

Investigation of the $\gamma \rightarrow \alpha$ Transformation in Fe-C-X Alloys Using Decarburization

Compiled by Hatem Zurob from works done at
McMaster University and LTPCM by:

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Chris Hutchinson, Yves Brechet and Armand
Beche.





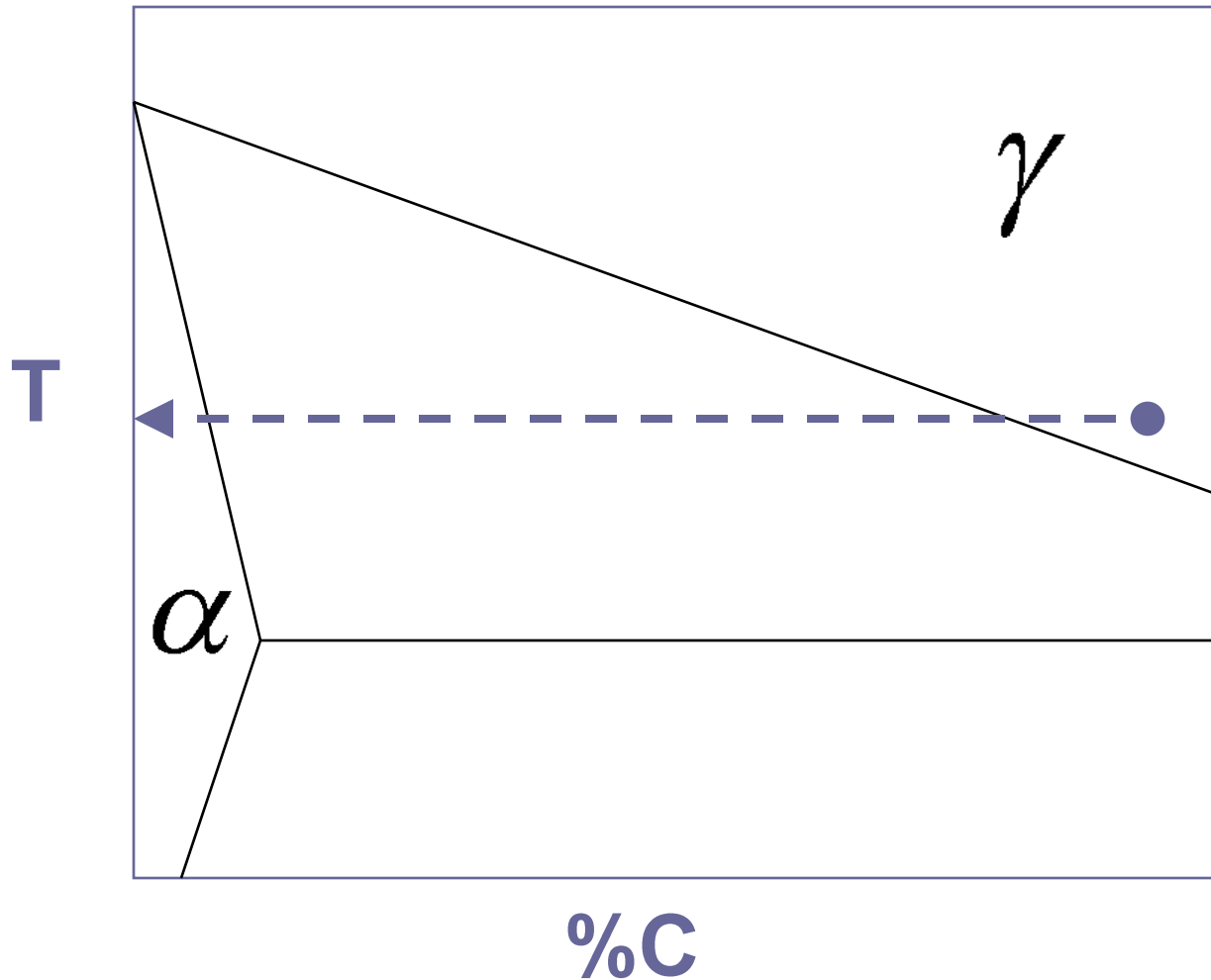
■ Decarburization Method.

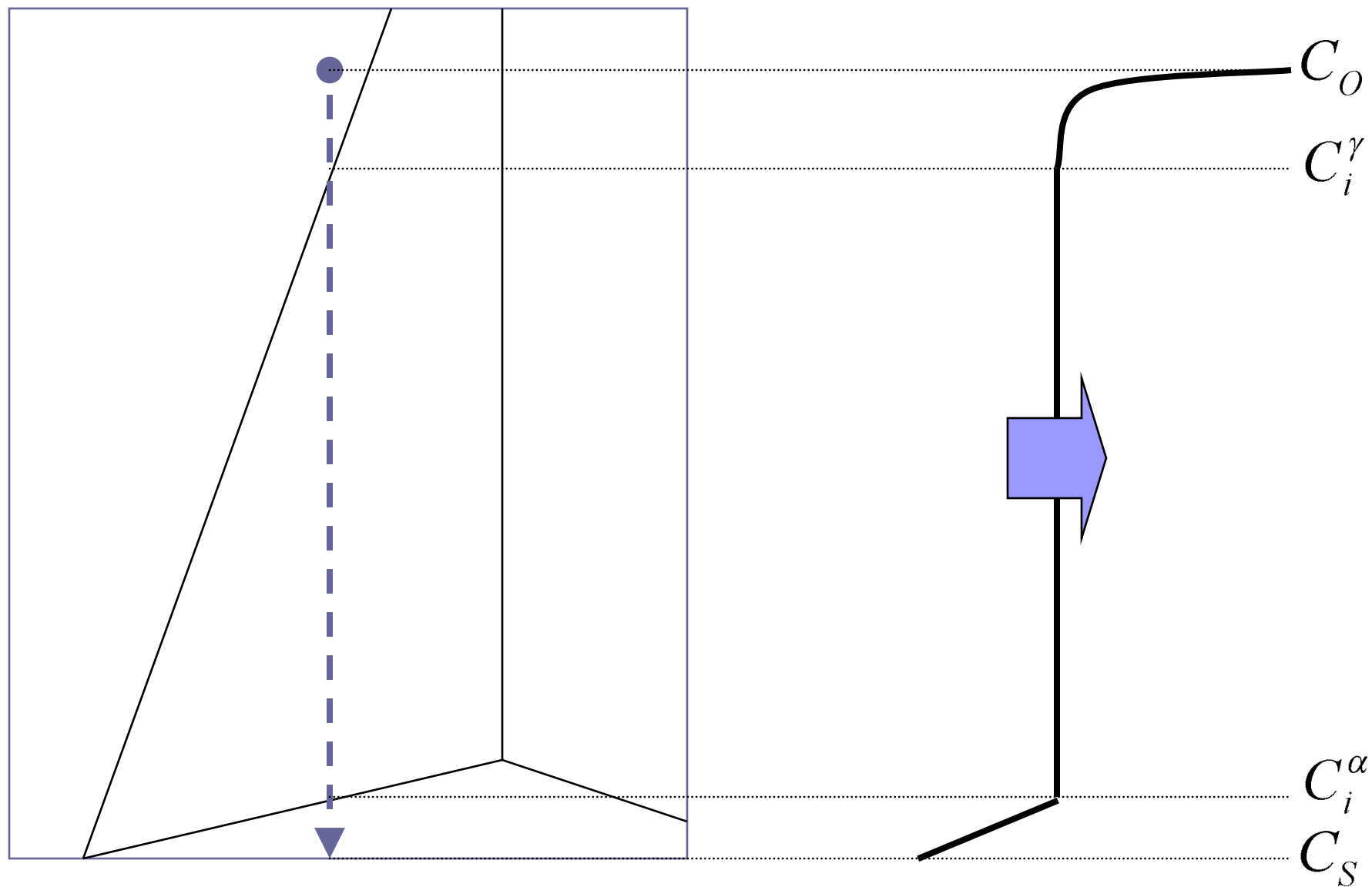
■ Summary of Results:

- ☐ Fe-C
- ☐ Fe-C-Ni
- ☐ Fe-C-Mn
- ☐ Fe-C-Cr
- ☐ Fe-C-Mo
- ☐ Fe-C-Si and Fe-C-Si-Mn
- ☐ Fe-C-Nb

■ Future Work.

1. The Decarburization Method





The rate of interface motion is given by:

