

Article

A Machine Learning Approach for the Non-Destructive Estimation of Leaf Area in Medicinal Orchid *Dendrobium nobile* L.

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Abstract: In this study, leaf area prediction models of *Dendrobium nobile*, were developed through machine learning (ML) techniques including multiple linear regression (MLR), support vector regression (SVR), gradient boosting regression (GBR), and artificial neural networks (ANNs). The best model was tested using the coefficient of determination (R^2), mean absolute errors (MAEs), and root mean square errors (RMSEs) and statistically confirmed through average rank (AR). Leaf images were captured through a smartphone and ImageJ was used to calculate the length (L), width (W), and leaf area (LA). Three orders of L, W, and their combinations were taken for model building. Multicollinearity status was checked using Variance Inflation Factor (VIF) and Tolerance (T). A total of 80% of the dataset and the remaining 20% were used for training and validation, respectively. KFold (K = 10) cross-validation checked the model overfit. GBR (R^2 , MAE and RMSE values ranged at 0.96, (0.82–0.91) and (1.10–1.11) cm²) in the testing phase was the best among the ML models. AR statistically confirms the outperformance of GBR, securing first rank and a frequency of 80% among the top ten ML models. Thus, GBR is the best model imparting its future utilization to estimate leaf area in *D. nobile*.

Keywords: leaf area; smartphone; ImageJ; *Dendrobium nobile*; Gradient Boosting Regression (GBR); Average Rank (AR)



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1. Introduction

Dendrobium nobile (Orchidaceae) is an endangered orchid species listed under Appendix II of CITES [1] and forms the largest vascular epiphytes having features of CAM plants [2]. It serves as an ornamental and medicinal food for humans [3]. Recently HPLC fingerprints showed a strong inhibiting effect on cancer cells, through increased bioactive compounds; signifying *D. nobile* as a functional herb for the market [4]. Leaf area is a necessary biometrical variable that helps not only in the computation of various physiological indices such as leaf area index (LAI), specific leaf area (SLA), net assimilation rate (NAR), specific leaf weight (SLW), leaf area duration (LAD) and various plant physiological mechanisms viz., photosynthesis, respiration, light interception, transpiration, etc. [5–7], but also water-related anatomical traits such as leaf density and vein density, thus, helping in water conservation and transport ultimately resulting in the ability of *Dendrobium* to cope in water stress situations in the environment [8]. Leaf area, shape, size, and number present per plant affects the source–sink relationship and ultimately the yield of the plant as green leaves act as a primary source of assimilates for petals in *Dendrobium* orchid [9,10]. Leaf area is also responsible for nutrient spray responsiveness in *Dendrobium* [11]. The leaf area trait, i.e., specific leaf area (SLA), is strongly influenced by phylogeny (k value > 1) indicating strong conservation in changes in this trait over evolution in *Dendrobium* [12].

Leaf area further can act as an important component of complex process-based plant growth models for the development of support decision-based systems for management of cultural practices in *Dendrobium*. Therefore, leaf area (LA) is a fundamental component of estimation of physiological, ecological, evolutionary, and anatomical traits that affect the survival, growth, and distribution of *Dendrobium*. Leaf area modeling enables an easy way of predicting LA through models which aid in a non-destructive estimation, rather than manual methods.

LA can be measured directly through techniques such as photographing [13], image analysis, blueprinting, use of expensive instruments including digital planimeter [14], scanning planimeter [15], etc. Major disadvantage in these direct methods lies in the removal of leaves, which poses a problem of not only for the experiments which are time-shared, but also leads to a decline in these valuable plant samples, in our case a rare and endangered orchid species. Other problems in these methods can be mentioned as requirements of time and labor, costly instruments, reduction in canopy affecting photosynthesis, growth, etc. The above bottleneck in LA estimation is successfully overcome through the various non-invasive technologies using very few measurements such as leaf length (L) and leaf width (W) or its different combination [16–19], constant leaf area term KA [20] or a correction factor [21]. These indirect LA estimation methods are performed in situ [22], low in cost, reliable, provide fast results without detachment of the leaves from the plants, and also excludes biological variations in the experiment [23]. In our study, an *Orchidaceae* family member, *D. nobile*, was assessed for LA estimation where conservation prioritization is essential and automated conservation assessments were recently performed through deep learning [24]; thus making it inevitable to use the non-destructive ML methods for development of LA prediction models.

Machine learning (ML) techniques have gained a lot of importance due to their non-requirement of explicit instructions, i.e., a data driven nature, and their reliable, non-destructive, fast, accurate, high throughput, user friendly and less laborious approach of solving real life problems. Recently, ML and modeling has become very popular in the field of LA estimation, as it allows the adjustments and improvements of the model for accurate prediction of LA in crops. The regression problem depicts the relationship between one continuous dependent variable and multiple independent variables. LA modeling in the present study underwent regression problem analysis as a continuous dependent variable leaf area (LA) was predicted by a numerical value using two independent variables leaf length (L) and leaf width (W). Multiple linear regression (MLR), the most common form of linear regression analysis, computes a weighted sum of the input features and a bias term, the intercept. Support vector regression (SVR) is a supervised ML algorithm to predict the dependent variable in the model. In principle, SVR is same as support vector machine (SVM) [25] which is capable of computing linear and non-linear regression problems. The gradient boosting regression (GBR) [26] algorithm is a decision tree-based ensemble technique for solving both linear and non-linear regression problems. The high flexibility to construct new base-learners which correlates maximally with the negative gradient of the loss function makes it a highly customizable data driven ML technique. Boosting algorithms are also very easy for various model design implementations, thus imparting gradient boosting machine applications in various fields of both practical and data mining and machine learning studies [27–29]. It is faster and has better model performance due to its principle of hypothesis boosting; the requirement of minimum data pre-processing has outperformed Artificial Neural Network (ANN) and SVM in many reports [30–32]. ANN techniques have been used in a limited number of crops which are summarized in the paper [33]. These models are self-adaptive, data driven, and non-linear in nature which helps to find the relationship between the predictor and predicted [34].

Performance metrics have been used for evaluation in multiple machine learning algorithms, but a unified and single standard metric is difficult to find. Another drawback of these metrics is since regression problem may have multiple values ranging from zero to infinity, merely a single metric cannot solve the problem of regression with respect to

the ground truth elements in the study. Thus, in our study three performance evaluation metrics were used to assess the results of regression viz., coefficient of determination (R -squared or R^2 [35]) which determines the proportion of variance in the dependent variable which can be predicted from the independent variables; mean absolute error (MAE [36]) depicting the quality of fit in terms of the distance of the regressor to the actual training points; and Root mean Square Error (RMSE) for the detection of outliers [37]. We also used another statistical ranking method Average Rank (AR) that follows Friedman's M statistic [38] to select the best model for *Dendrobium* orchid.

LA modeling has been reported for many crops including apple [39], walnut [13], apricot [40], onion [41], jatropha [42], cacao [43], cherry cultivars [14], chestnut [44], grapes [45,46], green and black peppers [33,47,48], ginger [49], medicinal and aromatic plants [50], niagara and grave vines [51], som [52], mango [53], tomato [54], cotton [55–57], multiple crops [58], kiwi fruit [59], durian [60], pecan [61], forest tree [62], and hazelnut [63]. Ornamentals received meagre attention; among them LA modeling was performed in zinnia [64], sunflower [65], rose [66], Euphorbia $\times$ lomi Thai hybrids [67], bedding plants [68], and bougainvillea [69]. The development of leaf dry weight and leaf area models in the four cultivars of *Phalaenopsis* orchids [70] was reported by considering the length and width of leaves and linear regression analysis, but a major limitation is the non-utilization of the progressively utilized ML techniques.

