

Kinetic Transitions in Fe-Mn-C Alloys

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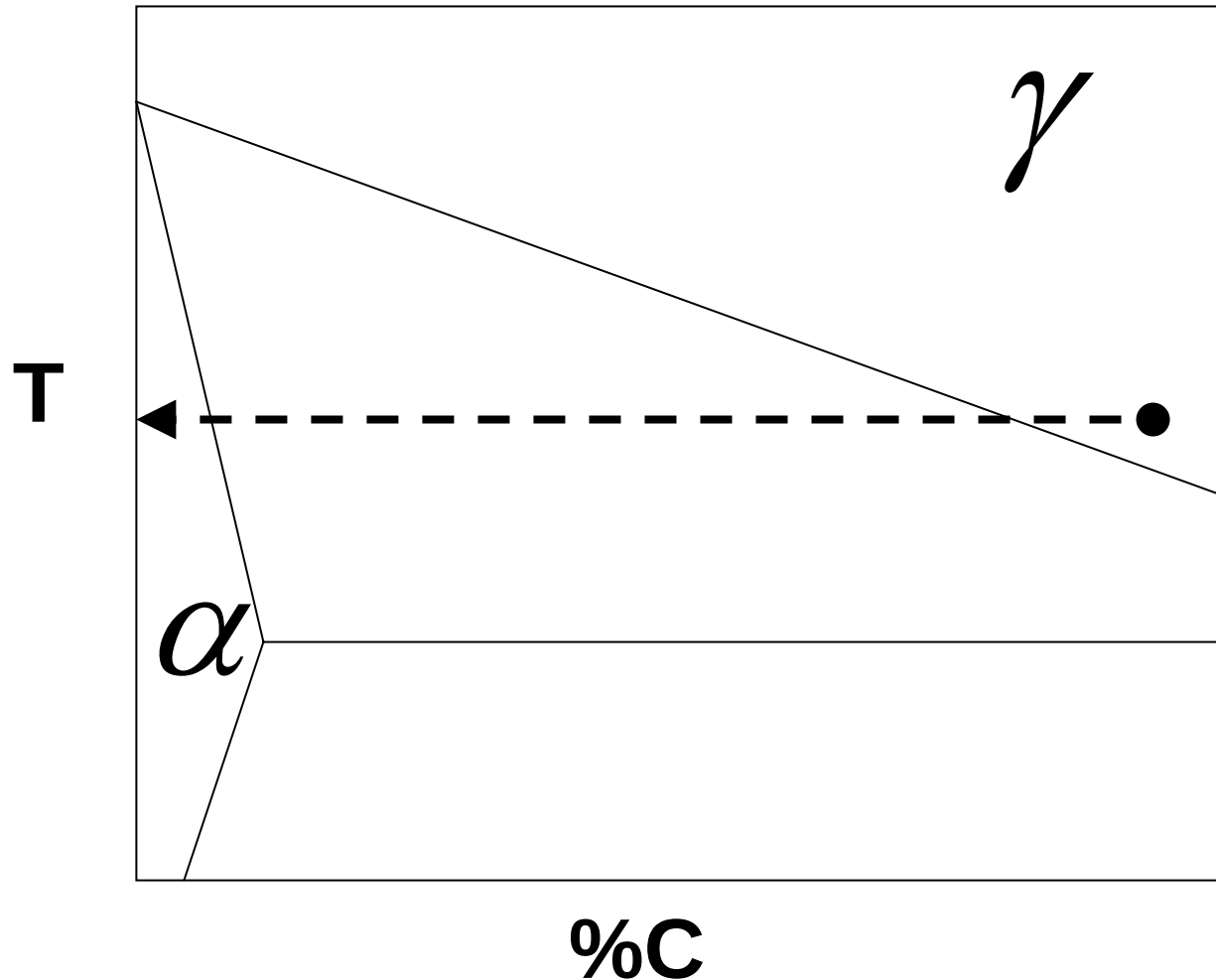


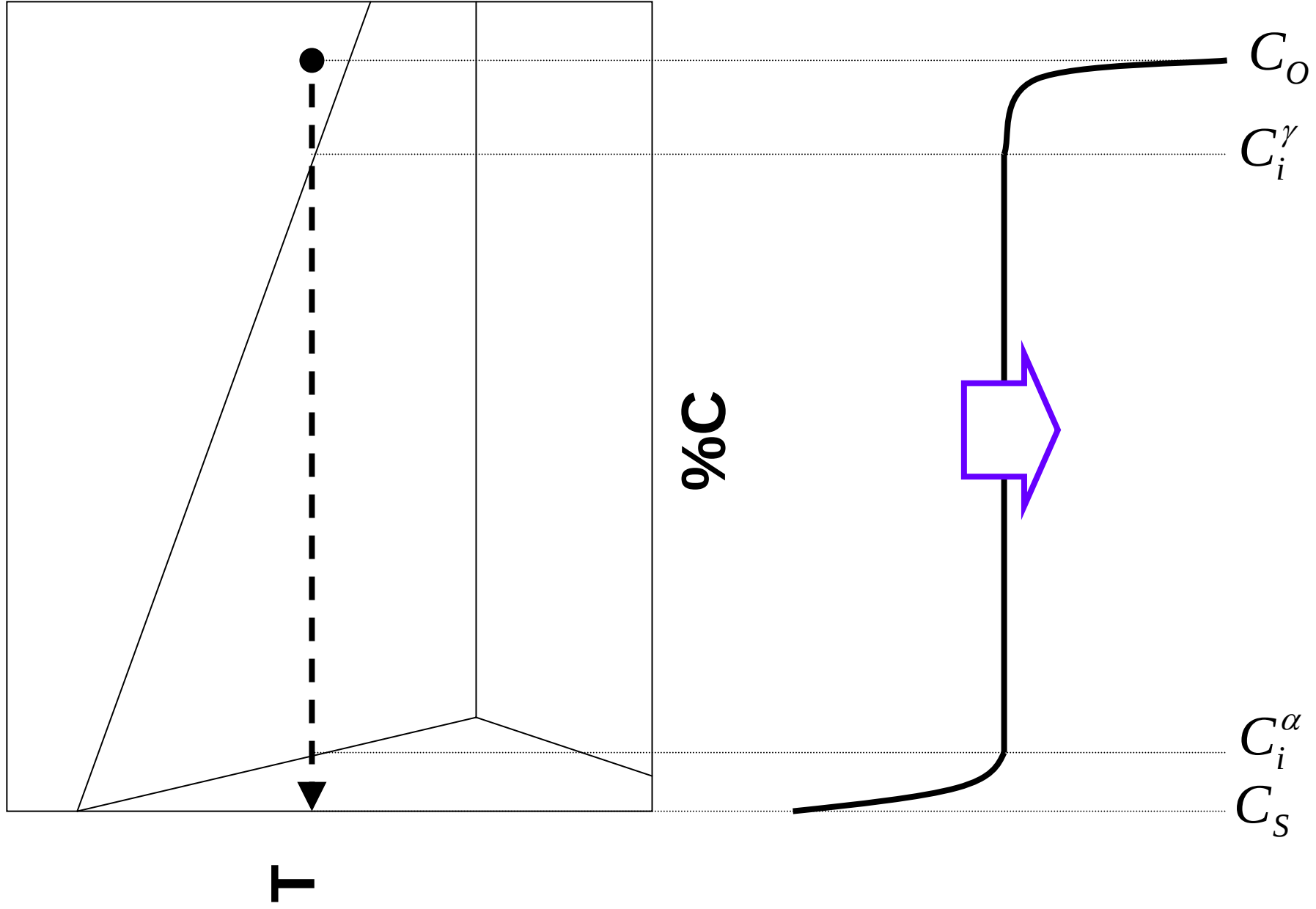
Outline

- Decarburization as a method of studying the $\gamma \rightarrow \alpha$ transformation.
- Summary of results on Fe-Ni-C.
- New results on Fe-Mn-C
 - Fe-1%Mn
 - Fe-0.5%Mn
 - Fe-2%Mn
- Possible Interpretations
- Conclusions

The Decarburization Approach:

> Background:





The rate of interface motion is given by:

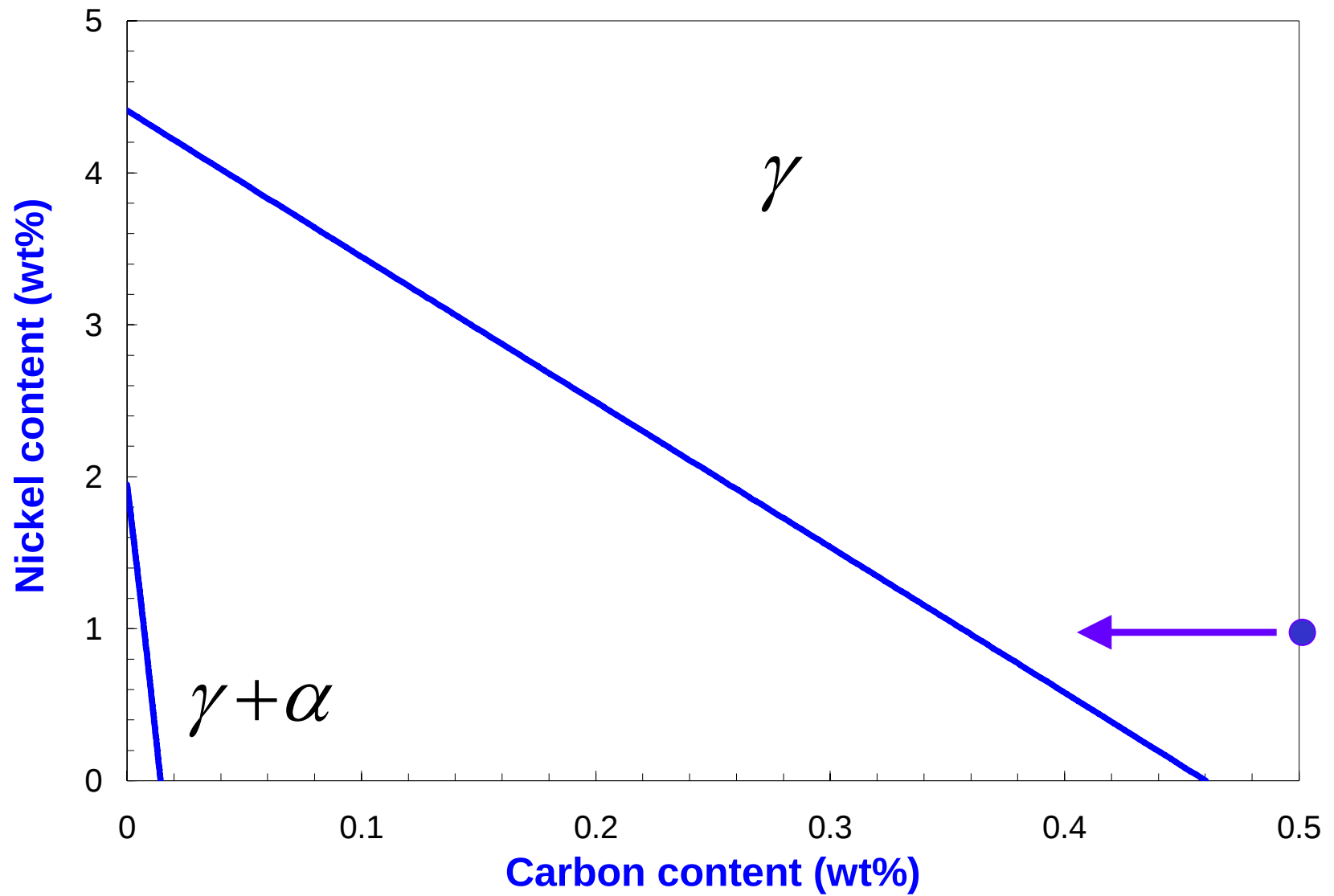
$$\frac{dz}{dt} = \frac{J_i^\alpha - J_i^\gamma}{C_i^\gamma - C_i^\alpha}$$

This differential equation has an analytical solution of the form:

