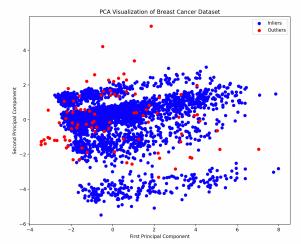
HW 3

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1. Preprocessing

First, I displayed basic characteristics of data like min, max, mean values, data types and missing values count.

Next, I proceeded with handling missing values using mean for numerical columns and mode for categorical columns. Before proceeding, I needed to transform categorical columns into numerical and for that I used OrdinalEncoding. Subsequently, I normalized the data using Z-score standardization and removed outliers using Local Outlier Factor method. This method removed 92 outliers. I then visualized the dataset by reducing the dimensionality to two using PCA:



2. Feature selection

To select the right features to train our models, I started with feature ranking using entropy. To be more specific I used sklearn's method called *mutual_info_classif*. The results showed a big drop in entropy for six columns which I decided to drop if their entropy was below 0.01.

```
HW3 /Users/aleksanderkapera/.pyenv/versions/3.12.1/bin/python "/Users/aleksanderkapera/School/Uni/NYU/Sem 3/PA/HW3/feature_selection.py" 
{'Survival Months': 0.13734327726086693, '6th Stage': 0.033974630379879356, 'N Stage': 0.02631538034934988, 'Reginol Node Positive': 0.021565849 
69569253, 'Progesterone Status': 0.017343852876732013, 'T Stage ': 0.017012253521944576, 'differentiate': 0.01536905688942472, 'Grade': 0.01320 
0536090228976, 'Estrogen Status': 0.012456459393355157, 'Marital Status': 0.007726107514020919, 'Tumor Size': 0.005848183255553252, 'A Stage': 0.005778163577640605, 'Age': 0.004611553289120218, 'Regional Node Examined': 0.0014706103187251962, 'Race': 0.0001994301537253662}

Features dropped due to low ranking: 
['Marital Status', 'Tumor Size', 'A Stage', 'Age', 'Regional Node Examined', 'Race']

Remaining features: ['T Stage ', 'N Stage', '6th Stage', 'differentiate', 'Grade', 'Estrogen Status', 'Progesterone Status', 'Reginol Node Positive', 'Survival Months', 'Status']
```

For feature selection, I experimented with forward and backward sequential feature selectors. Their results mostly overlapped and they only differed on three columns. I decided to stick with backward selection as it contained a feature that was individually ranked as the third most important, and the forward selector didn't select it.

3. Modeling

3.1. KNN

By implementing KNN, I achieved an accuracy of **87.95%**. I set k=3 based on a visual depiction of the dataset reduced to two dimensions where I visually identified 3 groups. The correctness of my KNN implementation I crosschecked with the one from sklearn library where I achieved similar results. Pros of KNN are that it's simple, understandable, and needs only one parameter (k) which can be easily optimized. Cons are that it's computationally intensive, and the performance degrades with the number of features (curse of dimensionality).

3.2. Naïve Bayes

Implementation with Naïve Bayes resulted in an accuracy of **81.49%.** The pros of Naïve Bayes are that it's simple and fast, can be trained on a low amount of training data, and performs well in multi-class classification. Cons are that it assumes all features are independent (in most cases they are not), Gaussian Naïve Bayes assumes normal distribution and may perform weakly if the data is distributed in an assumed way.

3.3. C4.5 Decision Tree

This model performed very well, scoring **86.96%** in terms of accuracy. Pros: because it's a decision tree, it's easy to visualize and explain the reasoning behind the prediction, and it can also handle both numerical and categorical data. Cons: it has tendencies to overfit, the decision tree can vary greatly when introduced to a new small set of training data, and tends to favor the majority class. Main hyperparameters: maximum depth of a tree.

3.4. Random Forest

This model achieved an accuracy of **87.06%.** Pros: because of multiple decision trees, this model is less susceptible to overfitting, it usually performs better than other forms of decision trees, and it doesn't bear any assumptions about data distribution. Cons: it is more computationally expensive (more trees to

train), it's more difficult to interpret the data (many trees), and it is sensitive to hyperparameter tuning. Main hyperparameters: number of trees in a forest, maximum depth of a tree.

3.5. Gradient Boosting

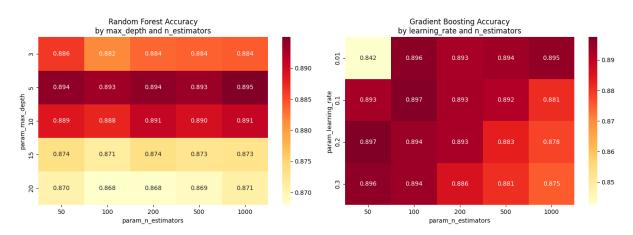
This model achieved an accuracy of **89.31%.** Pros: high performance, is flexible in terms of loss function or classification/regression. Cons: sensitive to hyperparameters, less interpretable, memory intensive. Main hyperparameters: number of stages (trees), learning rate, maximum depth of trees.

3.6. Neural Network

A simple neural network with one hidden layer containing 8 neurons achieved an accuracy of **89.81%.** Pros: flexible application, automatic feature selection, scalable. Cons: need large amounts of data to train, computationally intensive, requires many hyperparameters for tuning. Hyperparameters: network architecture (hidden layers and their sizes), learning rate, activation function, batch size.

4. Hyperparameter tuning

I chose to tune the hyperparameters of the Gradient boosting model as well as the Random forest. Below is the chart visualizing different values of hyperparameters and their respective performance. I used grid search to find the most optimal set of parameters. For Random forest, the hyperparameters that performed the best were maximum depth=5 and number of trees in the forest=1000 with accuracy=89.50%. In terms of Gradient Boosting, the best-performing combination of hyperparameters was learning rate=0.2 and number of gradient stages=50, achieving the accuracy of 89.75%.



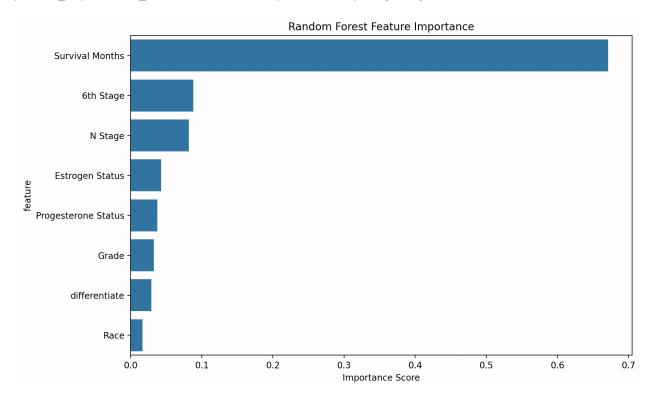
5. Results

The model that achieved the best accuracy score was neural network closely followed by, optimized gradient boosting and optimized random forest. The neural network model wasn't fine-tuned and, with more experimentation, is expected to yield even better results. Here is a table of all models:

Model	Accuracy
KNN	87.95%

Naïve Bayes	81.49%
Traite Bayes	01.1370
C4.5 Decision Tree	86.96%
Random Forest (not fine-tuned)	87.06%
Random Forest (fine-tuned)	89.50%
Gradient Boosting (not fine-tuned)	89.31%
Gradient Boosting (fine-tuned)	89.75%
Neural network	89.81%

To determine the most important feature for the random forest, I used built-in sklearn's parameter called *feature_importances_*. The most influential parameter, by a big margin, was survival months.



I then followed with the same approach for gradient boosting and received similar results. However, this time, survival months' importance was even more prevalent.

