Technical Communication

Biosci. Biotech. Res. Comm. 12(3): 741-747 (2019)



Significance of Accuracy Levels in Cancer Prediction using Machine Learning Techniques

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ABSTRACT

Across the world, any cancer becomes a calamity for a person who is suffering from it, mainly women are facing a real challenge when it comes to breast cancer. Breast cancer can be diagnosed at an early stage to overcome the consequences at a later stage. In the field of Computer Science, Machine Learning (ML) techniques are competent enough to diagnose the stages of cancer. ML techniques work upon the data which are collected from hospitals of suspected patients. There are various ML techniques which can build a model in order to diagnose cancer on the basis of finding accuracy level. In this paper, we have discussed the significance of accuracy level for predicting the cancer. In previous works, it has been observed that 100% accuracy is found on data analysis by some researchers. Although 100% accuracy must have given perfect prediction but it is observed that prediction was not so, sometimes it gives incorrect prediction also. So, prediction technique is scaled up with inclusion of more parameters precision, recall, F1- measure, Receiver Operating Characteristics (ROC) area and Area Under Curve (AUC) score.

KEY WORDS: ACCURACY, AUC, CANCER, F1-MEASURE, MACHINE LEARNING, PRECISION, RECALL, ROC

INTRODUCTION

Human body is made up of billions of cells and when cells start growing, becomes lumps and later on develops into tumor. Tumor is of two types: Malignant and Benign. Malignant is dangerous that causes another tumor and consequently cancer. Benign tumor usually does not cause cancer in short span of time. In this paper, there is a discussion of breast cancer which is due to internal or external disbalance of hormonal activities.

There are some traditional methods to diagnose and prognose the cancer but sometimes these methods take

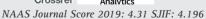
ARTICLE INFORMATION:

Corresponding Author: kumarajay7th@gmail.com Received 5th June, 2019

Accepted after revision 29th Aug, 2019 BBRC Print ISSN: 0974-6455

Online ISSN: 2321-4007 CODEN: USA BBRCBA





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Online Contents Available at: http://www.bbrc.in/

DOI: 10.21786/bbrc/12.3/29

long time to diagnose. A Machine Learning is a field of Artificial Intelligence (AI) in computer science which can implement many computational intelligent techniques for the fast and accurate prediction of cancer. ML techniques can precisely predict different type of tumors. It motivated us to work on cancer prediction technique and working to find alternative parameters for better prediction.

This paper comprises of four sections. In section 1, Introduction of cancer has been described. Section 2 summarizes the previous related work. In section 3, study on various machine learning techniques have been presented. Section 4 presents the various vital parameters including accuracy for breast cancer prediction. In section 5, breast cancer prediction performance metrics is discussed. In section 6, experiments and results are shown and in the last section, conclusion and discussion are explained.

Ahmad et al. (2013) compared 3 ML techniques viz. Decision Tree (C4.5), ANN and SVM on Iranian Centre for Breast Cancer datasets of 1189 patient and found the accuracy as 93.6%, 94.7% and 95.7%. Ali et al. (2019) proposed a prediction model using "big data" to explore feature selection and cross-validation in omics file datasets. Further described anti-drug drug response modelling and predicting their phenotypic responses. Asri et al. (2016) compared different ML algorithms: SVM, C4.5, NB and kNN for WBCD dataset which has 699 instances and 11 integer-valued attributes. Among all algorithms, SVM gave the highest accuracy 97.13% with lowest error rate conducted in WEKA data mining tool.

Bevilacqua et al. (2006) used IDEST novel approach based on ANN for WBCD datasets and found 98.6% accuracy. Boughorbel et al. (2017) focused on another metric known as Matthews Correlation Coefficient (MCC) to handle imbalance data using SVM and Bayes classifier. Burt et al. (2018) diagnosed a breast cancer with deep learning network using a system Computer-aided detection and diagnosis (CAD). It mainly looks for images captured by MRI, X-rays and compared with human expertise like radiologists, clinicians etc. Bychkov et al. (2018) took a sample of 420 images of colorectal cancer and apply deep learning outperformed AUC 0.69 which is better than AUC 0.58 and AUC 0.57 preformed by human expert and whole -slide level respectively.