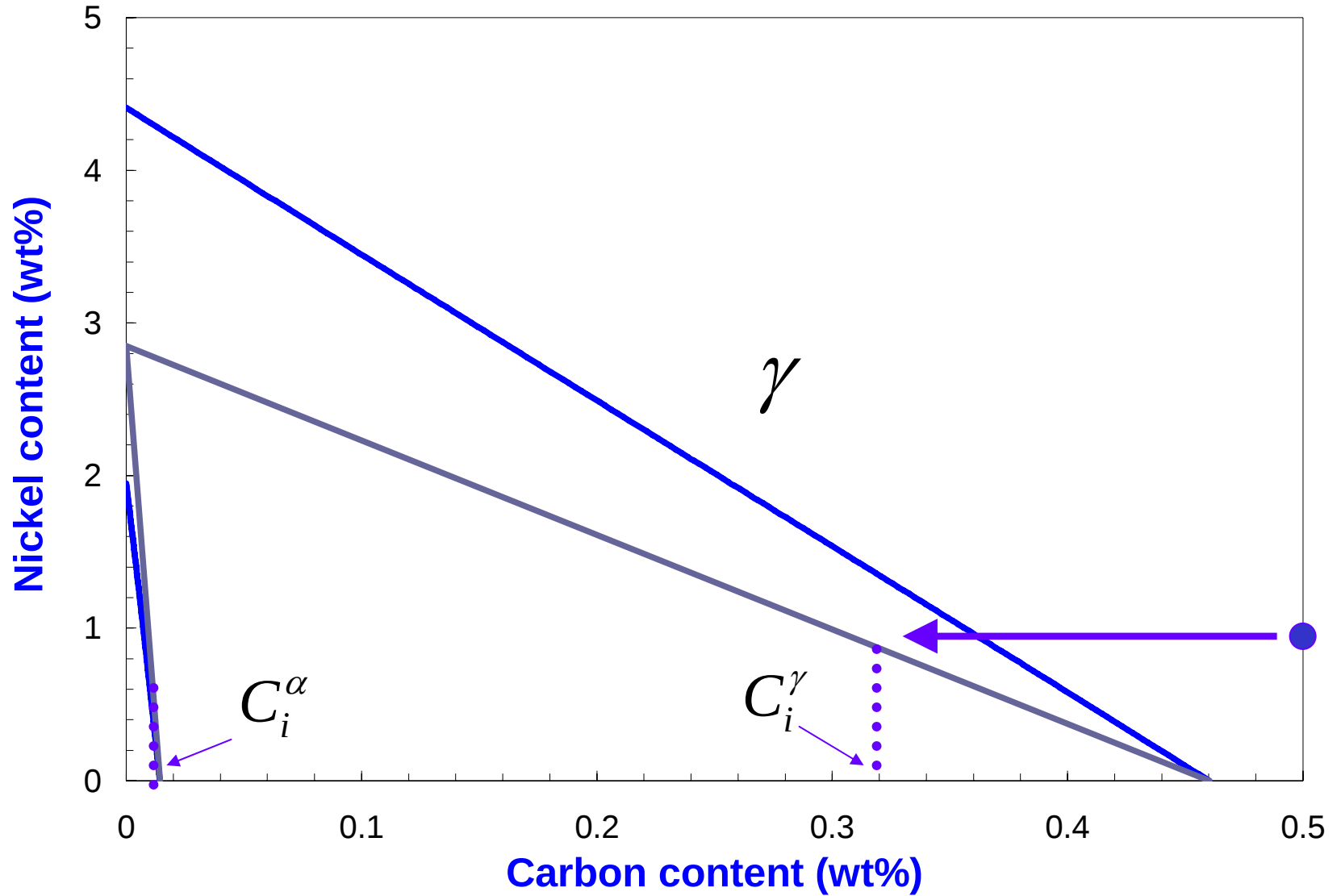
$$z = B\sqrt{t}$$

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

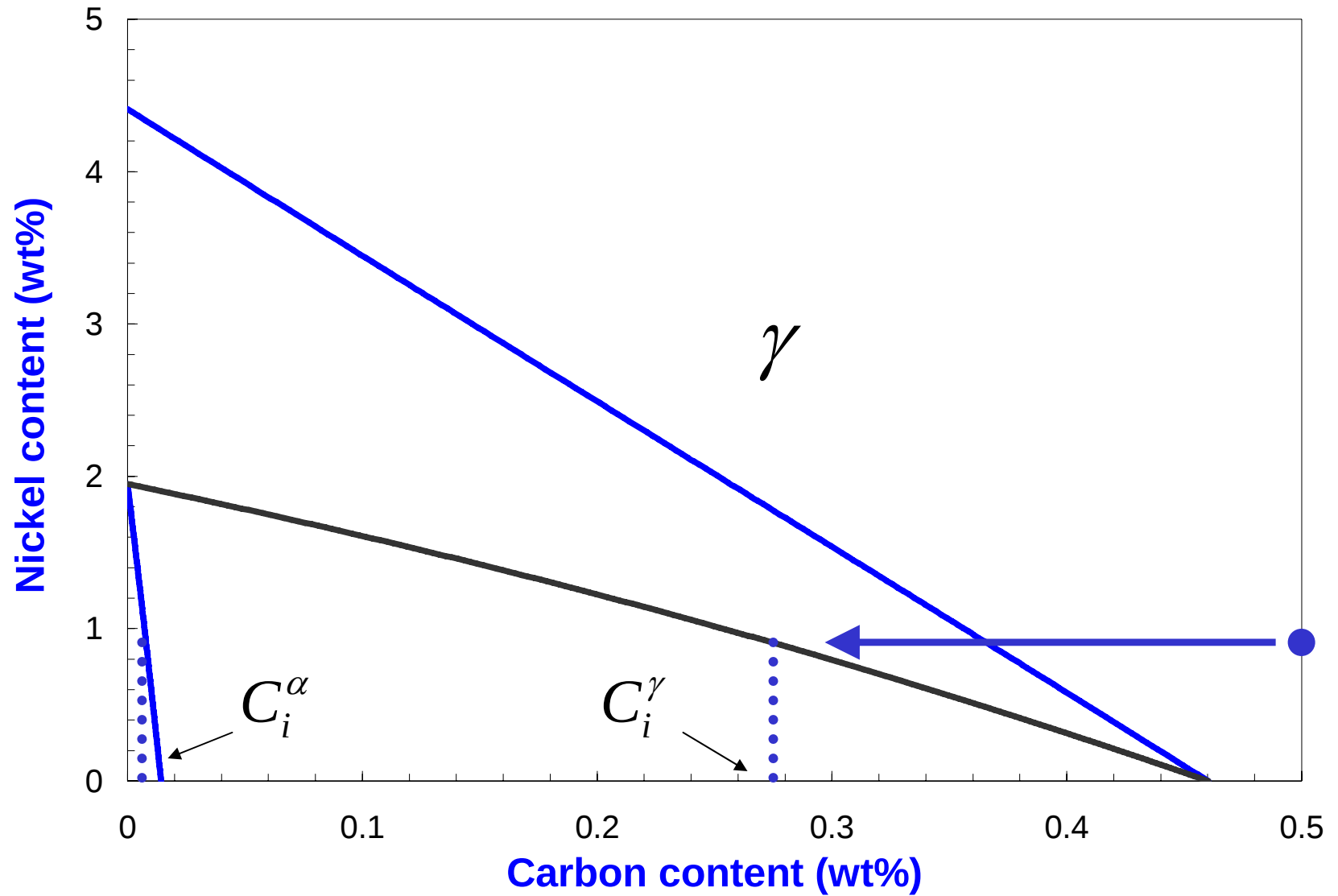
> Ternary Alloys.



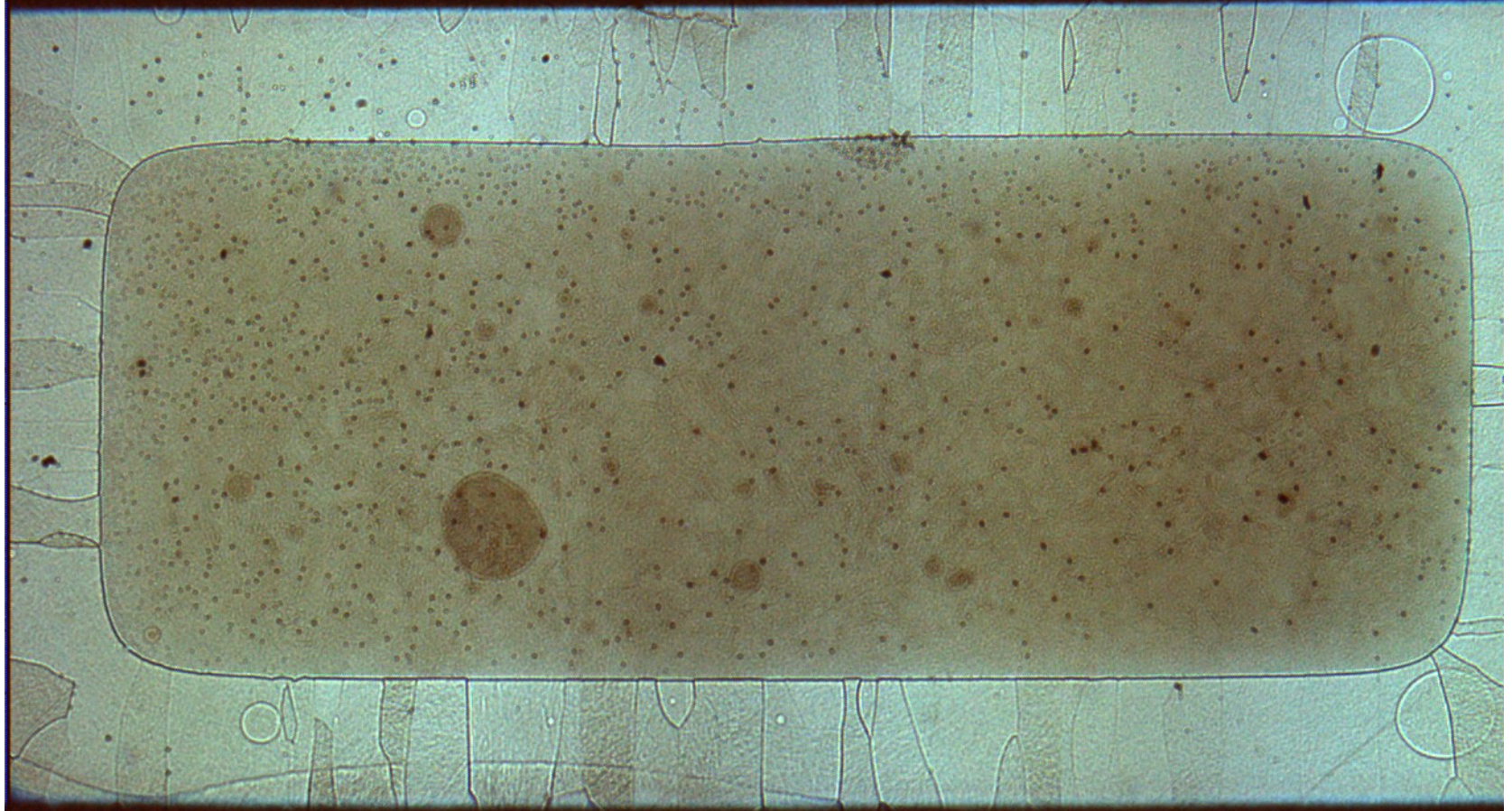
> Ternary Alloys: ParaEquilibrium Limit.



> Ternary Alloys: NPLE Limit.



1 mm

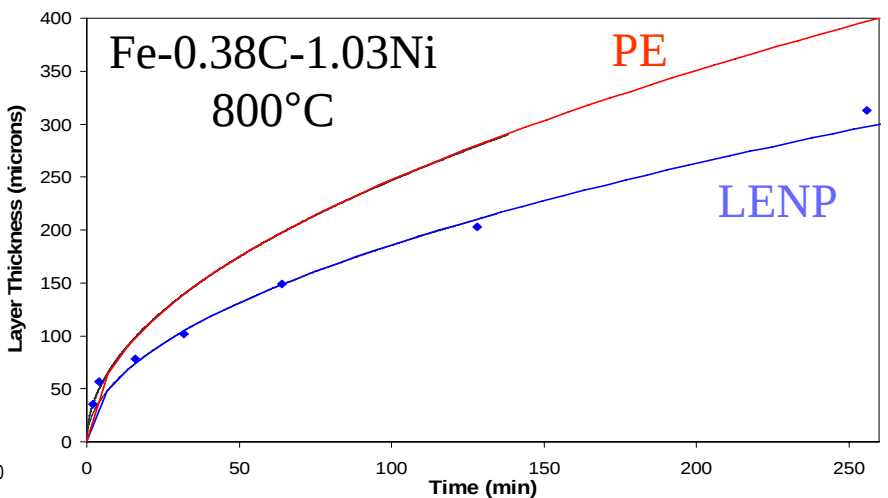
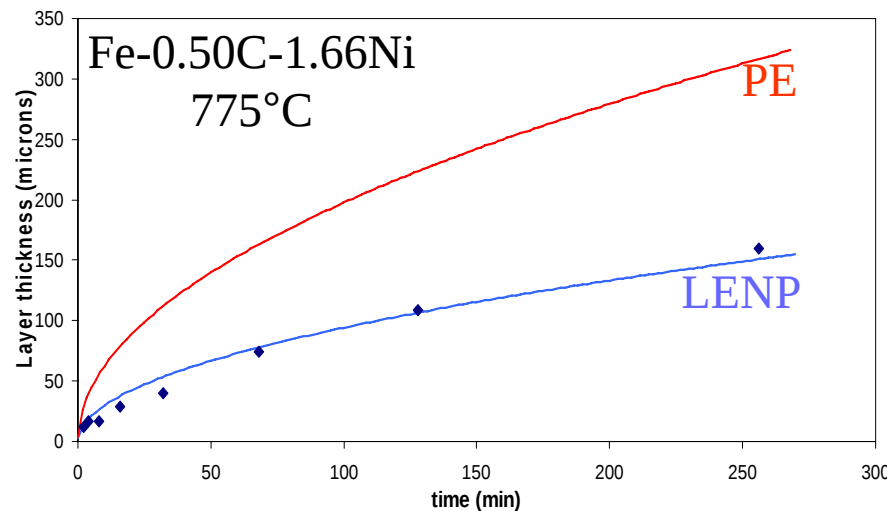
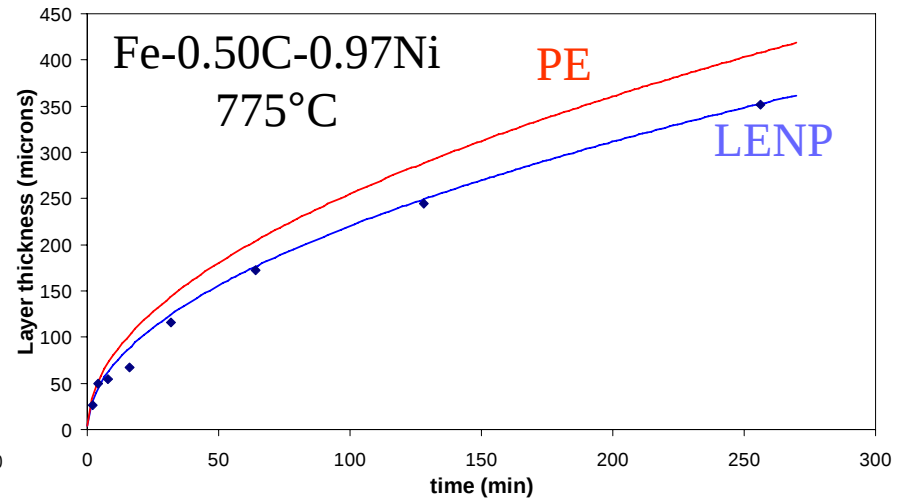
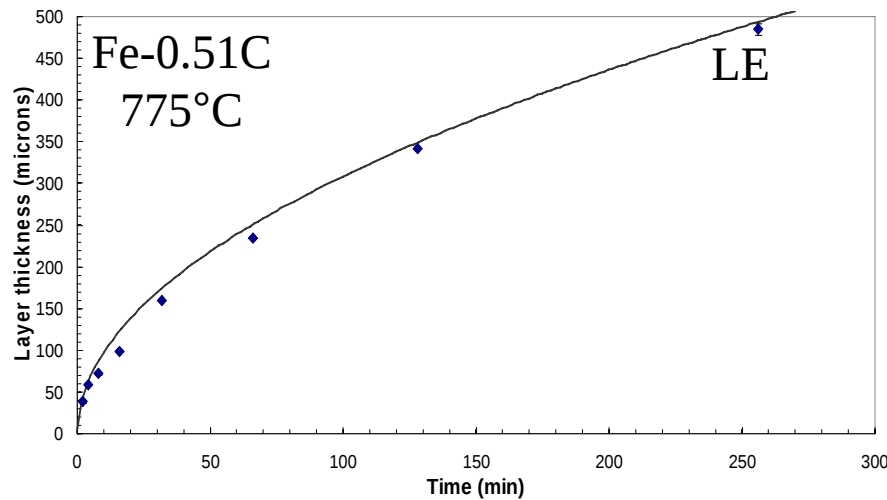


