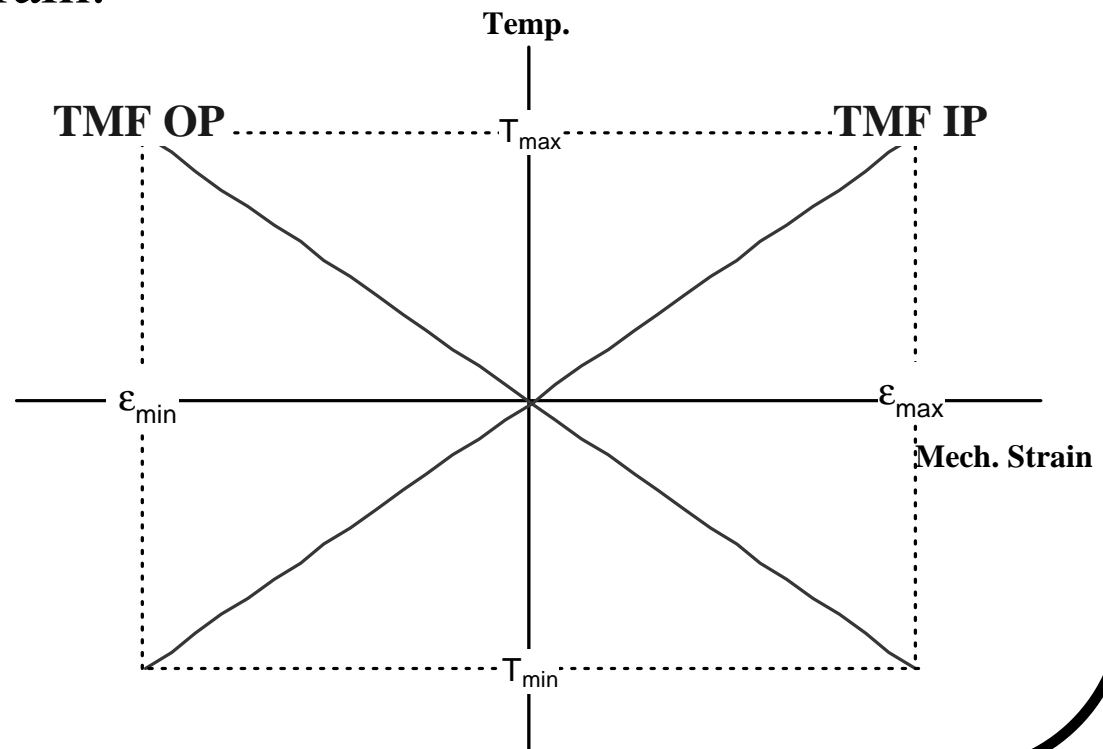
Thermo-Mechanical Fatigue

- Fatigue of materials subjected to simultaneously changing temperature and strain.

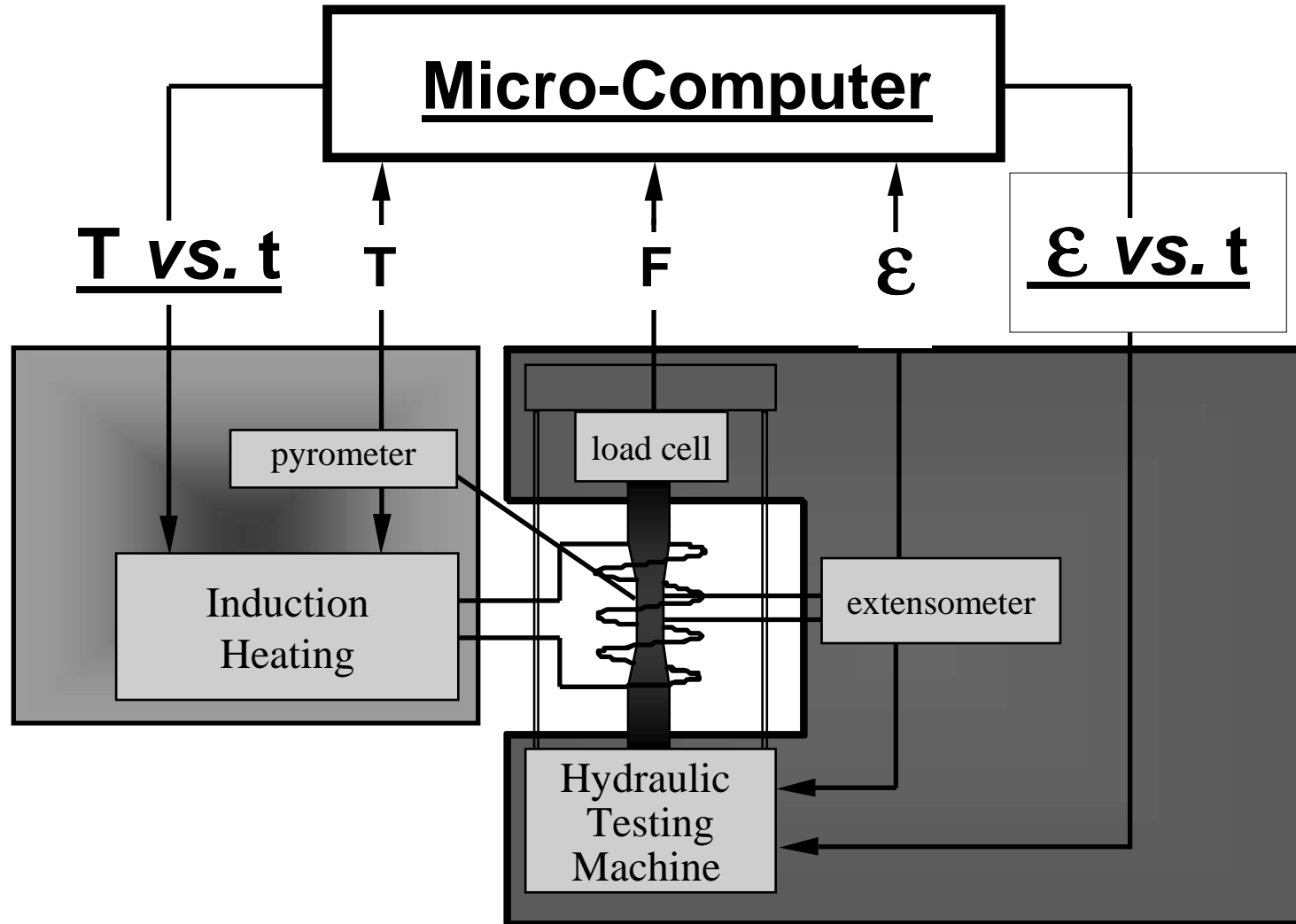
- $\epsilon_{\text{tot}} = \epsilon_{\text{th}} + \epsilon_{\text{mech}}$

- Terminology

- in-phase
- out-of-phase



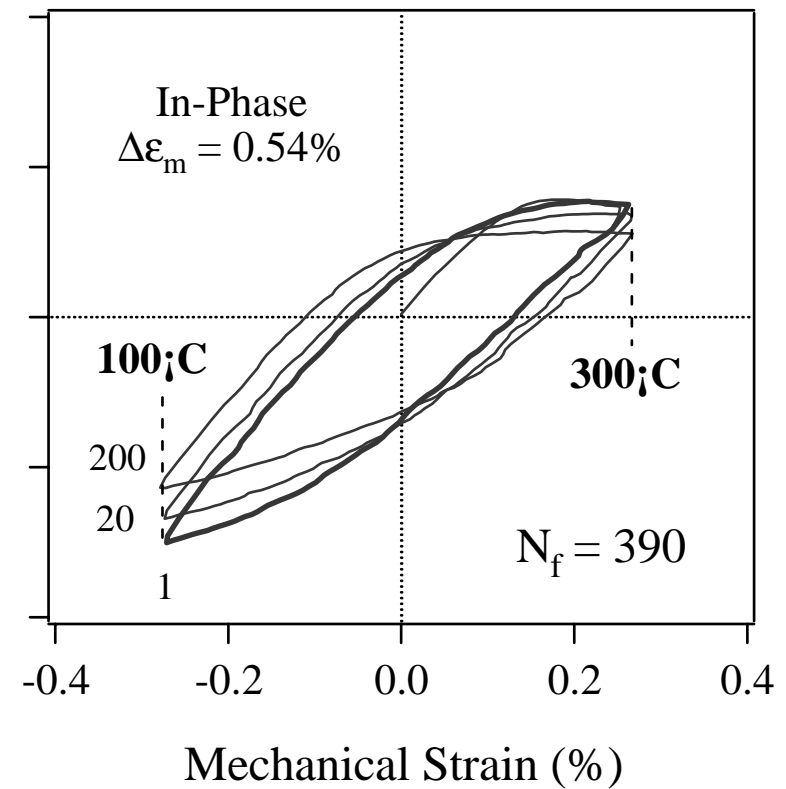
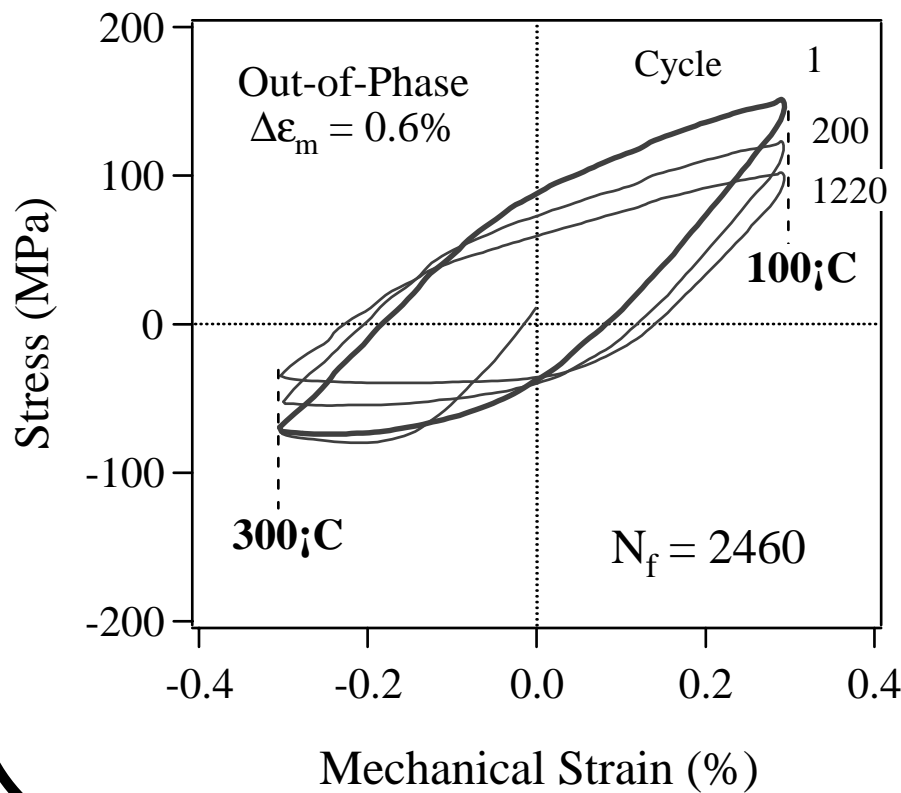
TMF Testing



