

BVRIT HYDERABAD College of Engineering for Women



Department of Information Technology

BREAST CANCER PROGNOSIS USING MACHINE LEARNING

Under the guidance of

Guide Name: Mr. B.Srinivasulu

Designation: Assistant Professor

Team- 09

R. Parimala (19WH1A1268)

P. Aditi Kiran(19WH1A1277)

Ch. Lakshmi Durga(19WH1A1295)

Raveena Yadlapalli (19WH1A1296)







- Summary of Stage 1
- Implementation of Experimental Design
- Execution video
- Analysis of Results, Discussion
- Conclusion & Future scope



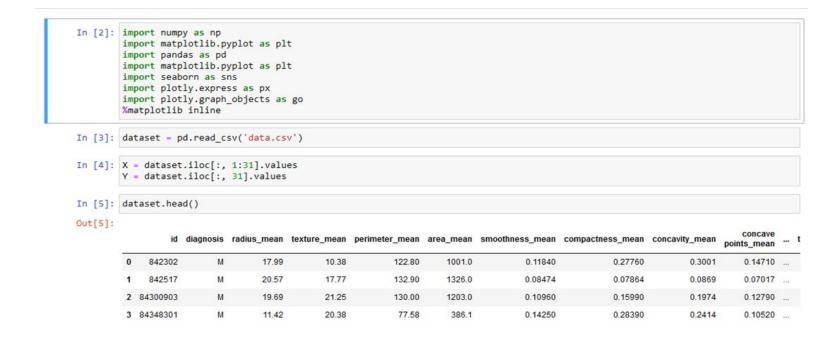


Summary of Stage 1

On the Wisconsin Breast Cancer Diagnostic dataset (WBCD) we applied five main algorithms which are: SVC, Random Forests, Logistic Regression, Decision Tree, K -NN, to calculate, compare and evaluate different results obtained based on confusion matrix, accuracy, sensitivity, precision, AUC to identify the best machine learning algorithm that are precise, reliable and find the higher accuracy. All algorithms have been programmed in Python using scikit-learn library in Jupyter notebook environment. After an accurate comparison between our models, we found that Support Vector Classifier achieved a higher Accuracy of 96.49%, efficiency of 97.2%, Precision of 97.5%, AUC of 96.6% and outperforms all other algorithms.

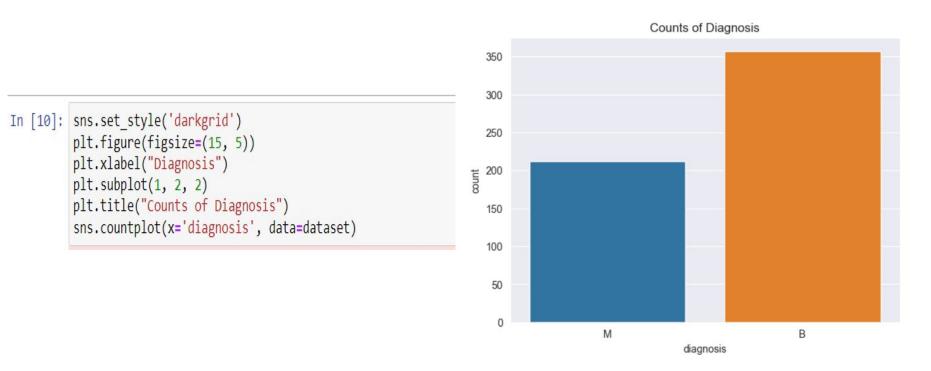
















```
models_list = {
    "LogisticRegression" : LogisticRegression(random_state=0),
    "K-NearestNeighbor" : KNeighborsClassifier(n_neighbors = 5, metric = 'minkowski', p = 2),
    "SVC" : SVC(kernel = 'rbf', random_state = 2,C = 2),
    "SVM" : SVC(kernel = 'linear', random_state = 0),
    "NaiveBayes" : GaussianNB(),
    "DecisionTreeClassifier" : DecisionTreeClassifier(criterion='entropy', random_state=0),
    "RandomForestClassifier" : RandomForestClassifier(n_estimators=10, criterion='entropy', random_state=0),
}
```





```
#train test splitting
from sklearn.model_selection import train_test_split
from sklearn.linear_model import LogisticRegression
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier
from sklearn.naive_bayes import GaussianNB
from sklearn.neighbors import KNeighborsClassifier
from sklearn.metrics import accuracy_score, confusion_matrix, f1_score
from sklearn.metrics import classification_report
from sklearn.model_selection import KFold
from sklearn.model_selection import cross_validate, cross_val_score
from sklearn.svm import SVC
from sklearn import metrics
```



