

Exploring Machine Learning Model Performance across Diverse Datasets: A Comparative Analysis

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Introduction

This report provides an analysis of using various machine learning algorithms on various datasets. The goal is to identify the best-use scenarios of model variants given the nature of input data. In this report, the authors analyze the

- Decision Tree, *provided by Polina Petrova*
- k Nearest Neighbors, *provided by Polina Petrova*
- Neural Network, *provided by Yan Mazheika*
- Random Forrest, *provided by Yan Mazheika*

against the handwritten digits, titanic survival, loan eligibility, and Parkinson's classification datasets.

It is our aim to explain the nature of the data, our models, and the performance of these classifiers. Throughout our report, we justify our algorithm choice for the given dataset and our choice of hyper-parameters for the tuning of the algorithm. These insights are for the reader's benefit; we also aim to provide insight into how to solve novel machine problems, which algorithms may work best, and how to adjust their hyper-parameters for optimal performance.

Approach

The authors have chosen to coordinate on every dataset. For the first two algorithms of every dataset, one of them comes from *Yan Mazheika* while the other comes from *Polina Petrova*. These algorithms may have been modified to handle a new type of data but their fundamental logic stays consistent from previous use cases.

We use cross-validation of $k=10$ folds when testing the performance of our models (except for kNN). We also transform the data into the $[0,1]$ range when using the neural network or kNN classifiers.

Model Performance on Diverse Datasets

Dataset:	Digits		Titanic		Loan		Parkinson's	
	Accuracy	F1-Score	Accuracy	F1-Score	Accuracy	F1-Score	Accuracy	F1-Score
k-NN	.8698	.9298	-	-	-	-	0.7526	0.8557
Decision Tree	-	-	.7023	.8234	.5354	.6860	-	-
Random Forrest	-	-	-	-	.7354	.6650	-	-
Neural Network	.9962	.9809	.9663	.9652	-	-	.9226	.8940
Network Forrest	-	-	-	-	-	-	-	-

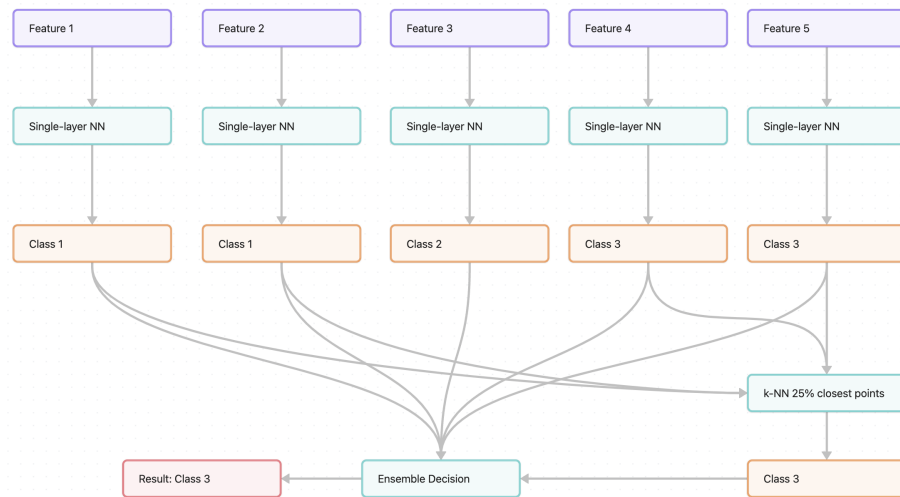
An Additional Variant

Consider a model variant which essentially performs logistic regression on each feature and then performs a vote on the output of each feature. The neural net architecture was modified to have a single neuron in one hidden layer to imitate logistic regression. We perform this logistic regression for each feature and then perform a vote. Afterwards, a k-NN variant gets to vote, which potentially breaks ties.

When each of the neural networks makes a vote, we take the vote or votes with the highest frequency. These represent the features which contributed to the majority vote. If the vote is evenly split, all of the features are used in the k-NN algorithm to break the tie. Otherwise, it's just the features that contributed to the majority vote. This k-NN variant performs the euclidean distance calculation to find the closest 25% of the training data to contribute to the vote.

The idea here is to rank the features by their importance in the dataset. If they contributed to the majority vote, we essentially confirm that they are important. In a way, it's like a decision tree in the way we prioritize the most important features. However, it's important to note that unlike a decision tree, we don't have access to the test dataset to confirm that they actually are. So essentially, we are estimating the best features to 'split' on rather than tuning them with the y-values.

We'll analyze this new variant over all of the datasets we look at. This graphic displays an example:



This is our attempt for extra credit #3 and #4. Because we analyze this variant over all of the datasets, this is also for extra credit #1.

