

Configuration Manual

MSc Research Project MSc Data Analytics

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MSc Project Submission Sheet

School of Computing

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Student ID:	X20255322	Year: 2022/2023				
Programme:	MSc Data Analytics	Year:	2022/2023			
Module:	Research Project					
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Date:	 E: Lilian Ifeoma Enwereobi X20255322 MSc Data Analytics Year: 2022/2023 Research Project Qurrat UI Ain 14-08-2022 Stoke prediction Using Model Comparison and Feature Selections 					
Project Title:	Stoke prediction Using Model Comparison and Feature Selections					

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Configuration Manual

Lilian Ifeoma Enwereobi Student ID:

1 Introduction

Python was implemented to create algorithms based on machine learning and Deep Learning techniques for Stroke Prediction Research. In this manual, the system's specs and usage are described in depth.

2 Technical Specifications

2.1 Hardware Required

System Software: Windows 10, 64 bits.

RAM: 12 GB

These are the only hardware-related criteria that have been met.

2.2 Software Required

You'll require to have Anaconda Jupyter Notebook installed to run the Python code. At the command prompt, type cd /some folder name to access the start-up folder. Enter jupyter notebook in the keyword search box to launch the Jupyter Notebook app. A fresh page or window in your browser will open with the user interface for the laptop.

Computation Syntax: Python

The most recent version of Anaconda, 5.0.0, as well as the Jupyter Notebook modifications it offers, are necessary since they let users define environment-specific kernels from the Jupyter Notebook interface. Designed to simplify package management and deployment, the anaconda is a package management system of the Python and R computational science programming languages.

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	Dataspell is an IDE for exploratory data enslysis and prototyping machine learning models. It combines the interactivity of Jupyter notebooks with the intelligent Python and R coding assistance of PyCherm	Run a cmd.exe terminal with your current environment from Nevigetor activeted	An extensible environment for interactive end reproducible computing, based on the Jupyter Notebook and Architecture.	Web-based, interactive computing notebook environment. Edit and run human-readable docs while describing the data analysis.	0.0.1 Run a Powershelt terminal with your current environment from Navigator activated	PyQt Gus that supports inline figures, proper multiline editing with syntax highlighting, graphical califips, and mo
	in one user-friendly environment.	Launch	Launch	Launch	Launch	Launch
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notebooks with eds of packages	Scientific Prython Development EnviRonment, Powerful Python IDE with advanced editing, interactive testing, debugging and introspection features	Kick-start your data science projects in seconds in a pre-configured environment. Enjoy coding assistance for Python, SQL, and R in Jupyter notebooks and benefit from no-code submations. Use Datafore	Deepnote is a notebook built for collaboration. Create notebooks in your browser, spin up your conda environment in seconds and share with a link.	IBM Watson Studio Cloud provides you the tools to enelyze end visualize data, to cleanse and shape data, to create and train mathine learning models. Prepare data and build models, using open source data	OCI Data Science offers a machine learning platform to build, trein, manage, and deploy your machine learning models on the cloud with your favorite open-source tools	1.9.0
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Fig 1: Anaconda Navigator Window

3. Collecting Dataset

The dataset can be obtained at https://www.kaggle.com/datasets/prosperchuks/health-dataset. There will be a file called Diabetes, Hypertension, and Stroke Prediction Dataset there, and downloading it only requires a single mouse click.

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(L)	Diabetes, Hypertension and Stroke Prediction	
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፠	About this file	IIII Stroke_data.csv
<>	This dataset has 17 feature variables and 1 target variable and its classes are balanced.	
	# Age = # Sex = # HighChol = # CholCheck 13-level age category (_AGEG5YR sec codebook) 1 = 18-24 9 = 60-64 13 = 80 or older patient's gender (1: male; 0: female). 0 = no high cholesterol 1 = high cholesterol 0 = no cholesterol in 5 years 1 = yes cholesterol check years	# BMI ol check Body N s s

Fig 2: Kaggle Dataset Repository

4. **Downloading data for Jupyter Notebook:** The dataset must first be saved on any system-available local drive. It was saved as an archive because it has 3 datasets, so I separated it to take the dataset I needed.

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Fig 2: My Download Drive held the dataset.

You can use the ANACONDA navigator to open the Jupyter Notebook version. The software used for the development of the notebook version makes use of the Python programming language.