Execution video









Analysis of Results

The accuracies of different algorithms are as follows:

| | model_name | score | accuracy_score | accuracy_percentage |
|---|------------------------|----------|----------------|---------------------|
| 5 | DecisionTreeClassifier | 1.000000 | 0.883041 | 88.30% |
| 6 | RandomForestClassifier | 0.994975 | 0.918129 | 91.81% |
| 2 | SVC | 0.962312 | 0.964912 | 96.49% |
| 1 | K-NearestNeighbor | 0.937186 | 0.964912 | 96.49% |
| 0 | LogisticRegression | 0.934673 | 0.959064 | 95.91% |
| 3 | SVM | 0.932161 | 0.959064 | 95.91% |
| 4 | NaiveBayes | 0.912060 | 0.929825 | 92.98% |
| | | | | |

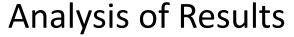




Analysis of Results

On comparing the accuracies of different algorithms, we observe that SVC (Support Vector Classifier) and KNN (k-nearest neighbors) shows best accuracy percentage, but we choose SVC because KNN is easier to interpret but can identify only a limited set of patterns.







```
import random
a = random.random()
meanR=round(random.uniform(0,30),3)
meanT=round(random.uniform(0,50),3)
meanS=round(random.uniform(0,1),5)
meanC=round(random.uniform(0,1),5)
meanSy=round(random.uniform(0,1),4)
meanF=round(random.uniform(0,1),5)
seR=round(random.uniform(0,2),4)
seT=round(random.uniform(0,3),4)
seS=round(random.uniform(0,1),6)
seC=round(random.uniform(0,1),6)
seSy=round(random.uniform(0,1),6)
seF=round(random.uniform(0,1),6)
input 1=[]
lst = map(lambda x : x[1], filter(lambda x : x[0].startswith('mean'), globals().items()))
for i in 1st:
    input 1.append(i)
lst1 = map(lambda x : x[1], filter(lambda x : x[0].startswith('se'), globals().items()))
for i in 1st1:
    input 1.append(i)
input array = np.asarray(input 1)
input reshaped = input array.reshape(1,-1)
predict = model.predict(input reshaped)
if (predict[0] < 0.5):</pre>
 print("breast cancer is malignant")
else:
  print("breast cancer is benign")
```

The result of the data given is that the cancer is benign.

breast cancer is benign







```
a = random.random()
meanR=round(random.uniform(0,30),3)
meanT=round(random.uniform(0,50),3)
meanS=round(random.uniform(0,1),5)
meanC=round(random.uniform(0,1),5)
meanSy=round(random.uniform(0,1),4)
meanF=round(random.uniform(0,1),5)
seR=round(random.uniform(0,2),4)
seT=round(random.uniform(0,3),4)
seS=round(random.uniform(0,1),6)
seC=round(random.uniform(0,1),6)
seSy=round(random.uniform(0,1),6)
seF=round(random.uniform(0,1),6)
input 2=[]
lst = map(lambda x : x[1], filter(lambda x : x[0].startswith('mean'), globals().items()))
for i in 1st:
    input 2.append(i)
lst1 = map(lambda x : x[1], filter(lambda x : x[0].startswith('se'), globals().items()))
for i in 1st1:
    input 2.append(i)
input array = np.asarray(input 2)
input_reshaped = input_array.reshape(1,-1)
predict = model.predict(input_reshaped)
if (predict[0] == 1):
  print("breast cancer is malignant and more chances to occur")
else:
   print("breast cancer is benign and more chances to occur")
```

Prediction of Recurrence module

The recurrence is predicted on dataset based on diagnosis and tumour size.

breast cancer is malignant and more chances to occur





Conclusion

- We have explored different prediction models and various evaluation methods.
- We measuring the performance of the models using real data.
- 5 popular classification algorithms have been used and after comparing these algorithms with different performance metrics we concluded that SVC is the best machine learning algorithm to predict breast cancer.

Future Scope

We discovered that SVC and KNN provide the most accurate results for predicting breast cancer. The accuracy of this work can be improved in the future by altering the currently used machine learning approaches or by creating new algorithms. The future of breast cancer prediction is machine learning, and machine learning models are improving daily thanks to the tremendous research being done in this area by researchers. Artificial intelligence and machine learning will revolutionise the medical sector in the next decades. The prediction of breast cancer will perform better than conventional pathology testing with the addition of cutting-edge technology like convolutional neural networks.





References

- [1] Usman Naseem, Junaid Rashid, "An Automatic Detection of Breast Cancer Diagnosis and Prognosis Based on Machine Learning Using Ensemble of Classifiers" 12 May 2022
- [2] Sharma, A. & Mishra, P. K. Performance analysis of machine learning based optimized feature selection approaches for breast cancer diagnosis. *Int. J. Inf. Tecnol.* 14, 1949–1960. (2022).
- [3] Ahmad, S. et al. A novel hybrid deep learning model for metastatic cancer detection. *Comput. Intell. Neurosci* (2022).
- [4]L Yang, B Fu, Y Li, Y Liu, W Huang, S Feng et al., "Prediction model of the response to neoadjuvant chemotherapy in cancers by a Naive Bayes algorithm", Computer methods and programs in biomedicine, vol. 192, pp. 105458, 2020.





THANK YOU