Digits Dataset

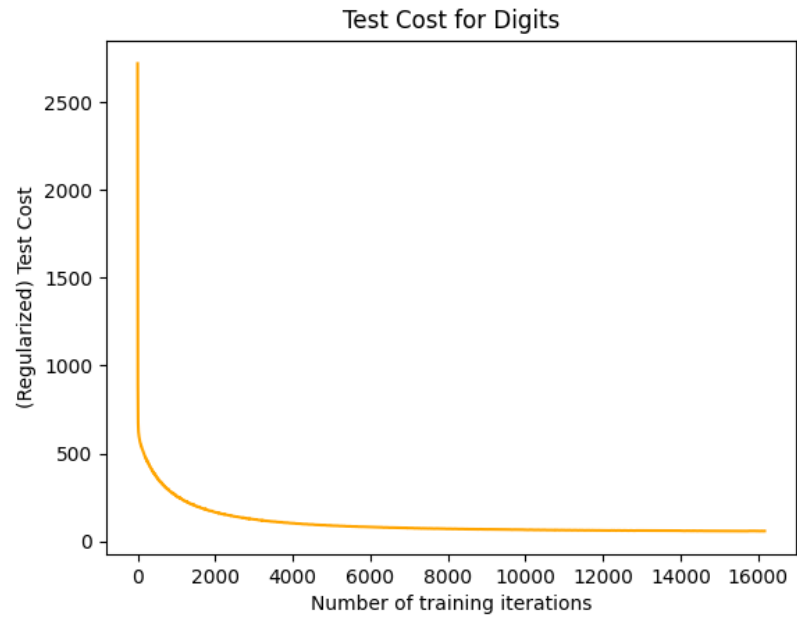
The digits dataset come's from the sci-kit learn library, available in Python. We have chosen to analyze the performance of a neural network classifier and the k-Nearest Neighbors classifier. Again, we mention that the neural network tuning was performed by *Yan Mazheika* while the kNN tuning was done by *Polina Petrova*.

We chose a neural network as one of our algorithms for this dataset because of the algorithm's ability to handle complex patterns and identify non-linear patterns in these data. Anecdotally, neural networks have shown impressive results in image classification tasks, making them a top contender for handwritten digit recognition, the focus of this dataset.

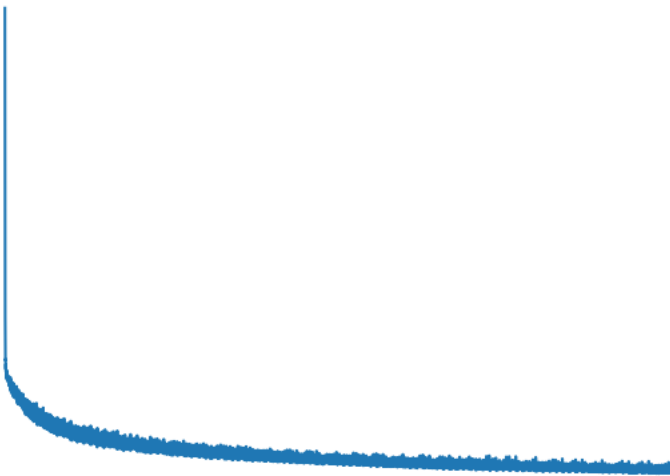
On the other hand, the k-NN algorithm is a simple yet effective classifier that makes predictions based on the closeness of past training instances in the feature space. We figured that similar pixel activations in the handwritten digits would translate into less distance in the 64-dimensional feature space. A concern we had when choosing this algorithm is precision loss; kNN performs poorly with a large number of features, which is 64 in our case. The distance measurement between two instances converges to zero as we add more features and normalize them. However, this wasn't a major problem in our analysis.

Performance Metrics

Neural Network

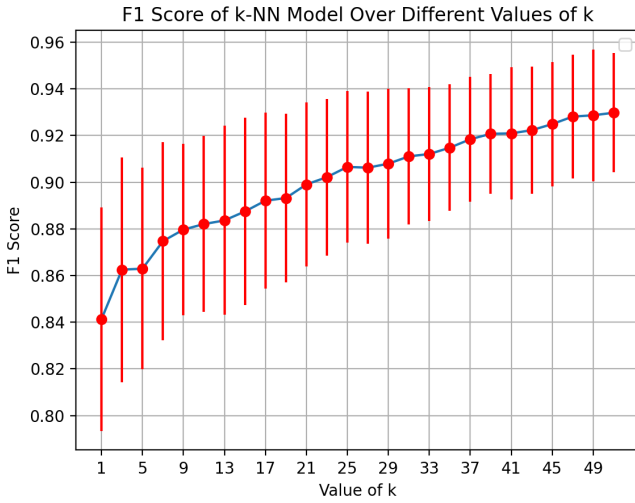
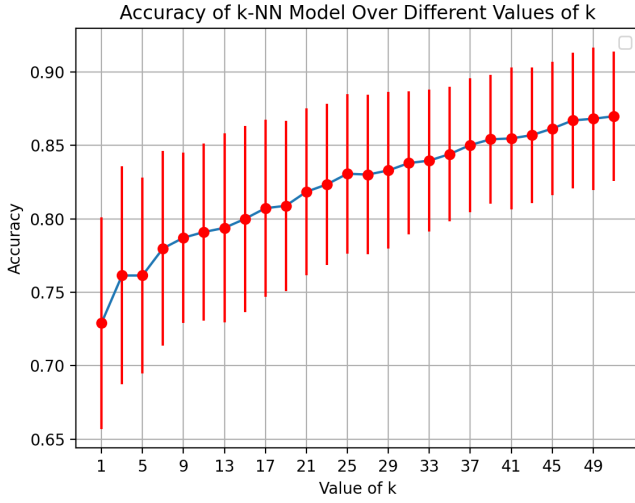