5. Package Installations and Library Importing

The following libraries were utilized in this study for data pre-processing, model construction, and model evaluation:

```
import warnings
import numpy as np
import pandas as pd
from sklearn.model_selection import train_test_split
import matplotlib.pyplot as plt
import seaborn as sns
from sklearn.metrics import confusion_matrix, accuracy_score, f1_score, classification_report, precision
import matplotlib.pyplot as plt
warnings.filterwarnings('ignore')
from tensorflow.keras.models import Sequential
from tensorflow.keras.layers import Dense
```

Fig 3: Imported Libraries and Packages

```
from sklearn.svm import SVC
from sklearn.ensemble import RandomForestClassifier
from sklearn.neighbors import KNeighborsClassifier
from boruta import BorutaPy
from lightgbm import LGBMClassifier
from sklearn.ensemble import GradientBoostingClassifier
from sklearn.ensemble import AdaBoostClassifier
from sklearn.preprocessing import StandardScaler,MinMaxScaler
from tensorflow.keras.wrappers.scikit_learn import KerasClassifier
from sklearn.model_selection import RandomizedSearchCV
from sklearn.metrics import accuracy_score , classification_report,ConfusionMatrixDisplay,precision_sco
from sklearn.model_selection import RandomizedSearchCV, cross_val_score, KFold
from sklearn.tree import DecisionTreeClassifier
from sklearn.neural_network import MLPClassifier
from sklearn.feature_selection import SelectKBest
from sklearn.feature_selection import chi2
from sklearn.preprocessing import QuantileTransformer
from sklearn.linear_model import LogisticRegression
from mlxtend.feature_selection import ExhaustiveFeatureSelector
from scipy.stats import yeojohnson
```

Fig 4: Imported libraries and Packages

To create frameworks for Stroke Prediction Using Model Comparison and Feature Selection, these libraries were imported into the Jupyter Notebook.

Figure 5 illustrates the process of installing the Boruta, TensorFlow, Keras, lightgbm, mlxtend, and Plotly packages. Installing this package is required to predict stroke.

pip install boruta
pip install lightgbm
pip install mLxtend
pip install tensorflow
pip istall keras
pip install plotly

Fig 5: Package Installation

As seen in Figure 6, the data is initially inserted through the Data frame and confirmed.

<pre>data = pd.read_csv("C:\\Users\\lilia\\Desktop\\stroke_data.csv")</pre>										
data										
	sex	age	hypertension	heart_disease	ever_married	work_type	Residence_type	avg_glucose_level	bmi	smoking_stat
0	1.0	63.0	0	1	1	4	1	228.69	36.6	
1	1.0	42.0	0	1	1	4	0	105.92	32.5	
2	0.0	61.0	0	0	1	4	1	171.23	34.4	
3	1.0	41.0	1	0	1	3	0	174.12	24.0	
4	1.0	85.0	0	0	1	4	1	186.21	29.0	
40905	1.0	38.0	0	0	0	4	1	120.94	29.7	
40906	0.0	53.0	0	0	1	4	0	77.66	40.8	
40907	1.0	32.0	0	0	1	2	0	231.95	33.2	
40908	1.0	42.0	0	0	1	3	0	216.38	34.5	
40909	1.0	35.0	0	0	0	4	0	95.01	28.0	
40910 rows × 11 columns										

Fig 6: Caption Data File

6. Checking For Zero Values in some Features and Replacing them with Median

```
#checking for 0 values in 4 columns
print(data[data['hypertension']==0].shape[0])
print(data[data['smoking_status']==0].shape[0])
print(data[data['age']==0].shape[0])
32162
35685
20921
23
# #replacing 0 values with median of that column
data['hypertension']=data['hypertension'].replace(0,data['hypertension'].mean())#normal distribution
data['heart_disease']=data['heart_disease'].replace(0,data['heart_disease'].mean())#normal distribution
data['smoking_status']=data['smoking_status'].replace(0,data['smoking_status'].median())#skewed distribution
```

data['age']=data['age'].replace(0,data['age'].median())#skewed distribution
Fig 7: Checking And Replacing the Zero Values

7. Exploratory Data Analysis

```
# Bivariate bar plot for categorical variables
# All data columns
feature_cols = data.columns
plt.figure(figsize = (30,50))
plt.suptitle('stroke by categorical features')
#subplots
for i in enumerate(feature_cols):
    plt.subplot(2,4, i[0]+1)
    x = sns.countplot(data=data, x=i[1], hue='stroke', palette = ['blue','crimson'])
    for z in x.patches:
        x.annotate('{:.1f}'.format((z.get_height()/data.shape[0])*100)+'%',(z.get_x()+0.25, z.get_height()/data.shape[0])*100)+'%',(z.get_x()+0.25, z.get_height()/data.shape[0])*100)+'%')
```

Fig 8: Code for Stroke Proportion Bar Plot

Figure 8 shows the percentage of people who have experienced a stroke compared to those who have not.