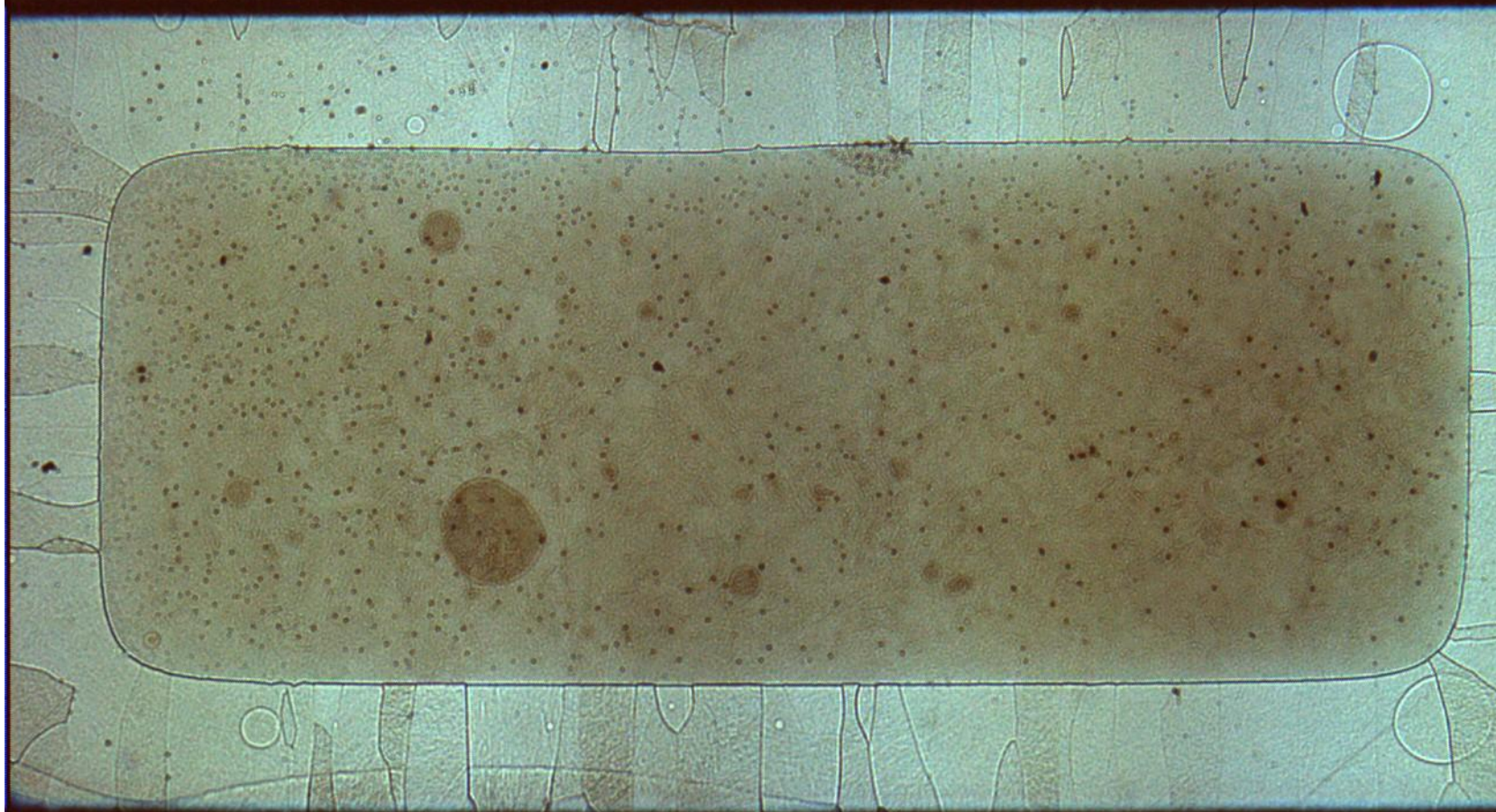
$$\frac{dz}{dt} = \frac{J_i^\alpha - J_i^\gamma}{C_i^\gamma - C_i^\alpha}$$

This differential equation has an analytical solution of the form:

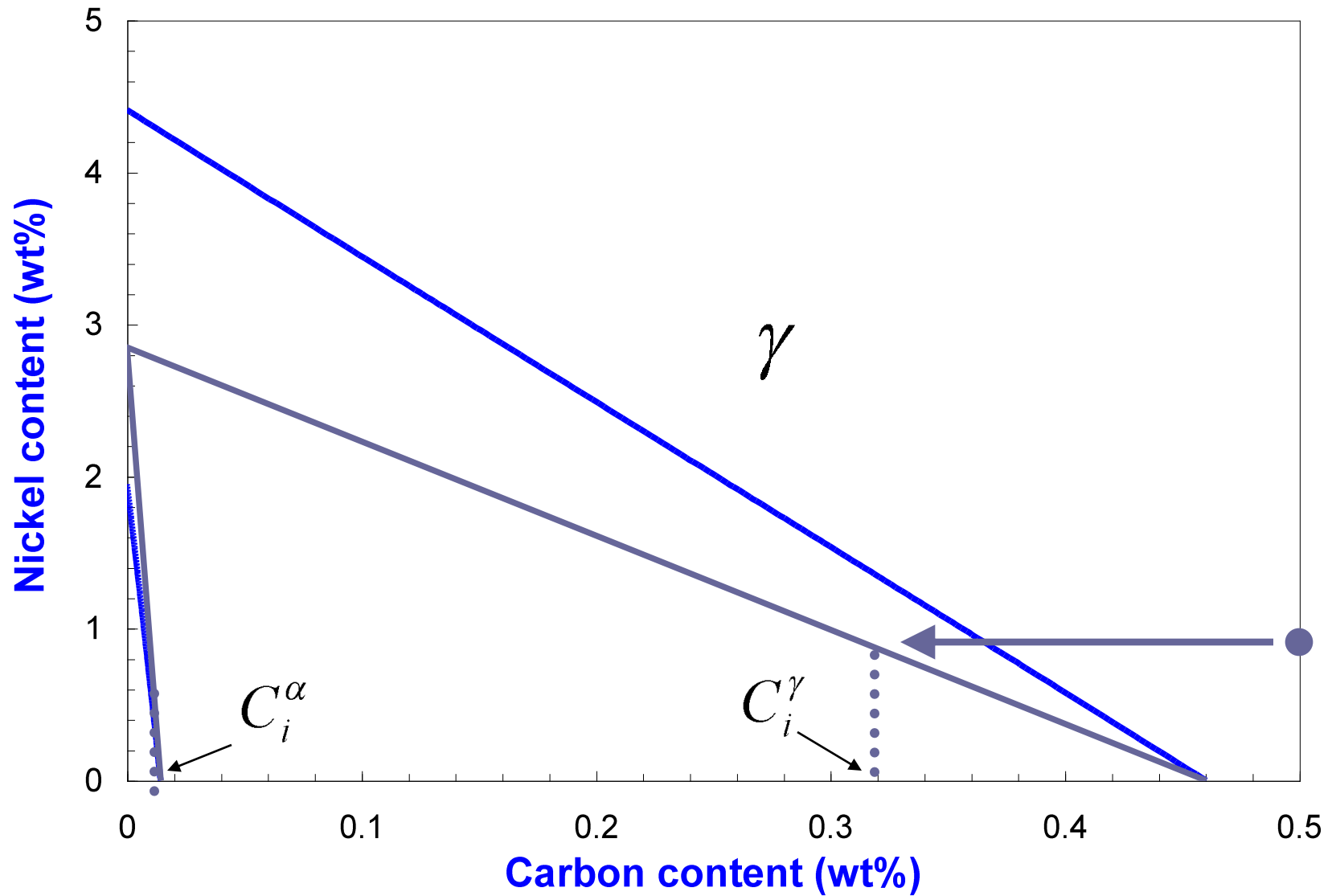
$$z = B\sqrt{t}$$

$$B = f(C_i^\alpha, C_i^\gamma, C_o)$$

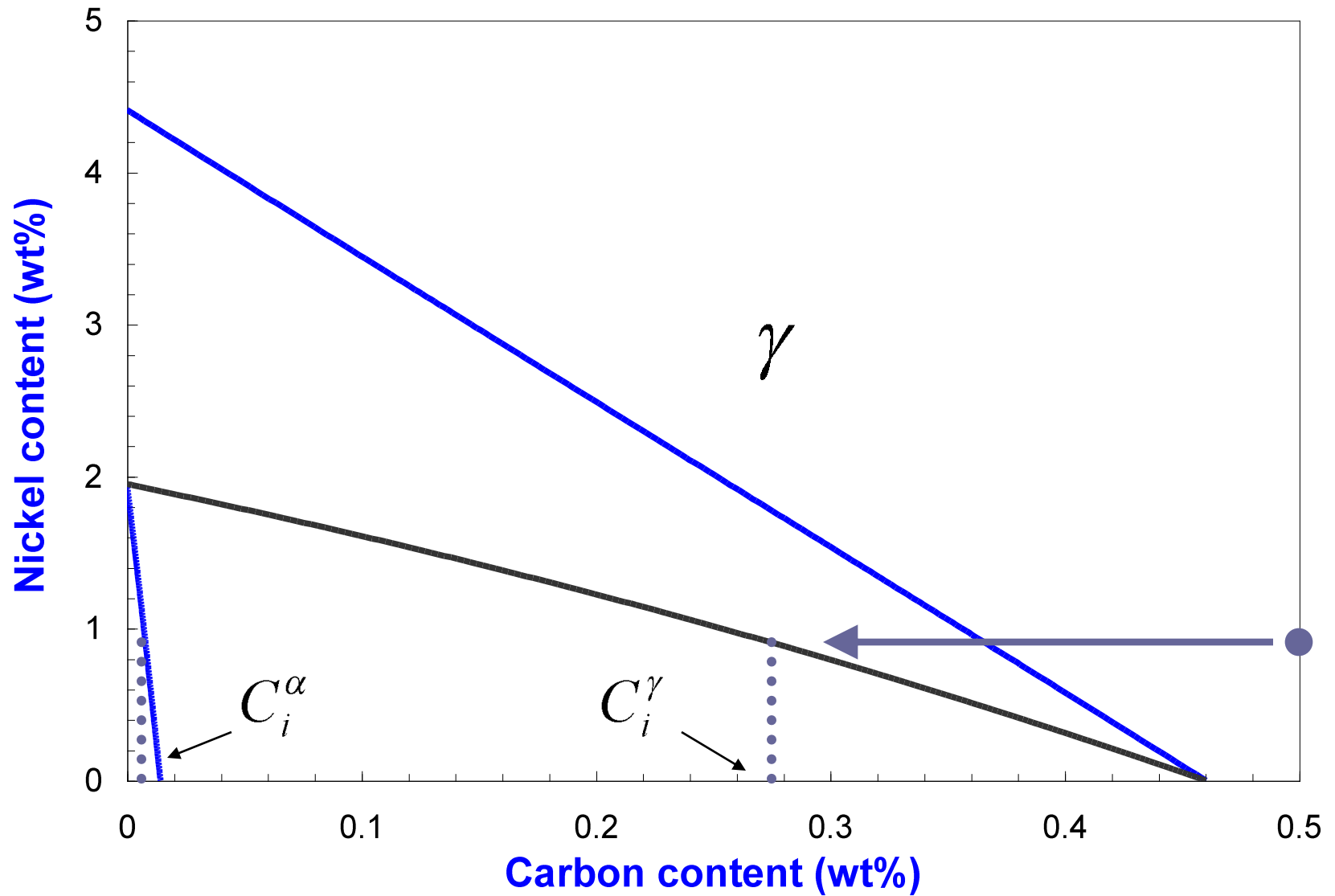
1 mm



➤ Ternary Alloys: ParaEquilibrium Limit.