Neural Network	
Learning Rate α	0.0500
Regularization λ	0.0100
Architecture	[64, 32, 32, 10]
Mean Accuracy	0.9962
Mean F1-score	0.9809
Mean Test Cost	59.3100



Training cost over 179,700 instances

k-NN Algorithm



k-NN	
Neighbors k	51
Mean Accuracy	0.8698
Mean F1-score	0.9298

Metrics for Preferred Model Variant

k	Test Accuracy	Test F-Score
1	0.7291	0.8413
3	0.7615	0.8625
5	0.7615	0.8629
7	0.7799	0.8748
9	0.7872	0.8797
11	0.7911	0.8821
13	0.7939	0.8836
15	0.8000	0.8875
17	0.8073	0.8921
19	0.8089	0.8932
21	0.8184	0.8991
23	0.8235	0.9022
25	0.8307	0.9066
27	0.8302	0.9062
29	0.8330	0.9079
31	0.8380	0.9111
33	0.8397	0.9121
35	0.8441	0.9148
37	0.8503	0.9184
39	0.8542	0.9208
41	0.8547	0.9209
43	0.8570	0.9223
45	0.8615	0.9249
47	0.8670	0.9281
49	0.8682	0.9287
51	0.8698	0.9298

Performance Metrics over various k

Tuning Hyper-parameters

For the neural network, we've chosen an α of 0.05 and a regularization constant of $\lambda = 0.01$. At first, with a $\lambda = 0$ and $\alpha = 0.50$, the model had good performance but a quick convergence in the training curve. We found increased performance and a slower convergence by decreasing α and increasing λ to a final accuracy of 99.62%. This indicates a misclassification of 6 total instances out of 1,797 total. The architecture was chosen because it had the lowest accuracy variance between each fold out of the models tested.

In our implementation of the k-NN algorithm, we tested different (integer) values of k in the range $[1, 52)$ for an ideal hyperparameter. We observed an initial jump in the evaluation metrics between $k = 1$ and $k = 2$ and a high variance. This points to the algorithm being sensitive to outliers at very small values of k . As the value of k increases, the values of the evaluation metrics increase, as well, and stabilise by $k = 51$ at 86.98% for accuracy and 92.98% for F1-score. Increasing this hyperparameter value also led to smaller variance, which is helpful when dealing with noisy datasets. In witnessing this convergence, we have chosen a final hyperparameter value for the k-NN algorithm to be $k = 51$. With a larger k value and smoother decision boundary, we mitigate the chance of overfitting our model to the data.

Analysis

Reporting on the final runs of the neural network and k-NN algorithms on the digits dataset, both models performed well on the testing instances. However, testing on the neural network resulted in higher accuracy and F1-score values as opposed to when testing on k-NN. Dealing with an 8x8 pixel image, k-NN is able to effectively measure distances in the feature space when making predictions. Although k-NN performs well on this dataset, it struggles in distinguishing between two or more digits that look alike when handwritten. For example, an analysis of its predictions made on an instance of class 0 may return a substantial number of predictions of class 9, which would be a factor in decreasing accuracy.

The neural network acts as a well-defined alternative to this issue, as it captures more complex patterns and data variability present in handwriting. Additionally, a higher neuron count facilitates learning these handwriting variations of numbers of the same class, making the network more reliable in predicting instances that look alike and a higher accuracy.

Titanic Dataset

We pre-processed some of the titanic dataset. At first, there was a 'Sex' column which was encoded into a 'Male' column indicating a 1 if the 'Sex' of a given instance contained 'male' and a 0 if an instance contained 'female'. There was also a 'Name' column which was removed entirely. At first, we attempted to process this by encoding every English letter into a number. This resulted in a very large data-frame that led to poor results after processing.

Afterwards, we tried to one-hot encode the first word of the 'Name' column which indicated the title of a person. Because this dataset represents a time when a person's title had greater significance, we figured that there would be differences in the survival rates between prestigious and non-prestigious titles (such as 'Rev' or 'Master' versus 'Mr' or 'Miss'). We also hypothesized that persons with the title 'Rev,' which indicates religious affiliation, would be less likely to survive as we thought that their role as a religious figure would influence them to stay on the ship for longer than others. Indeed, every reverend in the dataset did not survive.