Advanced ML techniques to predict leaf area are reported in a very limited number of crops and the existing LA models are mostly based on linear regression analysis. Until now, no other attempts have been reported for non-destructive estimation of LA through ML techniques in the *Orchidaceae* family, which constitutes the second largest (c. 28,000) flowering family in the world; it is the most widespread and accounts for 8% of angiosperm species diversity [71]. *Dendrobium*, in this study for LA modeling, constitutes the largest genus of *Orchidaceae*, over 1800 species [72]. To the best of our knowledge, only a linear regression-based model (LRM) and recently very few works on neural net (NN)-based models have been reported for LA prediction. Our study reports for the first time a non-linear regression-based model, i.e., support vector regression (SVR) and decision tree-based ensemble model gradient boosting regression (GBR) for LA modeling and its robustness was compared with the state-of-the-art ML models and a novel method of statistical ranking of the LA models based on average rank (AR) was proposed. Keeping this in mind, the objectives of this study were set as to (i) determine the individual leaf length (L), width (W) and the leaf area (LA) using the ImageJ software [73], to ultimately develop a leaf area model from the nine input combinations of L and W using different ML techniques viz., MLR, SVR, GBR, and the nine best selected ANN models (ii) to evaluate the model robustness through various statistical performance metrics (R^2 , MAE and RMSE) for selecting the best LA prediction model and (iii) to rank the ML models based on average rank (AR) ranking methodology in the orchid *D. nobile*.

2. Materials and Methods

2.1. Plant Material

The present study was conducted in the glass house located at ICAR-National Research Centre for Orchids (ICAR-NRCO), Pakyong, Sikkim, India (27.2267° N, 88.5877° E). The individual leaf images were collected from *D. nobile* orchid from 1.5 to 2 years of age. A mixture of stone or brick pieces, leaf mold, coconut husks, and semi rotten logs in the ratio of 1:1:1:1 in plastic pots of (5 to 6) inches was used as the potting media of *D. nobile*. The N:P:K composition as 20:10:10 @ 0.5% was sprayed on the 1 to 2 year old plants and potting materials at an interval of 15 days and other nutrients such as calcium nitrate @ 0.05%, iron sulphate @ 50 ppm, magnesium sulphate @ 0.1%, boric acid @ 50 ppm and zinc sulphate @ 50 ppm were sprayed at 60 days interval. Neem @ 3% and copper oxychloride @ 0.3% were sprayed once every fortnight for controlling sucking pests and foliar diseases, respectively. The number of average leaves in the plant was approximately 10 and image was captured by a smartphone (Samsung Galaxy J7 with 13 MP camera, resolution of 4128 $\times$ 3096 pixels,

autofocus, and LED flash) during 10:30 a.m. to 3:30 p.m. from July to August of 2020. The light intensity was recorded by lux meter (Lutron LX-101A, Delhi, India) every day at 12:00 p.m. regardless of the weather or cloud cover over the 4 weeks experimental period. The glass house ambient light intensity varied from 531 lux to 1413 lux. Direct bright sunlight was avoided by providing a green colored shade net of 50% over the glass house (Saveer Biotech Ltd., New Delhi, India). Ambient day and night temperature ranged to about 32 °C and 17 °C during the photoperiod and relative humidity varied from 60 to 80% inside the glass house.

2.2. Process Flow of Selection of Best Leaf Area Model in *D. nobile* through ML Techniques

Figure 1 depicts the whole workflow of the study conducted for the non-destructive estimation of leaf area using ML. Firstly, the individual leaves of the *D. nobile* were captured with a reference scale by a smartphone (Figure 2). Then the ImageJ (<https://imagej.nih.gov/ij/> (accessed on 27 April 2020)) software was used for determining the individual leaf length (L), width (W), and the leaf area (LA) of *D. nobile*. Three orders of independent variable viz., first order, second order, and third order polynomial functions of L and W were taken which can be denoted as L, L^2 , and L^3 and W, W^2 , and W^3 , respectively. L, L^2 , and L^3 were combined with W, W^2 , and W^3 , respectively, thus totalling to the number of input combinations as 3×3 , i.e., 9. These nine combinations of inputs were used for MLR, SVR, and GBR machine learning techniques for model building. For ANN, we used four types of architectures and considered the nine combinations of L and W, the total number of models were $4 \times 9 = 36$. Finally, the nine best ANN models were selected based on R^2 , MAE, and RMSE values. Thus, the above nine models from each of MLR, SVR, GBR, and ANN were compared together with their respective performance metrics. For the study, the original dataset was split into 80% training and 20% testing dataset for model development and its validation. To avoid the over fitting of the ML model KFold ($k = 10$) cross-validation [74] was performed using cross_val_score on the training dataset. The rest of the 20% holdout data were used for testing of the model. R^2 , MAE, and RMSE were used for ranking of the models. Selection of the best leaf area model was performed through higher R^2 values and lower MAE and RMSE values. Lastly, AR was used to statistically confirm the best performing models through ranks.

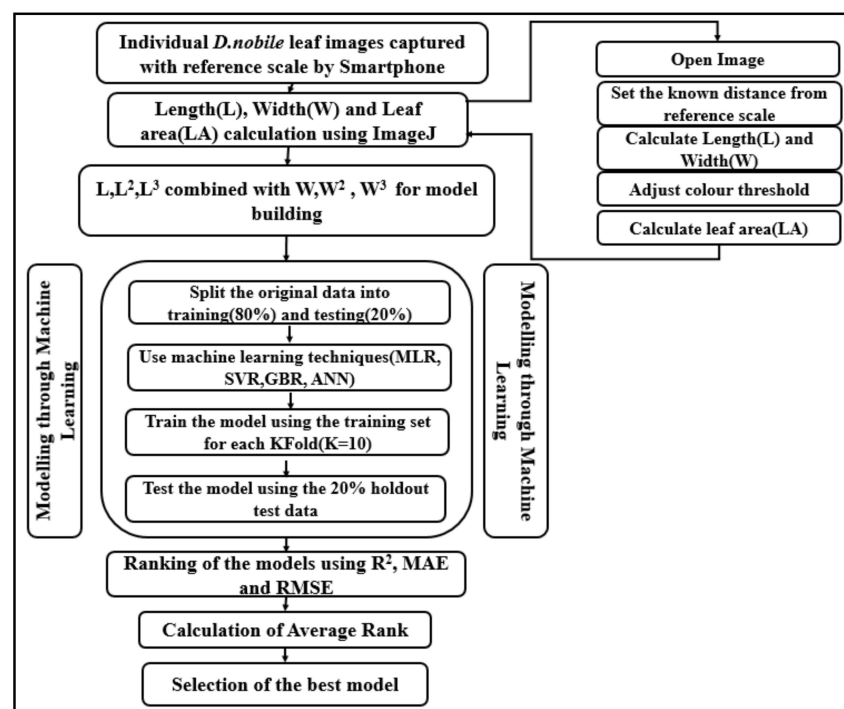


Figure 1. Flow diagram depicting the workflow of selection of best leaf area model through ML techniques in *D. nobile*.

