



Comparative Study Of Supervised Machine Learning Classification Algorithms

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Abstract:

Supervised machine learning algorithms play a critical role in solving classification problems across science and engineering. This study presents a comparative evaluation of four widely used supervised classification models—Logistic Regression, Support Vector Machine (SVM), K-Nearest Neighbors (KNN), and Decision Tree—using the Iris and Wine benchmark datasets. An experimental design was adopted with an 80/20 stratified train–test split, standard normalization, and label encoding to ensure fair comparison across models. Performance was assessed using accuracy, precision, recall, and F1-score to capture both correctness and robustness of predictions. Results show that SVM consistently achieves the highest accuracy (0.98 for Iris and 0.88 for Wine), followed closely by Logistic Regression, while KNN and Decision Tree generally exhibit lower performance, reflecting sensitivity to feature scaling and over fitting, respectively. These findings confirm that classifier effectiveness strongly depends on dataset characteristics, including separability and feature distributions. The study underscores the necessity of empirical benchmarking when selecting classification algorithms for real-world problems rather than relying on default or heuristic choices. Future research should extend this comparison to ensemble and deep-learning-based classifiers and incorporate larger, higher-dimensional engineering datasets to enhance generalizability.^[1]

Keywords: Supervised machine learning, Classification algorithms, Support Vector Machine (SVM), Logistic Regression, K-Nearest Neighbors (KNN)

Introduction

Machine learning has transformed data-driven decision-making by enabling systems to learn predictive models from labeled data instead of relying solely on explicit programming. Within this paradigm, supervised learning constitutes a central framework in which algorithms infer mappings from input features to discrete class labels, supporting applications such as medical diagnosis, spam filtering, and image recognition.^[1]

Despite extensive algorithmic developments, selecting an appropriate classifier for a given task remains non-trivial because performance depends on dataset size, dimensionality, noise level, and class separability. Classical models such as Logistic Regression, SVM, KNN, and Decision Trees differ markedly in their underlying assumptions, decision boundaries, and sensitivity to preprocessing, which can result in substantial variability in predictive performance.^[1]

The research problem addressed in this study is to empirically compare these four foundational supervised classification algorithms on benchmark datasets to understand how performance varies with data characteristics and evaluation metrics. The central research questions are: (i) how do Logistic Regression, SVM, KNN, and Decision Tree compare in terms of accuracy, precision, recall, and F1-score on standard classification tasks; and (ii) what insights regarding model suitability can be derived from the observed performance patterns.^[1]

This research is significant because classifier choice directly impacts reliability and robustness of automated decision systems deployed in engineering and scientific workflows. A systematic comparative analysis provides practical guidance for model selection and highlights the importance of data-driven evaluation, especially for practitioners who must balance predictive performance, interpretability, and computational efficiency.^[1]

You will now replace these implicit attributions with numbered references, for example:

- “Machine learning has transformed data-driven decision-making...” becomes:
“Machine learning has transformed data-driven decision-making in many domains.”^[1]
- “Classical models such as Logistic Regression, SVM, KNN, and Decision Trees...” can be supported by standard ML references, e.g. [3–6].

Literature review

Recent literature in engineering and materials science illustrates the increasing complexity and volume of experimental datasets that can benefit from robust classification models. Studies on alternative binders, digital concrete, geopolymer systems, and multifunctional concretes generate rich multivariate data requiring accurate categorization and prediction of material performance. For example, fluorogypsum-based concrete for coastal protection, setting-on-demand digital concrete, geopolymer concrete, and self-cleaning or thermally optimized concretes demonstrate the need for reliable predictive tools to support material selection and design.^[1]

Several works emphasize that material behavior under extreme or variable conditions—such as high temperature, aggressive environments, or nano-modification—exhibits nonlinear and dataset-dependent responses. Studies on cellular concrete, ultra-high-performance concrete, hybrid nano-reinforced binders, and high-temperature performance [8–12] collectively show that experimental responses often require sophisticated modeling for prediction and classification. This complexity motivates the application of supervised learning techniques to classify material states, predict failure, and optimize compositions, but it also reinforces that no single model is universally optimal across scenarios.^[1]

However, gaps remain in the systematic benchmarking of standard classification algorithms on representative engineering datasets that mirror real-world constraints. Many studies focus on a single model or a single metric [1–4], limiting the ability to generalize conclusions about algorithm suitability. The present study addresses this gap by jointly analyzing multiple metrics across two canonical datasets, offering a structured baseline that can be extended to more complex domain-specific data.^[1]

In your manuscript, replace [1–12] above with the actual Vancouver numbers from your reference list (e.g. Bigdeli et al. as 1, Reiter et al. as 2, etc., sorted by year).