➤ Ternary Alloys: LENP Limit.

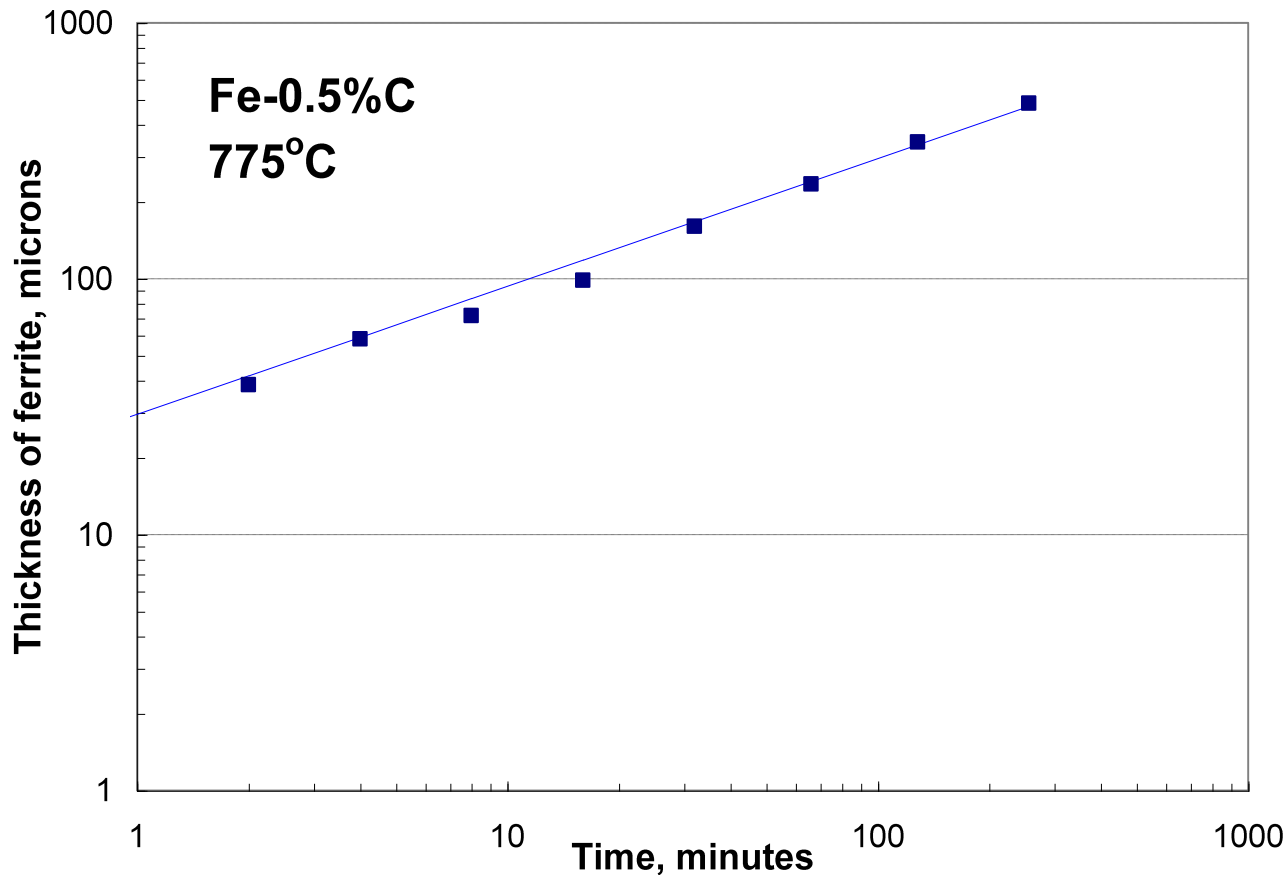




2. Summary of Results:

- 2.1. Fe-C (diffusion coefficient).
- 2.2. Fe-C-Ni (the LE-NP limit).
- 2.3. Fe-C-Mn (possible transition).
- 2.4. Fe-C-Cr (solute drag).
- 2.5. Fe-C-Mo (dissipation are not created equal).
- 2.6. Fe-C-Si and Fe-C-Si-Mn (combined effects).
- 2.7. Fe-C-Nb (precipitation).

2.1. Binary Fe-C System:

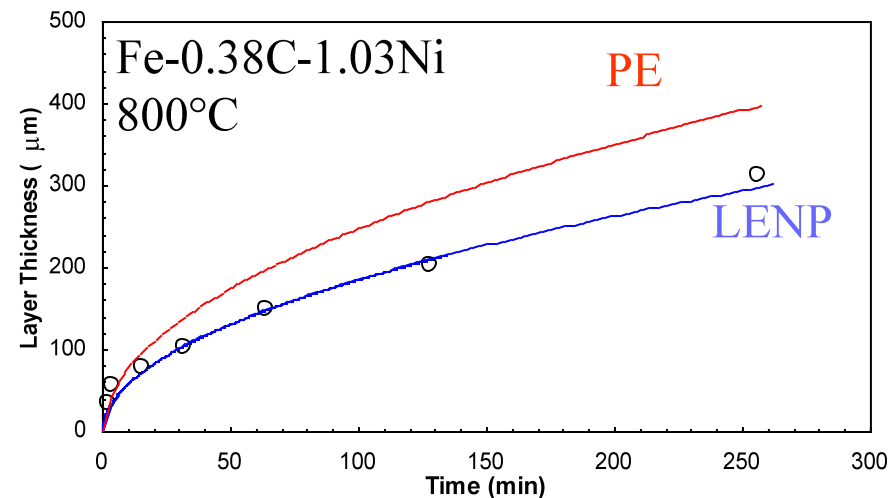
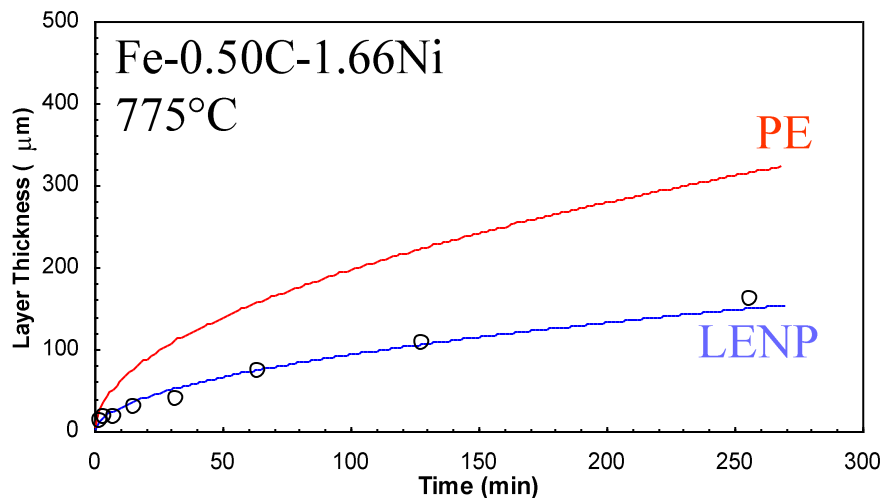
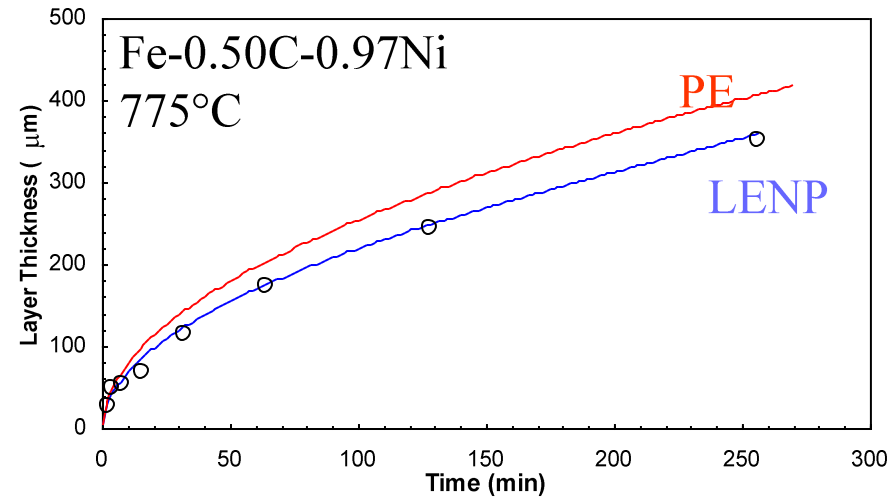
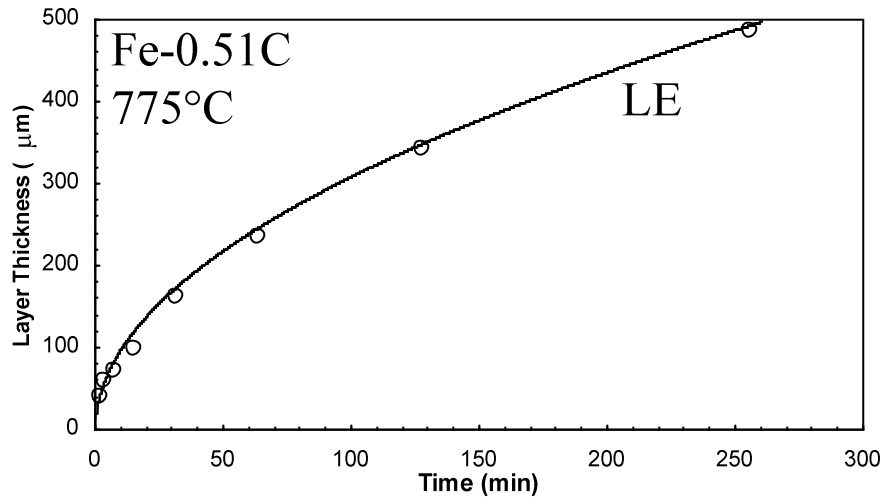