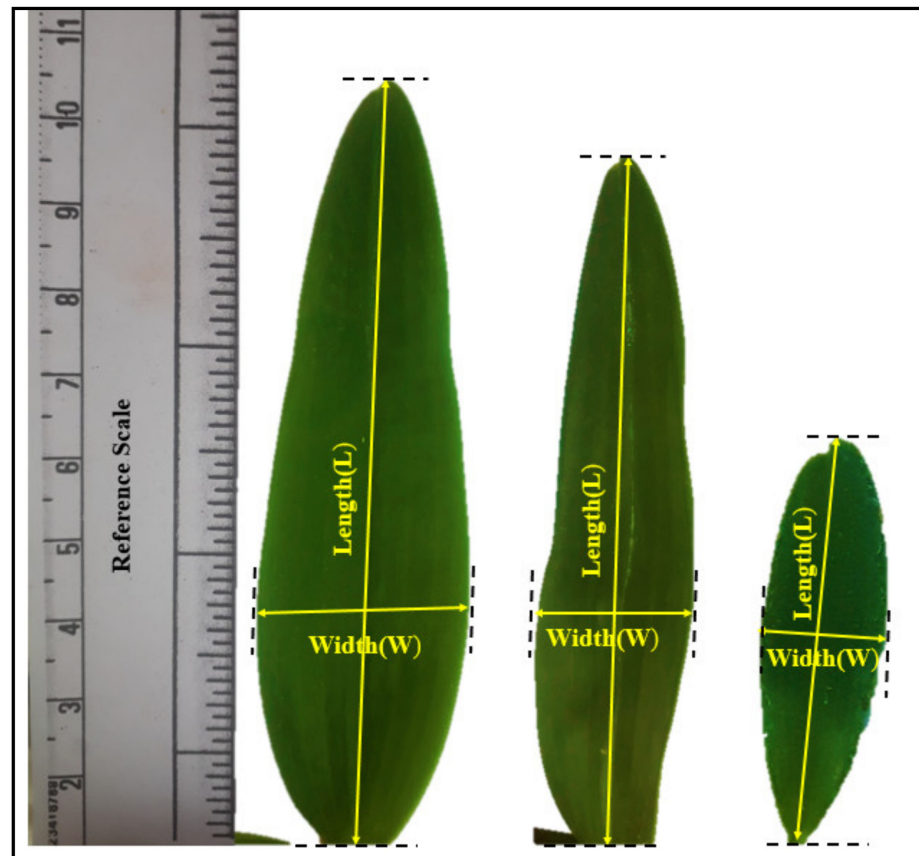


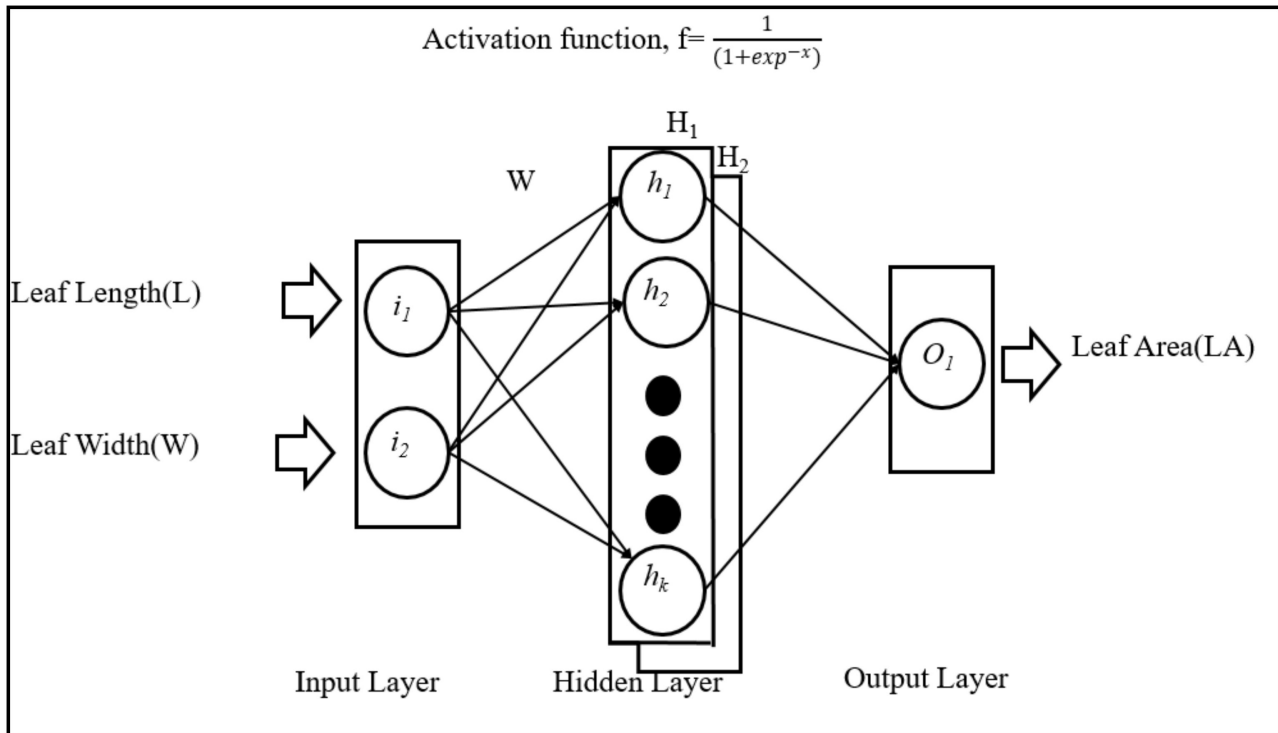
Figure 2. Images of individual leaf of *D. nobile* captured through smartphone with a reference scale for measurement of leaf length (L) and leaf width (W).

2.3. Dataset

A total of 1589 *D. nobile* leaf images were collected by Samsung Galaxy J7 smartphone for the study. The image was captured with a reference scale and the ImageJ software was used for estimation of the length, width, and area of the individual leaf [75]. The color threshold function of ImageJ was used to demarcate the captured leaf area from its background, thus automatically calculating the region of interest, i.e., measured leaf area (LA). The dataset contains one continuous dependent variable, i.e., leaf area (LA) and two independent variables—leaf length (L) and leaf width (W). The measurement of the leaf length (L) was taken from the tip of the lamina to the intersection point between the leaf and petiole. The leaf width (W) measurement was taken from end-to-end point between the broadest part of the lamina exactly perpendicular to the midrib of *D. nobile* lamina. All the variables were measured in cm scale. The values of three variables, length (L), width (W), and leaf area (LA) were stored in csv format for further model development processing and submitted to the dataset repository Mendeley data [<http://doi.org/10.17632/8tk2sc4ytg.1> (published on 28 June 2021)] and KRISHI, ICAR research data repository for knowledge management [<http://krishi.icar.gov.in/jspui/handle/123456789/71908> (published on 6 May 2022)] (Supplementary S1). Table 1 depicts the different combination of inputs of L and W along with their ML models. In this experiment three orders of L and W were taken, so L , L^2 , and L^3 were combined with W , W^2 , and W^3 . Each L and its higher order were combined with W and its higher order. The total number of input combination thus can be calculated as 3×3 , i.e., 9. These nine combinations of inputs were used for MLR, SVR, and GBR techniques. The higher orders of L and W were taken to observe their effect on model building. For ANN, four types of architectures were designed and the total number of models were $4 \times 9 = 36$. Finally, the nine best ANN models were selected based on R^2 , MAE, and RMSE values (Figure 3 and Table 2).

Table 1. Input combinations of L, W, and their variants with model names of MLR, SVR, and GBR.

Inputs	W	W ²	W ³
L	L, W (MLR1, SVR1, GBR1)	L, W ² (MLR2, SVR2, GBR2)	L, W ³ (MLR3, SVR3, GBR3)
L ²	L ² , W (MLR4, SVR4, GBR4)	L ² , W ² (MLR5, SVR5, GBR5)	L ² , W ³ (MLR6, SVR6, GBR6)
L ³	L ³ , W (MLR7, SVR7, GBR7)	L ³ , W ² (MLR8, SVR8, GBR8)	L ³ , W ³ (MLR9, SVR9, GBR9)

**Figure 3.** ANN network structures used in this study includes three types of layers. Input layer (with input neuron i_1 and i_2) taking values from leaf length (L) and leaf width (W), hidden layer (H_1, H_2) (with hidden neuron h_k , where $k = 3$ or 5 when architecture has single hidden layer and $k = 3$ when architecture has double hidden layer) and output layer (with output neuron O_1) were utilized to predict the leaf area (LA). Furthermore, $4 \times 9 = 36$ models of ANN were developed. W and f depict the weights and sigmoid activation function, respectively.**Table 2.** Combination of different inputs for ANN techniques with different ANN architectures.

Models	Input Variables	Output Variable	Layers (Input-Hidden-Output)
ANN1	L, W	LA	2-3-1
ANN2	L, W	LA	2-5-1
ANN3	L, W	LA	2-10-1
ANN4	L, W	LA	2-3-3-1
ANN5	L, W ²	LA	2-3-1
ANN6	L, W ²	LA	2-5-1
ANN7	L, W ²	LA	2-10-1
ANN8	L, W ²	LA	2-3-3-1
ANN9	L, W ³	LA	2-3-1
ANN10	L, W ³	LA	2-5-1
ANN11	L, W ³	LA	2-10-1
ANN12	L, W ³	LA	2-3-3-1
ANN13	L ² , W	LA	2-3-1
ANN14	L ² , W	LA	2-5-1

Table 2. Cont.