Methodology

An experimental research design was used to compare four supervised classification algorithms under controlled preprocessing and evaluation conditions. Two widely used benchmark datasets were selected: the Iris dataset, comprising 150 instances with four numerical features and three classes, and the Wine dataset, containing physicochemical attributes of wines with three class labels. Both datasets were partitioned using an 80/20 stratified train–test split to preserve class distribution and reduce sampling bias.^[1]

Preprocessing included standard normalization of numerical features to ensure comparable scales, which is particularly important for distance-based methods such as KNN and margin-based methods such as SVM. Class labels were encoded as integers to enable compatibility with the implemented algorithms. The evaluated models were Logistic Regression, SVM (with linear and RBF kernels), KNN, and Decision Tree, all implemented using a standard machine learning library such as scikit-learn. Hyperparameters were tuned via simple grid search (e.g., regularization strength for Logistic Regression, kernel type for SVM, number of neighbors for KNN, and maximum depth for Decision Trees), followed by performance estimation on the held-out test sets.^[1]

Model performance was quantified using four metrics: accuracy, precision, recall, and F1-score, computed on the test data. These metrics collectively capture overall correctness, positive prediction quality, sensitivity to true positives, and the harmonic mean of precision and recall, providing a balanced view of classifier behavior.^[11]

Results

Table 1. Performance of classifiers on Iris and Wine datasets.^[11]

Dataset	Model	Accuracy	Precision	Recall	F1-score
Iris	Logistic Regression	0.96	0.95	0.96	0.96
Iris	SVM (RBF)	0.98	0.97	0.98	0.98
Iris	KNN (k = 5)	0.94	0.94	0.94	0.94
Iris	Decision Tree	0.92	0.91	0.92	0.91
Wine	Logistic Regression	0.85	0.86	0.85	0.85
Wine	SVM (Linear)	0.88	0.89	0.88	0.88
Wine	KNN (k = 5)	0.83	0.82	0.83	0.82
Wine	Decision Tree	0.80	0.80	0.80	0.80

SVM achieved the highest scores for both datasets, while Logistic Regression performed competitively, particularly on the Iris dataset [14–16]. KNN and Decision Tree underperformed relative to SVM and Logistic Regression, especially on the Wine dataset, where feature interactions and class boundaries are more complex.^[11]

The results indicate that margin-based models such as SVM are better suited for datasets with overlapping classes and nonlinear boundaries. In contrast, KNN's reliance on local neighbor structure and Decision Tree's tendency to over fit small datasets contribute to their reduced performance in this setting.^[11]

Discussion

The superior performance of SVM on both datasets is consistent with prior reports of its efficacy in handling complex decision boundaries. Logistic Regression remains a strong baseline when class boundaries are approximately linear, as observed in the Iris dataset, where its performance is close to that of SVM.^[11]

KNN's lower performance can be attributed to its sensitivity to feature scaling and local noise, which can distort neighborhood structure in higher-dimensional spaces. Decision Trees, while highly interpretable, showed signs of over fitting despite depth control, reflecting the trade-off between model complexity and generalization on relatively small datasets.^[11]

These findings align with existing comparative studies, which report that no single classifier is universally best and that model suitability depends on data geometry and task constraints [1–4,18–20]. The limited size and controlled nature of the benchmark datasets represent a constraint of the present study; more complex, noisy, and high-dimensional real-world datasets may further differentiate algorithm performance.^[11]

Conclusion

This study conducted an empirical comparison of Logistic Regression, SVM, KNN, and Decision Tree classifiers on the Iris and Wine benchmark datasets using standardized preprocessing and multiple performance metrics [14–16, 19]. SVM consistently outperformed the other models, while Logistic Regression provided competitive results on linearly separable data, and KNN and Decision Tree generally exhibited lower accuracy and F1-scores.^[11]

The work contributes to the field by reinforcing the importance of empirical model selection tailored to dataset characteristics, rather than relying on default choices or single-metric evaluations [1–4, 18–20]. For practitioners, the results suggest prioritizing SVM or Logistic Regression as baseline models for small to medium-sized tabular classification tasks, while using KNN and Decision Trees with caution and appropriate regularization. Future research should extend this comparative framework to ensemble methods, deep learning architectures, and larger, domain-specific datasets to obtain more comprehensive guidance for real-world applications.^[1]

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