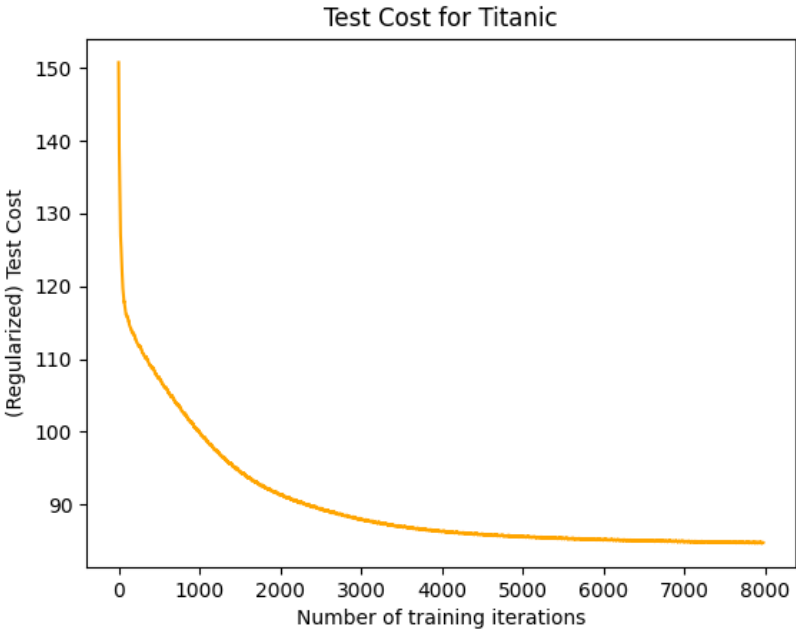
Empirically, we found that this didn't work well, which led to the removal of the 'Name' feature in the dataset. Perhaps this is because 'Pclass' contains the ticket class of a given passenger, which mitigates the role that the 'Name' column plays in indicating the socio-economic status of a passenger. Additionally, the added dimensionality of encoding around 8 titles could have mitigated the importance of the un-altered columns.

The choice of applying a neural network for this task can be justified by it's performance (speed) and ability to learn complex patterns. It's arguably the easiest model to tune with respect to the datasets this paper presents.

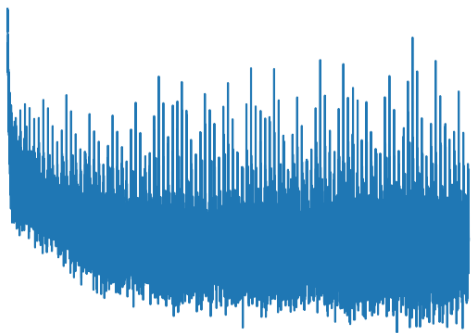
The choice of applying a decision tree for this task is natural as well. There are a number of more important features that a high-bias model can benefit from. These include passenger class, ticket fare, and gender. With three main features impacting survivability, we can benefit from a decision tree by quickly sub-setting the dataset on those features.

Performance Metrics

Neural Network



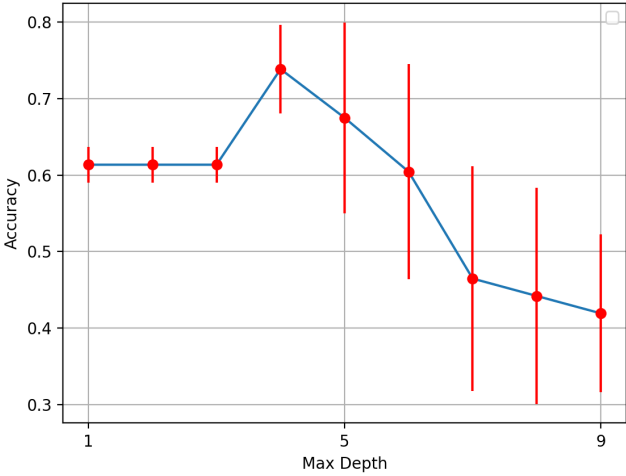
Neural Network	
Learning Rate α	0.0400
Regularization λ	0.0900
Architecture	[6, 4, 2]
Mean Accuracy	0.9663
Mean F1-score	0.9652
Mean Test Cost	84.8500



Training cost over 87,700 instances

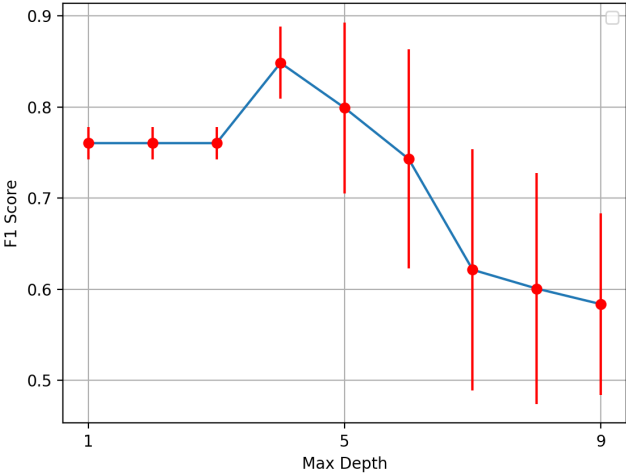
Decision Tree

Accuracy of Decision Tree Model Over Different Values of Max Depth



Maximum depth	Test Accuracy	Test F-score
1	0.7852	0.8793
2	0.6386	0.7774
3	0.5898	0.7372
4	0.7125	0.8291
5	0.6784	0.7920
6	0.6295	0.7540
7	0.5227	0.6619
8	0.4216	0.5802
9	0.4216	0.5802