Models	Input Variables	Output Variable	Layers (Input-Hidden-Output)
ANN15	L^2, W	LA	2-10-1
ANN16	L^2, W	LA	2-3-3-1
ANN17	L^2, W^2	LA	2-3-1
ANN18	L^2, W^2	LA	2-5-1
ANN19	L^2, W^2	LA	2-10-1
ANN20	L^2, W^2	LA	2-3-3-1
ANN21	L^2, W^3	LA	2-3-1
ANN22	L^2, W^3	LA	2-5-1
ANN23	L^2, W^3	LA	2-10-1
ANN24	L^2, W^3	LA	2-3-3-1
ANN25	L^3, W	LA	2-3-1
ANN26	L^3, W	LA	2-5-1
ANN27	L^3, W	LA	2-10-1
ANN28	L^3, W	LA	2-3-3-1
ANN29	L^3, W^2	LA	2-3-1
ANN30	L^3, W^2	LA	2-5-1
ANN31	L^3, W^2	LA	2-10-1
ANN32	L^3, W^2	LA	2-3-3-1
ANN33	L^3, W^3	LA	2-3-1
ANN34	L^3, W^3	LA	2-5-1
ANN35	L^3, W^3	LA	2-10-1
ANN36	L^3, W^3	LA	2-3-3-1

2.4. ML Methodologies Used for LA Prediction Modeling

2.4.1. Multiple Linear Regression Analysis (MLR) Models

Multiple linear regression (MLR) is a variant of simple linear regression with more than one explanatory or independent variable. It can capture linear relationships between independent variables and the dependent variable. It performs better in the absence of multicollinearity within the dependent variables and dependent and independent should be linearly correlated with each other [76]. The equation for MLR model estimation is as follows [77]:

$$y = \beta_0 + \beta_1 x_1 + \dots + \beta_2 x_2 + \dots + \beta_k x_k + \epsilon \quad (1)$$

where y is the dependent variable (LA), x_i is the i th independent variable (L or W), β_i is the polynomial coefficients of x_i , k is the number of independent variables, and ϵ is the possible variation form.

2.4.2. Support Vector Regression (SVR) Models

SVR is based on the Vapnik–Chervonenkis (VC) theory which in principle is structural risk minimization [78] and is an effective tool for estimating real-valued functions [79]. The advantages of SVR, which was harnessed in this study, is its highly effective performance for both the linear and non-linear regression dataset and not too large dataset which results in the use of Gaussian RBF kernel hyperparameter (for non-linear regression) thus mapping each training instance to an infinite-dimensional space [80]. The other usefulness of SVR can be summarized as its excellent capability of generalization with high accuracy prediction. Default hyperparameters used in `sklearn.svm.SVR` are kernel = 'rbf', degree = 3, gamma = 'scale', coef0 = 0.0, tol = 0.001, C = 1.0, epsilon = 0.1, shrinking = True, cache_size = 200, verbose = False, and max_iter = -1.

2.4.3. Gradient Boosting Regression (GBR) Models

GBR develops an additive model that serially adds predictors or new base learners to an ensemble and each one corrects its predecessors leading to the minimization of the loss function. In each iterative training process, the residual value in the current model

denotes the negative gradient of the loss function, and a new regression tree (predictor) is trained to fit the current residual and added subsequently to the previous model. For developing a GBM model, both the loss function and the function corresponding to the negative gradient must be specified. In our study, loss function for continuous response viz., least-square was used which usually neglects smaller deviations but refits larger ones. One of the advantages of GB models is that it provides the flexibility to use a variety of base learning models at the same time for designing a complex model for a particular solution. A model's regularization capability is of the utmost importance for model building from data because it prevents the overfit of the data [81]. Our study used shrinkage as a regularization technique which reduces unstable regression coefficients by shrinking it to zero which is normally used in ridge regression problems. The hyperparameters used in our study were *n_estimators* to control the ensemble training through the number of boosted trees to fit the training data; *min_samples* and *max_depth* to control the growth of decision trees; and learning rate to scale the contribution of each tree via a regularization technique known as shrinkage, the lower the value the better the generalization of the predictions will be. The value of the learning rate and *n_estimator* were adjusted in our experiment to 0.01 and 500, respectively, which performed better than the other combinations tried. The loss hyperparameter controls the cost function of GBR. Table 3 depicts the hyperparameters, their values and descriptions which were adjusted to improve the learning performances in GBR and ANN.

Table 3. Descriptions of hyperparameters tuned in GBR and ANN models.

GBR			ANN		
Hyper Parameters	Values	Descriptions	Hyper Parameters	Values	Descriptions
<i>n_estimators</i>	500	No of decision tree in the ensemble	<i>alpha</i>	0.001	Regularization parameter
<i>max_depth</i>	4	Maximum depth of the decision tree	<i>hidden_layer_sizes</i>	(3,); (5,); (10,); (3,3)	Four architecture 2-3-1, 2-5-1, 2-10-1, 2-3-3-1
<i>min_samples_split</i>	5	Minimum number if sample required to split an internal node	<i>max_iter</i>	1000	No. of iteration
<i>learning_rate</i>	0.01	Determine the impact of each tree on final outcome.	<i>activation</i>	'logistic'	Logistic sigmoid function $\frac{1}{(1+exp^{-x})}$
<i>loss</i>	'ls'	Least square loss function	<i>learning_rate</i>	'adaptive'	Keep learning rate constant as the initial learning rate

2.4.4. Artificial Neural Networks (ANNs) Models

ANNs are data-driven, non-linear, and self-adaptive techniques capable of establishing a relationship between the input data and target output data even when they have an unknown relationship [82]. The mathematical structures of ANN models have a resemblance with the neural system of the human brain [78]. They process data collectively through interconnected neurons depending on factors including thresholds, adjustable weights, and mathematical transformation functions of the images [83]. The ANN architecture used in our study comprised of three layers, where the input and output layers depict the independent and dependent variables, and the hidden layer (inter layer) deals with the computation of data and establishes interconnection between the input and output layers through the hidden neurons. The optimum neurons in the hidden layer were determined through the trial and error approach [84]. In our study, a single hidden layer was found

to provide better accuracy than the double hidden layers which is in accordance with the literature [61]. Figure 3 and Table 2 depict the different ANN models where nine types of inputs (as described in Table 1) were combined with four types of ANN architectures, i.e., input layer-hidden layer-output layer (2-3-1, 2-5-1, 2-10-1, 2-3-3-1) in total calculating to $4 \times 9 = 36$ models of ANN. This was performed to observe the effects of different hidden layers on ANN models leading to the selection of the nine best models based on the coefficient of determination (R^2), mean absolute errors (MAEs), and root mean square errors (RMSEs) values of the testing phase.

The learning rate was adaptive in nature, i.e., constant as the initial learning rate (0.001) as long as the training loss decreased and number of iterations were determined as 1000 epochs for convergence. To capture the non-linearity of the data an activation function, sigmoid, was used. The hyperparameters, their values and descriptions used for ANN model building are described in Table 3.

2.5. Feature Correlation Heatmap and Multicollinearity of the Independent Variables

A feature correlation heatmap of L, W, and its variants (L^2 , L^3 , W^2 , and W^3) and LA was generated through the matplotlib package of Python library. Pairwise Pearson's correlation coefficient, r [85] in the seaborn heatmap functions from seaborn module of Python was applied among the different combination of variables. The Variance Inflation Factor (VIF) and Tolerance values (T) were calculated to check the multicollinearity status of the two independent variables (L and W) and their variants (L^2 , L^3 , W^2 , and W^3). VIF values >10 or T value <0.10 denote the effect of multicollinearity on the estimation of the parameters of the model that lead to its non-reliability and for that problem at least one of the independent variables must be excluded from the prediction model [86].

The equations of Pairwise Pearson's correlation coefficient (r) and Variance Inflation Factor (VIF) [87] and Tolerance (T) [88] are shown below:

$$r = \frac{\sum_i^n (x_i - \bar{x})(y_i - \bar{y})}{\sum_i^n (x_i - \bar{x})^2 (y_i - \bar{y})^2} \quad (2)$$

where n = number of data points; i varies from 1 to n ; x and y are two independent variables; $\bar{x}$ and $\bar{y}$ are means of x and y , respectively.

$$\text{VIF} = \frac{1}{(1 - r^2)} \quad (3)$$

r is the correlation coefficient.

$$T = \frac{1}{\text{VIF}} \quad (4)$$

2.6. Programming Set-Up

The experiment was conducted on an HP Omen laptop computer with i7 9th generation processor (8 GB RAM, 4 GB GEFORCE GTX nVIDIA GPU, 1 TB HDD, 250 GB SSD Windows 10 operating system). The experiments were conducted in Pycharm Community Edition 2020.2 IDE using Python 3.7 version. For our study, we used many python packages including numpy, pandas, sklearn, matplotlib, etc.