Summary of Results on Fe-Ni-C

- ▣ Fe-Ni-C alloys seems to follow LE-NP kinetics at all the temperatures and compositions investigated

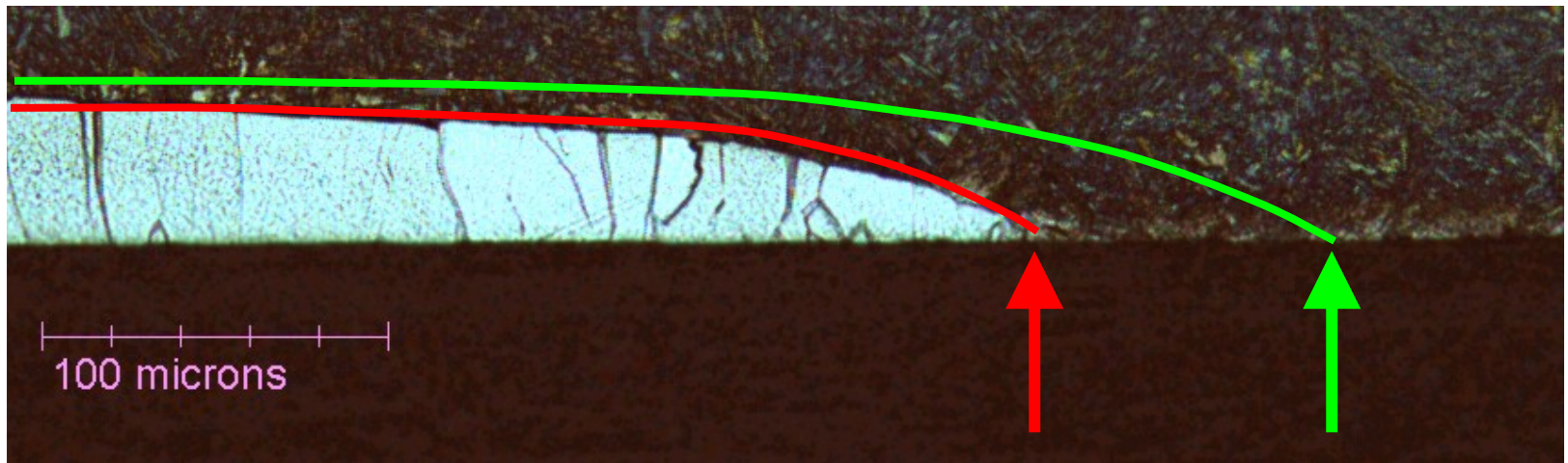
Fe-C-Ni: decarburization kinetics

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



Fe-C-Ni: diffusion couples

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



775°C for 4 min.

