

# MACHINE LEARNING IN HEALTHCARE: A HEAD-TO-HEAD COMPARISON OF DISEASE PREDICTION TECHNIQUES

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# AGENDA

- Problem Description
- Approach and Algorithms
- Implementation environment and code example
- Results
- Performance comparison
- Conclusion
- Contribution



# PROBLEM STATEMENT

- Dataset comprises of two CSV files containing a total of 133 columns each Out of these columns, 132 represent various symptoms that an individual may exhibit
- The last column denotes the predicted outcome or prognosis of the disease
- These symptoms are mapped to 42 diseases
- We had to train the model on training data and test it on testing data using different classification models.

# DATA PREPARATION



Removing Inconsistencies



We calculated the mean of each feature because in classification it's either 0 or 1, This also demonstrates that there is no further way to normalize the data



Vomiting and fatigue has a mean value of 0.404761905 and are considered very important characteristics, but when we remove them to see the change in accuracy, the output is different than expected. This means these functions are too general



After that we removed a feature 'fluid\_overload' which had value of mean as 0 , trained the model again, to conclude that no effect on the output

An abstract graphic on the left side of the slide, consisting of several overlapping, tilted rectangles and lines in white on a black background, creating a complex, layered geometric pattern.

# APPROACH AND ALGORITHMS USED

- Linear Regression
- Support Vector Machine (SVM)
- Decision Tree
- Random Forest
- XGBoost

# LINEAR REGRESSION

- Linear regression is a statistical method used to model the relationship between a dependent variable and one or more independent variables
- It assumes a linear relationship between the dependent variable and the independent variable(s)
- To predict the accuracy of the Prognosis, we use the Training dataset and Prognosis as the target variable

# CODE SNIPPET – LINEAR REGRESSION

## Linear Regression

```
✓ [398] from sklearn.linear_model import LinearRegression  
0s
```

```
✓ [399] lr = LinearRegression()  
0s
```

```
✓ [400] y_train = pd.get_dummies(y_train)  
0s
```

```
✓ [401] lr.fit(X_train,y_train)  
0s
```

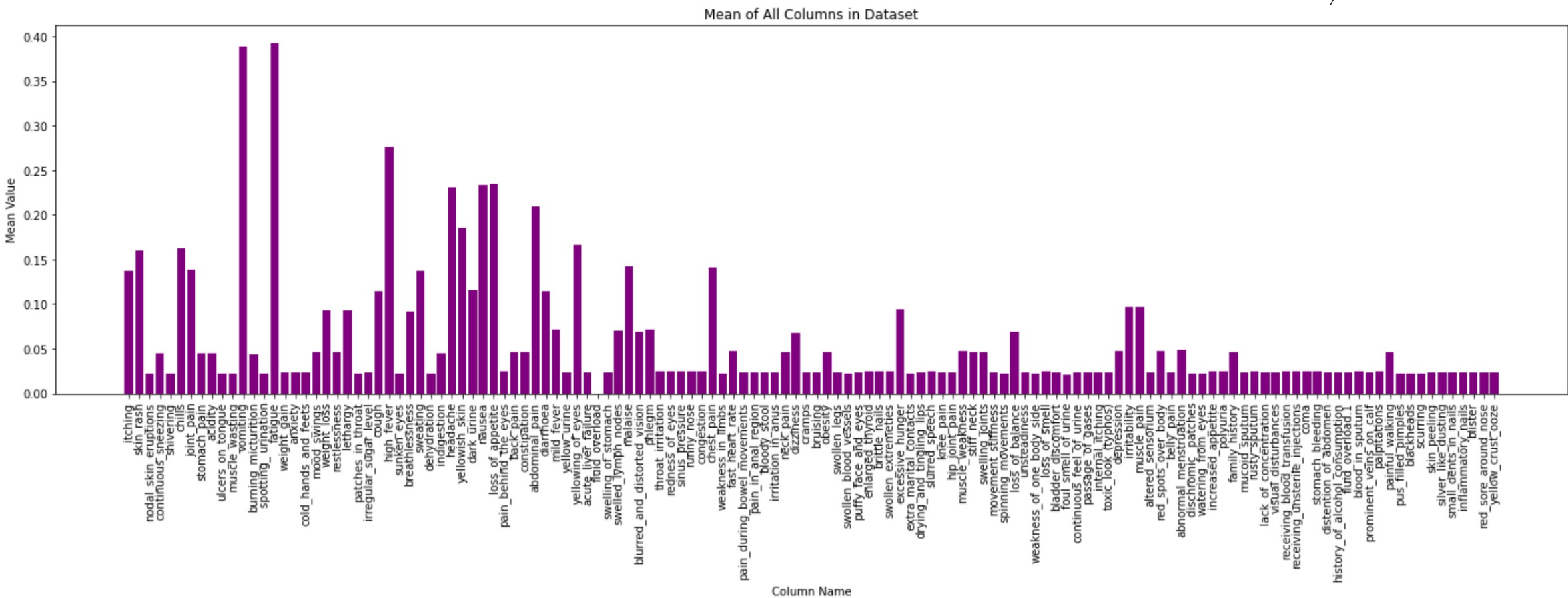
```
▼ LinearRegression  
LinearRegression()
```

```
✓ [402] y_test = pd.get_dummies(y_test)  
0s
```

```
✓ [403] a=lr.score(X_test,y_test)  
0s
```

```
✓ [404] a  
0s
```

```
0.9667341432134509
```