Chaurasia et al. (2014) explained data mining techniques to predict cancer when they applied RepTree (C4.5), RBF Network and Simple Logistics on 286 samples and found accuracy as 71.32%, 73.77% and 74.47% respectively. Coudray et al. (2018) used a model deep convolutional neural network (inception v3) on whole-slide images obtained from The Cancer Genome Atlas (TCGA) and obtained an improved AUC score from 0.733 to 0.856 in the detection of cancer subtype.

Elgedawy et al. (2017) applied 3 machine learning techniques: Naïve Bayes, SVM and RF. Out of them RF is the most appropriate and useful algorithm to give the best accuracy as 99.42% where SVM and NB produced 98.8% and 98.24% accuracy respectively. Huang MW et al. (2017) used SVM ensemble classifier along with boosting method and RBF kernel based SVM to predict the accuracy in cancer dataset. In case of small-scale dataset, GA+RBF SVM ensembled with boosting method and produced 98.28% accuracy whereas in large dataset, RBF SVM ensembled with boosting method produced 99.52% accuracy.

Nguyen et al. (2013) shown experiments on two datasets WBCD (diagnosis) & WBCP (prognosis) from Wisconsin Breast Cancer Dataset and claimed for 99.8% and 99.7% accuracy.Pirooznia et al. (2007) compared many ML techniques and found 100% accuracy when they applied SVM-RFE on 84 sample which is a less size of data.

Sahu et al. (2012) proposed a novel approach using PSO along with SVM and k-NN ML techniques applied on the size of 87 sample and found 100% accuracy. Sivakami et al. (2015) proposed a hybrid technique DT-SVM to forecast cancer prediction for the dataset of 699 instances of WBCD repository and obtained accuracy as 91%.

Steiner et al. (2018) shown a impact of deep learning assistance on lymph of breast cancer and found micro metastases in the images with a range from 0.02, 0.002, 0.018 and 0.0005. Lower the value of micro metastases like 0.0005, higher the accuracy in lymph node of breast cancer. Xiao et al. (2018) discussed a multi-model ensemble method based on deep learning to find the accuracy and effective of different classifier. The data were supplied in the form of gene expression. This method was tested on three public RNA-seq data sets.

MACHINE LEARNING TECHNIQUES

Machine Learning (ML) is a part of AI. ML is used to infer the knowledge from the behavior of data. There are many areas where ML can be applied. In this paper we are discussing the cancer related issues. So, ML uses the techniques to generalize the biological sample of a given datasets. Following popularly used ML techniques have been introduced briefly.

Decision Tree

A Decision Tree (AL-SALIHY et al. 2017, Yue et al. 2018, Ponnuraja et al. 2017) is a binary classifier used to take the decision on attributes of the dataset. It looks like the tree but it is an inverted tree.

Random Forest

A Random Forest (Okun et al. 2007, Nahid et al. 2017, Ghongade et al. 2018) is another classifier, based on

decision tree, which is a next step when multiple decision tree resides together.

Support Vector Machine

Support Vector Machine (SVM) is a very effective classifier that classify the feature's outcome in two categories with a hyperplane having distance between the samples. (Huang MW et al. 2017, Sweilam et al. 2010, Sewak et al. 2007)

k-Nearest Neighbor (kNN)

kNN (Pawlovsky 2017, Rodriguez et al. 2018, Meneses et al. 2019, Al-Hadidi et al. 2016) is an essential classifier which makes a group of similar patterns with dataset sample. In general, the value of k is randomly put from 1 to 5 or extends up to 11 depends on how many neighbors are needed for grouping. There are various distance measurement mathematical equations such as Euclidean, Manhattan, Minkowski, Chebyshev, Cosine Similarity, Cosine Distance to find the distance between the neighbor.

Naïve Bayes Classifier

This classifier is based on Bayes' theorem (Maysanjaya et al. 2018, Rashmi et al. 2015, Soria et al. 2008) which takes into consideration of independent feature of the data. Dependent feature creates more correlation effect in prediction.

SIGNIFICANCE OF LEVEL OF ACCURACY FOR BREAST CANCER PREDICTION

There are many ML techniques for cancer diagnosis and prognosis. Many researchers have calculated the level of accuracy in percentage using different ML techniques such as Random Forest, SVM, Naïve Bayes, Decision Tree (Nguyen et al. 2013, Ahmad et al. 2013, Chaurasia et al. 2014, Sivakami et al. 2015, Elgedawy et al. 2017) for cancer prediction as shown below in table 1 for breast cancer datasets.