F1 Score of Decision Tree Model Over Different Values of Max Depth



Decision Tree	
Maximum Depth	4
Mean Accuracy	0.7023
Mean F1-score	0.8234

Metrics for Preferred Model Variant

Tuning Hyper-parameters

The neural network was analyzed over various learning rates and regularization constants. We found that a learning rate of 0.04 and a regularization constant of 0.09 produced the best results. We found that a larger-than-usual regularization constant was necessary to reducing the variance in the model's predictions. At lower λ values, the model appeared to overfit the data, which was evident by the high variance in the training cost curve. A smaller architecture was chosen to reduce the complexity of the model as performance suffered with the presence of more than one hidden layer. This is likely due to the complexity of the dataset, which features rather arbitrary social factors.

Conducting multiple analyses of the decision tree algorithm on the titanic dataset with differing maximum tree depths, we have determined the preferred hyperparameter to be *max_depth* = 4. At this value, accuracy and F1-score reach a peak at 70.23% and 82.34%, respectively, after which both metrics decrease substantially. Deeper trees lead to overfitting and decrease in generalisation ability, which is evidenced by the sharp increase in variance on both accuracy and F-1 score after tree depth 4.

Analysis

Our decision tree struggled in providing reliable predictions on the titanic dataset. This is possibly due to the fact that the dataset contained numerical attributes such as *Ticket Fare* and *Age* that may be difficult to evaluate on this model. The decision nodes were split based on a numerical threshold that maximised information gain at each decision point. This approach is particularly challenging given the large range of numerical values present in the data. For example, there were passengers as old as 80 years-old on the ship, along with infants as young as 5 months-old. Similarly, the ticket fare values ranged from 0.0 to 512, showing a wide distribution of values. Although we employed feature normalisation, this variability in numerical attributes could have contributed to the difficulty in finding optimal splits during the tree construction process. Additionally, deeper trees tend to create more decision nodes based on numerical thresholds, which may struggle to effectively differentiate between instances with similar normalised values but different original scales.

On the other hand, employing a neural network on this dataset proved to be substantially more accurate in predicting instance classes. However, we observed a concerning trend in the neural network's performance metrics. While the accuracy on the test set was high, we witnessed a very high mean test cost, as well. Furthermore, the extreme variance in cost as the model processed more instances indicates that the neural network's predictions were inconsistent across different subsets of the dataset. This may be a consequence of class imbalance in the titanic dataset. The proportion of the class *Did Not Survive* was 61%, while for the class *Survived*, the proportion was only 38%. Our neural network model may be prioritising minimising errors on the majority class, leading to higher costs for misclassifying the minority class instances.

Loan Eligibility Dataset

The loan eligibility dataset is a binary classification problem. We have chosen to analyze the performance of a decision tree and a random forest classifier.

The choice of applying a decision tree to the problem was natural. Anecdotally, decision trees are usually taught in the context of loan eligibility problems. This is because decision trees are based on a series of if-else statements that are both easy to interpret and prioritize the most important features. In the context of this dataset, the presence of 11 total features makes some features more important than others, which will ideally be filtered for relevance by a decision tree.

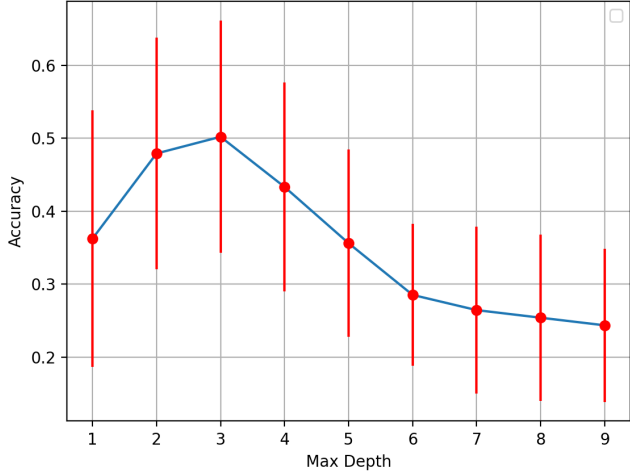
The choice of applying a random forest was also sensible. Random forests are ensembles of decision trees and aim to reduce the variance of output problems. With this algorithm, we can ideally expect a model less prone to over-fitting compared to a decision tree if the nature of the data is nuanced.