Linear regression of sklearn.linear_model, SVR of sklearn.svm, gradient boosting regressor of ensemble, and MLP regressor of sklearn.neural_network modules available in sklearn were used for implementation of MLR, SVR, GBR, and ANN ML techniques respectively. Training and testing was performed using train_test_split methods of sklearn.model_selection, mean_squared_error, mean_absolute_error, and r2_score of sklearn.metrics were used for calculation of the performance of the machine learning algorithm [89].

The equations are defined as under [77]:

$$R^2 = \frac{\sum_{i=1}^n (LA_{\text{mea}} - \overline{LA}_{\text{mea}})(LA_{\text{mea}} - \overline{LA}_{\text{est}})}{\sum_{i=1}^n (LA_{\text{mea}} - \overline{LA}_{\text{mea}})^2 \sum_{i=1}^n (LA_{\text{mea}} - \overline{LA}_{\text{est}})^2} \quad (5)$$

$$MAE = \frac{\sum_{i=1}^n |LA_{\text{mea}} - \overline{LA}_{\text{est}}|}{n} \quad (6)$$

$$RMSE = \sqrt{\frac{\sum_{i=1}^n (LA_{\text{mea}} - \overline{LA}_{\text{est}})^2}{n}} \quad (7)$$

where LA_{mea} = measured LA value; LA_{est} = estimated LA value; $(\overline{LA}_{\text{mea}})$ = average values of measured LA value; $(\overline{LA}_{\text{est}})$ = average values of estimated LA value; n = number of leaves used for training or testing phase.

2.7. Model Performance Evaluation

A model evaluation is an essential step in machine learning problems. Several methods are available in literature to evaluate machine learning methods. The statistical model performance criteria used in this study were: coefficient of determination (R^2), mean absolute errors (MAEs), and root mean square errors (RMSEs).

In general, R^2 , MAE, or RMSE cannot predict the departure patterns of observed values in the study against the predicted values. In the scatter plot an overlap of the reference line ($y = x$) and regression line of observed vs. predicted values were visualized. Departures of the regression line of observed vs. predicted values from the reference line suggests an over or underestimation [90].

2.8. Model Ranking Based on Average Rank (AR) Ranking Methodology

For each 36 ML models which includes the four ML techniques viz., MLR, SVR, GBR, and ANN, based on the performance metrics R^2 , MAE, and RMSE, ranks were assigned. The best algorithm was ranked as first (1st) and second best as second (2nd) and so on, both for the testing as well as training phase, so each model obtains six ranks, three from training (R^2 , MAE and RMSE) and three from testing (R^2 , MAE and RMSE). The average ranks were obtained by averaging the ranks for each model. The final rank was derived by arranging the average ranks in an ascending order.

The formula for the calculation of average rank (AR) is as follows [38]:

$$AR_{\text{model}_j} = \frac{\sum_{i=1}^6 r_i}{6} \quad (8)$$

where j varies from 1 to 36 for 36 ML models, r_i represents the i th rank of a model, i varies from 1 to 6

3. Results and Discussion

The LA prediction modeling study presented here is of importance to the physiologists, horticulturalist, and environmentalist as *D. nobile* is a rare and endangered epiphytic medicinal orchid listed officially in the Pharmacopoeia of the People's Republic of China (Chinese Pharmacopoeia Commission, 2015) for the Dendrobine content as an active ingredient category [91]. The low reproduction rate in the wild, slow growth, and poor regeneration ability may prove *D. nobile* to be vulnerable under the predicted scenario of climate change as suggested by MaxEnt ML models that their habitat will shrink from (1–10)% in the future [92]. Leaf area (LA) estimation is of the utmost importance for performing any physiological experiments investigating the growth characteristics for studying the impact of predicted climate change.

First, the descriptive statistics of *D. nobile* leaves was calculated, next feature correlation heatmap and multicollinearity status of the variables was assessed. A selection of nine best

performing ANN models based on R^2 , MAE, and RMSE was tested. Finally, a comparison of state-of-the-art ML models for estimating leaf area of *D. nobile* leaves based on R^2 , MAE, and RMSE was addressed and ranking of the models was performed statistically based on Average Rank (AR) ranking methodology.

3.1. Descriptive Statistics of *D. nobile* Leaves Used for LA Model Building

The descriptive statistics was studied and summarized in Table 4 where maximum length, minimum length, width, and leaf area as well as mean and standard deviation of those parameters of *D. nobile* leaves were described. Sampling of leaves for image capture was performed to cover a wide range of leaf sizes. Leaf length (L) ranged from 5.24 cm to 16.30 cm with the mean of 10.68 and standard deviation of 1.57. Width (W) ranged from 1.30 cm to 4.34 cm with the mean of 2.39 and standard deviation of 0.51. Leaf area (LA) ranged from 6.39 cm² to 43.902 cm² with the mean of 20.042 cm and standard deviation of 6.26.

Table 4. Descriptive statistics of *D. nobile* leaves.

Parameters	Maximum	Minimum	Mean and Standard Deviation
Leaf Length (L) cm	16.30	5.24	10.68 and 1.57
Leaf Width (W) cm	4.34	1.30	2.39 and 0.51
Leaf Area (LA) cm ²	43.90	6.39	20.04 and 6.26

3.2. Feature Correlation Heatmap between the Variables through Correlation Coefficient

The correlation heatmap illustrating the relationship between each variable and compared with every other variable is depicted in Figure 4. The following observations can be made from the heatmap generated for the two independent variables (L, L², L³, W, W², W³) and LA in our study:

- Leaf length (L) and width (W) are linearly least correlated ($r = 0.43$ – 0.49) to each other as depicted from the light green color;
- L, L², L³, and LA are nearly equally correlated ($r = 0.76$ – 0.78) to each other as depicted from the medium green color;
- Leaf width (W) shows more correlated to leaf area (LA) with correlation coefficient (r) values ranging between 0.87 and 0.9 than leaf length (L) 0.76–0.78 as depicted from the darker green color in the heatmap;
- Among the variants of leaf width, the best to worst correlation with the leaf area (LA) can be shown in the order of W ($r = 0.9$) = W² ($r = 0.9$) > W³ ($r = 0.87$);
- L, W, and their variants are strongly linearly correlated with leaf area (LA) where r ranges from 0.76 to 0.90 as depicted from dark green color in the heatmap.

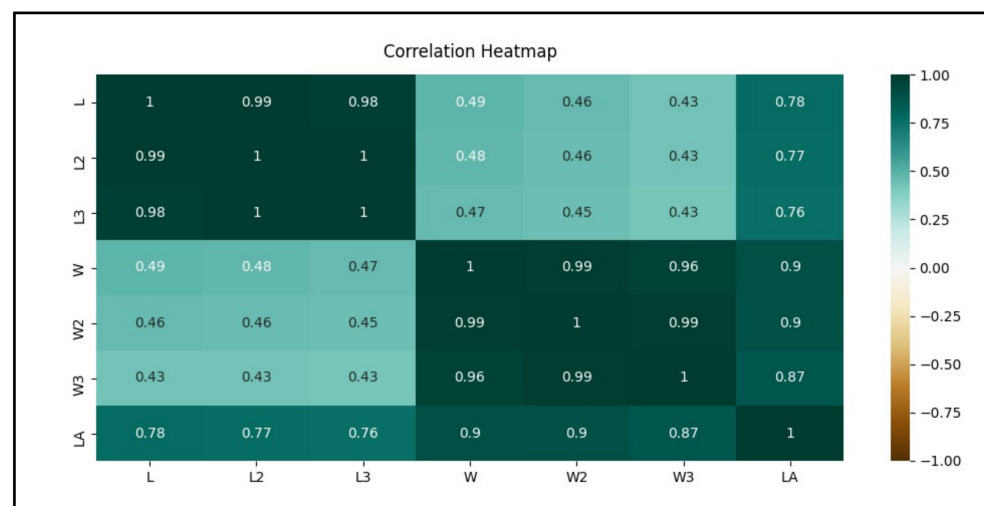


Figure 4. Correlation heatmap illustrating the relationship between each variable and compared with

every other variable. Dark green ($r = 1$) color indicates a positive correlation between the variables while white ($r = 0$) indicates no correlation between the variables and dark brown ($r = -1$) indicates a negative correlation. Pairwise Pearson's coefficient was used for correlation coefficient determination between the variables and seaborn module with matplotlib of Python was used for generation of the heatmap.