7.1. Analysis of All the Features

```
# All data columns
feature_cols = data.columns
plt.figure(figsize=(25, 40))
# Loop for subplots
for i in range(len(feature_cols)):
    plt.subplot(9, 5, i+1)
    plt.title(feature_cols[i])
    plt.xticks(rotation=90)
    plt.hist(data[feature_cols[i]], color="blue")
plt.tight_layout()
```

Fig 9: Code for Histograms of all the Features

8. Implementation and Evaluation

8.1. Feature Selection

Three feature selection was used to select important features by eliminating the features that are relevant to stroke predictions.

Boruta Feature selection model

```
# Fit Boruta on the dataset
poruta_selector = BorutaPy(model, n_estimators='auto', verbose=2, random_state=42)
poruta_selector.fit(X.values, y.values)
# Check the selected features
selected_features = X.columns[boruta_selector.support_]
# Print the feature importances
feat_importances = pd.Series(boruta_selector.ranking_, index=X.columns)
plt.figure(figsize=(8, 6))
feat_importances.nsmallest(6).plot(kind='barh')
plt.show()
```

Fig 10: Code for implementing Boruta.

SelectKBest Feature Selection Model

```
#apply SelectKBest class to extract top 5 best features #Do this before quantile transformation
# Initialize the SelectKBest class for feature selection
best_features = SelectKBest(score_func=chi2, k=5)
# Fit the feature selector on the data
fit = best_features.fit(X_transformed, y)
# Get the scores and feature names
feature_scores = pd.DataFrame(fit.scores_)
feature_names = pd.DataFrame(X.columns)
# Concatenate the feature scores and names
feature_scores_df = pd.concat([feature_names, feature_scores], axis=1)
feature_scores_df.columns = ['Specs', 'Score']
# Print the top 5 best features
print(feature_scores_df.nlargest(5, 'Score'))
```

Fig 11: Code for Implementing SelectKBest.

Exhaustive Feature Selection Model

Fig 12: Code for Implementing Exhaustive Model

All the models were hyper-tuned with RandomSearchCV and these are the codes we used on each model and their estimators

Gradient Boosting Classifier (XGB)

```
param grid xgb = {
    'n_estimators': [50, 100, 200],
'learning_rate': [0.1, 0.01, 0.001],
    'max_depth': [3, 5, 7],
    'subsample': [0.8, 0.9, 1.0],
}
# Create the RandomizedSearchCV objects for each model
random_search_xgb = RandomizedSearchCV(
    estimator=GradientBoostingClassifier(),
    param_distributions=param_grid_xgb,
    n_iter=10,
    scoring='accuracy',
    cv=5,
    verbose=3,
    random_state=42,
    n_jobs=-1
)
```

Fig 13: XGB Hyperparameter Tuning.

Figure 13 shows the code for utilizing RandomSearchCV and 10-fold cross-validation to discover the optimal parameters for XGB. The n estimator for XGB is 100.

AdaBoosting Classifier (Adaboost)

```
param_grid_ada = {
    'n estimators': [50, 100, 200],
    'learning_rate': [0.1, 0.01, 0.001]
3
from sklearn.ensemble import AdaBoostClassifier
from sklearn.tree import DecisionTreeClassifier
# Define the parameter grid for the base estimator (DecisionTreeClassifier)
param_grid_base = {
    'max_depth': [1, 2, 3],
'min_samples_split': [2, 5, 10],
    'min_samples_leaf': [1, 2, 3]
3
# Create an instance of the base estimator
base estimator = DecisionTreeClassifier()
# Perform RandomizedSearchCV for AdaBoostClassifier
random_search_ada = RandomizedSearchCV(
    AdaBoostClassifier(base_estimator=base_estimator),
    param_grid_ada,
    cv=5
)
random_search_ada.fit(X_train, y_train)
```

Fig 13: Adaboost Hyperparameter Tuning.