# Visualizing The Score with and without “fluid\_overload”

- Indicating no change to the accuracy

## Linear Regression

```
✓ [398] from sklearn.linear_model import LinearRegression
```

```
✓ [399] lr = LinearRegression()
```

```
✓ [400] y_train = pd.get_dummies(y_train)
```

```
✓ [401] lr.fit(X_train,y_train)
```

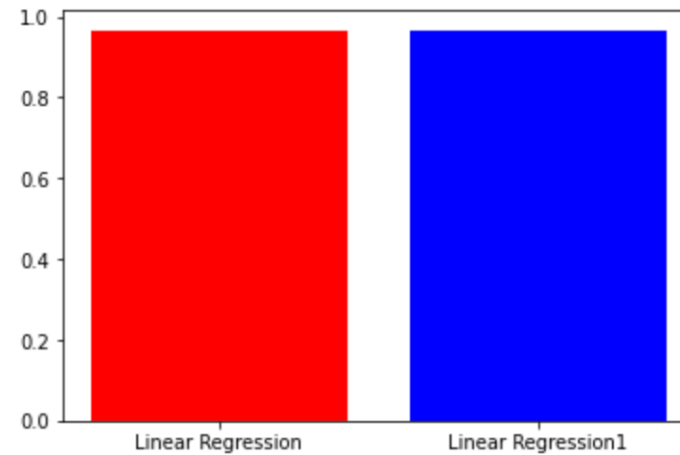
```
▼ LinearRegression  
LinearRegression()
```

```
✓ [402] y_test = pd.get_dummies(y_test)
```

```
✓ [403] a=lr.score(X_test,y_test)
```

```
✓ [404] a
```

```
0.9667341432134509
```



```
[410] X_train = X_train.drop('fluid_overload',axis='columns')  
X_test = X_test.drop('fluid_overload',axis='columns')
```

```
from sklearn.linear_model import LinearRegression  
lrl = LinearRegression()  
y_train = pd.get_dummies(y_train)  
lrl.fit(X_train,y_train)  
y_test = pd.get_dummies(y_test)  
b=lrl.score(X_test,y_test)
```

```
[412] b
```

```
0.9667234006397065
```

# SUPPORT VECTOR MACHINES (SVM)

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- Support vector machines (SVMs) are a type of supervised learning algorithm used for classification and regression
- SVMs work by finding a hyperplane that best separates classes in the data
- Kernel functions are used to transform the data into a higher-dimensional space where it can be linearly separable
- To predict the accuracy of the Prognosis, we use the Training dataset and Prognosis as the target variable

# CODE SNIPPET – SVM

## RBF

```
✓ [250] from sklearn.svm import SVC
```

```
✓ [251] rbf_model = SVC(kernel='rbf')
```

```
✓ [252] rbf_model.fit(X_train, y_train)
```

```
▼ SVC  
SVC()
```

```
✓ [253] rbf_model.score(X_test, y_test)
```

```
1.0
```