3.3. Multicollinearity Status of the Two Independent Variables (L and W)

In our study, the VIF ranged from (1.22 to 1.31) and T ranged from (0.76 to 0.82) for independent variables L, W, and their variants (Table 5) which are less than the non-acceptable VIF values, i.e., >10 and higher than T value < 0.10 to state the effect of multicollinearity on the estimation of the parameters of the model. Thus, the independent variables and the variants can provide accurate and reliable model building results overcoming the effects of multicollinearity.

Table 5. The values of VIF and T of L, W, and their input combinations.

Methods	Input Combinations								
	L, W	L, W ²	L, W ³	L ² , W	L ² , W ²	L ² , W ³	L ³ , W	L ³ , W ²	L ³ , W ³
Variance Inflation Factor (VIF)	1.31	1.27	1.23	1.30	1.27	1.23	1.28	1.25	1.22
Tolerance (T)	0.76	0.78	0.81	0.77	0.79	0.81	0.78	0.80	0.82

3.4. Selection of Nine Best Performing ANN Models Based on R², MAE and RMSE

From 36 models of ANN, nine models were selected based on nine different input combinations as mentioned in Table 2. One best model was selected from every input combination. ANN1–ANN4 came under the LW combination of inputs. ANN2 showed the highest R² value (0.96) and lowest MAE (0.86) and RMSE (1.13 cm²) in the testing phase. ANN5–ANN8 came under the LW² combination of inputs. ANN6 showed highest R² value (0.96) and lowest MAE (0.84) and RMSE (1.18 cm²) in the testing phase. ANN9–ANN12 came under LW³ combination of inputs. ANN10 showed highest R² value (0.97) and lowest MAE (0.87) and RMSE (1.18 cm²) in the testing phase. ANN13–ANN16 came under L²W combination of inputs. ANN15 showed highest R² value (0.97) and lowest MAE (0.86) and RMSE (1.11 cm²) in the testing phase. ANN17–ANN20 came under the L²W² combination of inputs. ANN18 shows the highest R² value (0.96) and lowest MAE (0.77) and RMSE (1.11 cm²) in the testing phase. ANN21–ANN24 came under the L²W³ combination of inputs. ANN22 showed the highest R² value (0.97) and lowest MAE (0.86) and RMSE (1.19 cm²) in the testing phase. ANN25–ANN28 came under the L³W combination of inputs. ANN26 showed the highest R² value (0.88) and lowest MAE (1.35) and RMSE (2.06 cm²) in the testing phase. ANN29–ANN32 came under the L³W² combination of inputs. ANN31 showed the highest R² value (0.94) and lowest MAE (1.01) and RMSE (1.47 cm²) in the testing phase. ANN33–ANN36 came under the L³W² combination of inputs. ANN35 showed the highest R² value (0.96) and lowest MAE (0.97) and RMSE (1.31 cm²) in the testing phase. Finally, the selected nine best performing ANN models which were to be compared with the other three models viz., MLR, SVR, and GBR includes ANN2, ANN6, ANN10, ANN15, ANN18, ANN22, ANN26, ANN31, and ANN35. Table 6 enumerates the performance metrics (R², MAE, and RMSE) of 36 ANN models.

Table 6. Comparative analysis of 36 ANN models with different inputs, i.e., LW, LW², LW³, L²W, L²W², L²W³, L³W, L³W², L³W³ for selection of 9 best models for each input combinations based on the statistical performance metrics (R², MAE and RMSE) of models.

Inputs	Model	Training			Testing		
		R ²	MAE	RMSE (cm ²)	R ²	MAE	RMSE (cm ²)
LW	ANN1	0.96	0.93	1.29	0.96	0.89	1.21
LW	ANN2	0.96	0.87	1.27	0.96	0.86	1.13
LW	ANN3	0.96	0.86	1.26	0.96	0.93	1.32
LW	ANN4	0.85	1.03	2.51	0.96	0.98	1.32
LW ²	ANN5	0.96	0.96	1.30	0.96	0.88	1.23
LW²	ANN6	0.96	0.90	1.25	0.96	0.84	1.18
LW ²	ANN7	0.96	0.91	1.28	0.96	0.84	1.21
LW ²	ANN8	0.93	1.13	1.50	0.94	1.18	1.62
LW ³	ANN9	0.96	0.92	1.28	0.95	0.91	1.35
LW³	ANN10	0.96	0.90	1.27	0.97	0.87	1.18
LW ³	ANN11	0.96	0.90	1.23	0.96	0.91	1.29
LW ³	ANN12	0.95	1.12	1.63	0.96	0.88	1.20
L ² W	ANN13	0.94	1.00	1.55	0.94	1.12	1.60
L ² W	ANN14	0.95	0.96	1.48	0.96	0.92	1.40
L²W	ANN15	0.95	0.94	1.41	0.97	0.86	1.11
L ² W	ANN16	0.92	2.28	5.17	0.95	1.07	1.45
L ² W ²	ANN17	0.95	1.00	1.44	0.96	0.91	1.34
L²W²	ANN18	0.94	0.93	1.36	0.97	0.77	1.11
L ² W ²	ANN19	0.95	0.94	1.39	0.96	0.89	1.24
L ² W ²	ANN20	0.74	1.60	4.05	0.00	4.62	5.84
L ² W ³	ANN21	0.94	1.03	1.47	0.95	0.88	1.32
L²W³	ANN22	0.95	0.97	1.39	0.97	0.86	1.19
L ² W ³	ANN23	0.95	0.90	1.32	0.96	0.92	1.34
L ² W ³	ANN24	0.83	1.67	3.19	0.95	0.93	1.26
L ³ W	ANN25	0.47	2.73	3.74	0.88	1.42	2.06
L³W	ANN26	0.73	2.77	3.78	0.88	1.35	2.06
L ³ W	ANN27	0.70	2.46	4.32	0.88	1.38	2.17
L ³ W	ANN28	0.39	3.19	4.41	0.83	1.61	2.68
L ³ W ²	ANN29	0.86	1.63	3.46	0.93	1.07	1.61
L ³ W ²	ANN30	0.81	1.36	2.15	0.89	1.13	1.97
L³W²	ANN31	0.88	1.41	2.22	0.94	1.01	1.47
L ³ W ²	ANN32	0.75	1.90	3.02	0.88	1.54	2.06
L ³ W ³	ANN33	0.87	2.04	2.32	0.85	1.58	2.45
L ³ W ³	ANN34	0.77	1.57	2.54	0.86	1.41	2.38
L³W³	ANN35	0.88	1.29	2.34	0.96	0.97	1.31
L ³ W ³	ANN36	0.85	1.92	3.58	0.89	1.38	1.99

3.5. Comparisons of Different ML Models for Estimating Leaf Area of *D. nobile* Leaves

A sum of 63 different ML models was developed for estimating the LA of *D. nobile* leaves. Input combinations are described in Table 1, which forms the various ML models: MLR (1–9), SVR (1–9), and GBR (1–9), each has nine models using the various input combinations. ANN also used these input combinations along with different network architectures which are described in Table 2. Among the 36 ANN models (Table 6) developed for each input combination, the best nine ANN models were taken based on higher R² and lower MAE and RMSE values. Therefore, for LW, LW², LW³, L²W, L²W², L²W³, L³W, L³W², and L³W³ input combinations ANN2 (2-5-1), ANN6 (2-5-1), ANN10 (2-5-1), ANN15 (2-10-1), ANN18 (2-5-1), ANN22 (2-5-1), ANN26 (2-5-1), ANN31 (2-10-1), and ANN35 (2-10-1) models, respectively, were compared with other models, i.e., MLR, SVR, and GBR. Table 7 depicts the LA models developed using different input combinations of L and W of *D. nobile* leaves using MLR (MLR1-MLR9) machine learning techniques.