We had to pre-process this dataset as well. The provided file contained a `Loan_ID` container which the models would misinterpret as a feature and end up memorizing the target class as a function of the `Loan_ID`. Therefore, we removed this column from the dataset. To successfully process the dataset through the random forest algorithm, the classes had to be encoded as numbers but none of the features were one-hot encoded. This is because the random forest algorithm can handle categorical data without the need for this. For instance, the `graduate` feature was encoded as 0 for 'No' and 1 for 'Yes'.

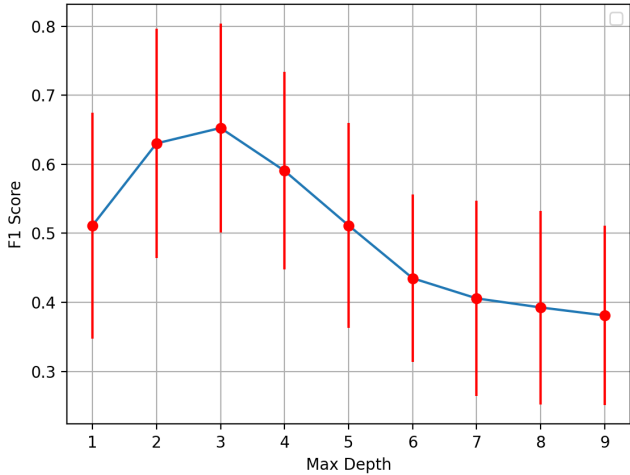
Performance Metrics

Decision Tree

Accuracy of Decision Tree Model Over Different Values of Max Depth



F1 Score of Decision Tree Model Over Different Values of Max Depth

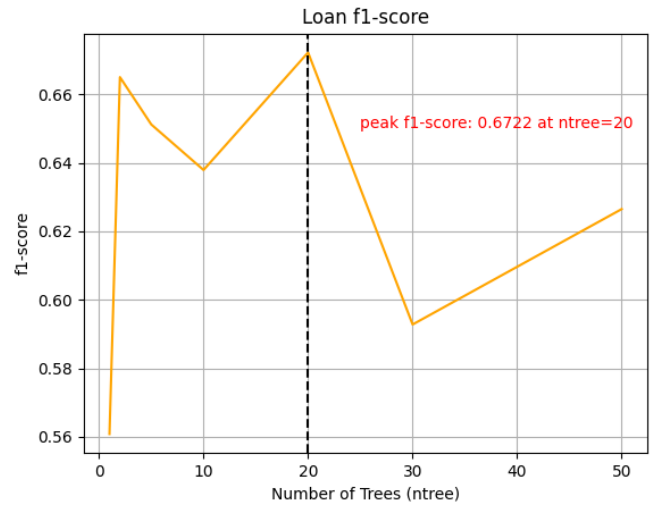
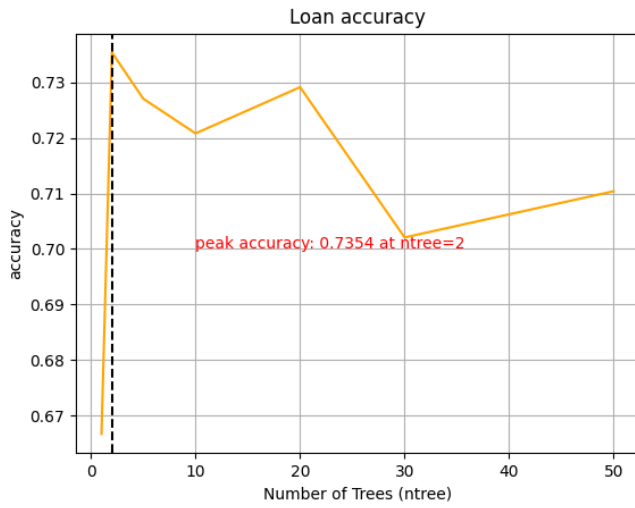


Maximum depth	Test Accuracy	Test F-score
1	0.8083	0.8930
2	0.6458	0.7657
3	0.4708	0.6357
4	0.3583	0.5139
5	0.3104	0.4606
6	0.3042	0.4609
11	0.2437	0.3894
16	0.2437	0.3894
21	0.2437	0.3894
26	0.2437	0.3894
31	0.2437	0.3894
36	0.2437	0.3894
41	0.2437	0.3894
46	0.2437	0.3894
51	0.2437	0.3894
56	0.2437	0.3894
61	0.2437	0.3894
66	0.2437	0.3894
71	0.2437	0.3894
76	0.2437	0.3894
81	0.2437	0.3894
86	0.2437	0.3894
91	0.2437	0.3894
96	0.2437	0.3894

Decision Tree	
Maximum Depth	2
Mean Accuracy	0.5354
Mean F1-score	0.6860

Metrics for Preferred Model Variant

Random Forrest



accuracy	precision	recall	f1-score	ntree
0.6667	0.5729	0.5493	0.5609	1
0.7354	0.6875	0.6440	0.6650	2
0.7271	0.6755	0.6286	0.6512	5
0.7208	0.7436	0.5585	0.6379	10
0.7292	0.8348	0.5627	0.6722	20
0.7021	0.6756	0.5281	0.5928	30
0.7104	0.7538	0.5360	0.6265	50