Figure 13 shows the code for utilizing RandomSearchCV and 10-fold cross-validation to discover the optimal parameters for the Adaboost classifier. The n_estimators for the Adaboost classifier is 50.

Light Gradient Boosting Classifier (LightGBM)

```
param_grid_lgbm = {
      'boosting_type': ['gbdt', 'dart'],
      'num_leaves': [31, 63, 127],
     'learning_rate': [0.1, 0.01, 0.001],
      'n_estimators': [100, 200, 300],
     'max_depth': [5, 10, 15],
'min_child_samples': [20, 50, 100],
'subsample': [0.8, 0.9, 1.0],
'colsample_bytree': [0.8, 0.9, 1.0]
3
random_search_lgbm = RandomizedSearchCV(
     estimator=LGBMClassifier(),
     param_distributions=param_grid_lgbm,
     n_iter=10,
     scoring='accuracy',
     cv=5.
     verbose=3.
     random_state=42,
     n jobs=-1
)
random_search_lgbm.fit(X_train, y_train)
best_params_lgbm = random_search_lgbm.best_params_
best_lgbm_model = LGBMClassifier(**best_params_lgbm)
cv_scores_lgbm = cross_val_score(best_lgbm_model, X_train, y_train, cv=k_fold, scoring='accuracy')
average_cv_score_lgbm = np.mean(cv_scores_lgbm)
print("LightGBM - Best Hyperparameters:", best_params_lgbm)
print("LightGBM - Average Cross-validation Score:", average_cv_score_lgbm)
```

Fig 14: LightGBM Hyperparameter tuning.

Figure 14 shows the code for utilizing RandomSearchCV and 10-fold cross-validation to discover the optimal parameters for the LightGBM classifier. The n_estimators for the LightGBM classifier is 300.

Random Forest (RF)

```
param_grid_rf = {
     'n_estimators': [100, 200, 300], # Number of trees in the forest
     'criterion': ['gini', 'entropy'], # Split quality criterion
'max_depth': [None, 5, 10], # Maximum depth of the tree
     'min_samples_split': [2, 5, 10], # Minimum number of samples required to split an internal node
     'min_samples_leaf': [1, 2, 4], # Minimum number of samples required to be at a leaf node
'max_features': ['auto', 'sqrt', 'log2'], # Number of features to consider at each split
     'bootstrap': [True, False] # Whether bootstrap samples are used when building trees
}
random_search_rf = RandomizedSearchCV(
    estimator=RandomForestClassifier();
     param_distributions=param_grid_rf,
     n iter=10.
     scoring='accuracy',
     cv=5,
     verbose=3.
     random_state=42,
     n_jobs=-1
)
random_search_rf.fit(X_train, y_train)
```

Fig 15: Random Forest Hyperparameter Tuning

Figure 15 shows the code for utilizing RandomSearchCV and 10-fold cross-validation to discover the optimal parameters for the Random Forest classifier. The n_estimators for the Random Forest classifier is 200.

Artificial Neural Networks (ANN)

```
ann_classifier = KerasClassifier(build_fn=create_model)

param_grid_ann = {
    'hidden_layer_sizes': [(32,), (64,), (128,)],
    'activation': ['relu', 'tanh'],
    'solver': ['adam'],
    'learning_rate_init': [0.001, 0.01, 0.1],
    'alpha': [0.0001, 0.001, 0.01]
}

random_search_ann = RandomizedSearchCV(MLPClassifier(), param_grid_ann, cv=5)
random_search_ann.fit(X_train, y_train)
```

Fig 16: Artificial Neural Network Hyperparameter Tuning

Figure 15 shows the code for utilizing RandomSearchCV and 10-fold cross-validation to discover the optimal parameters for the Artificial Neural Networks.

9. Conclusion

You can adhere to the following steps to carry out the complete piece of code in Jupyter successfully. The system's 12 gigabytes of random-access memory guarantee that the code will run more swiftly and without hiccups.

10. References

Dritsas, E. and Trigka, M., 2022. Stroke risk prediction with machine learning techniques. Sensors, 22(13), p.4670. Sailasya, G. and Kumari, G.L.A., 2021. Analyzing the performance

of stroke prediction using ML classification algorithms. International Journal of Advanced Computer Science and Applications, 12(6)