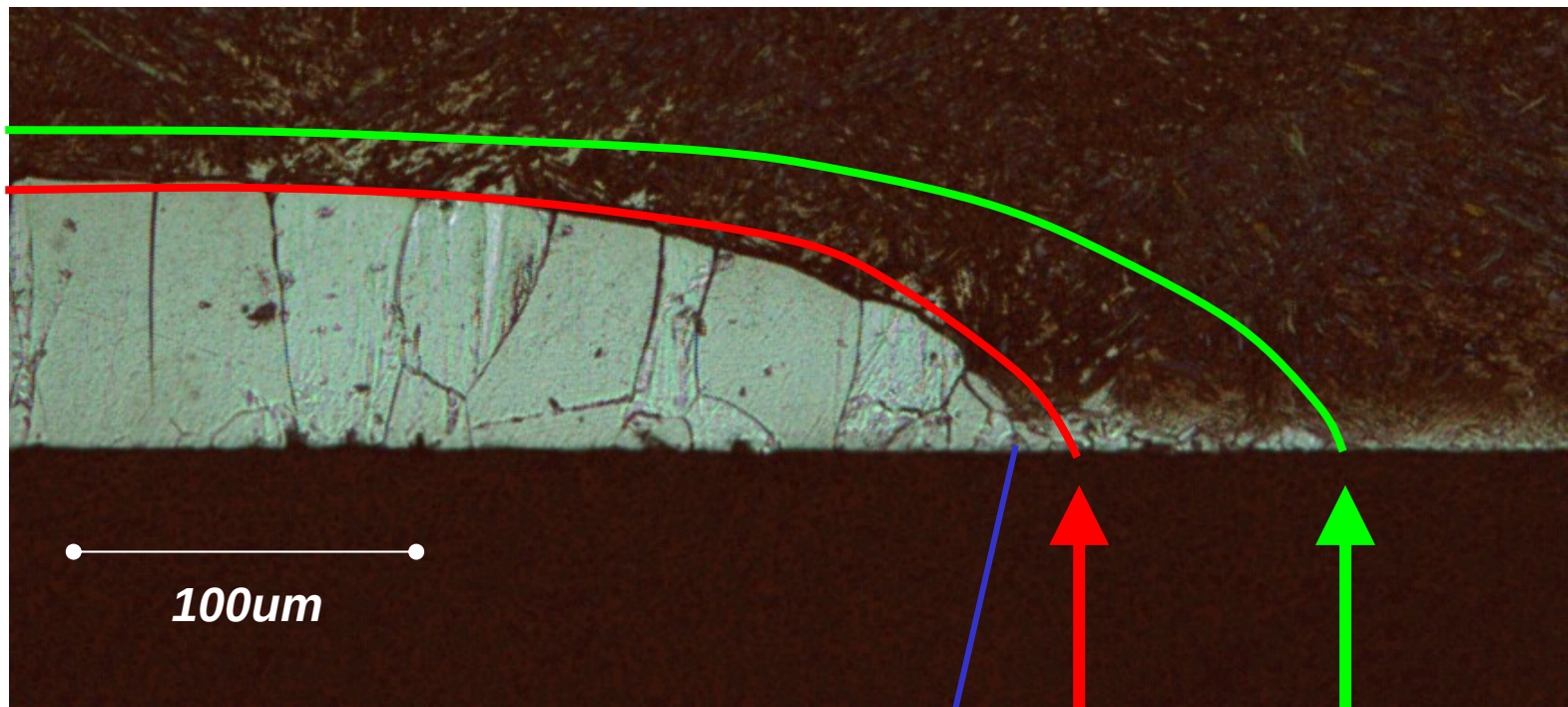
1.95%

1.89 +/- 0.33

2.85%

Fe-C-Ni: diffusion couples

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



775°C for 16 min

1.95%

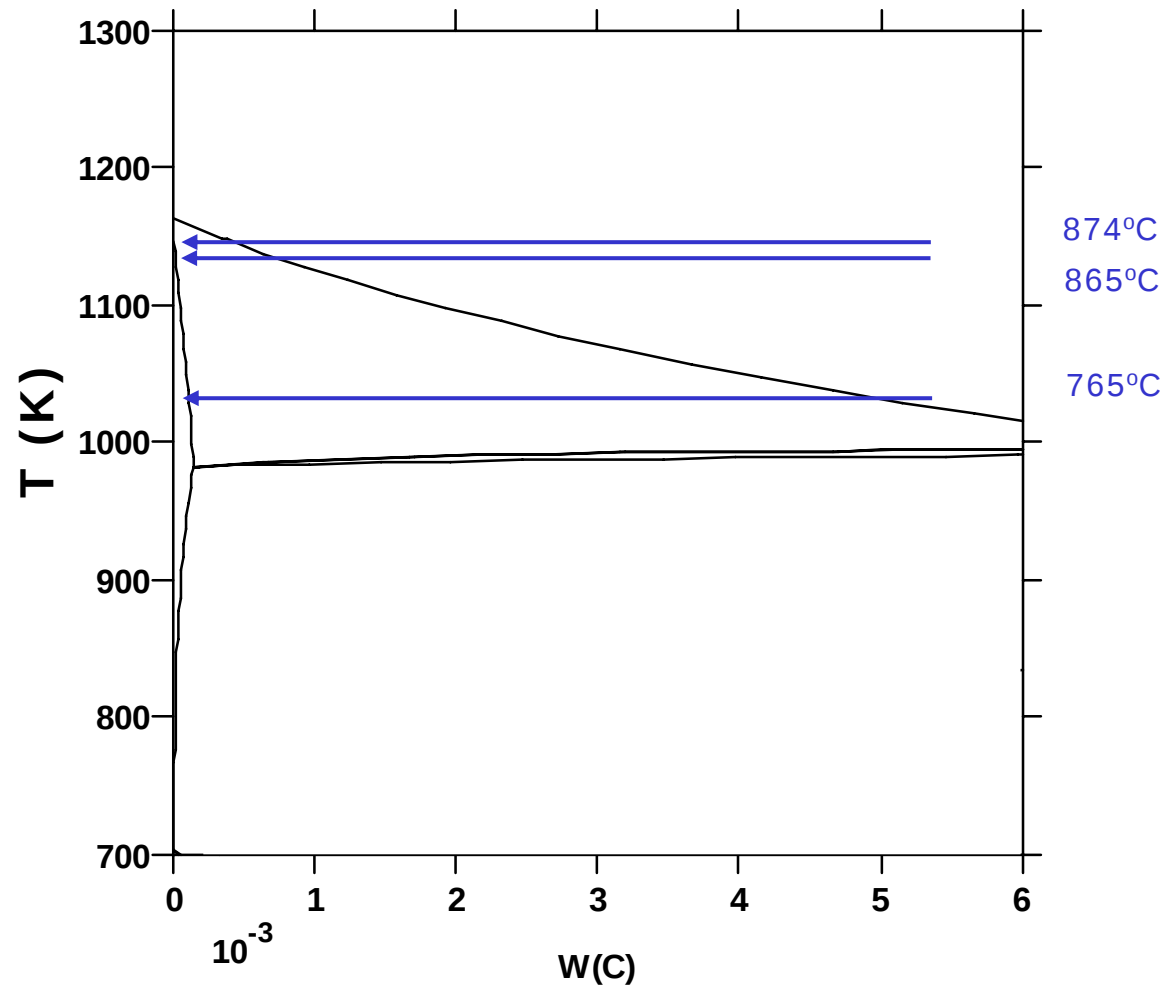
2.85%

1.85+/-0.33

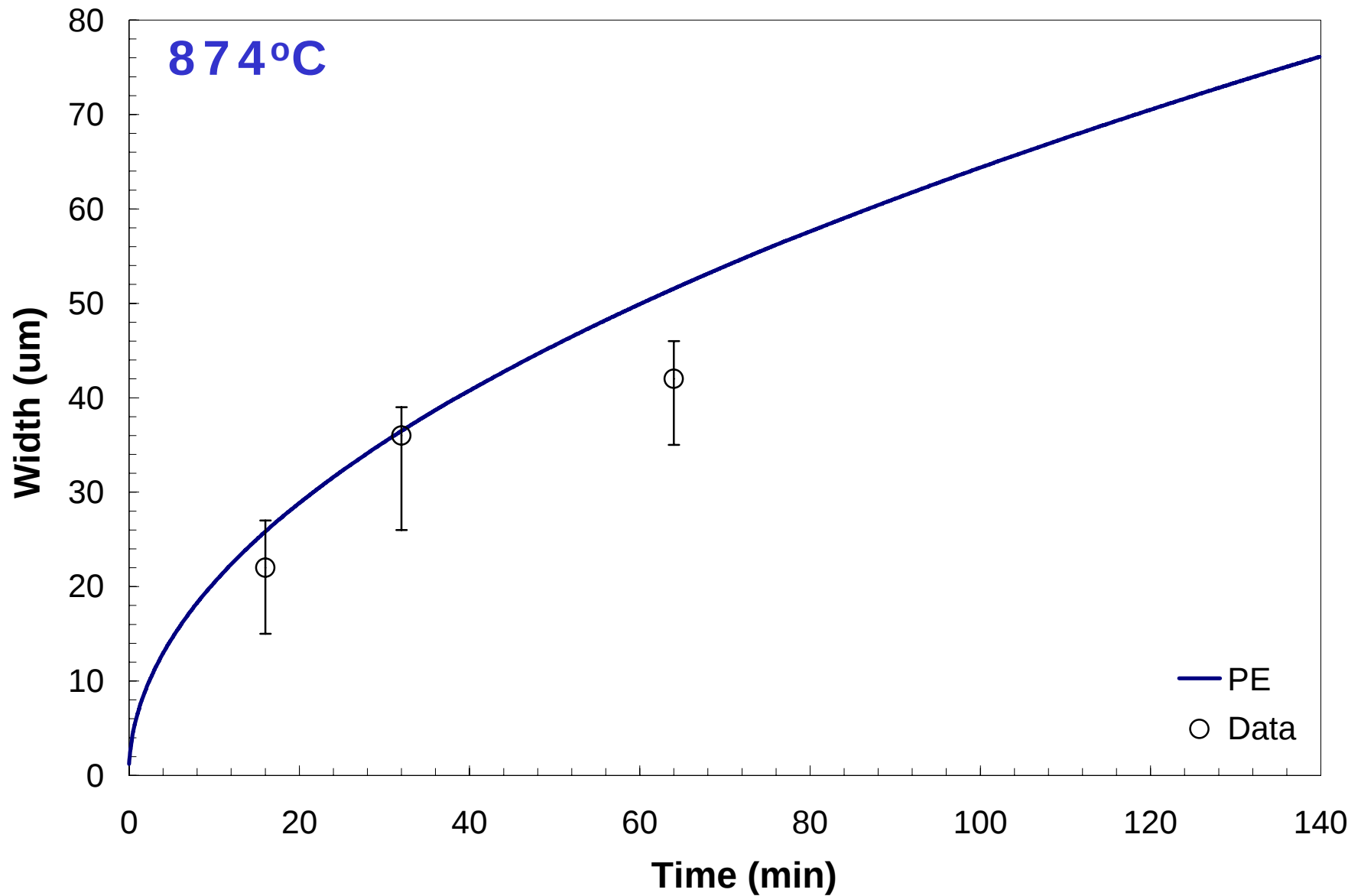
New Results on Fe-Mn-C

- ▣ Fe-0.5%Mn-C: Paraequilibrium at high Temperatures, unclear at low T.
- ▣ Fe-1%Mn-C: Definite transition from LENP to PE as the temperature increases.
- ▣ Fe-2%Mn-C: Ferrite forms above the LENP limit, but at a rate smaller than that predicted by PE.

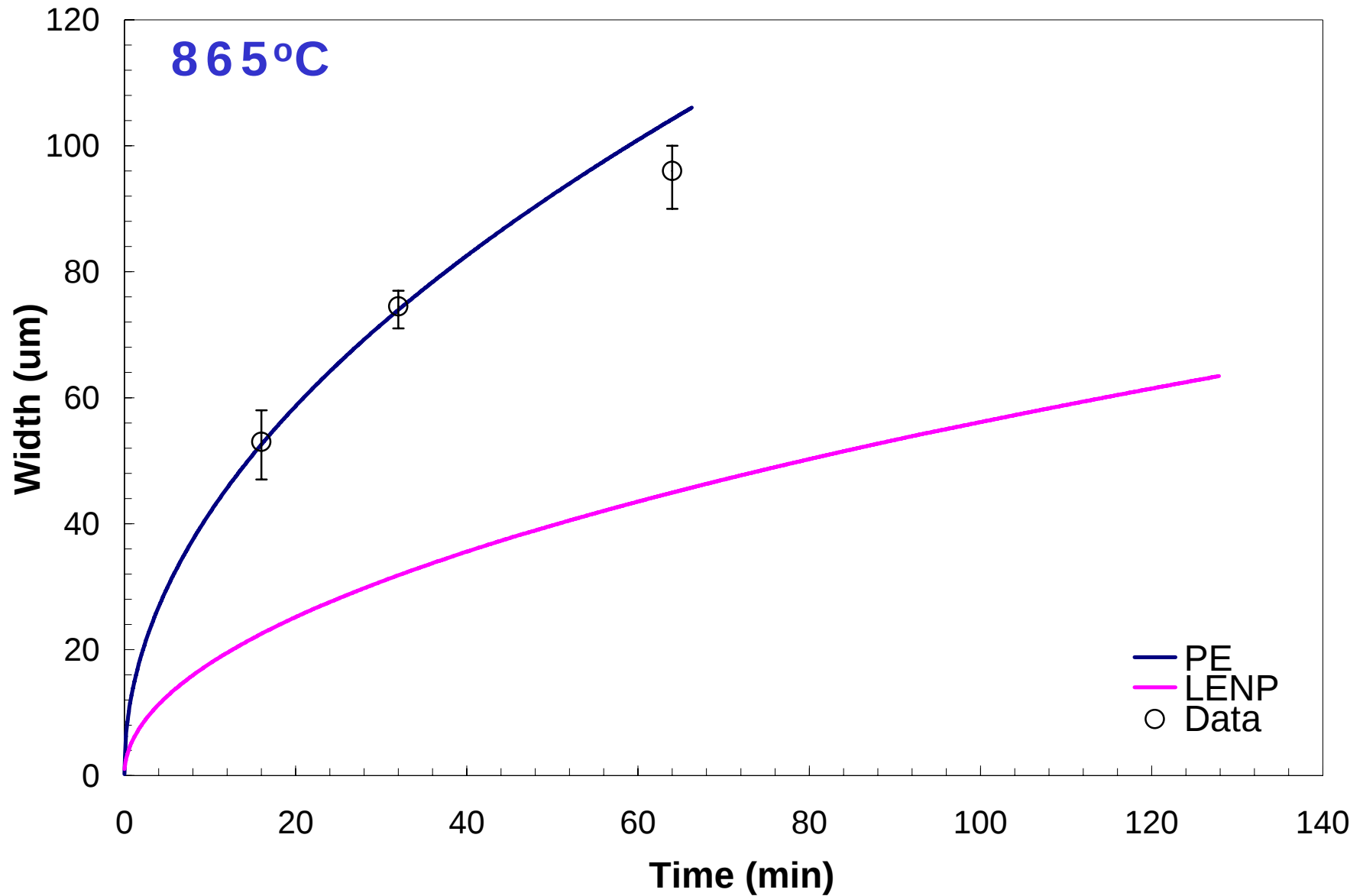
Fe-0.5%Mn-C



Fe-C-Mn: 0.48 % Mn-0.54 % C



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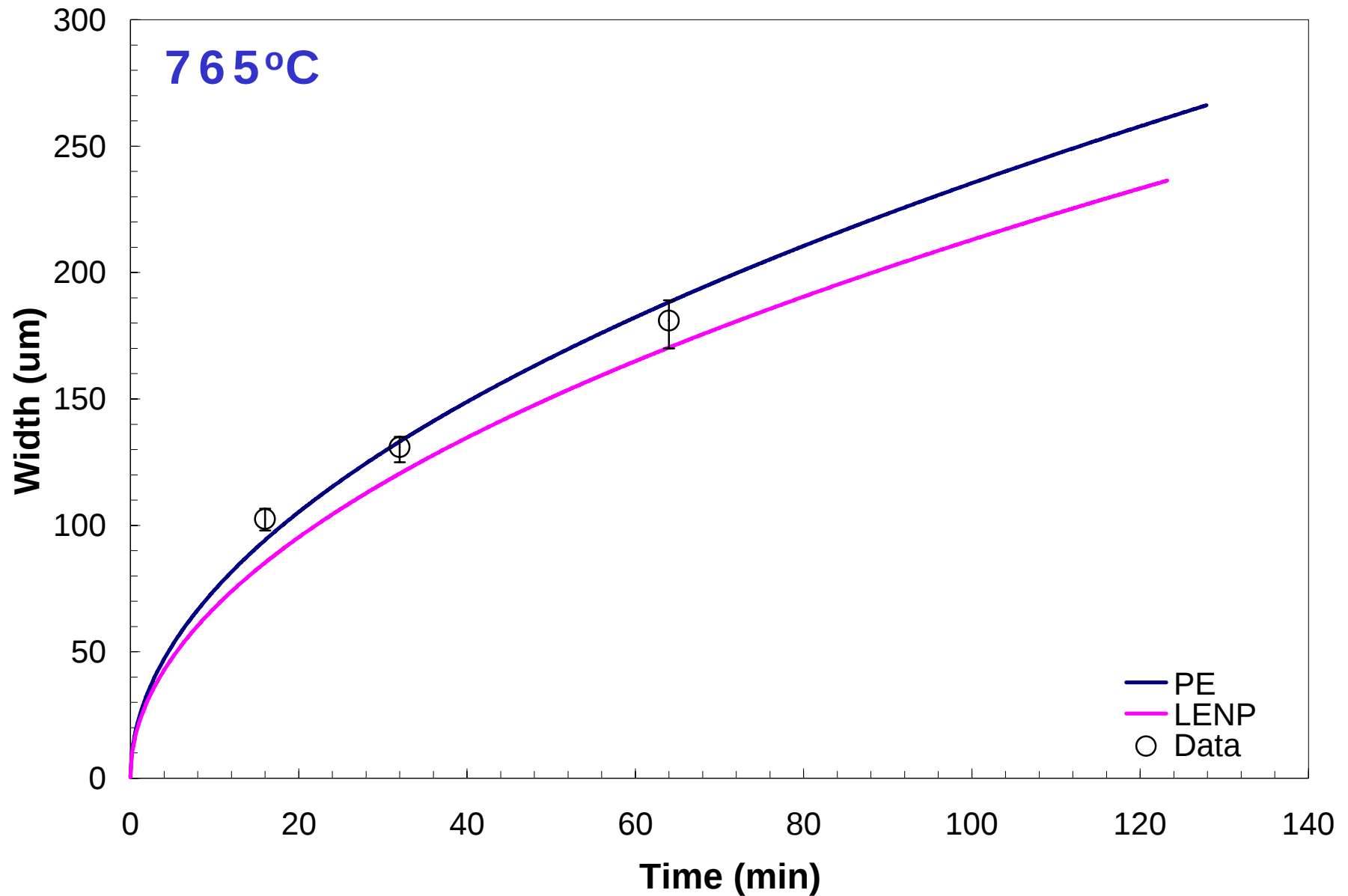


