Final Report:

**Supervised linear regression machine learning model to predict the primary dendritic arm spacing**

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# Problem Statement

During the solidification of metallic systems, the microstructure grows in treelike dendritic shape and the morphology of dendrites has a strong effect on the mechanical properties. Spacing between the dendrite arms provides information on segregation patterns and distribution of precipitates between the dendritic arms, which influences the mechanical properties. There are two well-known theoretical models that relate the primary dendritic arm spacing (PDAS) to solidification conditions, namely pulling velocity (V) and temperature gradient (G), and material properties like liquid diffusivity (, the freezing range (, Gibbs‐Thomson coefficient (, and equilibrium partition coefficient (. The two models, named Hunt-Burden (HB) [1] and Kurz-Fisher (KF) [2], are different in the way the material properties are accounted:

HB model   
 . KF model

Despite the previous attempts, there are still important issues to develop a predictive model of PDAS for a wide range of solidification rates. In rapid solidification cases, such as in laser additive manufacturing of materials, the analytical relations and models for PDAS have significant deviations from computational predications and experimental data for the same solidification condition [3]. In this project we want to develop more reliable analytical models for prediction of PDAS in a wide range of solidification rates. For this, a supervised linear regression machine learning model will be used to produce a new relation between PDAS and the solidification rate (pulling velocity and temperature gradient in the model) that works for different material systems.

# Data Generating

PDAS data for the PDAS versus solidification rate and for different materials systems is generated for 7 alloy systems, with various solidification velocity range (from 10‐4 m/s to 0.12 m/s), temperature gradient range (from 106 K/m to 107 K/m) of 10 different alloys to generate enough data points for machine learning analysis. Obtaining quantitative prediction of the characteristics of solidification process with phase field simulation approach requires additional knowledge of crystal melt properties, such as anisotropic crystal-melt interfacial free energy and anisotropic kinetic coefficient. A major limiting factor is the scarcity of available experimental data pertaining to crystal-melt interfacial properties. Integrated atomistic and phase-field simulations will be performed in order to investigate the solidification of binary alloys.

In a separate study I have performed simulation study to obtain the computational data on PDAS for 7 different alloy systems (250 data points totally), for a wide velocity and temperature gradient range (given in table 1). We have also calculated all the material properties needed for each alloy systems ().

Table 1- Summary of the computational PDAS and the range of velocity (V), temperature gradient (G), and the number of datapoints for each alloy system.

|  |  |  |  |
| --- | --- | --- | --- |
| Alloy | V (m/s) | G (K/m) | Number of data points |
| Ti-3.4 at% Ni | 0.01-0.09 | 107, 8×106, 7×106,6×106, 5×106 | 45 |
| Ti-7.1 at% Ni | 0.01-0.09 | 107, 8×106,7×106,6×106, 5×106 | 42 |
| Ti-10.4 at% Ni | 0.0001-0.02 | 107, 5×106 | 22 |
| Mg-9 at% Al | 0.005-0.12 | 5×106,3×106, 1×106 | 21 |
| Al-6 at% Cu | 0.001-0.06 | 107, 8×106, 7×106, 6×106, 5×106 | 47 |
| Al-8 at% Cu | 0.01-0.08 | 107, 8×106,7×106,6×106, 5×106 | 43 |
| Al-10 at% Cu | 0.008-0.005 | 107, 8×106,7×106,6×106 | 30 |

**Data Wrangling**

It is important to obtain a clear visual view of the correlation between PDAS and the independent variables (material properties, V, and G). Visualization in this section helped us make sure we do not have any outliers in the PDAS data. The expected descending correlation between PDAS and V is more clear than the correlation between PDAs and G. However, as we have said previously, we need to check the logarithmic form of PDAS. The results do not present a specific trend in for G , mat\_HB, mat\_Kf in comparison with the

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Figure 1 Variations of PDAS with V, G, material Properties.

In order to use a regression model, we will take the logarithm from both sides of the PDAS equation for both HB and KF models, while keeping the coefficients on ln(V), ln(G), ln(mat\_HB), and ln(mat\_KF) as unknown in the linear regression models.

Then the data from computational PDAS will input into the linear regression models to obtain A, α, β, and γ.

We find a strong correlation (negative) between PDAS\_ln and V\_ln, also PDAS\_ln and G\_ln. both HB and KF material properties also have strong but positive effects on PDAS\_ln. This contradicts our original models (HB and KF) which showed G playing a more dominant role on the PDAS calculations.

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Figure 2 Heatmap graph based on Pearson correlation.

**Modeling**

*Parametric ML models- Linear Regression Models*

Chart, scatter chart

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Figure 3 The actual Ln(PDAS) versus the model predictions using HB and KF material properties.

*Nonparametric ML models*

We also wanted to test different non parametric ML models. For this we performed hyperparameter tuning for k-nearest neighbor (KNN), random forest (RF), decision tree (DT) models. With the hyperparameter branching, we found the optimum values for each model’s parameters which minimized the MSE. Figure 4 presents the performance of both parametric and non-parametric ML models. In general, the non-parametric ML models show a much better performance in comparison to the parametric linear regression models. Even the difference between the regular linear regression models with Ridge or Lasso was not dramatically different. Among all the models the DT showed the best performance (0.0328).

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Figure 4 the comparison of parametric vs nonparametric ML models based on computational training dataset.

However, in order to validate our models, we checked tested the models using the experimental data and the results, as shown in figure 5, presented an opposite trend. The parametric linear regression models have a much better performance (MSE= 0.239) in comparison to RF (MSE=2.23), DT(MSE=1.68) and KNN (MSE=1.62) models. This can be due to the fact that i) Despite the inaccuracy of analytical model’s coefficient, the general equation is correct, and that’s the reason the regression models which follows the same equational forms fits bettwe with all the data ii) the dataset we had was very small so, the performance of the model for nonparametric models was high because the model was overfitted. Despite the performance of parametric models were

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Figure 5 the comparison of parametric vs nonparametric ML models based on experimental training dataset.

**Conclusion**

1- We obtain a very important conclusion that in contract to analytical models, where temperature gradient plays much important role in the PDAS estimations, the velocity and temperature play almost equal roles in PDAS correlation!

2- In addition we compared different parametric and non parametric models. Despite the better performance of paranetric models on the computational data, when we compared them with the experimental predictions, they presented a much poorer performance and therefore, the parametric linear regression models works the best.

3- As we observed by the cross validation, the optimized alpha values for Ridge and Lasso regression models were very small. Therefore, we were already expecting the predictions of linear regression be close to Ridge and lasso models. As expected, the prediction of these models are very close to each other both for the experiment and computational data!

**Suggestion**

We suggest the fellow researchers to use the modified HB and KF models instead of the classical analytical ones:

Modified HB model

Modified KF model

**References:**

[1] M.H. Burden, J.D. Hunt, Cellular and dendritic growth. II, Journal of Crystal Growth, 22 (1974) 109-116.

[2] W. Kurz, D.J. Fisher, Dendrite growth at the limit of stability: tip radius and spacing, Acta Metallurgica, 29 (1981) 11-20.

[3] W. Xiao, S. Li, C. Wang, Y. Shi, J. Mazumder, H. Xing, L. Song, Multi-scale simulation of dendrite growth for direct energy deposition of nickel-based superalloys, Materials & Design, 164 (2019) 107553.