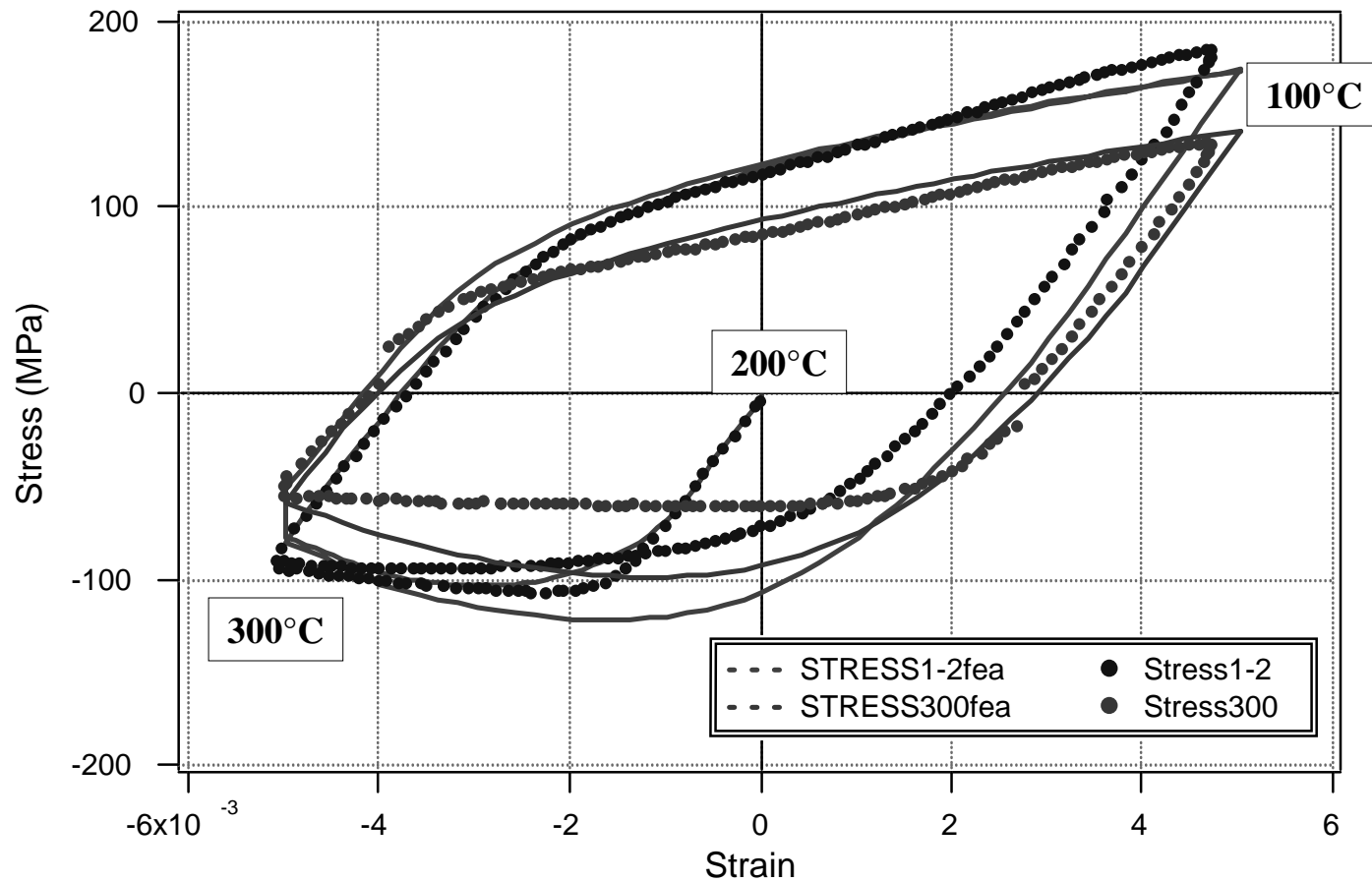
Experimental Procedures

- Isothermal LCF
 - 20°C, 150°C, 250°C and 300°C
 - $2 \times 10^{-1} \text{ s}^{-1}$, $4 \times 10^{-3} \text{ s}^{-1}$ and $5 \times 10^{-5} \text{ s}^{-1}$
- Thermo-Mechanical Fatigue
 - 100–300°C — $5 \times 10^{-5} \text{ s}^{-1}$