Performance metrics across various n-tree

Tuning Hyper-parameters

Similar to the implementation of the decision tree model for the titanic dataset, the performance of the model diminished with maximum depth values above 2. Thus, we chose the hyperparameter *max_depth* = 2 to represent this model. Furthermore, we found that repeated testing with maximum depth values of 2 and 3 produced similar evaluation values. The decision to select *max_depth* = 2 stemmed from the oscillation in performance metrics during repeated testing on *max_depth* = 3, which produced greater instability than the former hyperparameter value.

We found that the random forest model performed best with 2 ensemble trees. Perhaps other values were optimal such as *ntree* = 20, as seen by the graph above, but the variant with the highest accuracy was preferred despite the peak f1-score being at *ntree* = 20 and a relatively high accuracy as well. The stopping condition of each tree in the random forrest was the same: a minimal split entropy of 0.21. This was chosen through the method of moments, simply the value that produced the best results at 5 decision trees. Alternative stopping conditions that were tested with cross validation include the minimal split entropy, ranging from 0.01 to 0.27 in 10 equal step sizes, the maximum depth of each decision tree, ranging from 3 to 22 in 10 equal steps, and the proportion of the dataset required in each branch to perform a split, which ranged from 1% to 27% in 10 equal steps.

Analysis

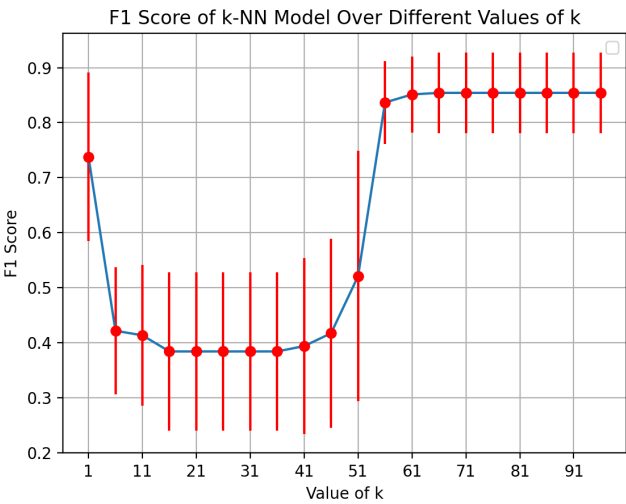
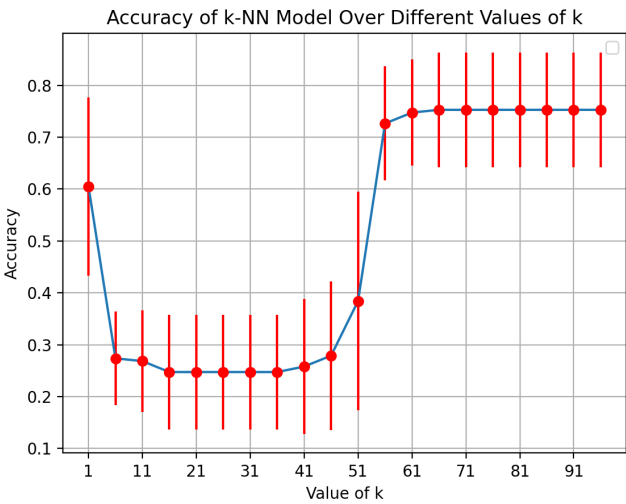
For the loan dataset, we employed the decision tree and random forest algorithms. It is crucial to note that the random forest model used entropy as a stopping criterion, while the decision tree model used tree depth as its stopping criterion. This may be a factor in why the random forest classifier performed better on both accuracy and F1-score than the decision tree classifier. That being said, both models converged to poor local optimums, indicating a lapse in classification ability for tree-based algorithms on this dataset.

It'd be difficult to compare the effect of adding an extra decision tree to result in the random forrest model because the algorithms are tuned to different stopping conditions. However, the higher performance of the random forrest model suggests that an ensemble is beneficial, given the nature of the data. This suggests that the data is nuanced and requires greater complexity to capture less-subtle patterns. Indeed, the relatively high performance of the random forest at *ntree* = 20 suggests that this is the case and that this alternative model variant could also be acceptable. The relatively similar precision and recall scores for the 2-tree ensemble suggests that the model is classifying false positives and negatives at similar rates. This is interesting since 67% of the dataset is of class 'Y'. This potentially explains the precision and recall metrics of the model at *ntree* = 20, where the model is classifying false negatives at a higher rate than false positives.