Fe-C-Mn: 0.40 % Mn-0.38 % C

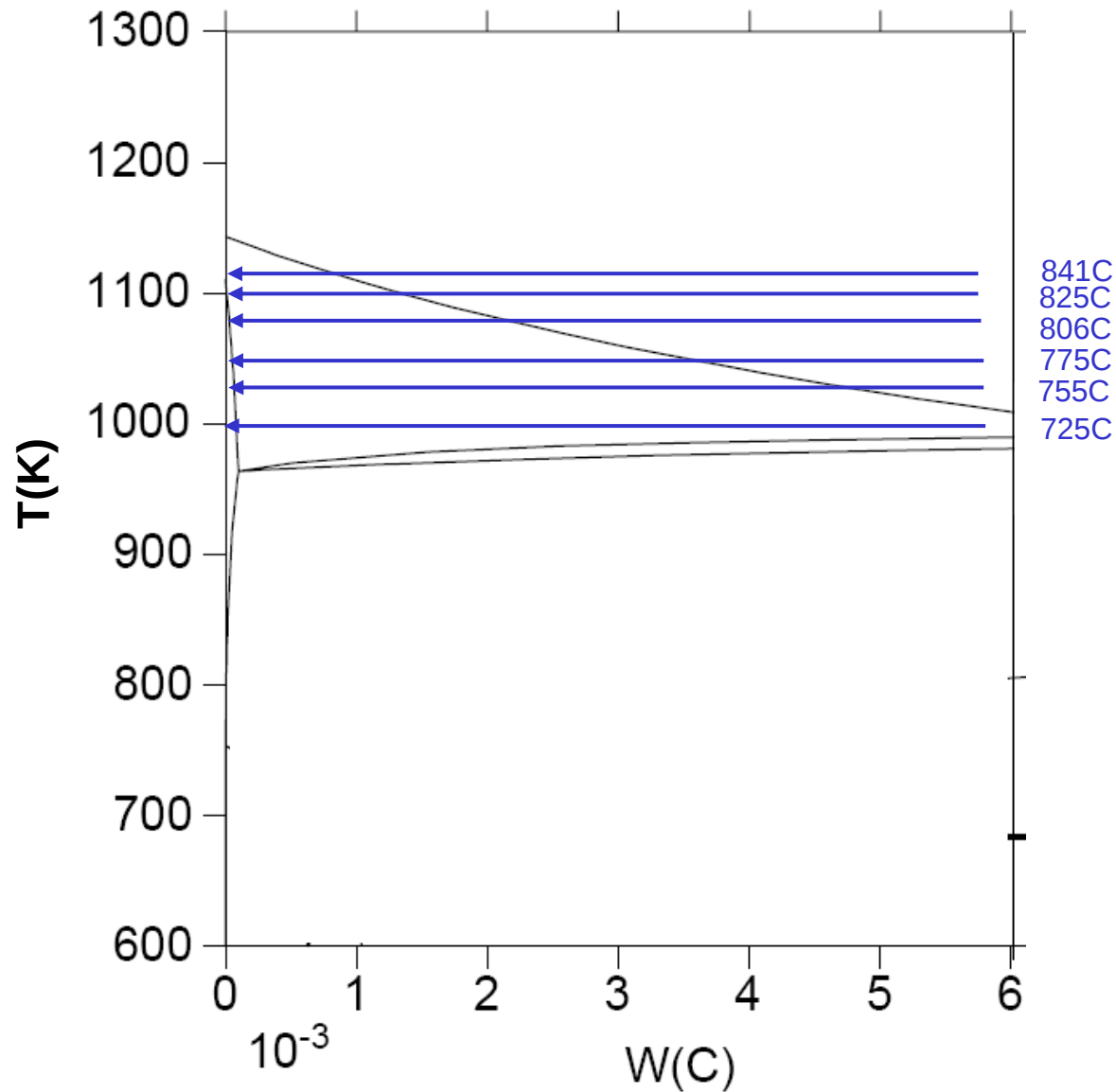
838°C, 32min, PE

810°C, 32min, PE

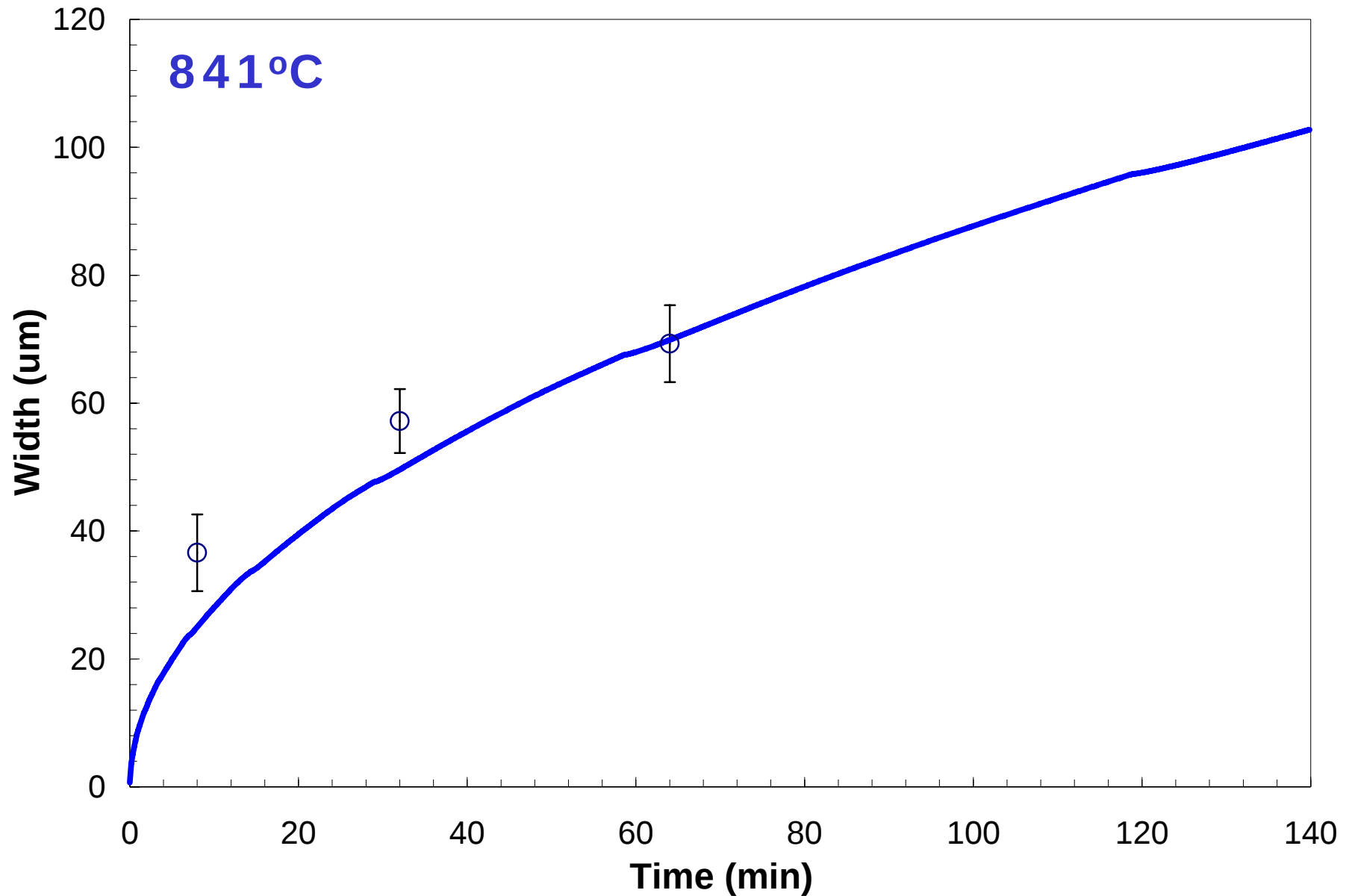
Fe-C-Mn: 0.48 % Mn-0.54 % C



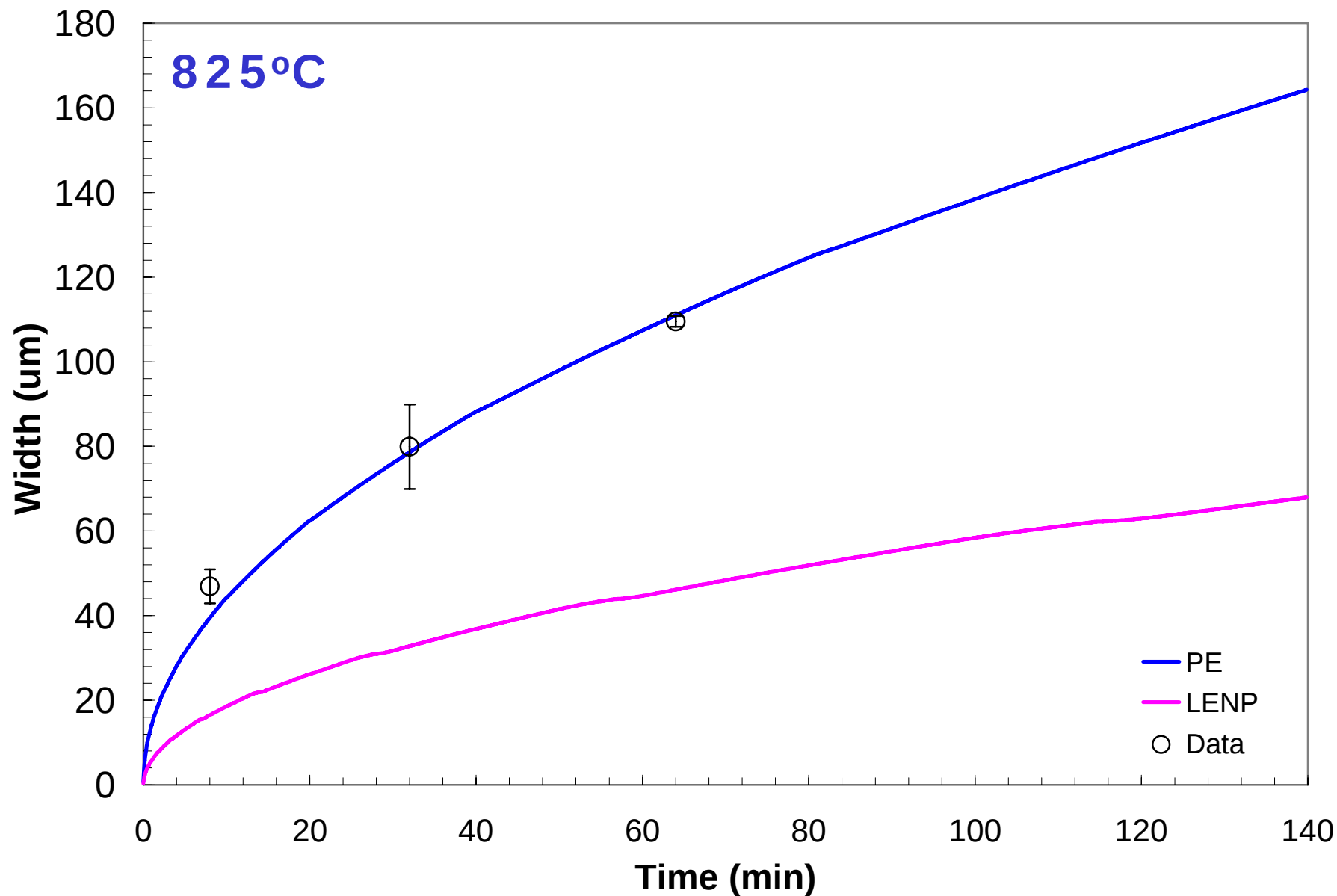
Fe-1%Mn-C



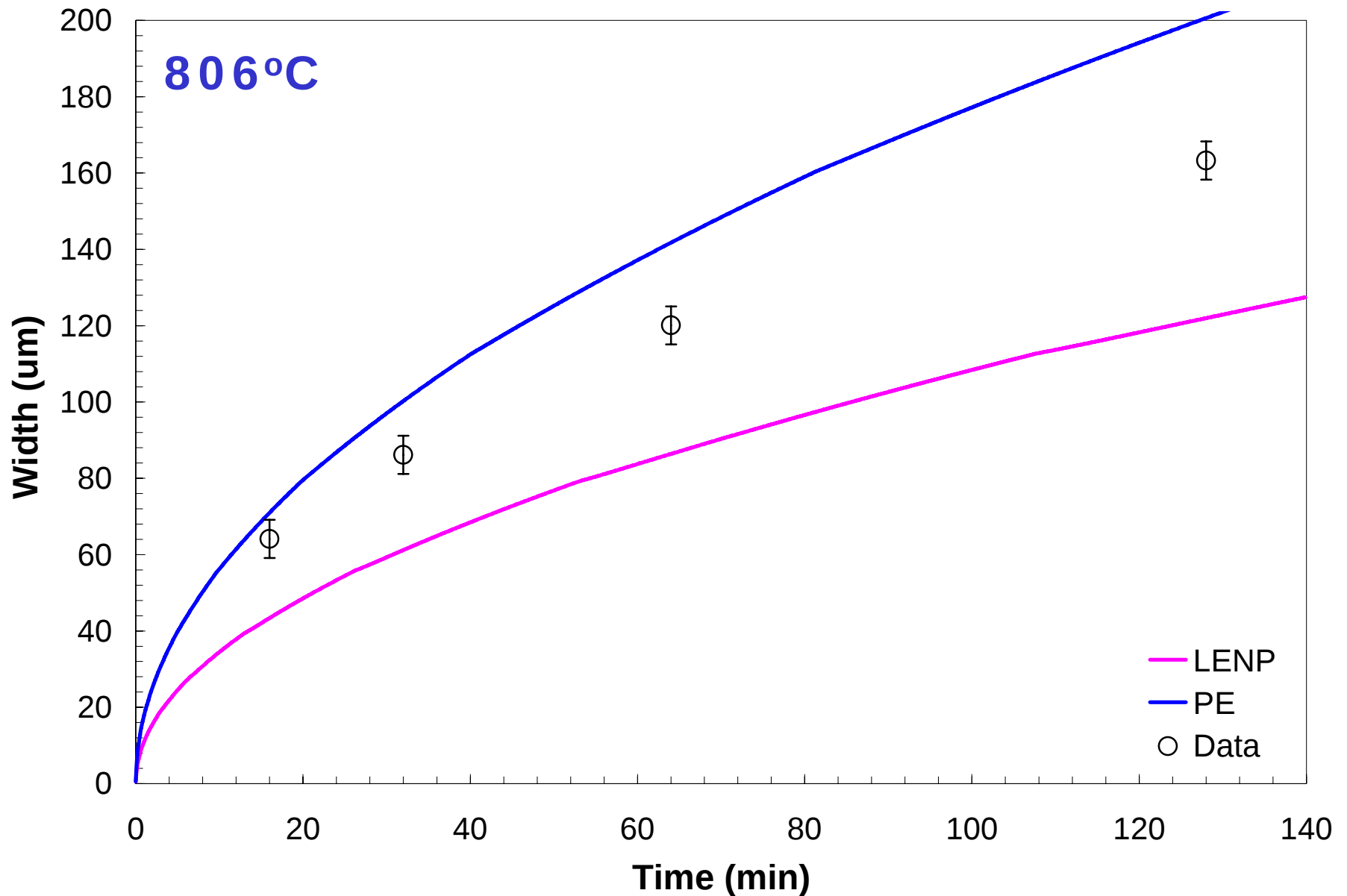