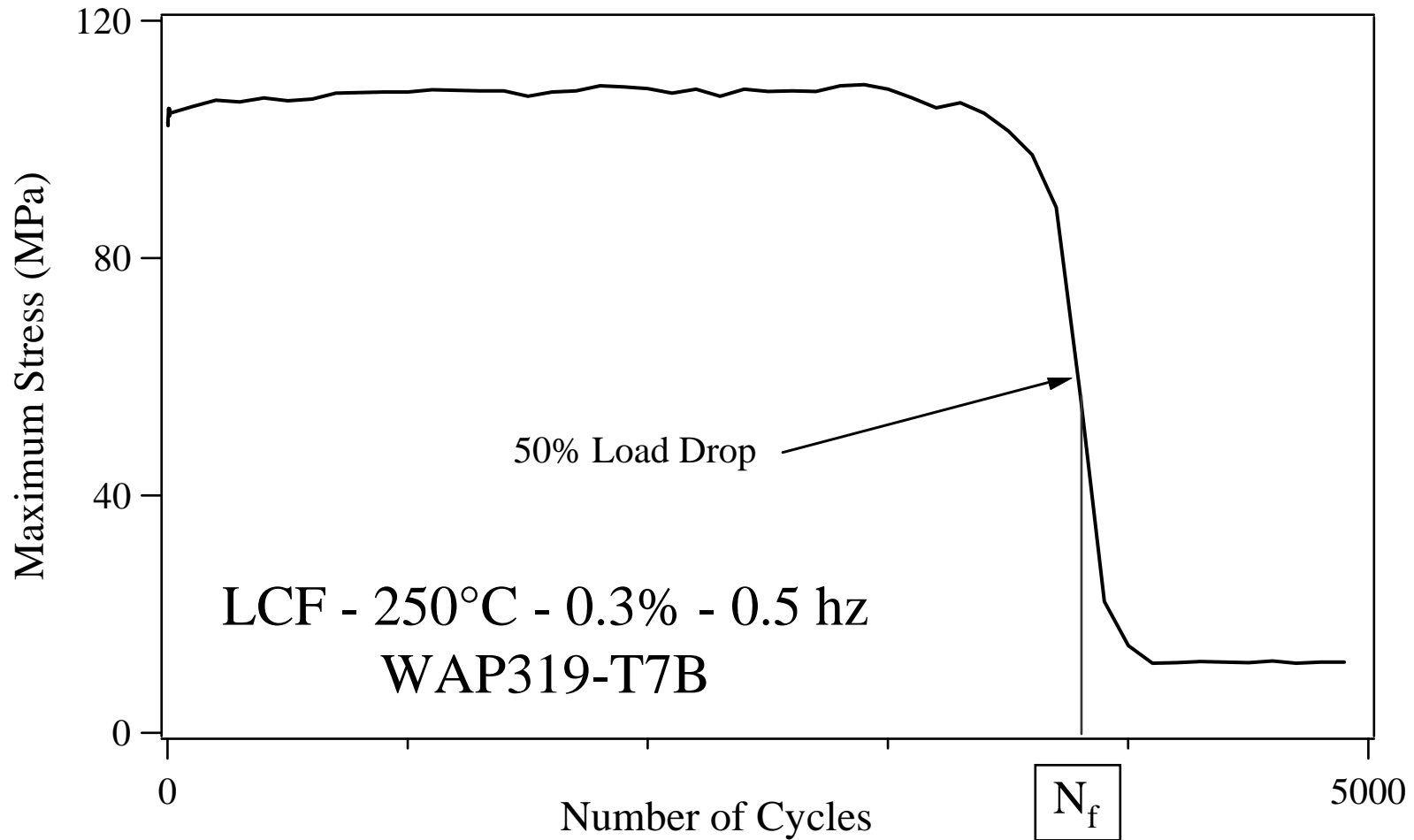
TMF Loops



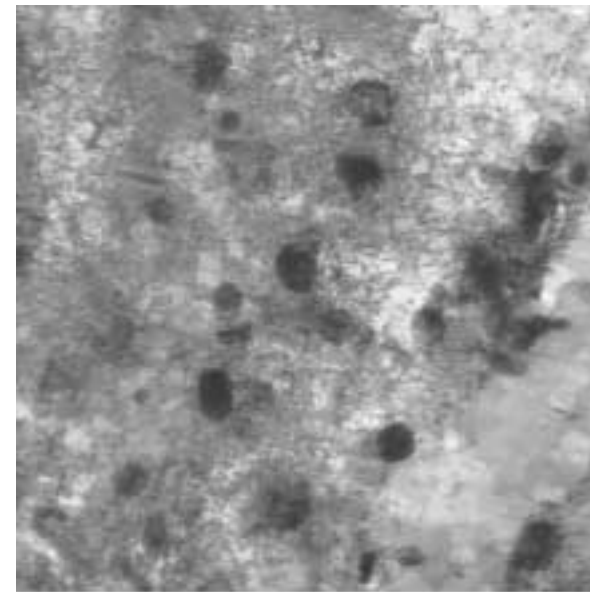
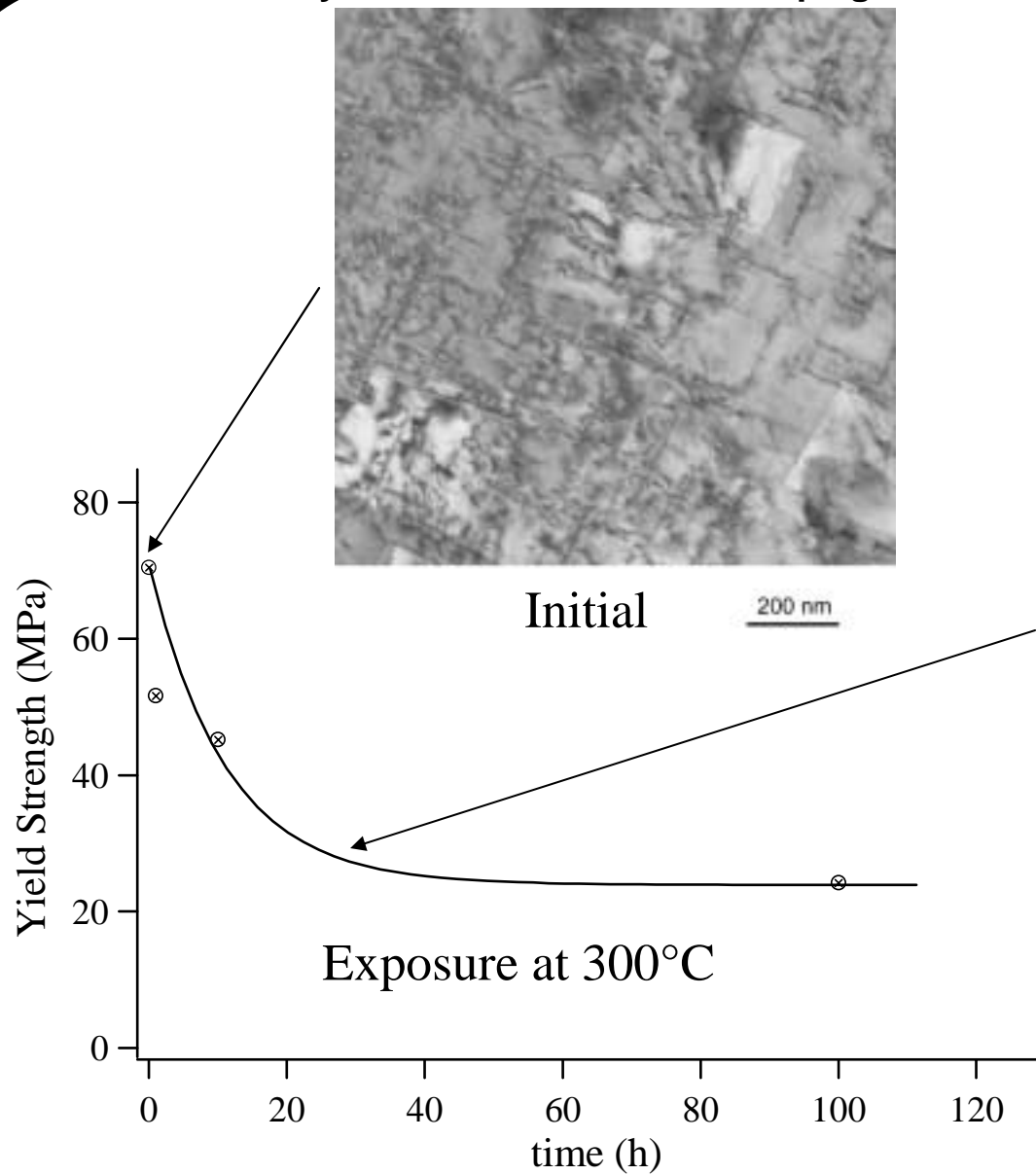
TMF OP 100-300°C 1.0%