	Table 1. ML Techniques' Accuracy level for prediction of Breast Cancer								
Type of	Type of Cancer: Breast Cancer								
S. No.	ML Technique	ML Technique Sample Accuracy							
1	Random Forest	699	99.82%						
2	DT-SVM	699	91%						
3	Random Forest		99.24%						
	SVM	699	98.8%						
	Naïve Bayes		98.24%						
4	Decision Tree (C4.5)		93.6%						
	ANN	1189	94.7%						
	SVM		95.7%						
5	RepTree (C4.5)		71.32%						
	Radial Basis Function Network	286	73.77%						
	Simple Logistic		74.47%						

Table 2. Level of accuracy of ML Techniques for small data size							
S. No.	ML Technique	Types of Cancer	Sample	Accuracy			
1	SVM-RFE	Breast Cancer	84	100%			
2	PSO-KNN PSO-SVM	Breast Cancer	97	100%			

On the basis of above results, it is not easy to say that a particular Machine Learning technique is fit suitably for the diagnosis of breast cancer for a particular dataset on the basis of level of accuracy only because 100% accuracy comes in underfit condition where drawn conclusion of cancer prediction can't be correct.

Following table 2 (Pirooznia et al. 2007, Rajeshwari et al. 2011, Sahu et al. 2012, Gunavathi et al. 2014) shows 100% accuracy level using some other ensemble Machine Learning techniques on selected datasets where 100% accuracy is not predicting cancer correctly.

Table 3. Performance parameter metrics for dataset *BCWD11										
Algorithm	Confusion Matrix Components				Performance Parameters (*BCWD11)					
	True Positive (TP)	True Negative (TN)	False Positive (FP)	False Negative (FN)	Accuracy	Precision	Recall	F-Measure	MCC	ROC Area
Naïve Bayes	436	235	22	6	95.99%	96.2%	96%	96%	0.914	98.6%
SVM (SMO)	445	231	13	10	96.7%	96.7%	96.7%	96.7%	0.927	96.5%
KNN (IBK)	443	222	15	19	95.13%	95.1%	95.1%	95.1%	0.892	94.5%
Decision Tree (J48)	438	223	20	18	94.56%	94.6%	94.6%	94.6%	0.880	95.5%
Random Forest	444	230	14	11	96.42%	96.4%	96.4%	96.4%	0.921	99%

Table 4. Performance parameter metrics for dataset **WBCD32										
Algorithm	Con	Confusion Matrix Components				Performance Parameters (**WBCD32)				
	True Positive (TP)	True Negative (TN)	False Positive (FP)	False Negative (FN)	Accuracy	Precision	Recall	F-Measure	MCC	ROC Area
Naïve Bayes	190	337	22	20	92.61%	92.6%	92.6%	92.6%	0.842	97.6%
SVM (SMO)	201	356	11	1	97.89%	97.9%	97.9%	97.9%	0.955	97.3%
KNN (IBK)	200	347	12	10	96.13%	96.1%	96.1%	96.1%	0.917	95.6%
Decision Tree (J48)	194	335	18	22	92.97%	93%	93%	93%	0.85	92.3%
Random Fores	196	350	16	7	95.95%	96%	96%	95.9%	0.913	99.1%

BREAST CANCER PREDICTION METRICS

The basis of performance parameter metrics is confusion matrix and then metrics such as accuracy, precision, recall, f-measure, Mathew's Correlation Coefficient (MCC), is calculated (Kourou et al. 2015, Baker 2003, Yang et al. 2017, Tilaki 2013).

CONFUSION MATRIX

A Confusion matrix is a summary of prediction results on a classification problem. In it, the number of correct and incorrect predictions are summarized with count values and broken down by each class, the concept is shown in table 5.

On the basis of above table, various metrics parameter has been defined below. Accuracy (Barlow et al. 2004, Tharwat 2018) is the number of correct predictions divided by total number of predictions made. Mathematically accuracy (Acc) is given by the following formula

$$Acc = \frac{TP + TN}{TP + TN + FP + FN}$$

Sometimes accuracy yields the same value with respect to multiple classifiers and this makes more complication with Error Rate (ERR) or misclassification rate (Jensen et al. 2010, Barlow et al. 2004, Tharwat 2018). Error Rate can be calculated as below

$$ERR = 1 - Acc = 1 - \frac{TP + TN}{TP + TN + FP + FN}$$

Table 5. Confusion Matrix							
Actual							
		Valid	Not Valid				
Predicted	Accept (Recurrence)	True Positive (TP)	False Positive (FP)				
	Reject (No Recurrence)	False Negative (FN)	True Negative (TN)				

Sometimes because of accuracy paradox, accuracy is not sufficient to find the best model. Improving the accuracy by reducing the error is not appropriate. Therefore, ROC Curve and AUC score are better option to use for prediction instead of using the only parameter accuracy.