T (°C)	D/D _A
775	1
806	0.8
825	0.8
850	0.8

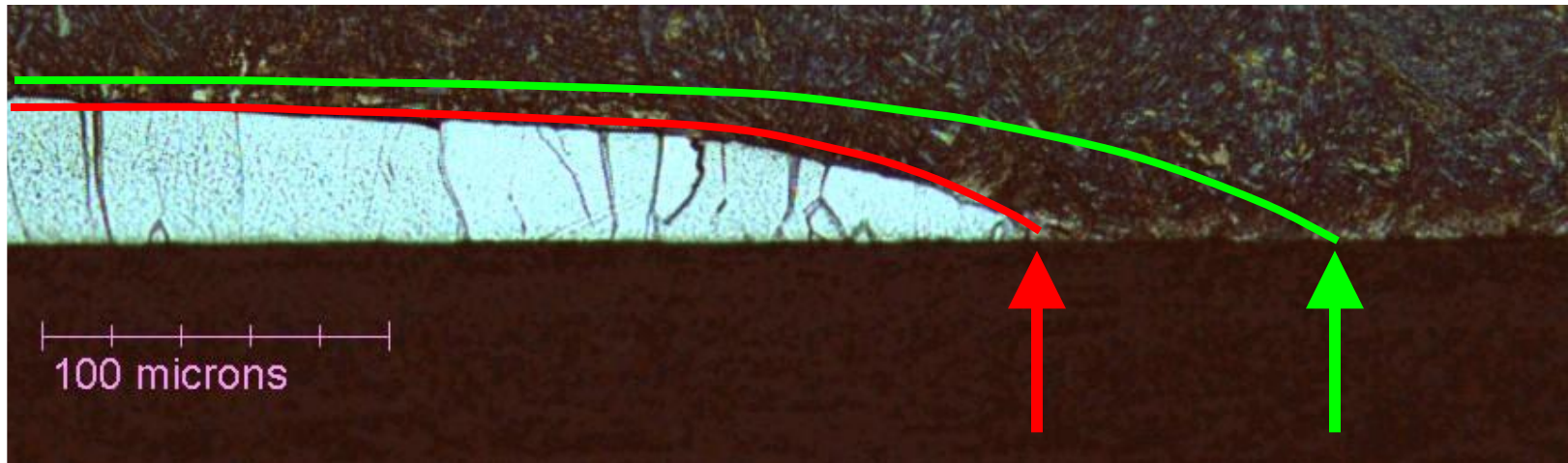
$$D_C^\alpha = 2.63 \times 10^{-10} m^2 / s \quad vs. \quad D_C^\alpha = 2.68 \times 10^{-10} m^2 / s$$

2.2. Fe-C-Ni Decarburization Experiments (ideal case)

Phillion, Zurob, Hutchinson, Guo, Malakhov, Nakano and Purdy, *Metall Trans.*, 35A, 1237-1242, 2004.



Fe-Ni diffusion couple: 775°C for 4 min.



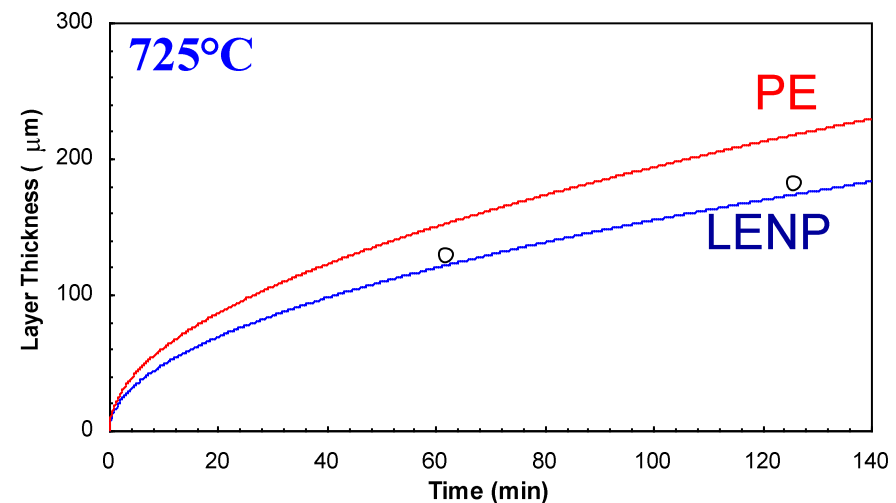
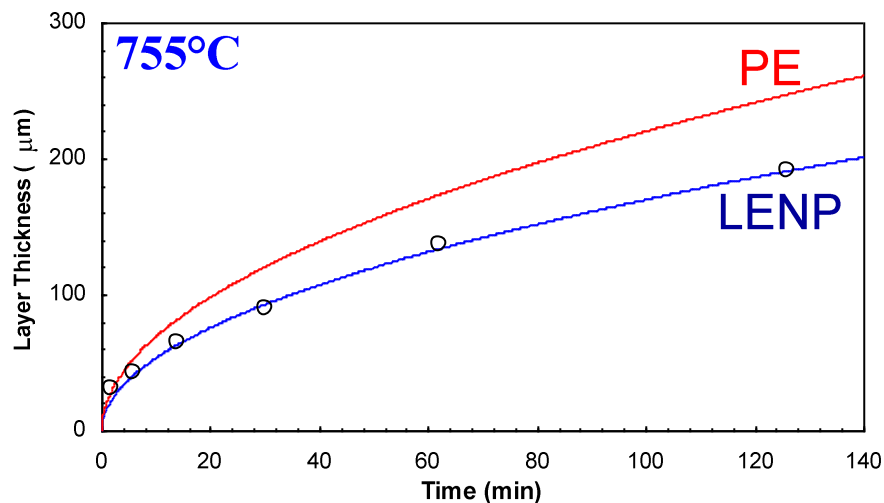
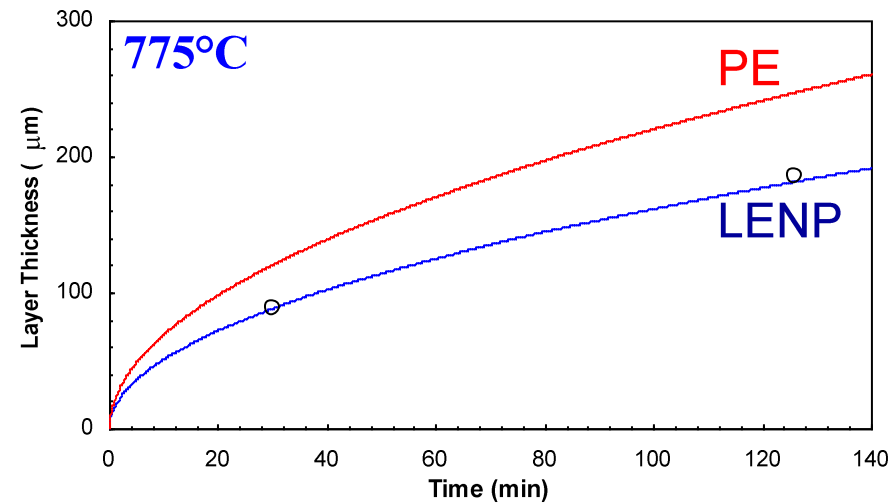
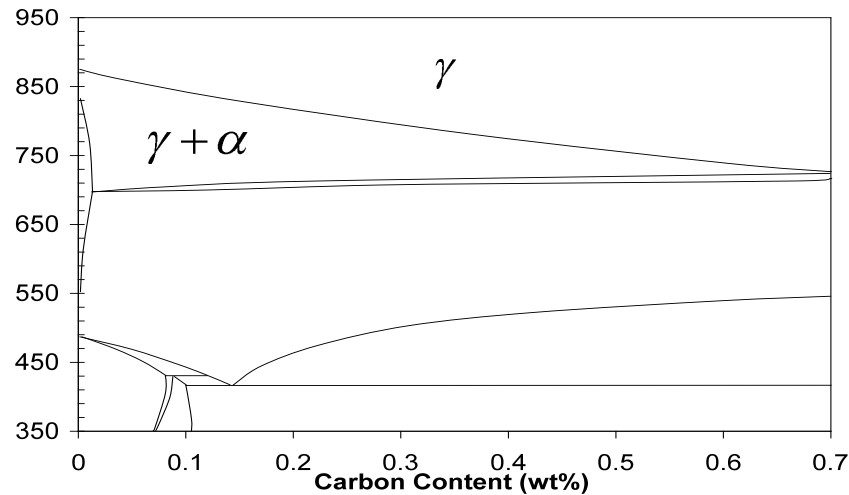
1.95%