Fatigue Life Criterion

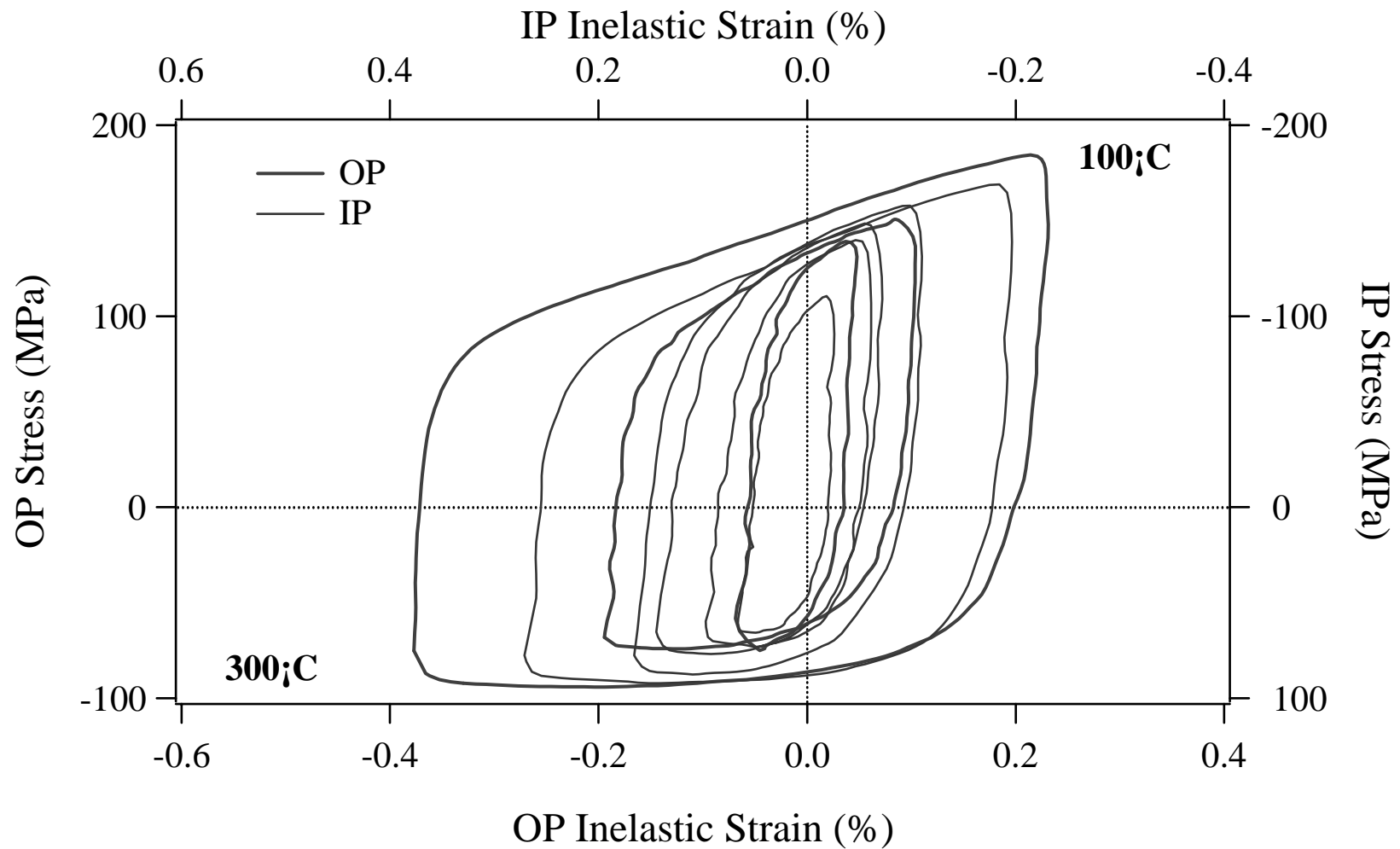


Precipitate Coarsening

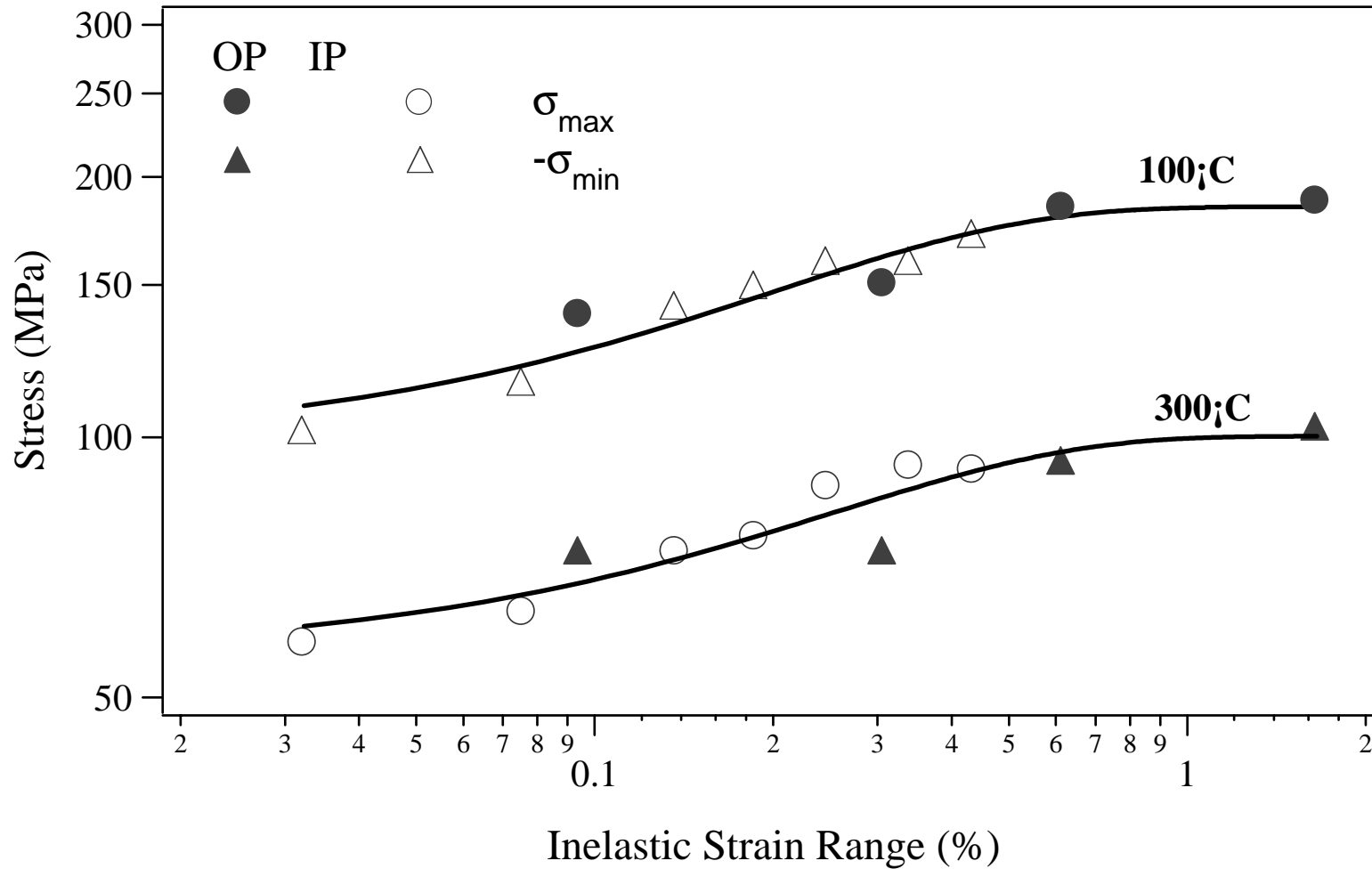


45000 cycles / ~25 hours
(300°C, $\Delta\epsilon_m = 0.2\%$).

