General Linear Regression of Superconducting Critical Temperature on Superconductor Materials

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Abstract—This paper analyzes the relationship between various material properties and the critical temperature (T_c) of superconducting materials. The main objective is to determine if a minimal set of these properties can be used to predict T_c through various general linear models. We will explore several derived metrics and use general linear regression to identify the most significant factors. The findings will demonstrate how these metrics are linearly related to the critical temperature.

Index Terms—Superconductors, Critical Temperature, General Linear Model, Gaussian, Gamma, Tweedie, Predictive Modeling

I. Introduction

This study investigates the linear relationship between a set of predictive metrics and the critical temperature (T_c) of various materials. We aim to determine if a linear model can effectively predict T_c and, if so, to identify the most suitable statistical distribution family for this relationship. The data, including several highly correlated derived metrics, will be analyzed to answer these questions.

We will begin by detailing the characteristics of the dataset and address any data anomalies. This will be followed by an explanation of the preprocessing methods used to prepare the data for our Generalized Linear Models (GLM). The analysis will incorporate insights from both the training and validation datasets to build and evaluate the predictive models.

Finally, we will present which GLM distribution family provides the best fit, as determined by its lowest error rate, for predicting the critical temperature of the materials under consideration.

II. ANALYSIS

A. Data

This study utilizes a dataset on superconducting materials, sourced from reference [1]. The dataset contains information about the chemical composition and critical temperature (T_c) for various materials.

- 1) Data Structure and Features: The dataset includes eight derived chemical features for each material's chemical formula:
 - Atomic Mass
 - · Atomic Radius

- Density, Electron Affinity
- First Ionization Energy (FIE)
- Fusion Heat
- Thermal Conductivity
- Valence

These eight features are summarized across ten statistical metrics: mean, weighted mean (wtd_mean), geometric mean (gmean), weighted geometric mean (wtd_gmean), entropy, weighted entropy (wtd_entropy), range, weighted range (wtd_range), standard deviation (std), and weighted standard deviation (wtd_std). This results in a total of 80 potential features (8 features × 10 statistical metrics).

Additionally, the dataset includes as an additional feature the **Number of Elements** (number_of_elements) in each material's chemical formula.

2) **Data Handling**: The units of measurement were not provided with the dataset, so all features were treated as **dimensionless numerical values**. A key characteristic of this data is the **high correlation** among many of the features, particularly between a feature and its weighted counterpart (e.g. mean and wtd_mean), which is a known aspect of this type of materials science data.

B. Methods

C. Analysis

D. Results

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III. CONCLUSION(S)

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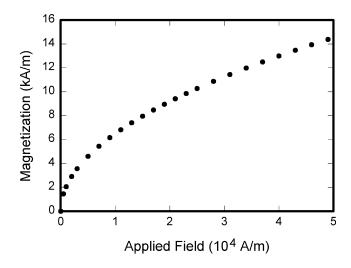


Fig. 1. Example of a figure caption.

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APPENDIX

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