Table 7. Developed LA models for three orders of L and W input combinations of *D. nobile* leaves using MLR techniques.

Model Input Combinations	Models
MLR1	$LA = -18.72 + L \times 1.74 + W \times 8.41$
MLR2	$LA = -9.01 + L \times 1.79 + W^2 \times 1.65$
MLR3	$LA = -6.76 + L \times 1.92 + W^3 \times 0.40$
MLR4	$LA = -9.56 + L^2 \times 0.08 + W^2 \times 8.34$
MLR5	$LA = 0.28 + L^2 \times 0.75 + W^2 \times 0.98$
MLR6	$LA = 3.37 + L^2 \times 0.089 + W^3 \times 0.40$
MLR7	$LA = -6.6 + L^3 \times 0.00049 + W \times 8.45$
MLR8	$LA = 3.5 + L \times 1.79 + W^2 \times 1.65$
MLR9	$LA = 6.8 + L^3 \times 0.0054 + W^3 \times 0.40$

A comparison of the performance statistics of MLR, SVR, GBR, and ANN models used for LA estimation was performed based on testing and training results using R^2 , MAE, and RMSE values which are enumerated in Table 8. The R^2 value in the training phase varied from (0.95–0.97), 0.94, 0.96, and (0.73–0.96) in case of MLR, SVR, GBR, and ANN, respectively. Similarly, the MAE values in the training phase varied from (0.84–1.11), (0.87–0.93), 0.86 and (0.87–2.77), for MLR, SVR, GBR, and ANN, respectively. RMSE values in the training phase ranged from (1.15–1.42) cm^2 , (1.47–1.58) cm^2 , 1.18 cm^2 and (1.25–3.78) cm^2 , for MLR, SVR, GBR, and ANN, respectively. In the testing phase, the R^2 values ranged from (0.94–0.96), 0.96, 0.96, and (0.88–0.97) for MLR, SVR, GBR, and ANN, respectively. Similarly, the MAE values in the testing phase varied from (0.85–1.12), (0.83–0.86), 0.82 and (0.77–1.35), for MLR, SVR, GBR, and ANN, respectively. RMSE values ranged from (1.13–1.46) cm^2 , (1.21–1.28) cm^2 , 1.11 cm^2 and (1.03–2.06) cm^2 , for MLR, SVR, GBR, and ANN, respectively. It was observed in our study that R^2 , MAE, and RMSE values among the different input combinations in the training as well as in the testing dataset in GBR models showed nearly indifferent performance values. GBR showed a stable high R^2 values and low MAE and RMSE values in the dataset studied (Table 8) and outperformed other ML methods studied as also evident from the AR ranking methodology (Table 9). This can be attributed due to GBR models ensembling individual models which can have characteristics of being both weak and overfitting nature but the final ensembled model leads to overcome these problems, also it requires lower data pre-processing leading to low error of prediction and high stability [93,94]. No reports of GBR for LA prediction modeling in plants is known to date. However, the outperformance of GBR over other ML methods are noticed in other fields. A recent study reported that gradient boosting regression algorithm (GBR) and random forest (RF) was employed for predicting and analyzing the net ecosystem carbon exchange (NEE) at UK-Gri, based on the flux data and meteorological data. GBR outperformed three state-of-the-art ML regression prediction models viz., stochastic gradient descent, support vector machine, and Bayesian ridge as GBR allows to tune a sufficient number of hyperparameters [31]. Another study depicted a model built on the gradient boosting regression tree (GBRT) with a limited sample size of data for predicting complex battery dynamics and its lifespan; having many extracted features was shown to outperform other ML algorithms. They suggested the hyperparameters viz., learning rate, maximum number of splits, and number of trees as the key to such performance. A 5-fold cross-validation technique was also applied to GBRT to prevent overfit [32]. Similar tuning of the hyperparameters (n_estimator, maximum_depth, min_sample_split and learning rate) and a 10-fold cross-validation technique were applied in our leaf area modeling study and may have resulted in the best performance of GBR in our study. Again, an experiment was performed for predicting the return temperature over other ML models of district heating systems in Tianjin, China on a decision tree-based ensemble algorithm gradient boosting (GB) and it can be seen that GB outperforms other ML models without the requirement of complex feature transformation and they highlighted the importance of n_estimator in lowering the RMSE values of the tree-based

model GB and RF [95]. A plot of RMSE of GBR with $n_estimator$ tuned for our study is drawn in Figure 5. In our study RMSE equals to 2.78 when $n_estimator$ is 100 and there is a decreasing trend upon increment of the $n_estimator$. RMSE is reduced to around 1.21 and 1.12 when $n_estimator$ is 300 and 400, respectively. Eventually, the RMSE stabilizes to 1.1 at $n_estimator$ 500 and achieves asymptote until it reaches 1000. This revealed the general principle that an improvement in the performance of the serial framework-based model, i.e., GBR model may be achieved (through decreased RMSE) upon increasing the $n_estimator$ number [95]. Similarly, another report supporting GBR performance over other ML methods revealed that gradient boosting models achieved the best performance outperforming random forest and SVM for identifying the relation extraction of medications to the adverse effects of drugs, [96]. A recent report on the outperformance of GBR over ANN can be seen in the prediction of risk analysis of offshore platform integrity for its subsequent use or reuse for alternative energy applications. They studied the effect of stressors on the Remaining Useful Life (RUL). The performance metrics in the study viz., R^2 , MAE, MSE, and RMSE depicted a slightly higher value of R^2 for GBR and a lower value for MAE, MSE, and RMSE than ANN model [30]. The results of these studies on the recent literatures were in general agreement of our observation highlighting our result of the best performance of GBR models in LA prediction in *D. nobile* orchid over the three ML models MLR, SVR, and ANN.

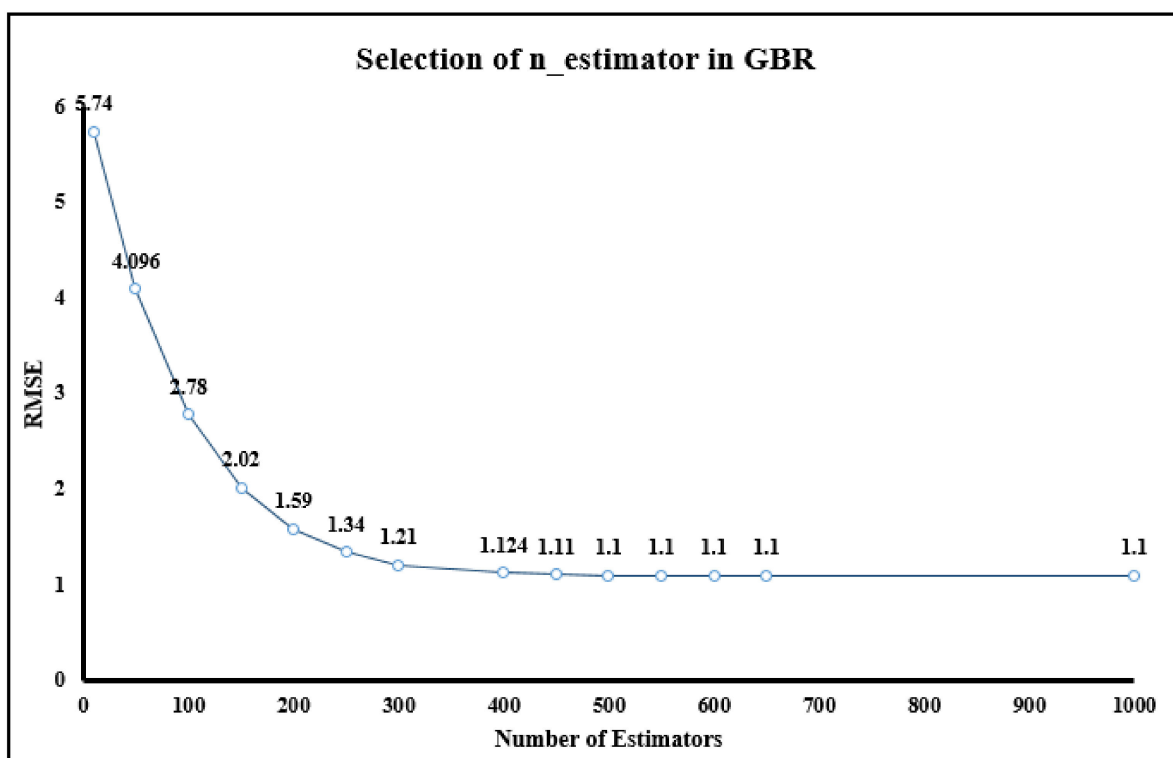


Figure 5. RMSE of GBR with different numbers of $n_estimator$.