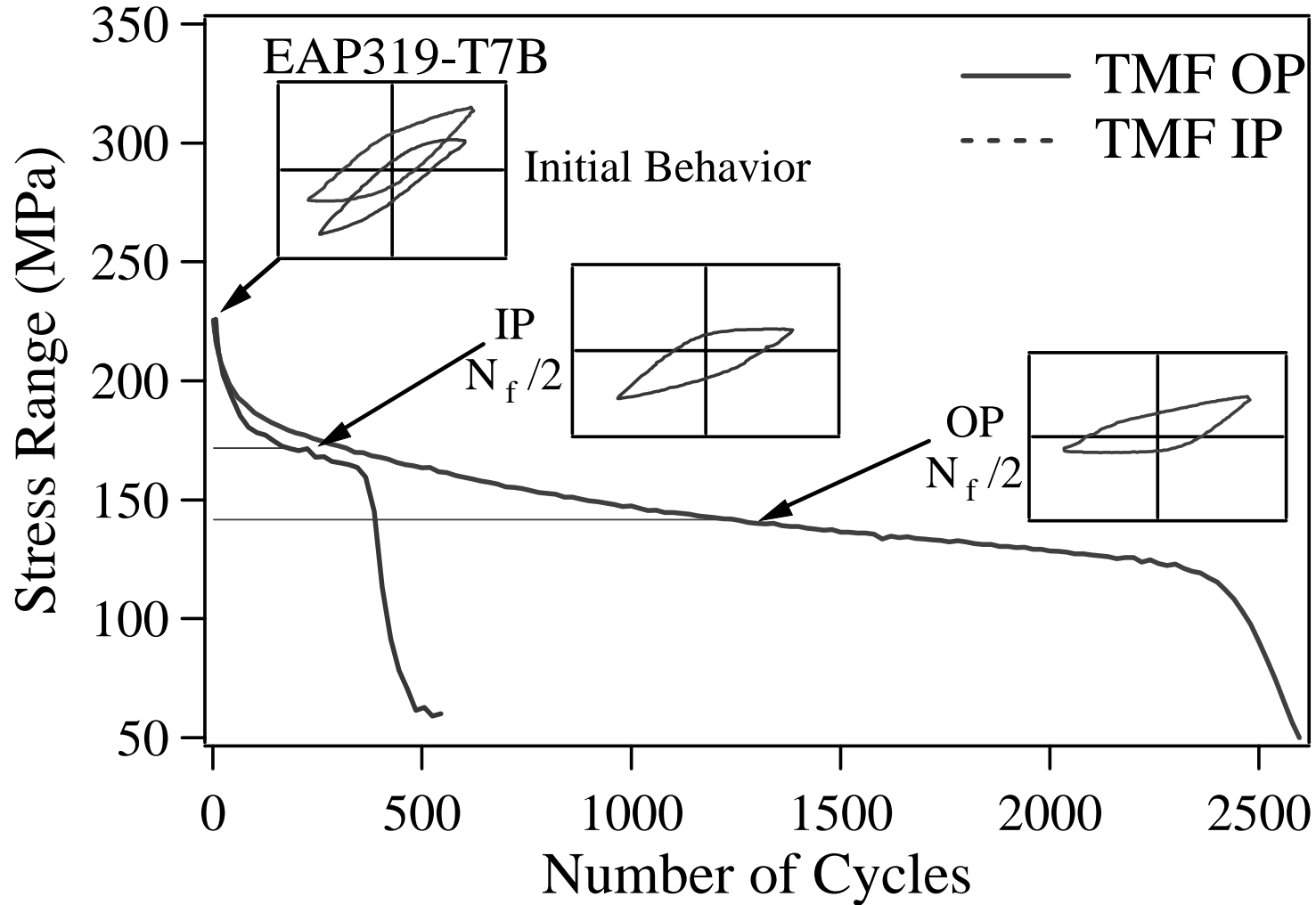
TMF Loops



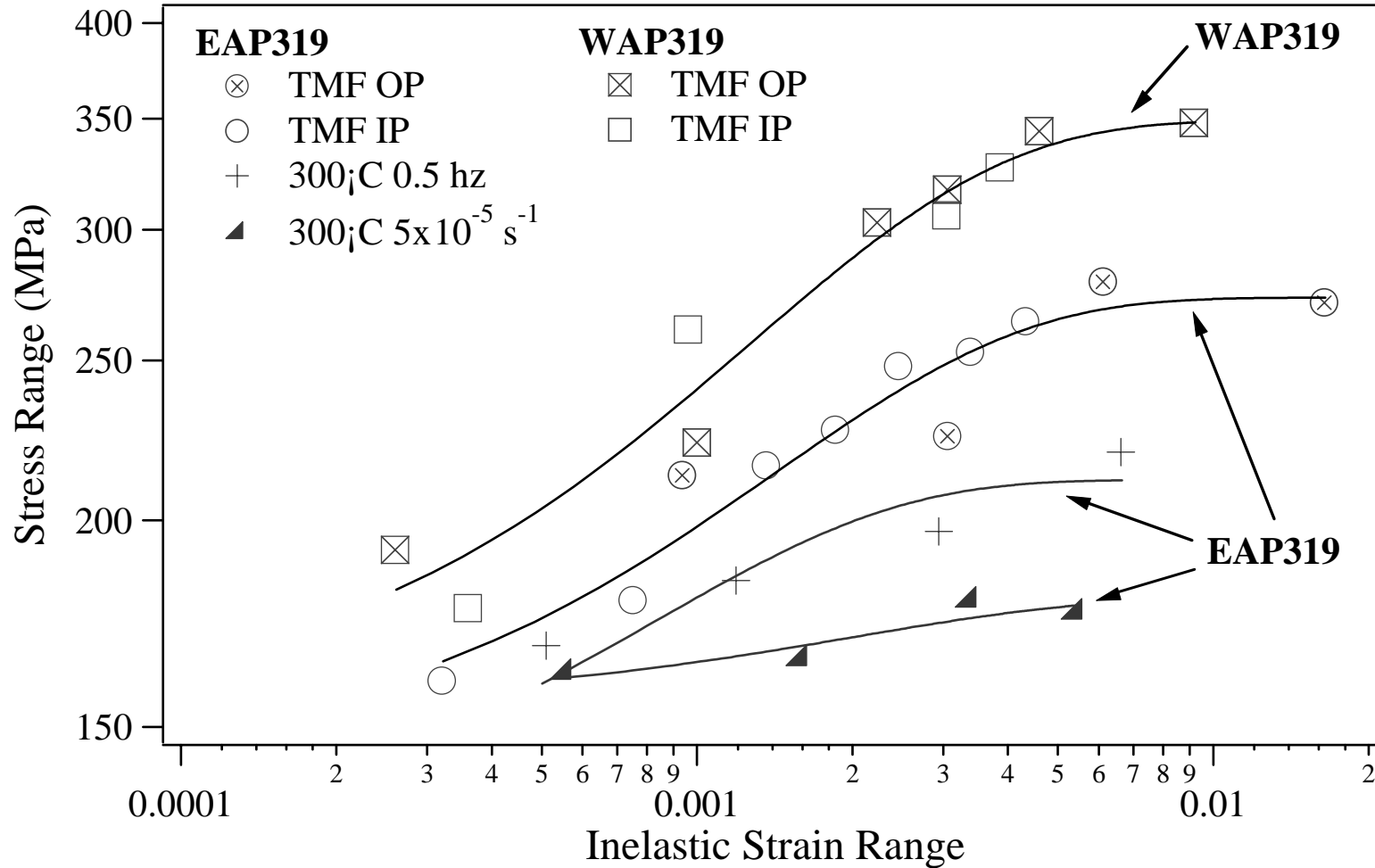
TMF – Peak Stresses



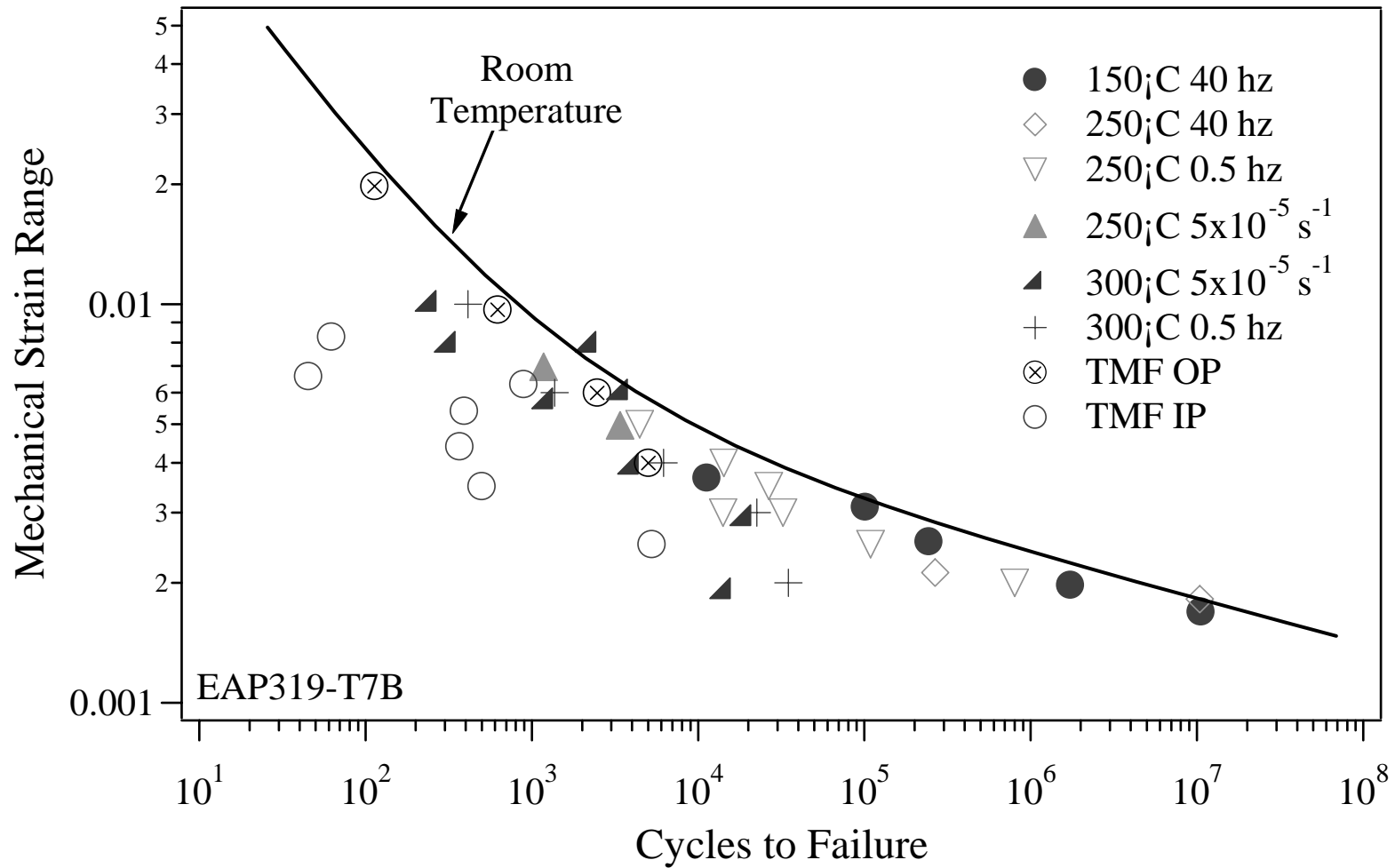
TMF – Stress Range Evolution



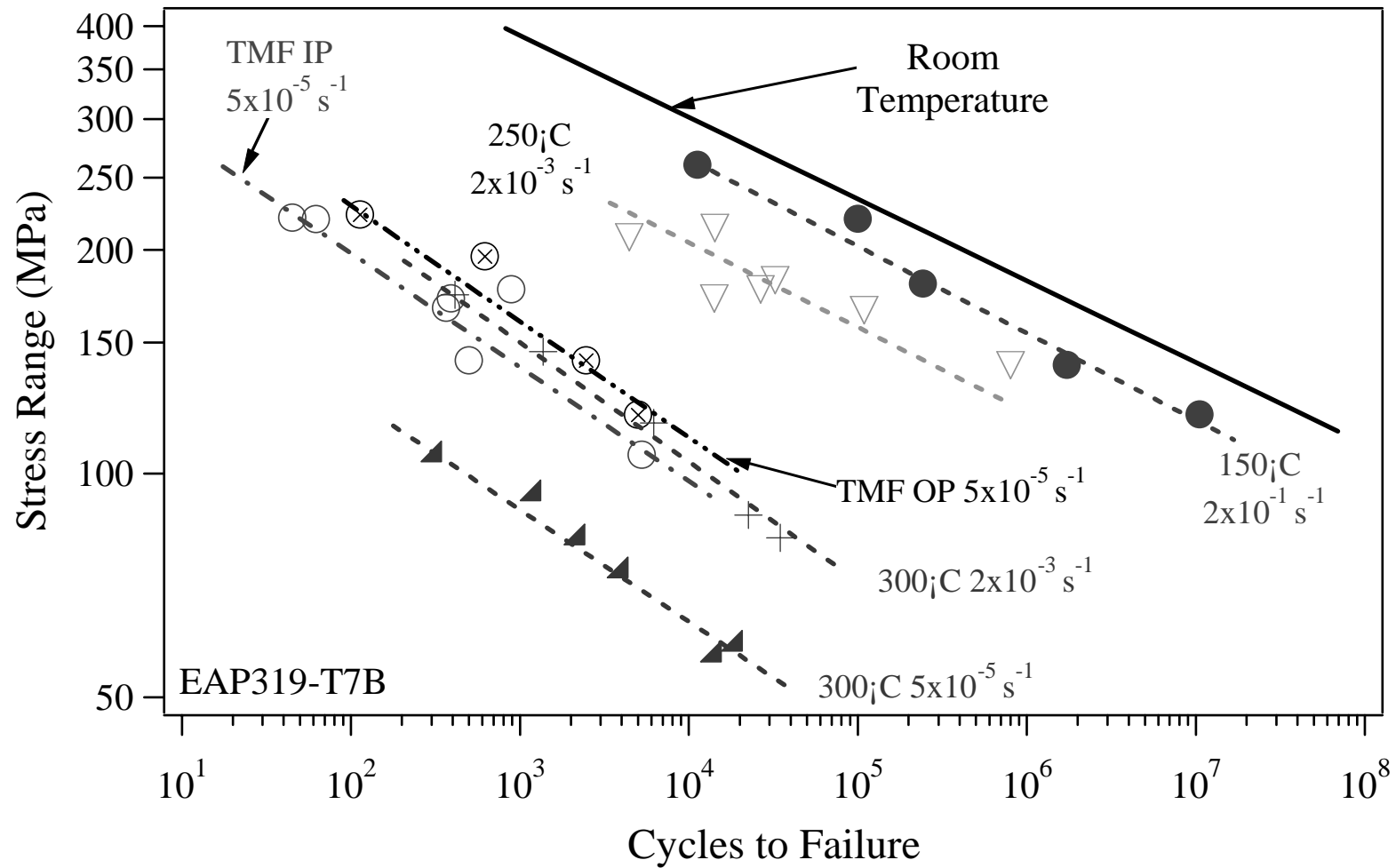