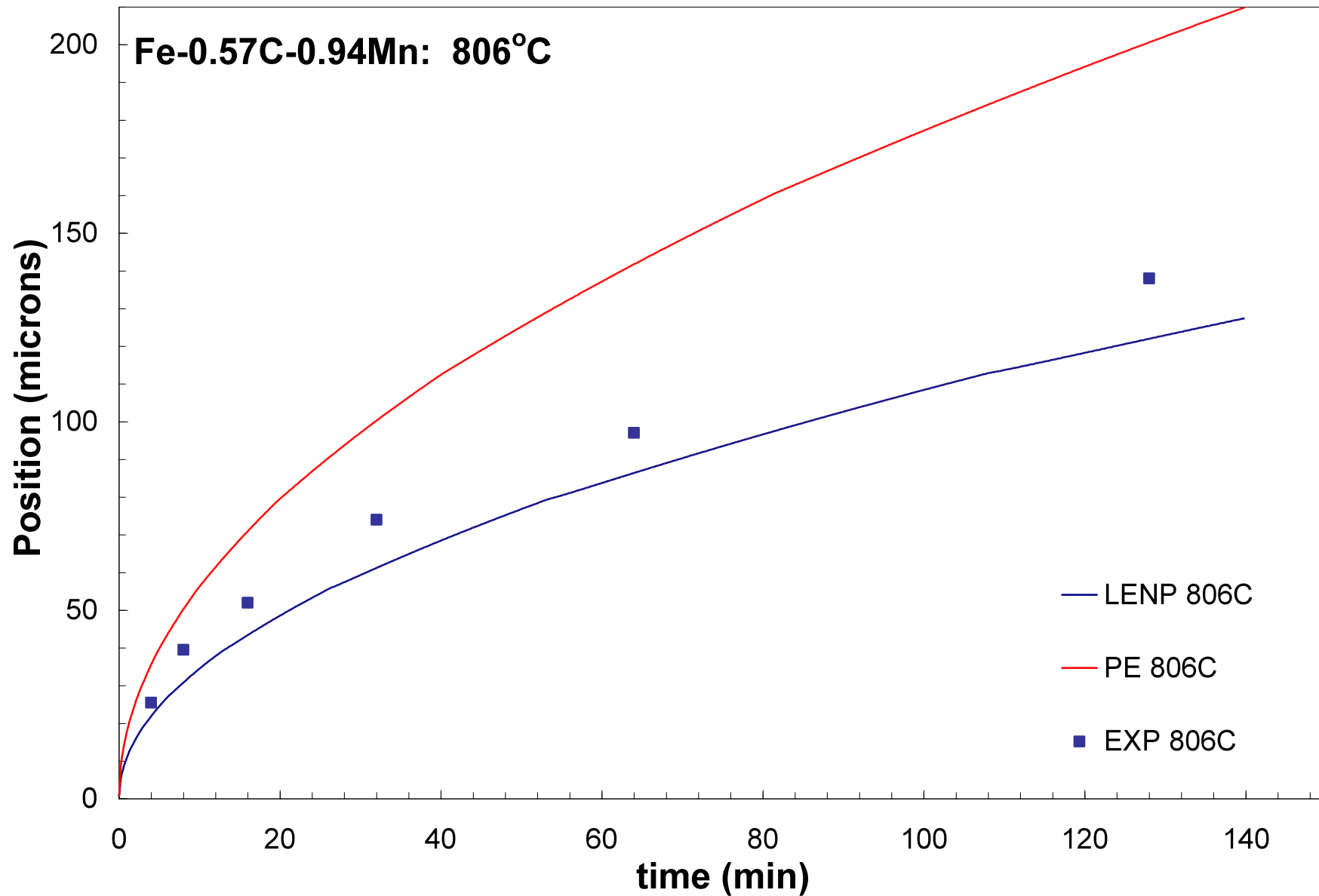
2.85%

1.89 +/- 0.33

2.3. Fe-C-Mn Decarburization Experiments

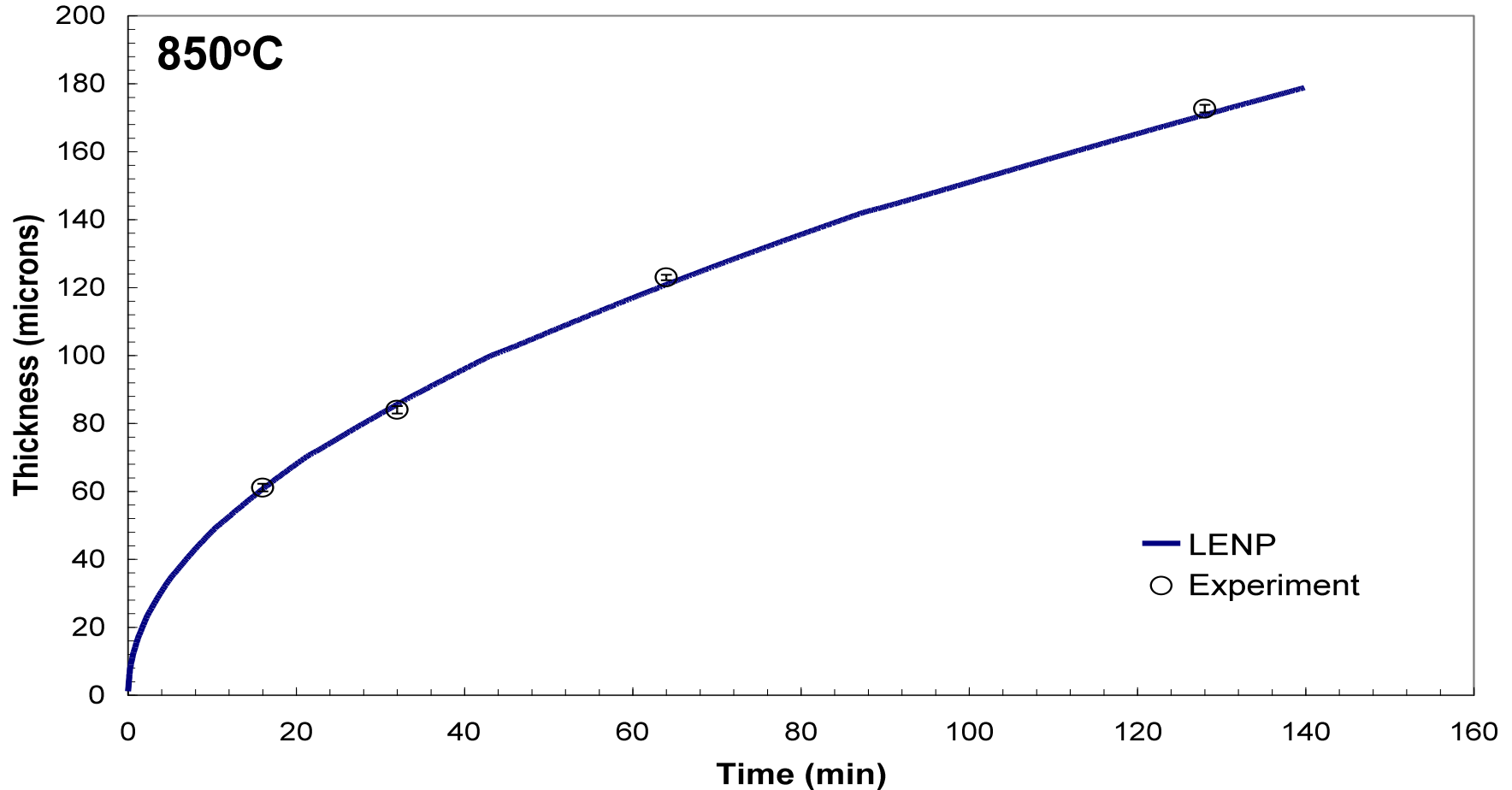
Fe-0.57C-0.94Mn (wt. %)





2.4. Fe-C-Cr Decarburization Experiments

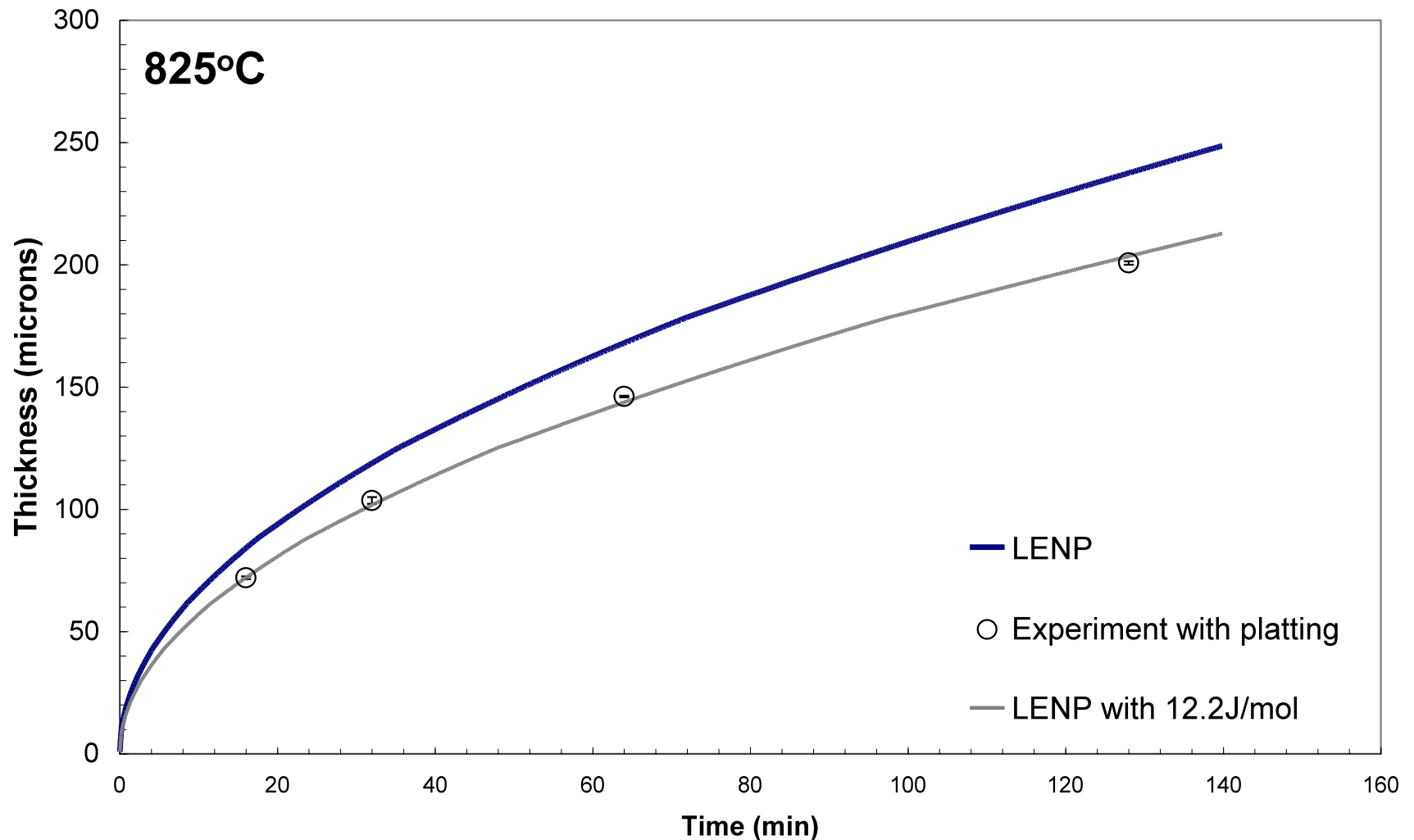
Fe-0.58C-1.99Cr (wt. %)