Oxford Parkinson's Disease Dataset

Performance Metrics

k-NN Algorithm

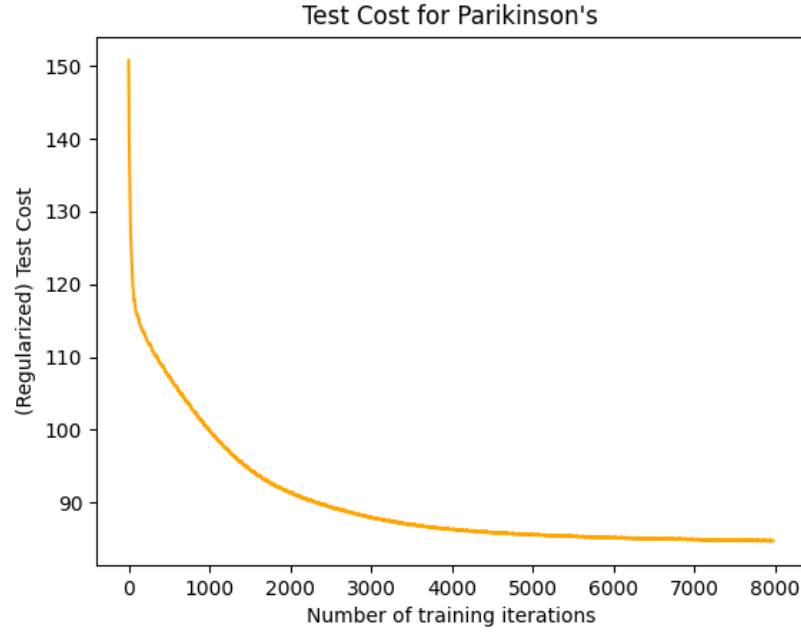


k	Test Accuracy	Test F-Score
1	0.5947	0.7367
6	0.2789	0.4319
11	0.2789	0.4319
16	0.2474	0.3921
21	0.2474	0.3893
26	0.2474	0.3893
31	0.2474	0.3893
36	0.2474	0.3893
41	0.2579	0.4030
46	0.4211	0.5625
51	0.6000	0.7320
56	0.7211	0.8350
61	0.7526	0.8561
66	0.7526	0.8561
71	0.7526	0.8561
76	0.7526	0.8561
81	0.7526	0.8561
86	0.7526	0.8561
91	0.7526	0.8561
96	0.7526	0.8561
101	0.7526	0.8561

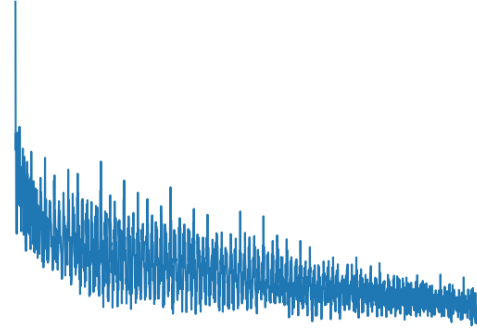
k-NN	
Neighbors <i>k</i>	61
Mean Accuracy	0.7526
Mean F1-score	0.8557

Metrics for Preferred Model Variant

Neural Network



	Neural Network
Learning Rate α	0.1000
Regularization λ	0.0001
Architecture	[22, 12, 2]
Mean Accuracy	0.9226
Mean F1-score	0.8940
Mean Test Cost	13.4800



Training Cost Over 19,500 Instances

Tuning Hyper-parameters

.7538 accuracy of random classifier.

The k-NN algorithm witnessed a large dip in performance between k values of 1 and 56. The high variance observed in predictions at these values may be indicative of the algorithm's sensitivity to local fluctuations in the training data. At lower values of k , the algorithm tends to overfit to the noise in the training data, resulting in poor generalisation performance on the test set, which is what we witnessed with our metrics dipping below 26% for accuracy and below 40% for F1-score. At $k = 61$, the algorithm finally converged to an optimum at 75.26% accuracy and 85.57% F1-score.

As this was the first k value to produce stable results, we selected it for our hyperparameter, looking to mitigate the overfitting phenomenon observed with smaller values of k .

Analysis

With regard to this dataset, we found that both of our classifiers exhibited high variance, indicating their sensitivity to fluctuations in the training data. In numerical datasets, such as this one, individual data points may exhibit significant variability, especially if they represent continuous variables with a wide range of values. For the k-NN algorithm, variance spiked between $k = 1$ and $k = 56$, representative of the overfitting issue mentioned previously. This increase in variance suggests that the algorithm's performance deteriorated as it tried to adapt to noise in the training data. However, once k hit 61, variance decreased substantially, indicating that considering a larger number of neighbors allowed the algorithm to smooth out the impact of individual data points and consider the dataset comprehensively.

A similar spike in variance was observed in the neural network algorithm. Initially, as the model processed the data, we noticed a substantial fluctuation in the cost function across different instances, indicating high sensitivity to variations in the training data. However, the algorithm converged to a more consistent behaviour the closer it got to processing 19,500 instances, reaching a mean test cost of 13.48. It may be helpful to increase the complexity of the neural network architecture to capture this data variance and learn more intricate patterns and relationships within it.