## Linear

```
✓ [254] linear_model = SVC(kernel='linear')
```

```
✓ [255] linear_model.fit(X_train, y_train)
```

```
▼ SVC  
SVC(kernel='linear')
```

```
✓ [256] linear_model.score(X_test, y_test)
```

```
1.0
```

## Sigmoid

```
✓ [257] sigmoid_model = SVC(kernel='sigmoid')
```

```
✓ [258] sigmoid_model.fit(X_train, y_train)
```

```
▼ SVC  
SVC(kernel='sigmoid')
```

```
✓ [259] sigmoid_model.score(X_test, y_test)
```

```
1.0
```

# DECISION TREE

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- A Decision Tree is a machine learning algorithm used for both regression and classification tasks. It is a simple and intuitive model that builds a tree-like structure of decisions and their possible consequences.
- The basic idea behind Decision Trees is to recursively partition the data into subsets that are as homogeneous as possible with respect to the target variable.
- Use decision tree algorithm to predict how the independent variables affects the target variable
- To predict the accuracy of the Prognosis, we use the Training dataset and Prognosis as the target variable

# CODE SNIPPET – DECISION TREE

✓ [231] `from sklearn import tree`  
0s

✓ [232] `model=tree.DecisionTreeClassifier()`  
0s

✓ [233] `model.fit(X_train,y_train)`  
0s



▼ DecisionTreeClassifier  
DecisionTreeClassifier()

✓ [234] `model.score(X_test,y_test)`  
0s

0.9761904761904762

# RANDOM FOREST

- Random Forest is a machine learning algorithm used for both regression and classification tasks. It is an ensemble method that combines multiple decision trees to make a prediction.
- The basic idea behind Random Forest is to create a forest of decision trees, each of which is trained on a random subset of the training data and a random subset of the features. The final prediction is made by aggregating the predictions of all the individual trees.
- To predict the accuracy of the Prognosis, we use the Training dataset and Prognosis as the target variable

# CODE SNIPPET – RANDOM FOREST

```
[240] from sklearn.ensemble import RandomForestClassifier
```

```
[241] model=RandomForestClassifier(n_estimators=3) #n_estimators is number of decision trees
```

```
▶ model.fit(X_train,y_train)
```



▼ RandomForestClassifier

RandomForestClassifier(n\_estimators=3)

```
[243] model.score(X_test,y_test)
```

0.9761904761904762

# XGBOOST

- XGBoost( Extreme Gradient Boosting), is an ensemble machine learning algorithm used for classification and regression tasks
- It uses a gradient boosting framework and employs multiple decision trees to make predictions
- XGBoost is known for its speed, accuracy, and ability to handle large datasets with high dimensionality



# CODE SNIPPET - XGBOOST

## Visualization

```
✓ [357] from xgboost import XGBRegressor  
s      import xgboost as xgb
```

```
✓ [358] xgb = xgb.XGBRegressor()  
s
```

```
✓ [359] xgb.fit(X_train,y_train)  
fs      xgBoostPrediction = xgb.predict(X_test)
```

```
✓ [360] xgb.score(X_test,y_test)  
s
```

```
☞ 0.9204344520364687
```

# PERFORMANCE EVALUATION

Linear Regression : 0.966

SVM: 1

Decision Tree: 0.971

Random Forest: 0.976

XGBoost: 0.92

# CONCLUSION



As per the comparative analysis of using these algorithms on the symptom-based classification problem, it is concluded that SVM fits the best as the accuracy score is 1



SVM outperformed Decision Tree in terms of classification accuracy, precision, and recall. SVM's ability to handle high-dimensional data and find the optimal hyperplane to separate the classes in the feature space made it a suitable choice for this problem.



Decision Tree had the advantage of being easier to interpret and visualize, but it lacked the accuracy and robustness of SVM

# CONTRIBUTION

Soham Shah – 100%

Arundhati Pathrikar – 100%

Mamtha Shetty – 100 %

Shantanu Pingale – 100%

A series of white, overlapping geometric lines and polygons on a black background, located on the left side of the slide.

# THANK YOU