Fe-C-Mn: 0.94 % Mn-0.57 % C



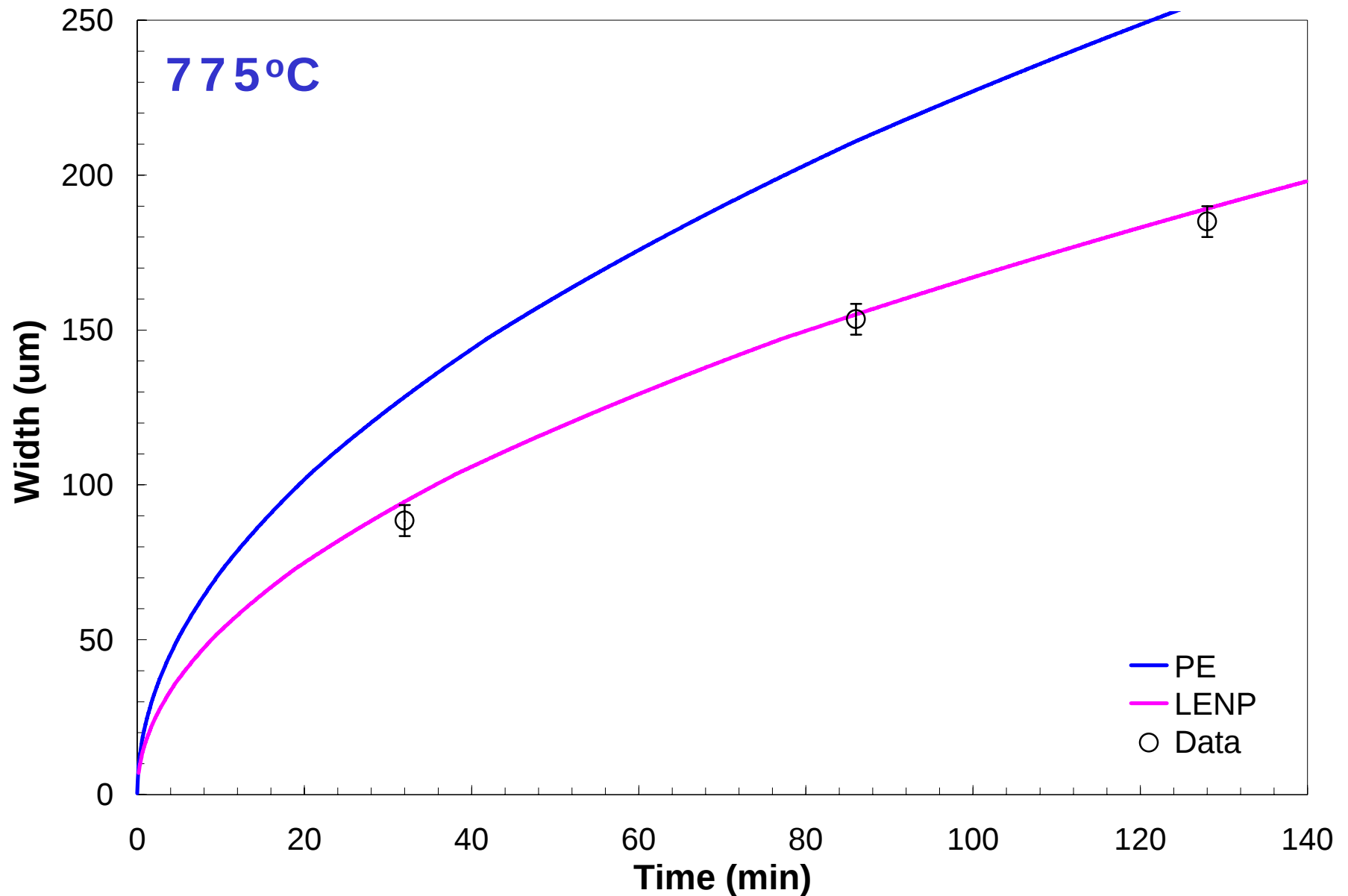
Fe-C-Mn: 0.94 % Mn-0.57 % C



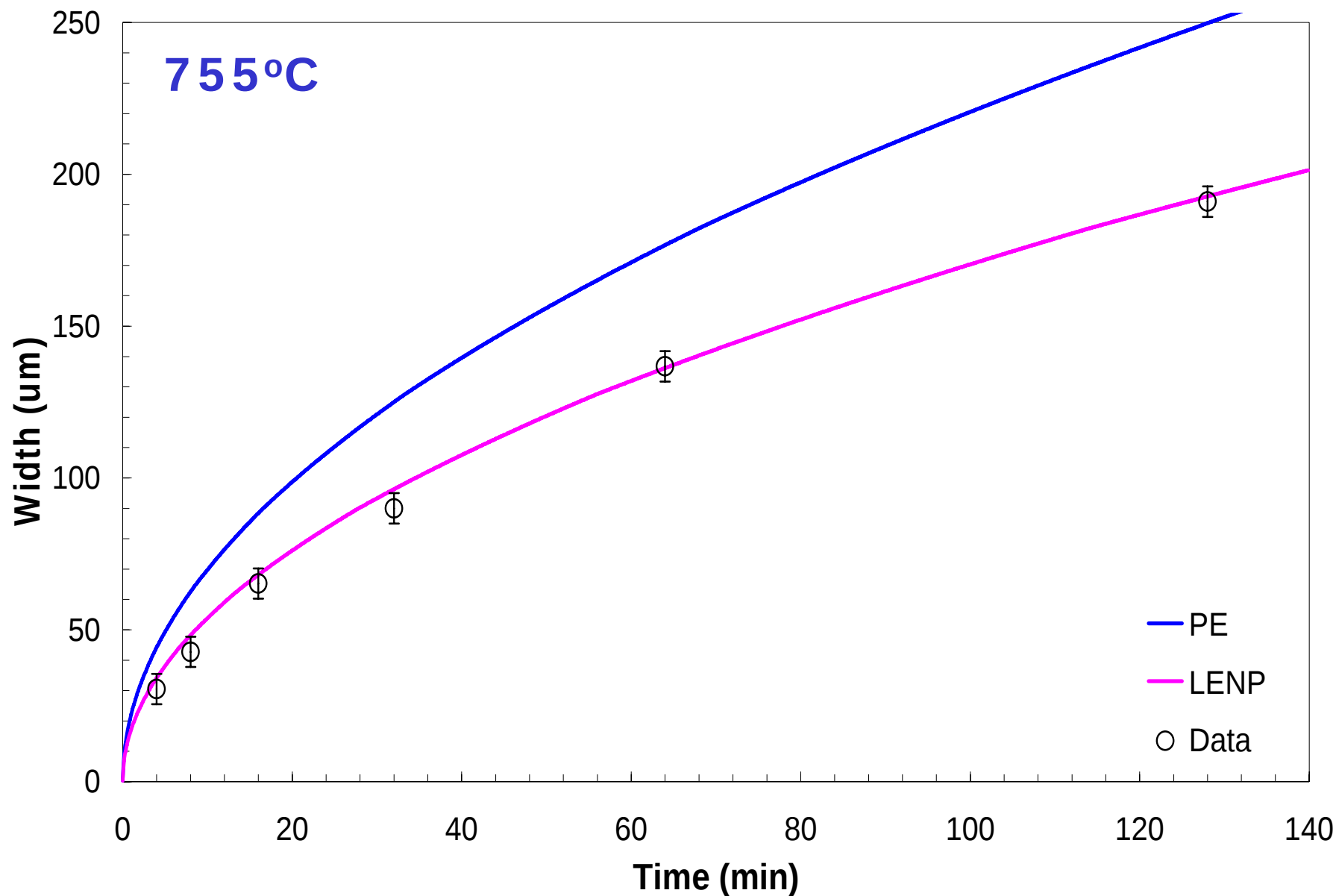
Fe-C-Mn: 0.94 % Mn-0.57 % C



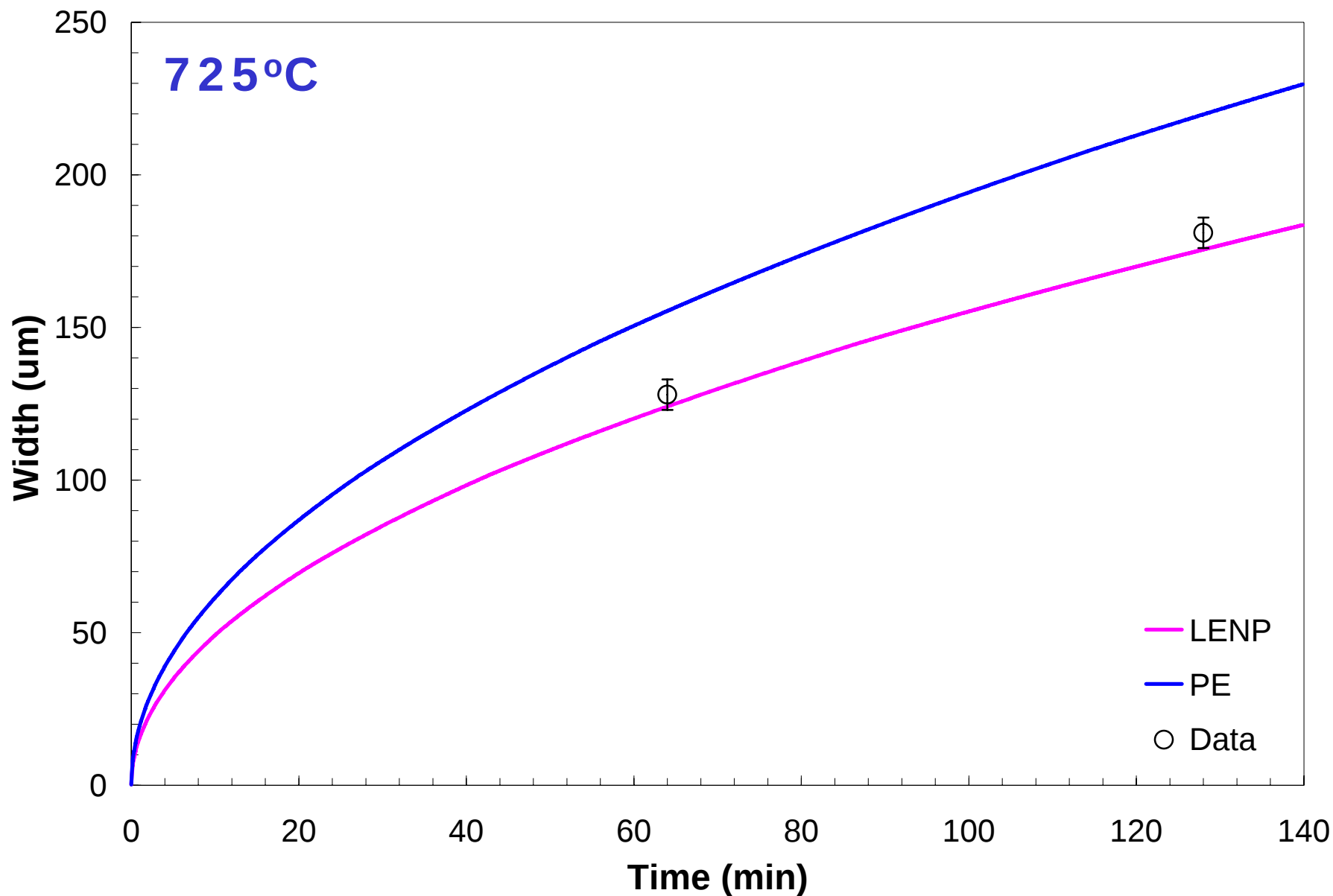
Fe-C-Mn: 0.94 % Mn-0.57 % C



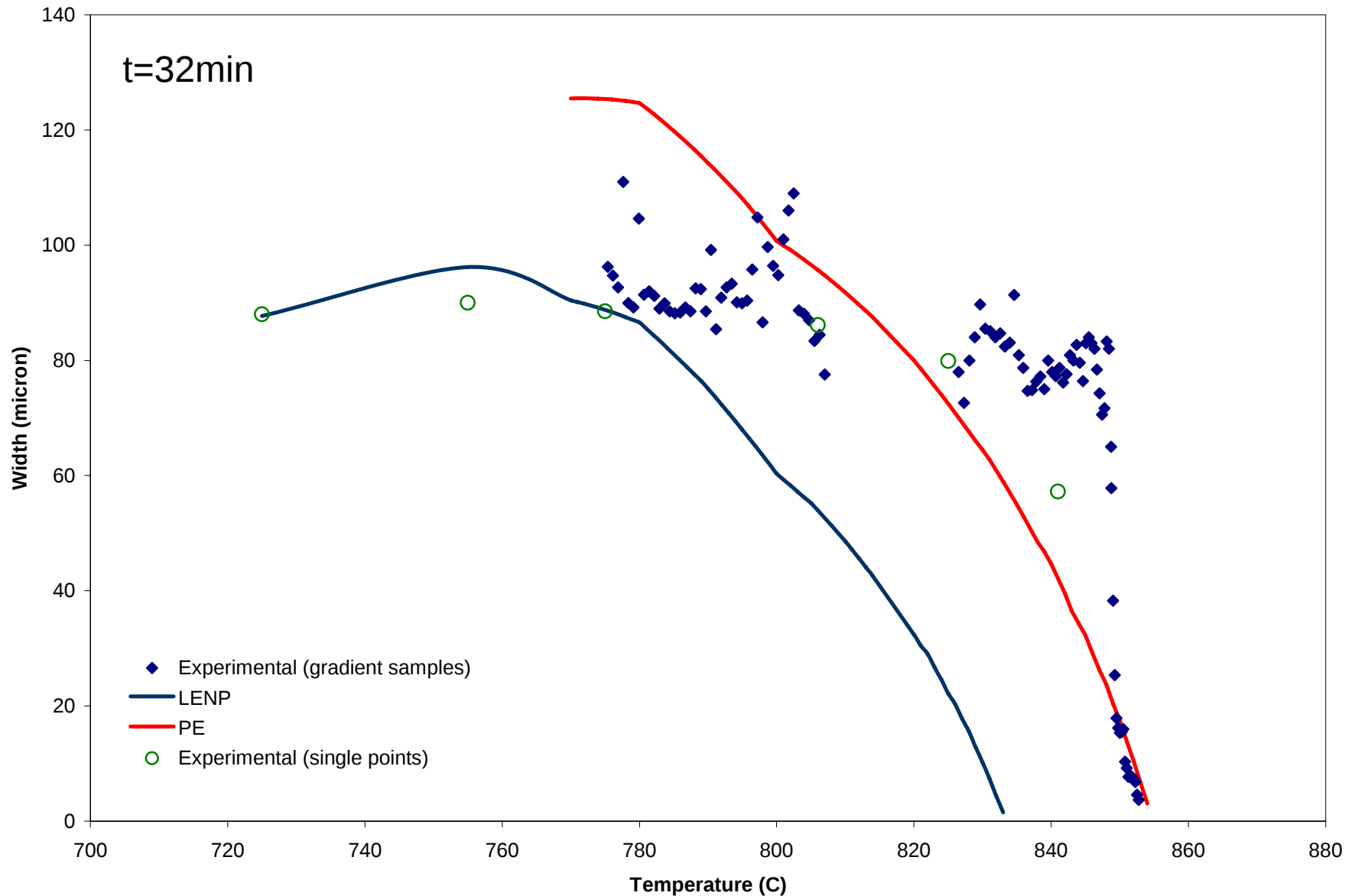