Cyclic Stress-Strain Curves



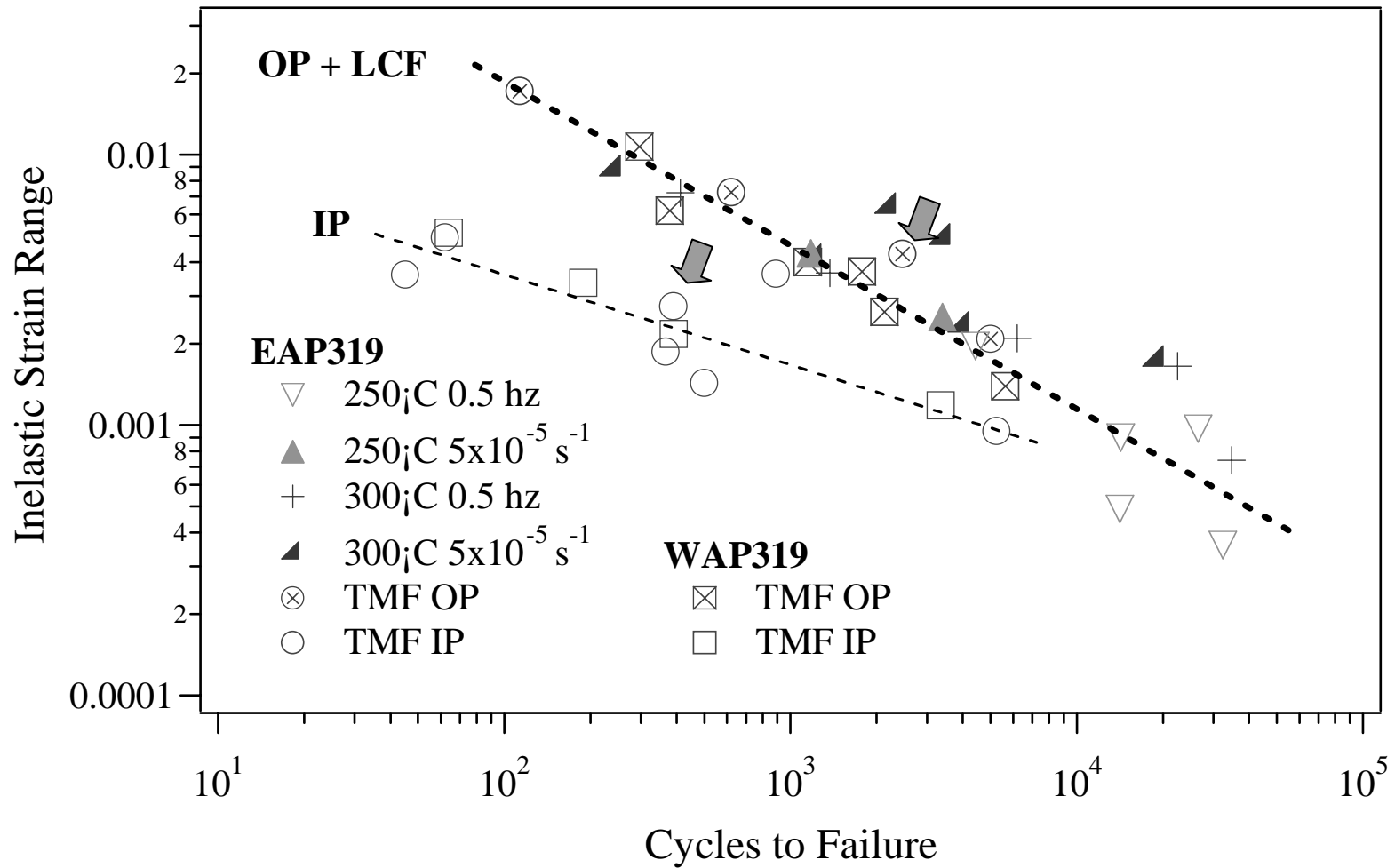
TMF Life



TMF Life



TMF Life



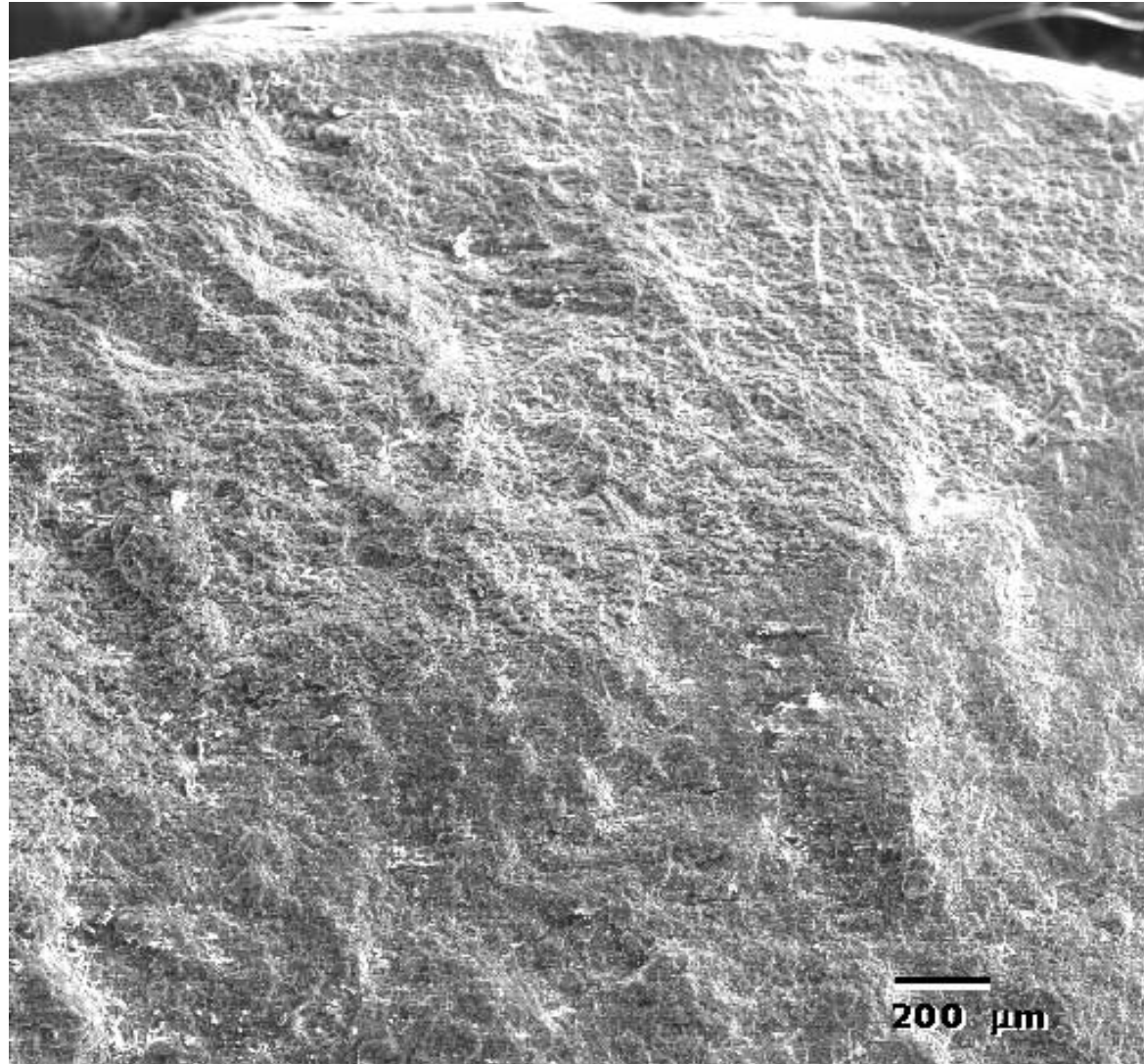
EAP319-T7B

TMF-OP

100–300°C

$\Delta\varepsilon_m = 0.6\%$

$N_f = 2460$ c.