A. Beche, senior thesis.

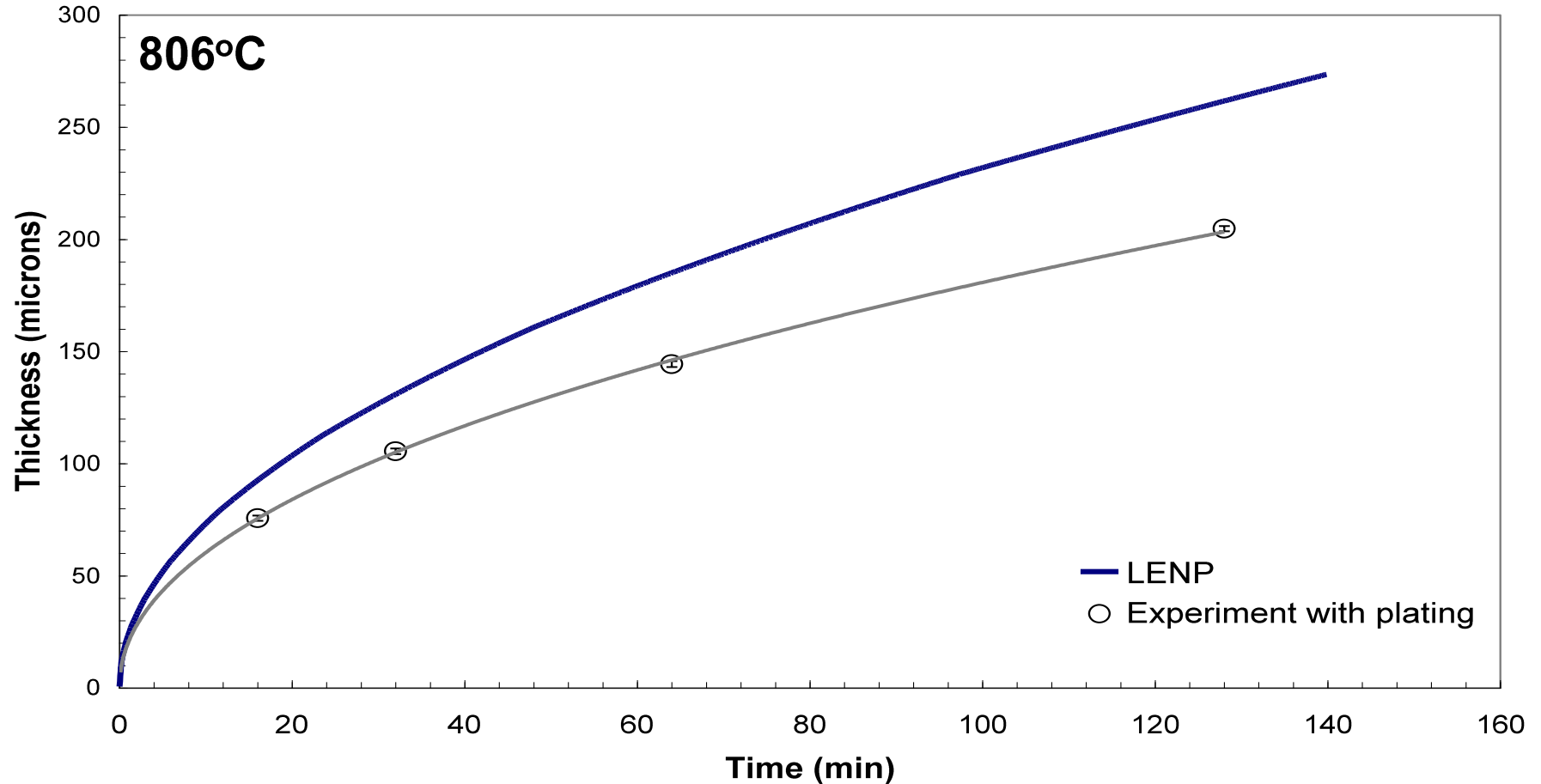
2.4. Fe-C-Cr Decarburization Experiments

Fe-0.58C-1.99Cr (wt. %)

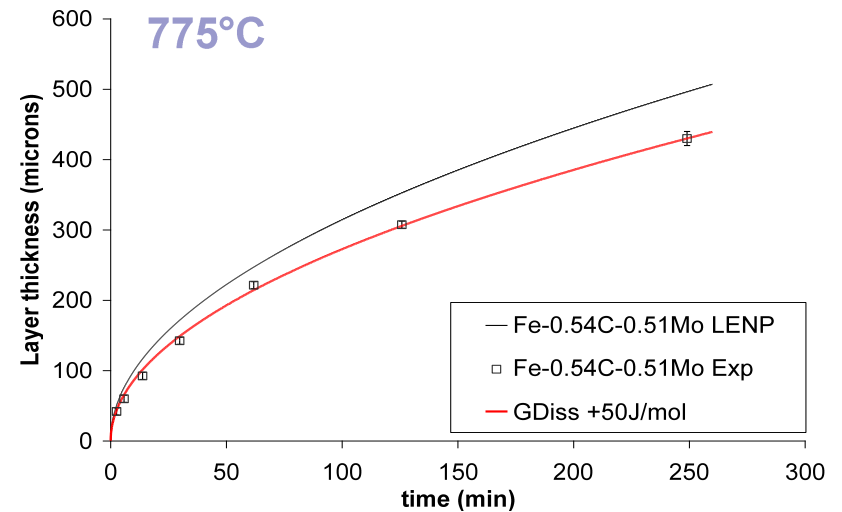
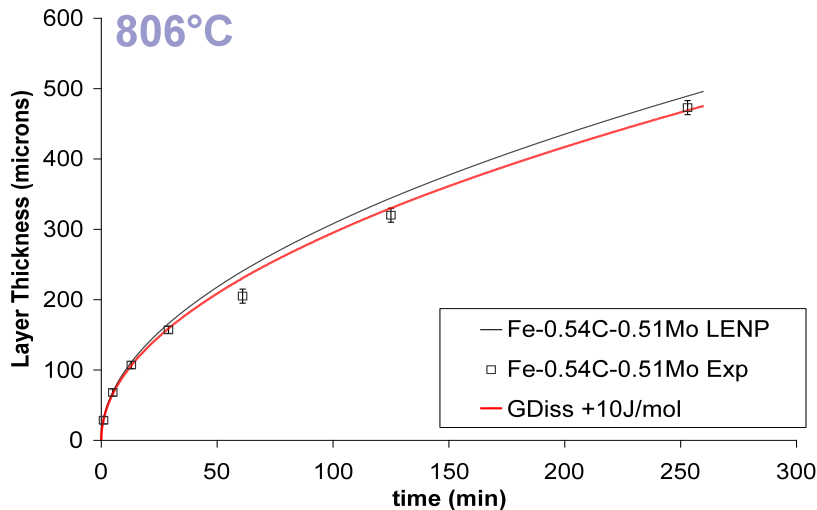
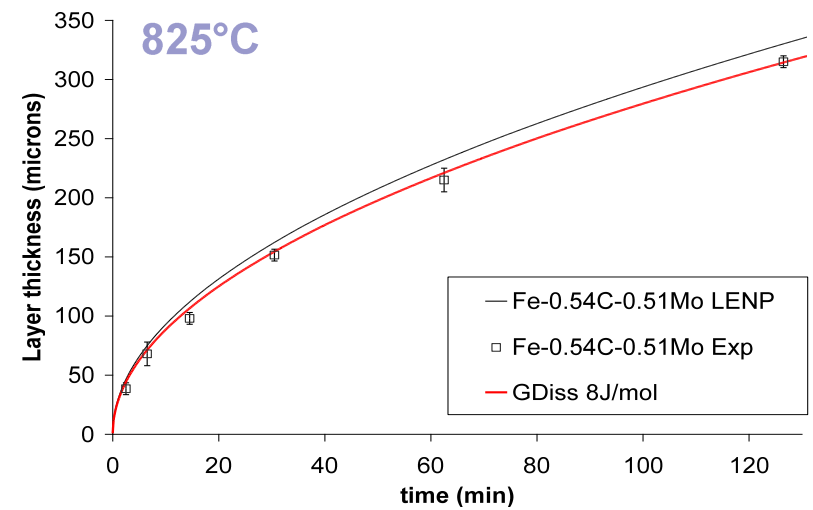
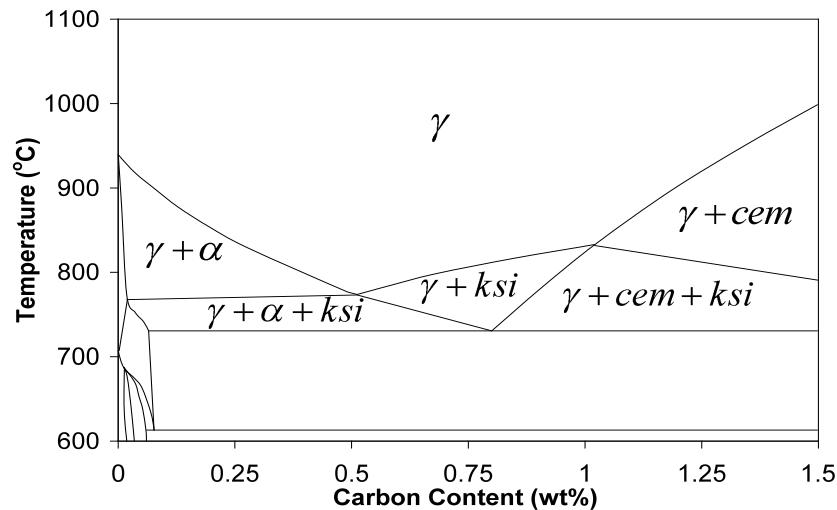