Fe-C-Mn: 0.94 % Mn-0.57 % C



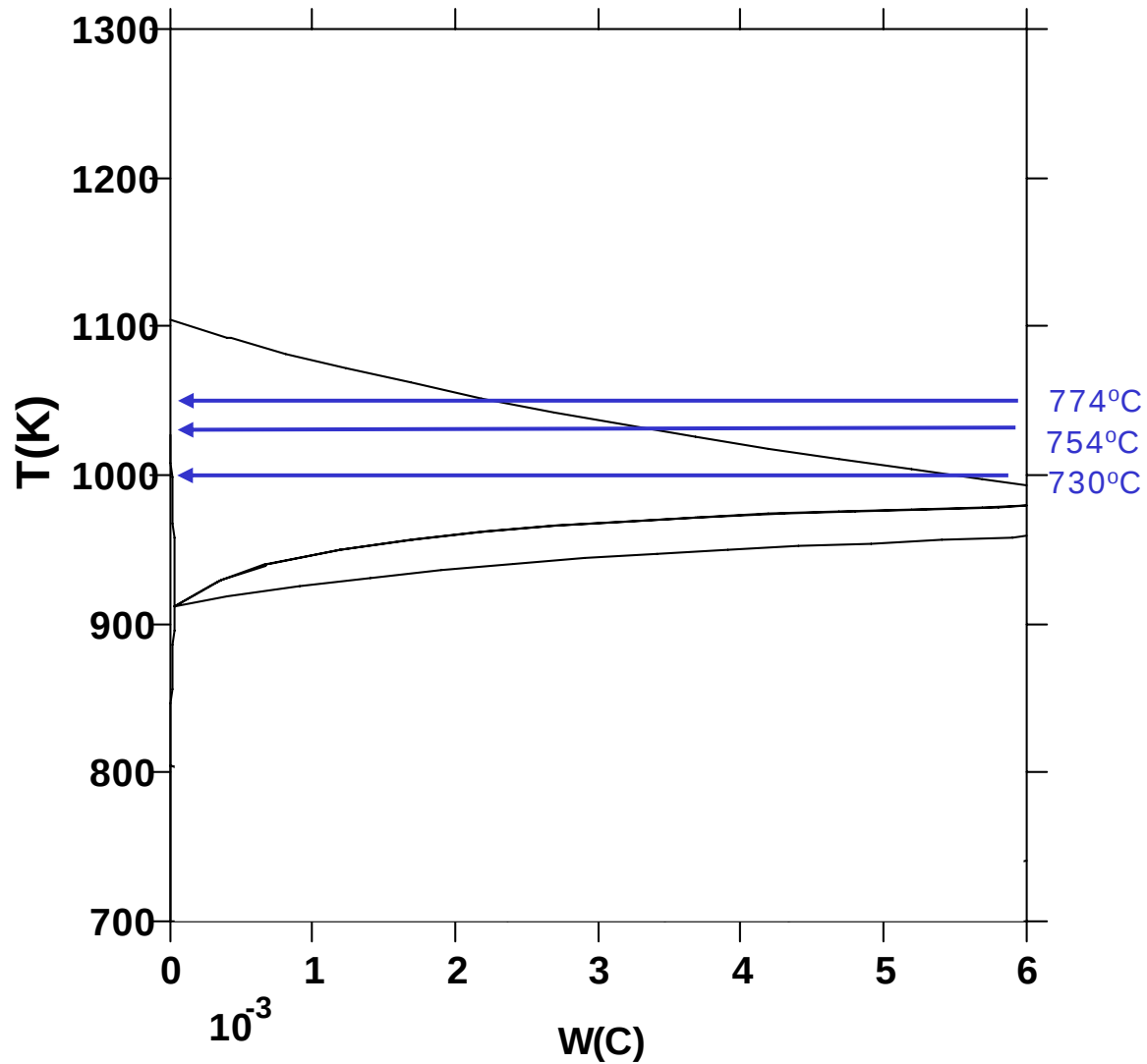
Fe-C-Mn: 0.94 % Mn-0.57 % C



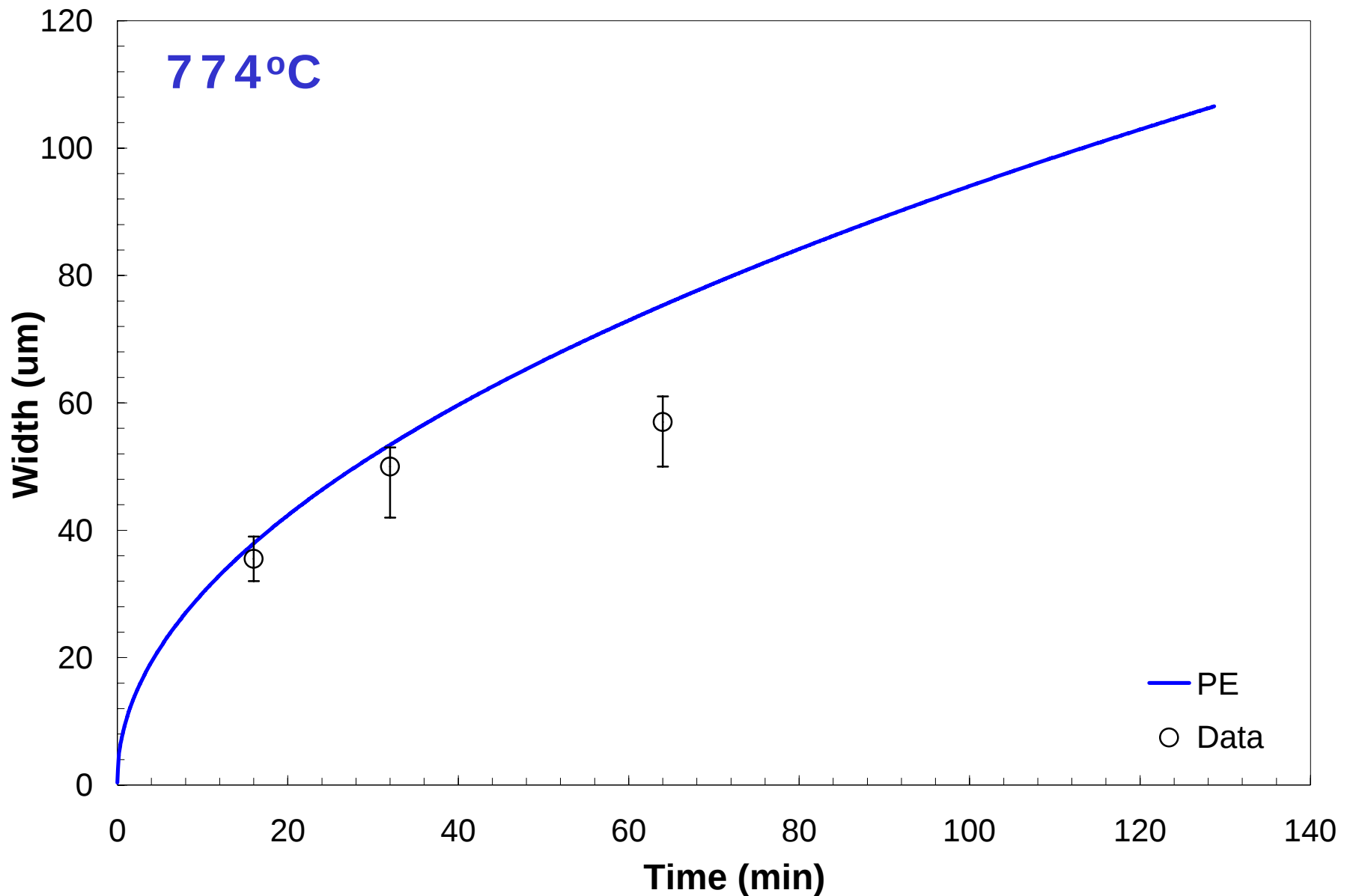
Blue dots show experimental data obtained by decarburizing a long sample in a temperature gradient for 32 min.