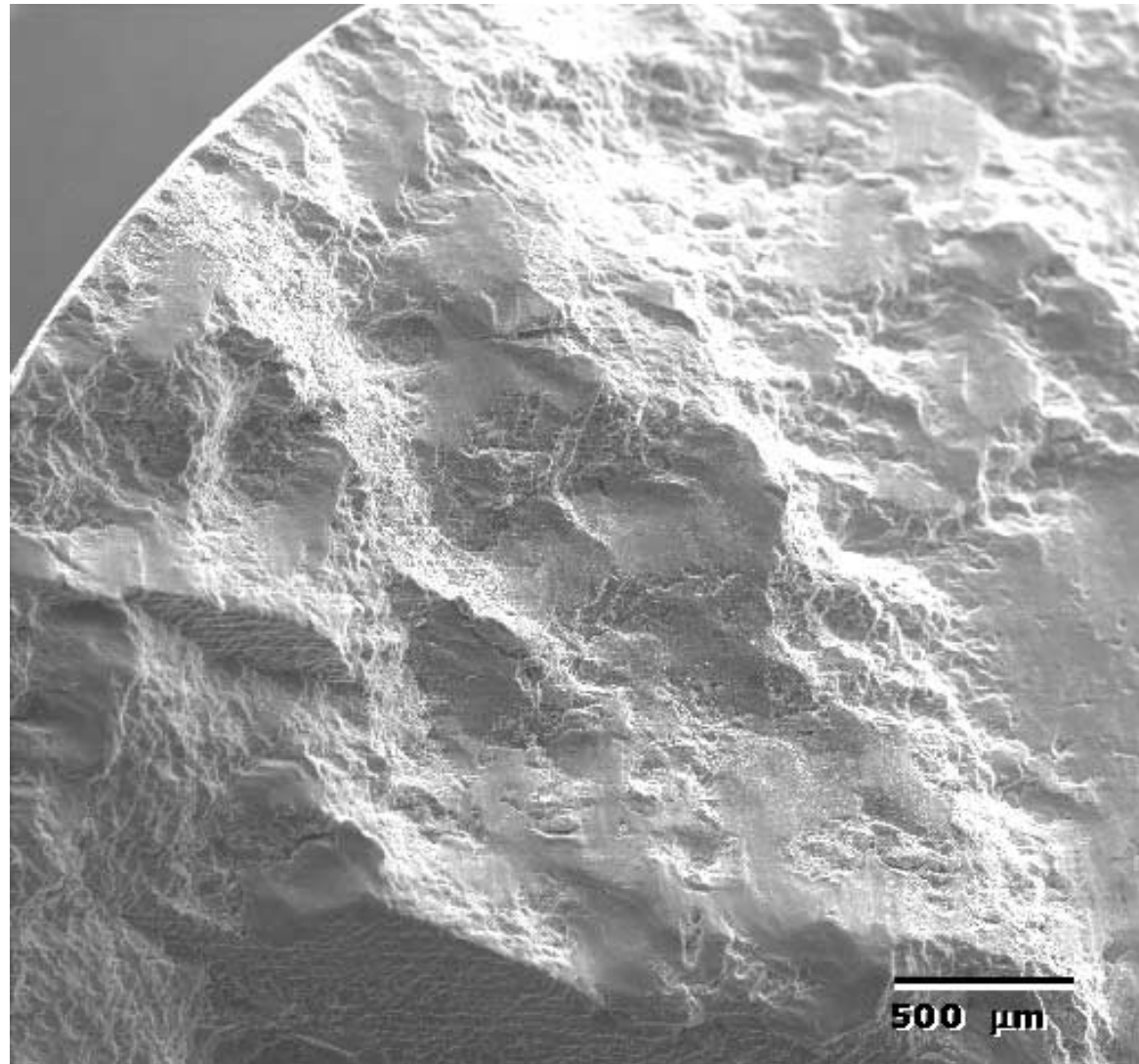
EAP319-T7B

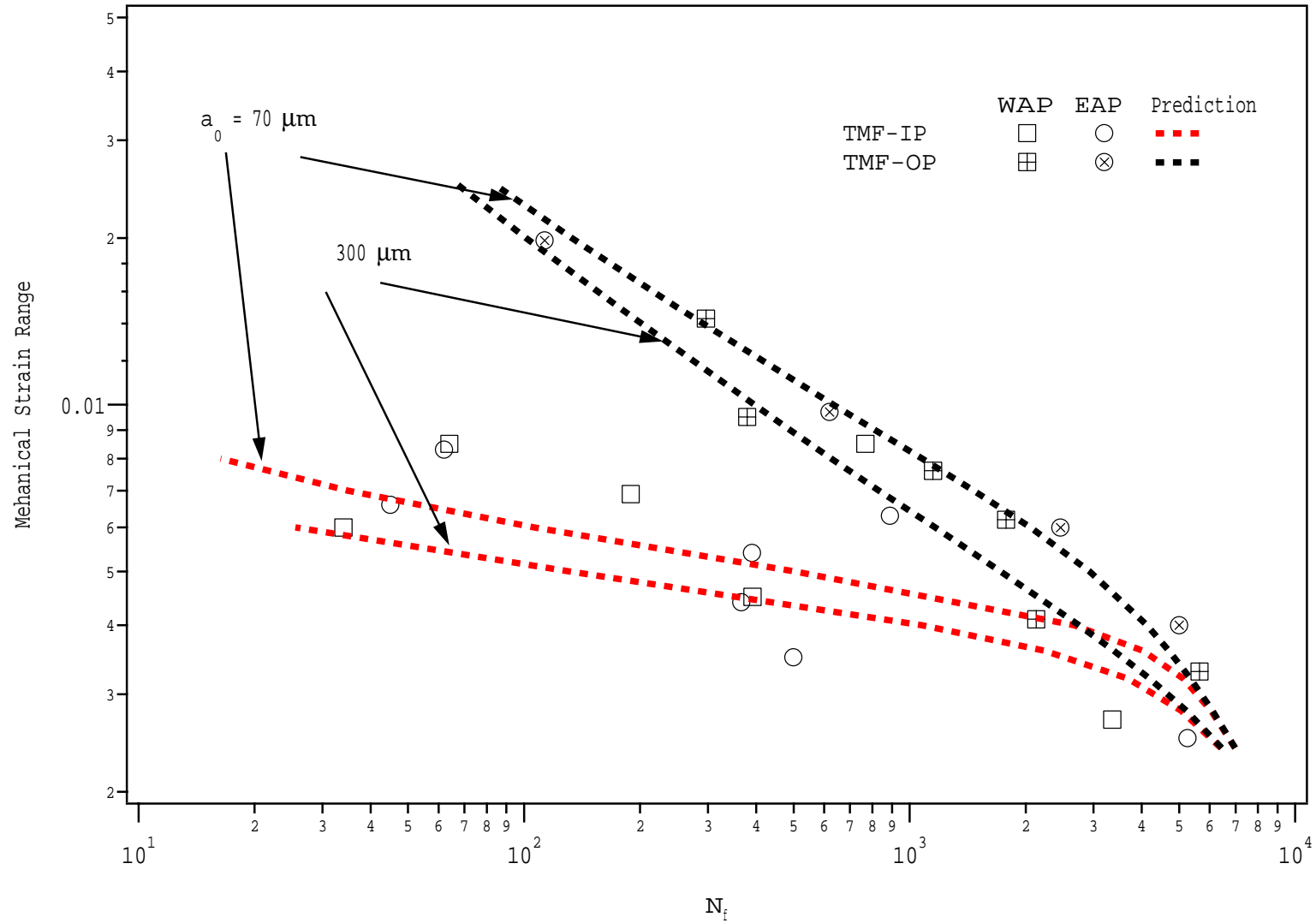
TMF-IP

100–300°C

$\Delta\varepsilon_m = 0.54\%$

$N_f = 390$ c.




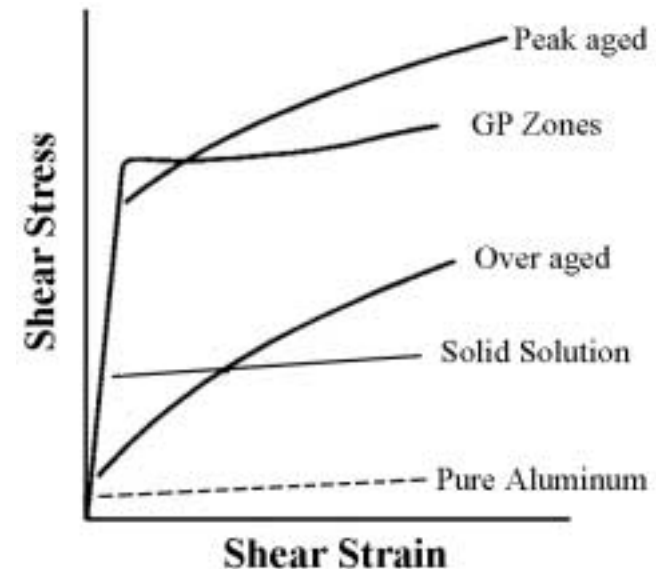


Summary

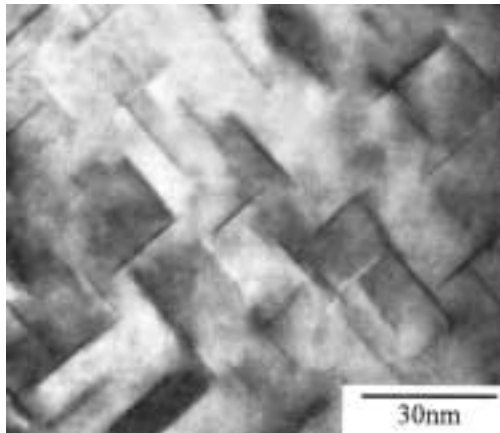
1. TMF stress-strain behavior is identical for both IP and OP loading conditions. TMF-IP lives are shorter than TMF-OP (based on the mechanical or inelastic strain range) lives.
2. Creep damage dominates for TMF-IP loading and in the high strain range regime.
3. The secondary alloy (EAP319) is softer than the primary alloy (WAP319), but TMF lives are very similar.

Aluminum-Copper Alloys

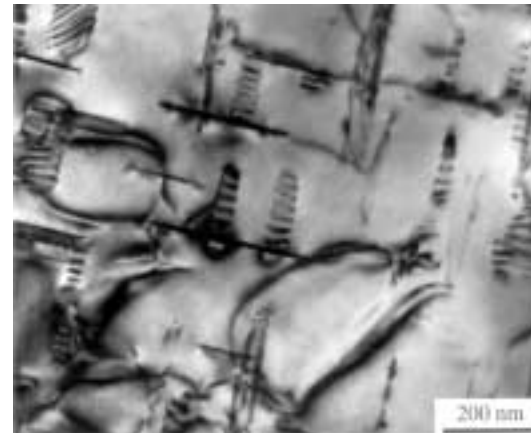
- Precipitate-dislocation interactions
 - Anisotropy on plastic flow behavior (Hosford & Zeisloft '72, Bate *et al.* '81, Barlat & Liu '98, Choi & Barlat '99)
 - Bauschinger effect (Abel & Ham '66, Moan & Embury '79, Wilson '65)
- Coherent particles - GP zones and θ'' (Price and Kelly '64)
 - Higher yield stress than Al shearing of particles
 - Comparable work hardening rates and deformation to Al 
- Semi-coherent - θ' (P & K '64, Russell & Ashby '70)
 - High yield stress and high work hardening rates
- Incoherent particles - θ (P & K '64, R & A '70)
 - Low initial yield stress
 - Highest rates of work hardening



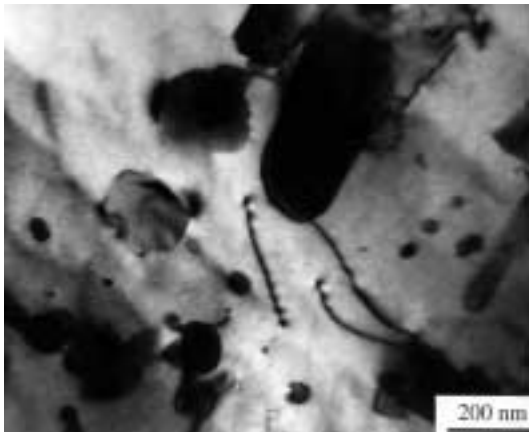
Precipitate Development



GP zones *

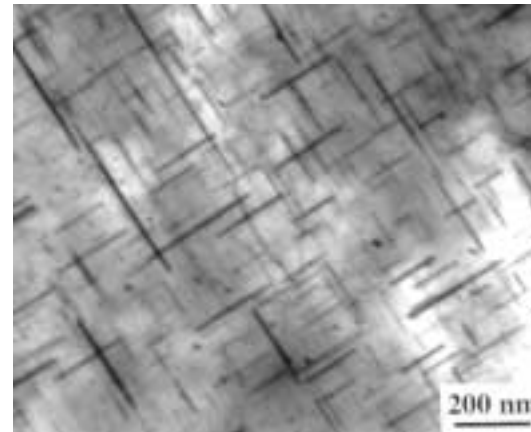
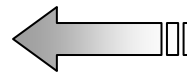