2.4. Fe-C-Cr Decarburization Experiments

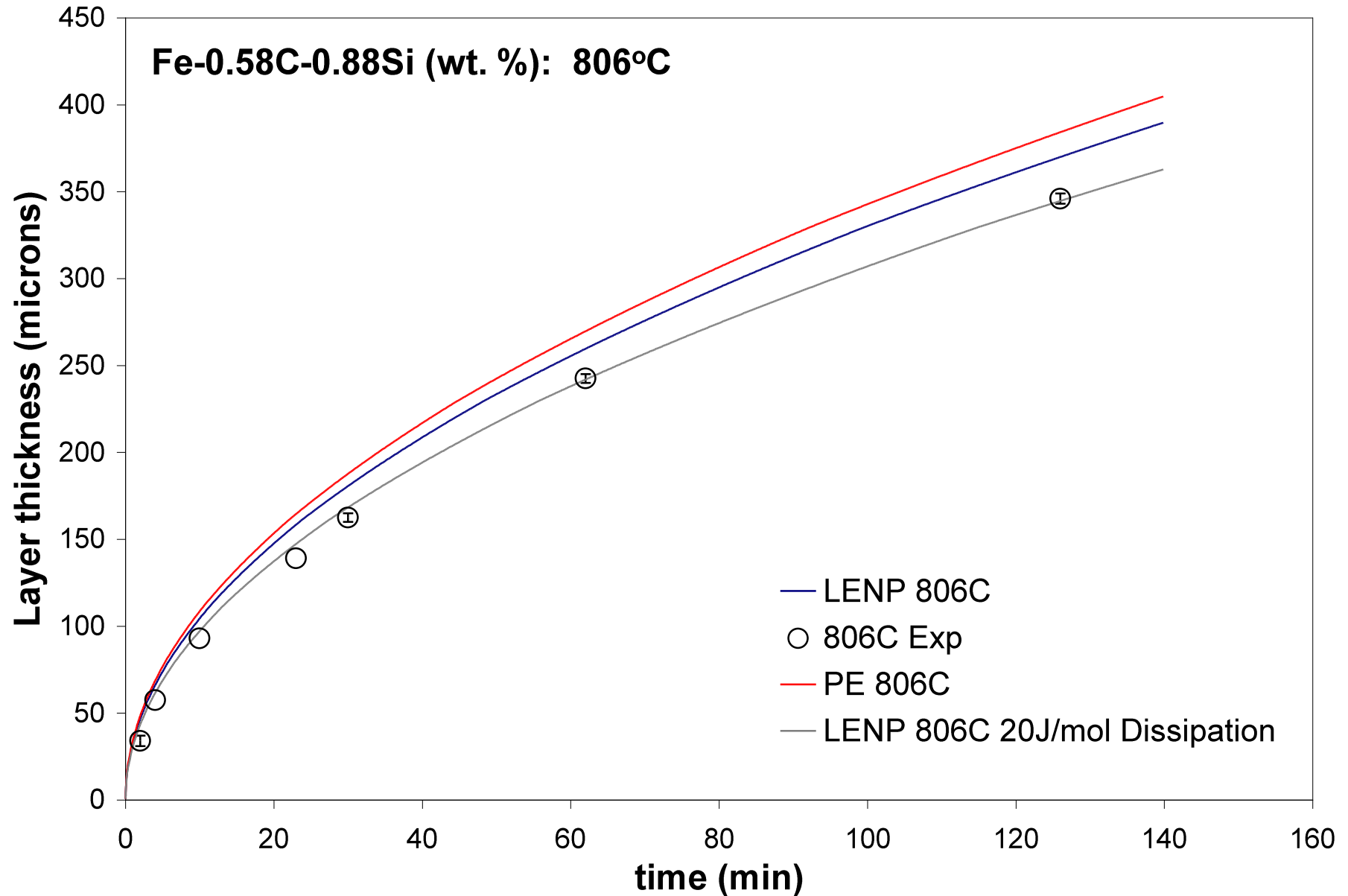
Fe-0.58C-1.99Cr (wt. %)



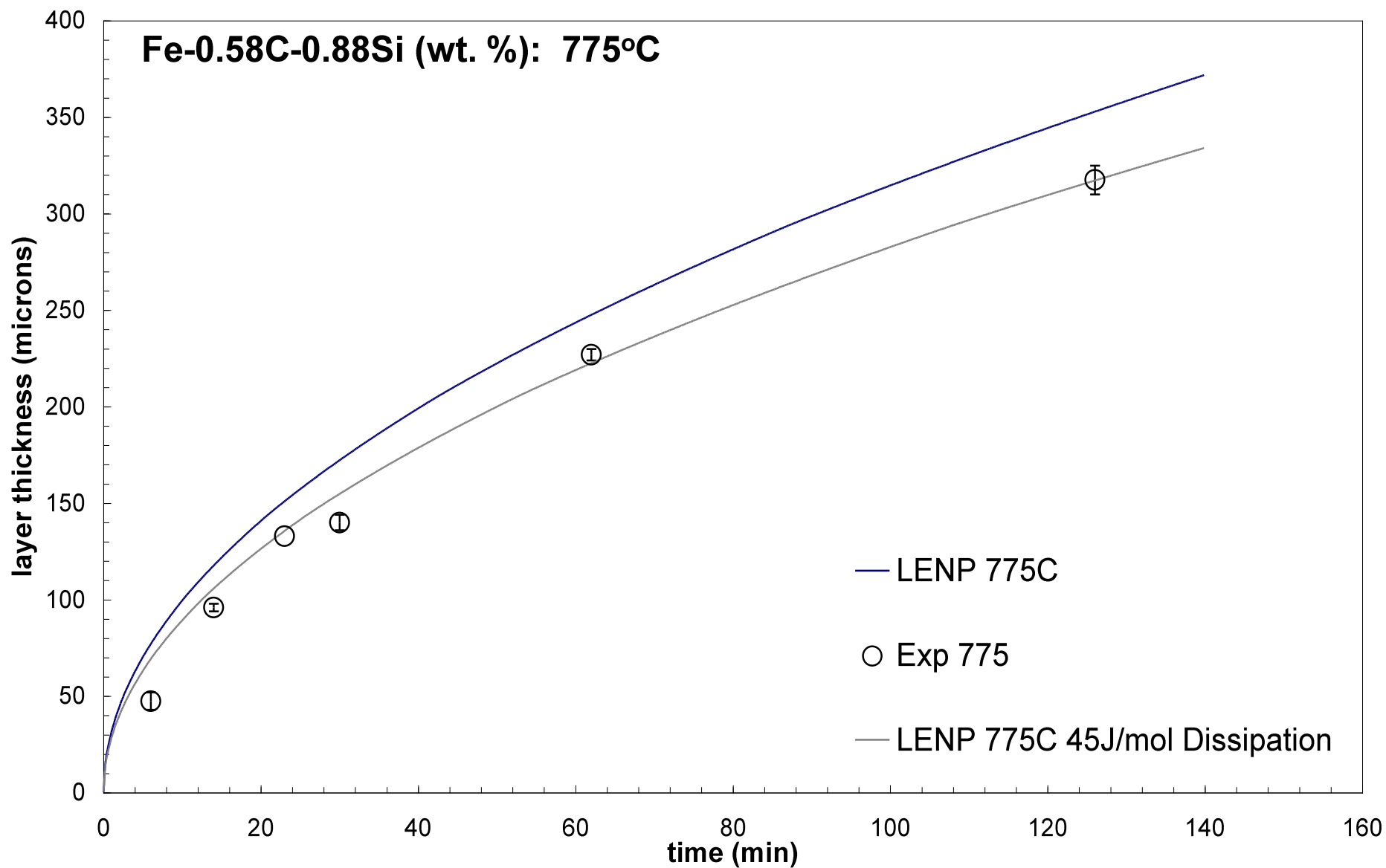
2.5. Fe-C-Mo Decarburization Experiments