In this paper other additional parameter such as ROC and AUC metrics have been calculated to find the actual best suitable fit model.

ROC AND AREA UNDER CURVE (AUC)

ROC (Baker 2003, Yang et al. 2017, Tilaki 2013) curve demonstrates the tradeoff between the true positive fraction and false positive fraction to evaluate the positivity. AUC (Yang et al. 2017, Tilaki 2013) is a measure of the model's performance which is based on the ROC curve. This curve plots two parameters: True Positive Rate & False Positive Rate. Both the parameters are defined below.

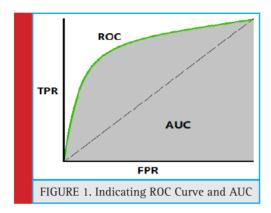
True Positive Rate (TPR), termed as sensitivity also, is the fraction of positives correctly classified divided by total positives and is defined as below

True Positive Rate,
$$TPR = \frac{TP}{TP+FN}$$

False Positive Rate (FPR) is the fraction of negative incorrectly classified divided by total negatives and is defined as below

False Positive Rate,
$$FPR = \frac{FP}{FP+TN}$$

AUC is an effective and combined measure of TPR and FPR that describes the inherent validity of diagnostic tests (Kouruo et al. 2015). In below figure 1, FPR and TPR are represented by x-axis and y-axis respectively. The ROC indicates the curve of value ranging from 0 to 1 whereas AUC shows the area under curve. Dashed-line partitions True values and False value. True values lie above the dashed line and False values lie below the line.



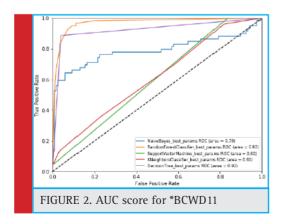
RESULTS AND DISCUSSION

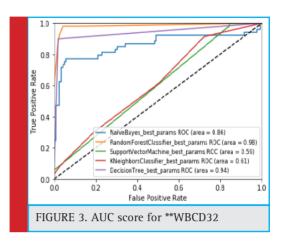
The experimental setup is designed under the environment of Windows 10 operating System, Python 3.x version and smart IDE Spyder which is a unit of Anaconda distribution. The dataset has been used from Wisconsin Breast Cancer Dataset having 10 major features in *BCWD11 and 30 major features in **WBCD32 (Kumar et al. 2019). The respective performance measurement metrics is calculated on the basis of confusion matrix given in table 3 & table 4 respectively.

On the basis of TPR and FPR, the ROC curve is plotted for all machine learning classifiers to obtain AUC score. The following graph is showing the AUC score in figure 3 below.

On the basis of fact that higher the value of AUC near to 1, the model is considered to be the best. In figure 2, it is observed that the best model for said dataset is Random Forest classifier, has the highest value of AUC i.e. 0.97.In similar fashion, another experiment has been performed for **WBCD32 and performance of Random Forest classifier is again observed as the best model for AUC score 0.98 shown in figure 3 below.

This paper discussed the confusion matrix and performance parameters useful for prediction. Parameters are accuracy, precision, F-measure, MCC and ROC area.





Based on all the parameters, the outcome of this paper was to find out the AUC score whereas in previous works, only accuracy was calculated.

We focused to locate the best fit model for selected breast cancer dataset. In general, it is observed that only accuracy is used for cancer prediction. But accuracy does not give perfect prediction. It is identified that AUC score is significant to consider for correct prediction of breast cancer instead of the only parameter 'accuracy'. Random Forest classifier found fit perfectly for the used dataset among all other four classifiers viz Naïve Bayes, SVM, kNN and Decision Tree. There are also few more considerable measurement metrics such as Youden's index (YI), Discriminant power (DP), Balanced classification rate (BCR), Optimization precision (OP), Jaccard (or Tanimoto similarity coefficient), and many more. In future, we intend to do feature engineering for better prediction of breast cancer.

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