Peak aged, θ'



Very over aged, coarse θ

*Sato & Takahashi, 1983



Over aged, θ' & fine θ

Limitations of current models

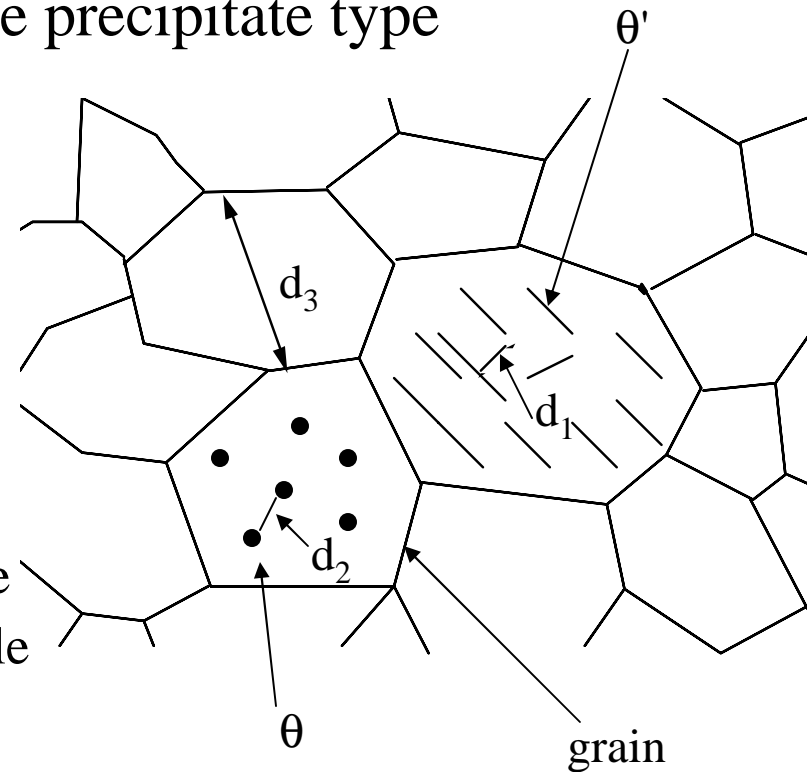
- No implicit consideration of aging treatment.
 - Models were developed for one specific aging treatment
 - Peak aged, θ'
- No inclusion of length scale
 - Volume fraction, precipitate size, mean free path etc. with aging treatment
- Empirical hardening models with a microstructural basis.

Proposed Hardening Law

- Single crystal formulation - one precipitate type

$$\dot{\gamma} = \left[\frac{K_o \alpha^2 \mu^2 b}{2(\tau - \tau_o) d} + \theta_o \left(\frac{\tau_s - \tau}{\tau_s - \tau_o} \right) \right] \sum_k |\dot{\gamma}^k|$$

- Polycrystal formulation
 - More than one type of precipitate
 - Incorporate grain size length scale



$$\dot{\gamma} = \left[\frac{\mu^2 b}{2(\tau - \tau_o)} \left(\frac{\alpha_1^2 K'_1}{d_1} + \frac{\alpha_2^2 K'_2}{d_2} + \frac{\alpha_3^2 K'_3}{d_3} \right) + \theta_o \left(\frac{\tau_s - \tau}{\tau_s - \tau_{o3}} \right) \right] \sum_k |\dot{\gamma}^k|$$

Constitutive Equations

- Relate stress and strain rate at single crystal and polycrystal level.

$$\dot{\epsilon}_i^n = \left\{ \dot{\gamma}^p \sum_{s=1}^S \frac{m_i^s m_j^s}{\tau_c^s} \left(\frac{m_k^s \sigma'_k}{\tau_c^s} \right)^{n-1} \right\} \sigma'_j \quad \text{where } n = 1, 5$$

- Can be written in pseudo-linear form.

$$\dot{\epsilon}_i^n = M_{ij}^{c(\text{sec})}(\sigma') \sigma'_j$$

- Assume overall polycrystal response described by law similar to that of single crystal.

$$\dot{E}_i^{in} = M_{ij}^{(\text{sec})}(\bar{\Sigma}') \bar{\Sigma}'_j$$

Hardening with Precipitates

- Start with dislocation evolution equation

$$\dot{\rho} = \sum_k \left[\frac{K_o}{db} + k_1 \sqrt{\rho} - k_2 \rho \right] |\dot{\gamma}_k|$$

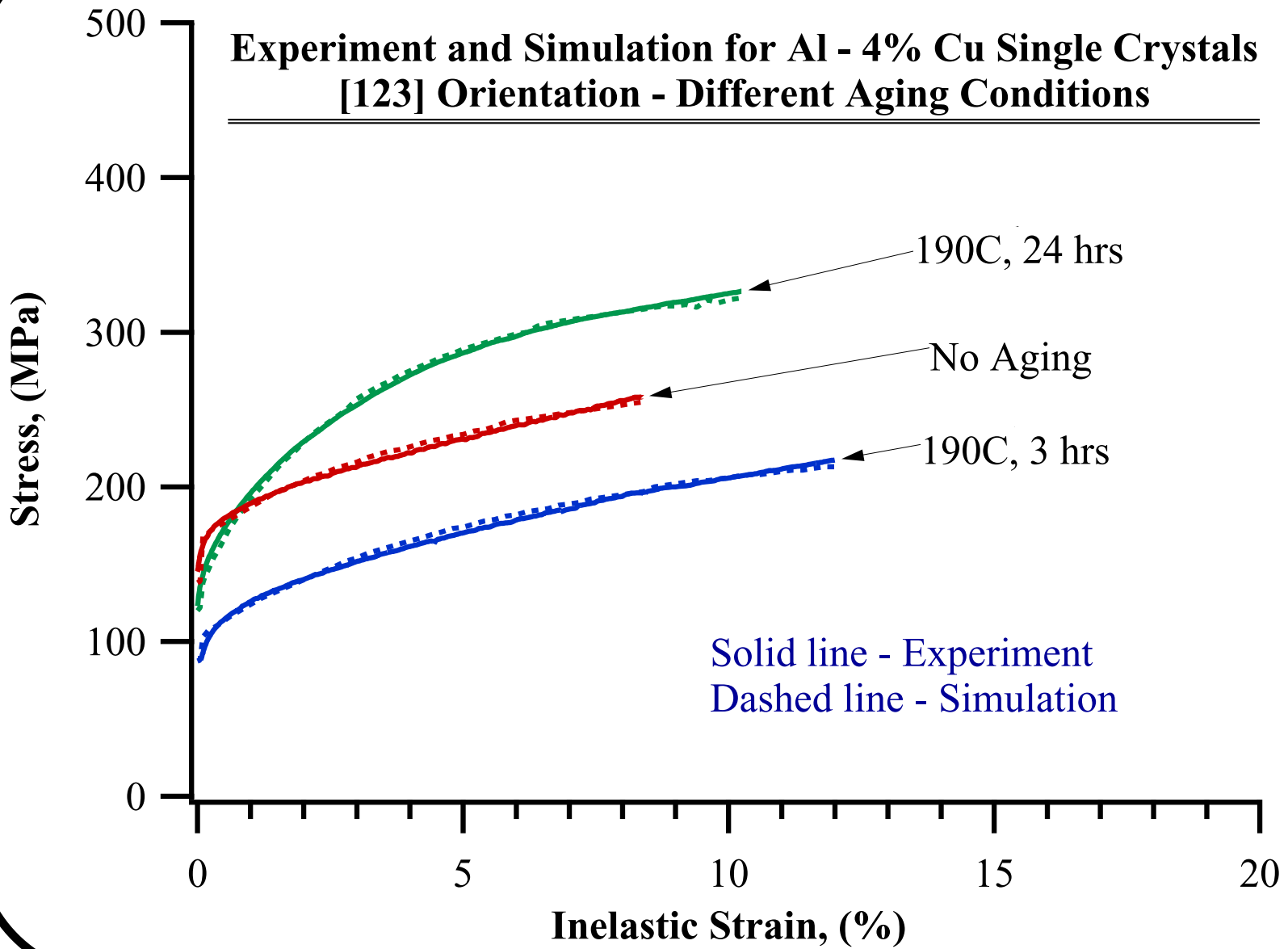
Geometric storage term due to boundaries / obstacles

Dynamic recovery of dislocations

- Combine with the Bailey-Hirsch relationship for flow stress.

$$\tau = \tau_o + \alpha \mu b \sqrt{\rho}$$

**Experiment and Simulation for Al - 4% Cu Single Crystals
[123] Orientation - Different Aging Conditions**



Solid line - Experiment
Dashed line - Simulation

Summary

- The hardening law including the effects of precipitates on the deformation behavior of binary Al - Cu alloys is physically based and accounts for precipitate size, orientation, and mean free path.
- The model incorporates hardening law and predicts single crystal behavior of pure Al and Al-Cu alloys, it also predicts polycrystalline experiments from knowledge of single crystal behavior.