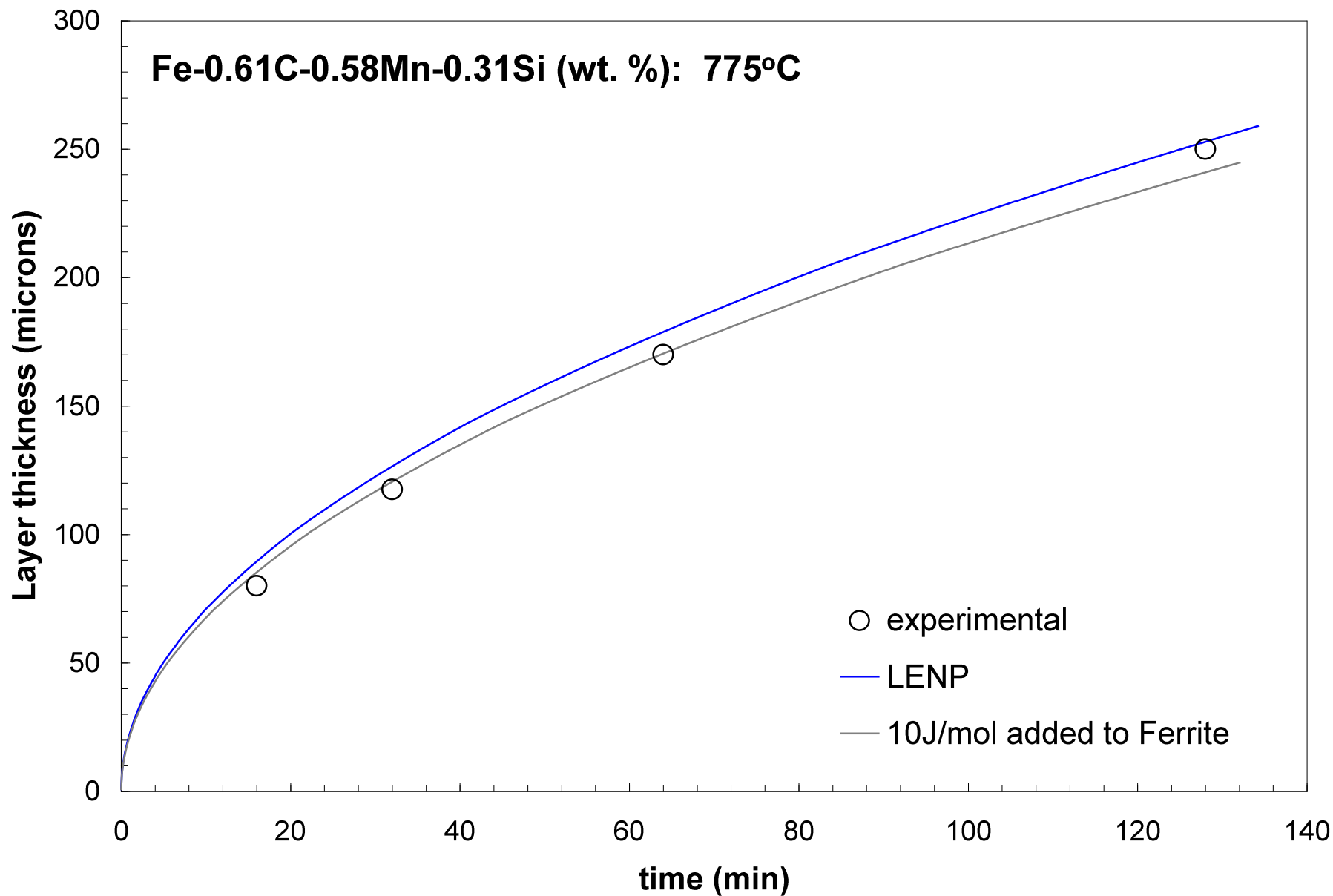
2.6. Fe-C-Si Decarburization Experiments



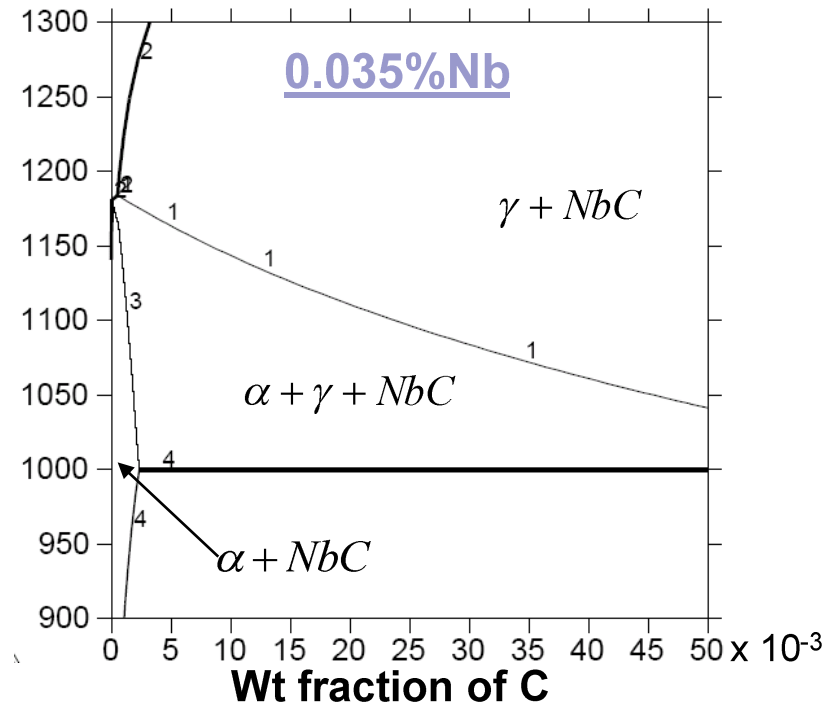
Fe-0.58C-0.88Si (wt. %): 775°C



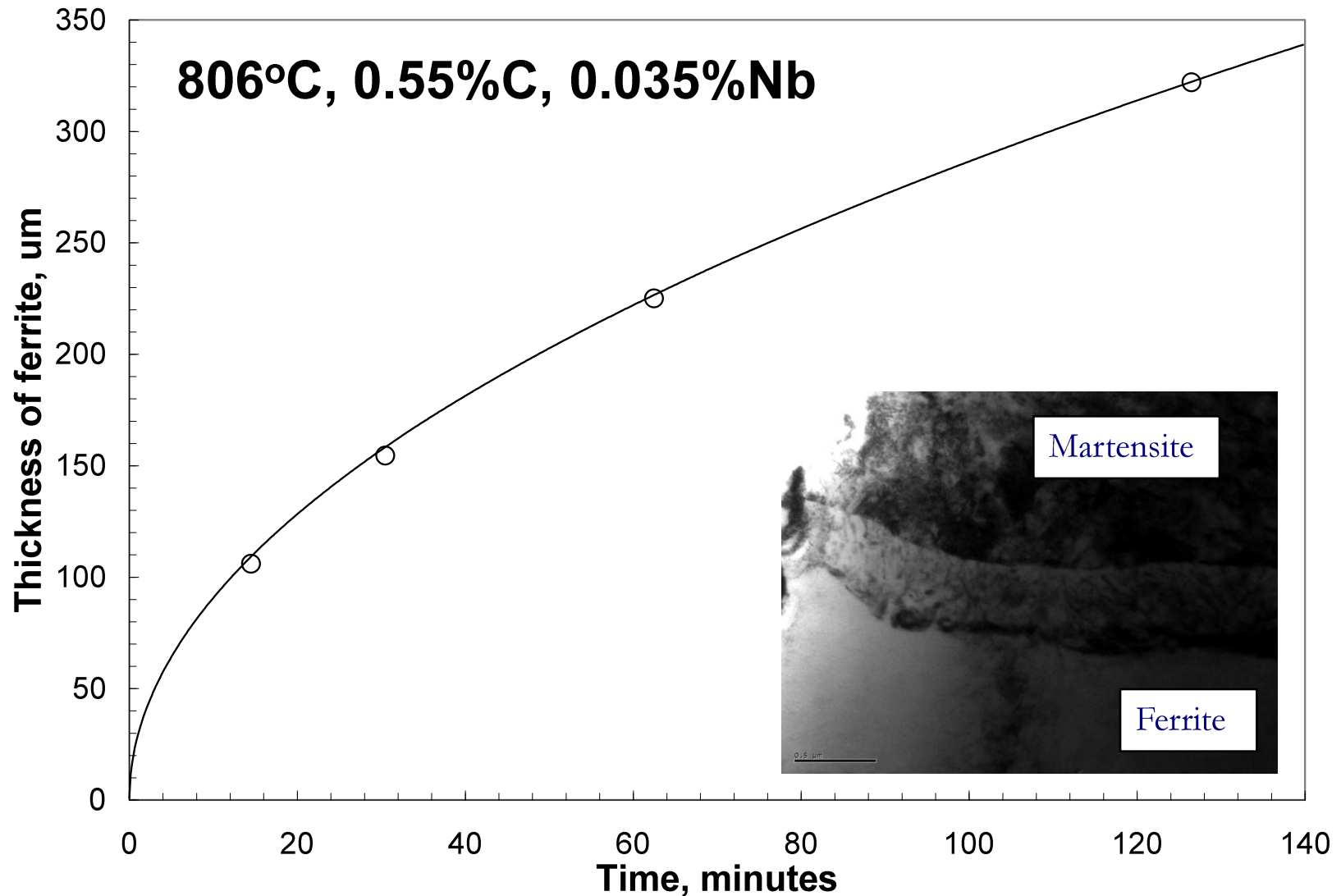
Fe-C-Si-Mn Decarburization Experiments



2.7. Fe-C-Nb Decarburization Experiments



2.7. Fe-C-Nb Decarburization Experiments



Summary

Alloy (wt. %)	Treat. Temp (°C)	Interfacial Compositions	Exp. observed excess dissipation	
			(J/mole)	(MPa)
Fe-0.50C-0.97Ni	775	LENP	-	-
Fe-0.50C-1.66Ni	775	LENP	-	-
Fe-0.38C-1.03Ni	800	LENP	-	-
Fe-57C-0.94Mn	806	LENP+	-	-
	775	LENP	-	-
	755	LENP	-	-
	725	LENP	-	-
Fe-0.54C-0.51Mo	825	LENP-	8	1.1
	806	LENP-	10	1.4
	775	LENP-	50	7.4
Fe-0.58C-0.88Si	806	LENP-	20	2.9
	775	LENP-	45	6.4
Fe-0.61C- 0.58Mn-0.31Si	775	LENP-	10	1.4
Fe-0.56C-1.97Cr	850	LENP	-	-
	825	LENP-	12	1.7
	806	LENP-	30	4.2

3. Future Work:

- Search for Transition in Fe-Mn-C at high temperatures.
- More detailed investigation of the quaternary Fe-C-Mn-Si.
- One more temperature in Fe-C-Cr.
 - Question: If we use 775°C, will precipitation be a problem?
- Denitriding of Fe-N-Cr down to 675°C.
 - Question: will denitriding be controlled by N diffusion? How can we check?