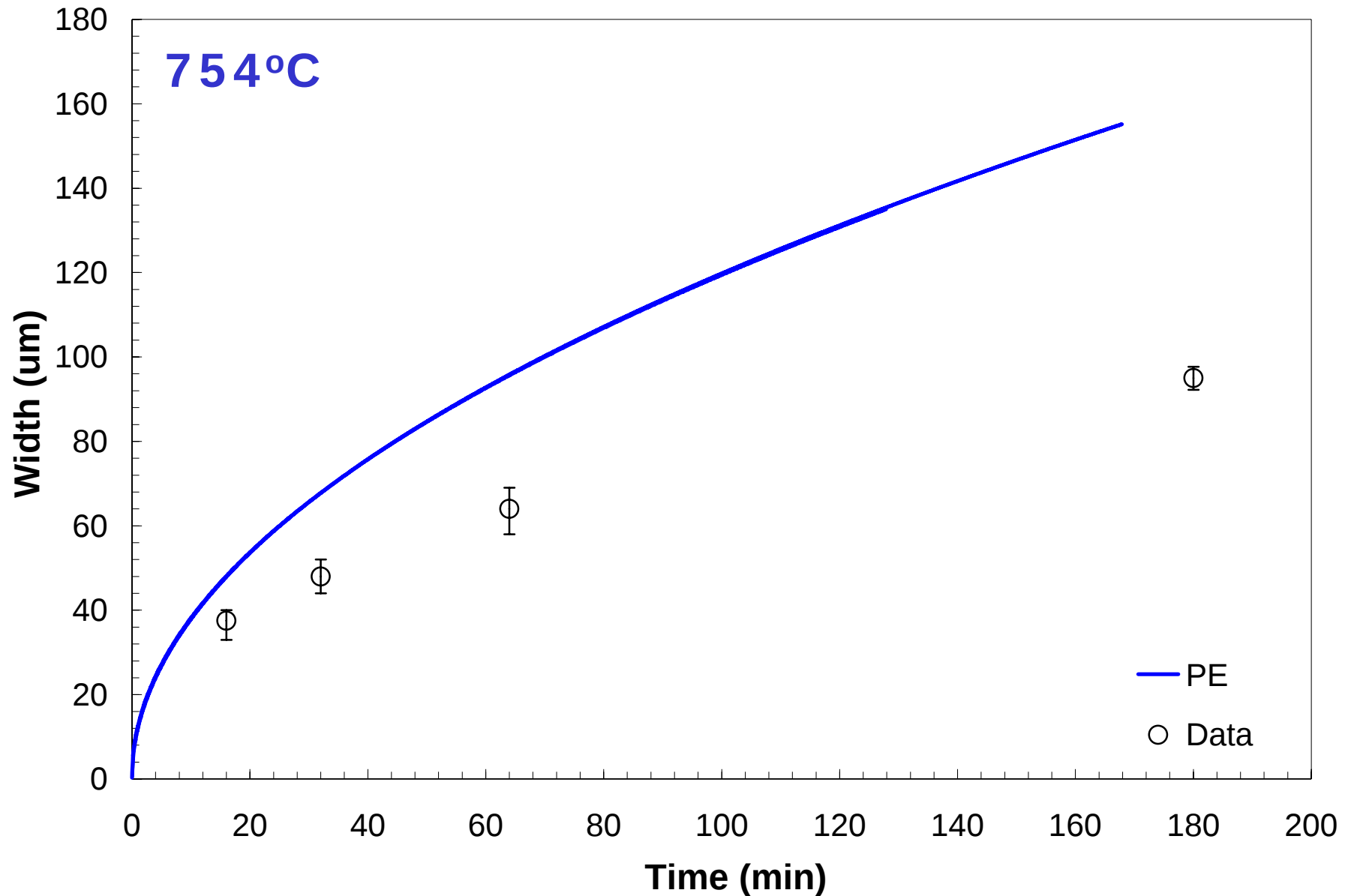
Fe-2%Mn-C



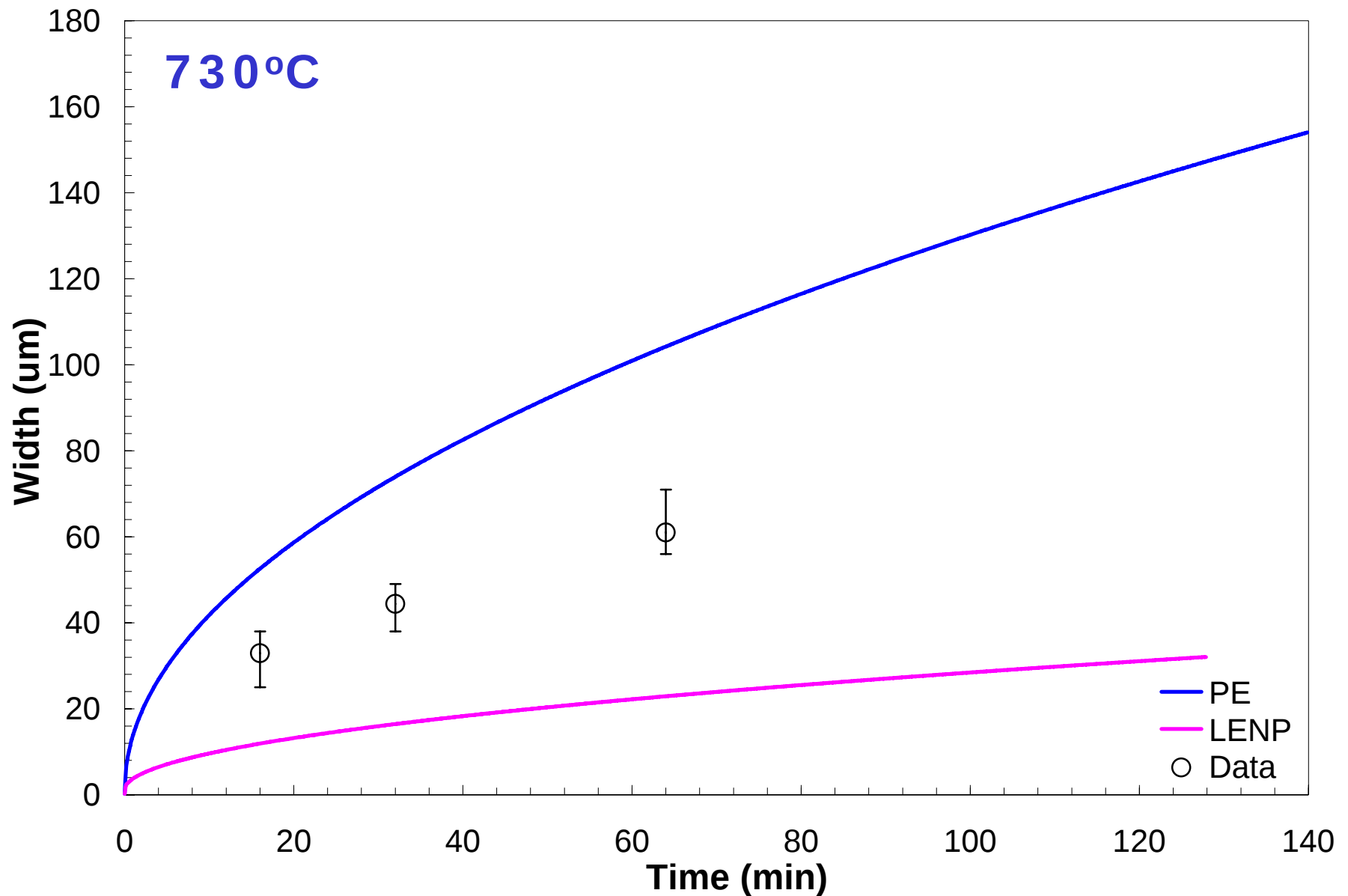
Fe-C-Mn: 2 % Mn-0.6 % C



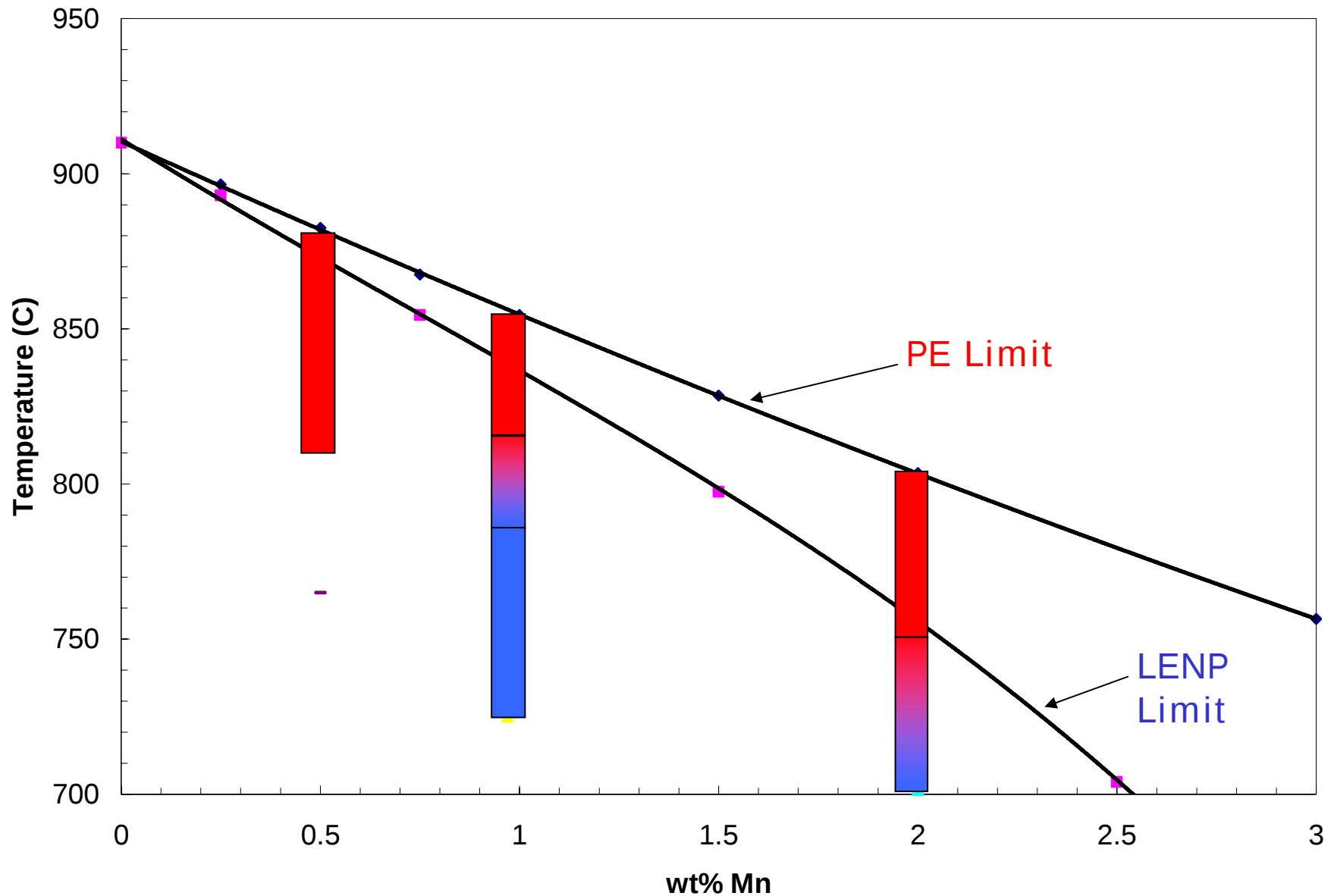
Fe-C-Mn: 2 % Mn-0.6 % C



Fe-C-Mn: 2 % Mn-0.6 % C



Fe-C-Mn: T vs. Mn content map

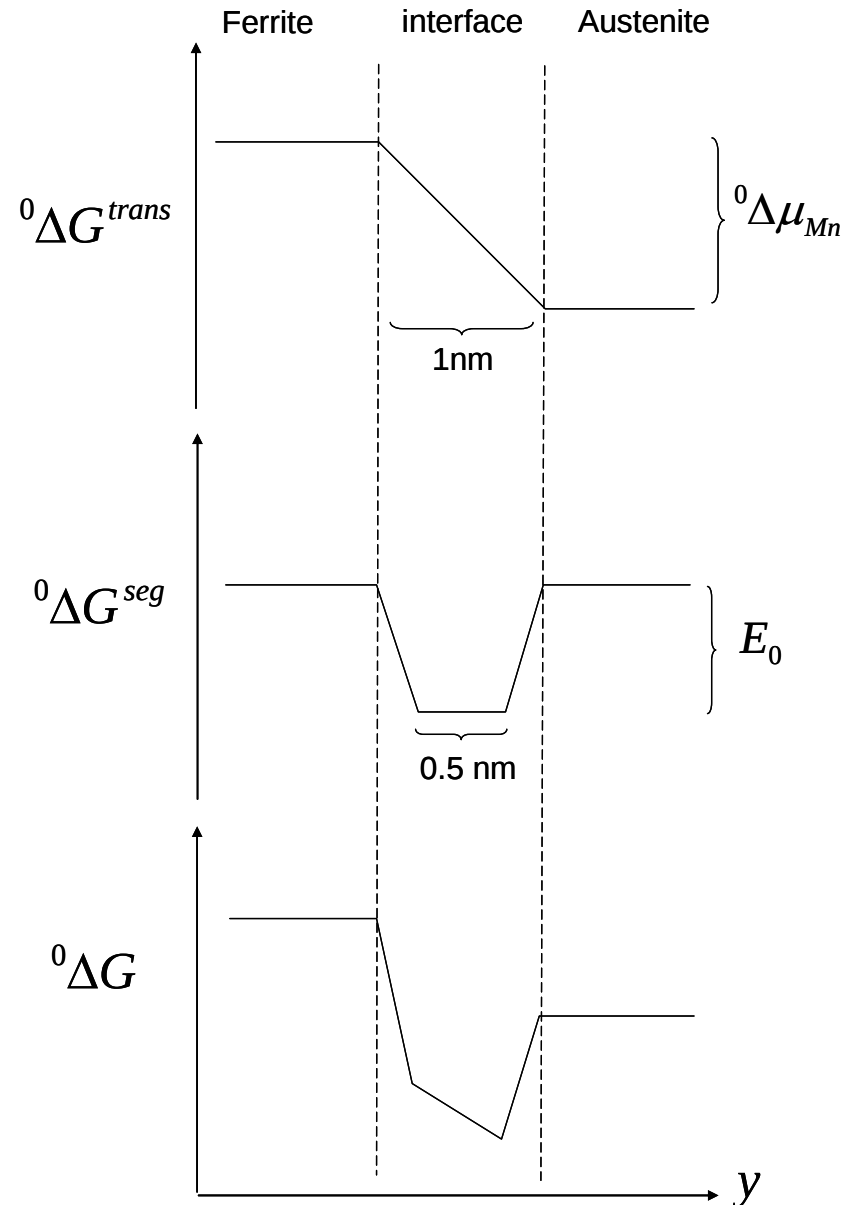


Possible Interpretations

- We need an interpretation that accounts for:
 - Difference between the behaviors of Fe-Mn-C and Fe-Ni-C.
 - Accounts for the transition between PE and LENP.
 - Explains long-lived intermediate states between LENP and PE.

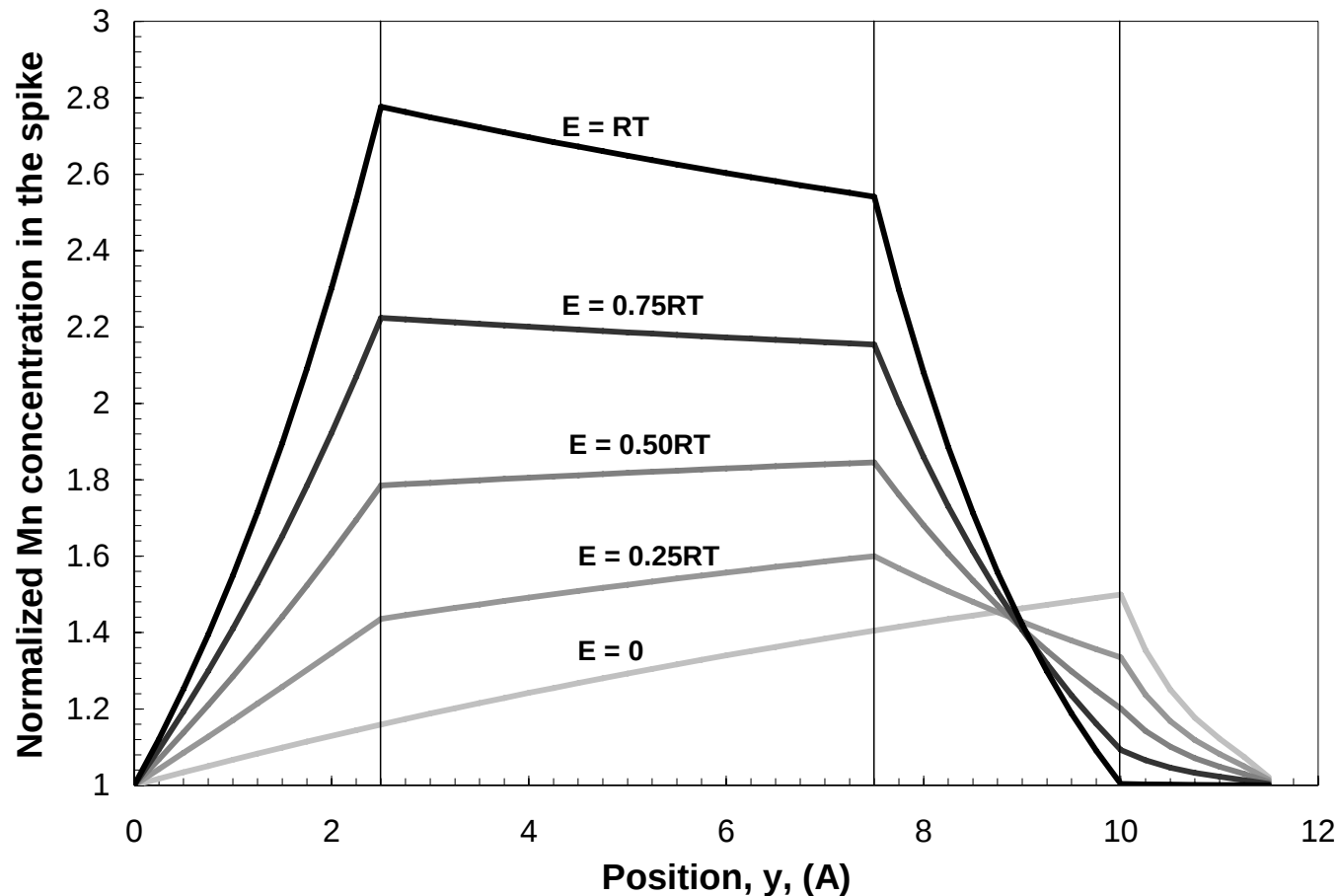
Possible Interpretations

□ Important Role of Segregation:



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□ Important Role of Segregation:



Challenges:

- Temperature dependence is problematic:
 - As T increases, segregation should become weaker. At the same time D become larger. Both factors should favor LENP at high T .
- Concentration dependence is also difficult to explain.
- Solute Drag
 - Segregation would lead to the development of solute drag.

Conclusions:

- ▣ Kinetic transitions are observed in the Fe-Mn-C system during decarburization.
- ▣ It appears that segregation to the interface plays an important role in determining the ferrite growth kinetics.
- ▣ Additional modeling is needed to rationalize the data.

Acknowledgements

- Dr. T. Furuhashi is gratefully acknowledged for providing the Fe-2%Mn-0.6% Alloy.
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