In our study, it can be seen that within the 10 best models MLR5 and MLR4 secured second and seventh position, respectively, based on AR ranking methodology (Table 9). Suitability criteria of MLR to produce good prediction accuracy are dependent on the factors which include the absence of multicollinearity between the predictors (measured by VIF and T), strong linear correlation between the predictors and dependent variable, and finally imposing the cross-validation methodology to prevent the over fit of the models. The results depicted from VIF and T values of our experiment show that L, W, and their variants are linearly independent to each other, i.e., there is an absence of multicollinearity. From the feature correlation heatmap it can be inferred that L, W, and their variants are strongly linearly correlated with leaf area (LA) where r ranges from 0.76 to 0.90. The

10-fold cross-validation was also performed to prevent overfitting of the models. Although the MLR model for LA prediction in previous studies did not outperform ANN or other decision tree-based models, there are reports in other fields in agreement with our results of outperformance over ANN and other complex models. A study by [97] showed MLR outperformed random forest based on the VIF values and cross-validation techniques for predicting Soil Organic Carbon stocks. Similar results were seen in another report [98] of the outperformance of MLR over other more complex modeling approaches. The reports of better prediction performance of MLR over ANN were further inferred by [99].

Table 8. Comparative statistical performance analysis of MLR, SVR, GBR, and ANN with different inputs, i.e., LW, LW², LW³, L²W, L²W², L²W³, L³W, L³W², L³W³ studied for LA estimation of *D. nobile*.

Inputs	Model	Training			Testing		
		R ²	MAE	RMSE (cm ²)	R ²	MAE	RMSE (cm ²)
LW	ANN2	0.96	0.87	1.27	0.96	0.86	1.13
LW	GBR1	0.96	0.86	1.18	0.96	0.82	1.11
LW	SVR1	0.94	0.87	1.47	0.96	0.83	1.21
LW	MLR1	0.96	0.90	1.2	0.96	0.9	1.24
LW ²	ANN6	0.96	0.90	1.25	0.96	0.84	1.18
LW ²	GBR2	0.96	0.86	1.18	0.96	0.82	1.11
LW ²	SVR2	0.94	0.89	1.52	0.96	0.84	1.23
LW ²	MLR2	0.97	0.87	1.15	0.96	0.94	1.22
LW ³	ANN10	0.96	0.90	1.27	0.97	0.87	1.18
LW ³	GBR3	0.96	0.86	1.18	0.96	0.82	1.11
LW ³	SVR3	0.94	0.93	1.58	0.96	0.86	1.25
LW ³	MLR3	0.95	1.05	1.34	0.95	1.06	1.42
L ² W	ANN15	0.95	0.94	1.41	0.97	0.86	1.11
L ² W	GBR4	0.96	0.86	1.18	0.96	0.82	1.10
L ² W	SVR4	0.94	0.88	1.48	0.96	0.83	1.21
L ² W	MLR4	0.96	0.86	1.15	0.96	0.86	1.13
L ² W ²	ANN18	0.95	0.93	1.36	0.97	0.77	1.03
L ² W ²	GBR5	0.96	0.86	1.18	0.96	0.82	1.10
L ² W ²	SVR5	0.94	0.89	1.53	0.96	0.85	1.24
L ² W ²	MLR5	0.97	0.84	1.16	0.96	0.85	1.13
L ² W ³	ANN22	0.95	0.97	1.39	0.97	0.86	1.19
L ² W ³	GBR6	0.96	0.86	1.18	0.96	0.82	1.10
L ² W ³	SVR6	0.94	0.92	1.57	0.96	0.86	1.26
L ² W ³	MLR6	0.95	1.04	1.36	0.95	1.06	1.36
L ³ W	ANN26	0.73	2.77	3.78	0.88	1.35	2.06
L ³ W	GBR7	0.96	0.86	1.18	0.96	0.82	1.10
L ³ W	SVR7	0.94	0.88	1.49	0.96	0.85	1.25
L ³ W	MLR7	0.96	0.89	1.17	0.95	0.88	1.22
L ³ W ²	ANN31	0.88	1.41	2.22	0.94	1.01	1.47
L ³ W ²	GBR8	0.96	0.86	1.18	0.96	0.91	1.10
L ³ W ²	SVR8	0.94	0.89	1.53	0.96	0.85	1.26
L ³ W ²	MLR8	0.96	0.88	1.17	0.96	0.88	1.2
L ³ W ³	ANN35	0.91	1.29	2.34	0.91	0.97	1.31
L ³ W ³	GBR9	0.96	0.86	1.18	0.96	0.91	1.10
L ³ W ³	SVR9	0.94	0.92	1.56	0.96	0.86	1.28
L ³ W ³	MLR9	0.95	1.11	1.42	0.94	1.12	1.46

Table 9. Comparative statistical performance analysis of MLR, SVR, ANN, and GBR based on average rank (AR) ranking methodology.

Models	Ranks Based on Training Results			Ranks Based on Testing Results			Average Rank (AR)	Final Rank
	R ²	MAE	RMSE (cm ²)	R ²	MAE	RMSE (cm ²)		
GBR7	7	4	8	11	2	2	5.67	1
MLR5	1	1	3	7	13	12	6.17	2
GBR6	8	5	9	12	3	3	6.67	3
GBR5	9	6	10	13	4	4	7.67	4
GBR4	11	8	11	14	5	5	9.00	5
GBR9	3	2	6	9	28	6	9.00	6
MLR4	10	7	1	5	19	13	9.17	7
GBR8	5	3	7	10	29	7	10.17	8
GBR3	12	9	12	15	6	8	10.33	9
GBR2	14	10	13	16	7	9	11.50	10
ANN18	22	27	20	2	1	1	12.17	11
MLR2	2	12	2	6	30	22	12.33	12
MLR8	4	15	4	8	25	18	12.33	13
GBR1	17	11	14	17	8	10	12.83	14
ANN10	13	22	18	1	24	16	15.67	15
ANN6	15	23	16	19	11	15	16.50	16
ANN2	18	13	17	20	20	14	17.00	17
MLR7	6	18	5	30	26	21	17.67	18
ANN15	23	29	23	4	18	11	18.00	19
ANN22	21	30	22	3	17	17	18.33	20
SVR1	33	14	25	21	9	19	20.17	21
MLR1	16	24	15	18	27	25	20.83	22
SVR4	30	17	26	22	10	20	20.83	23
SVR7	27	16	27	23	14	26	22.17	24
SVR2	32	21	28	24	12	23	23.33	25
SVR8	26	19	29	25	15	28	23.67	26
SVR5	29	20	30	26	16	24	24.17	27
SVR9	25	25	31	27	21	30	26.50	28
SVR6	28	26	32	28	22	29	27.50	29
MLR6	20	31	21	32	34	32	28.33	30
SVR3	31	28	33	29	23	27	28.50	31
MLR3	24	32	19	31	33	33	28.67	32
MLR9	19	33	24	33	35	34	29.67	33
ANN35	34	34	35	35	31	31	33.33	34
ANN31	35	35	34	34	32	35	34.17	35
ANN26	36	36	36	36	36	36	36.00	36

Previous studies on leaf area modeling by [34,47,54,60,100,101] depicted that ANN outperformed MLR. On the contrary, in our study ANN could not perform better in other input combinations than MLR as well as GBR except in L²W² input combinations of testing results where, ANN18 with (2-5-10) architecture secured a rank of eleventh based on AR ranking (Table 9) (R² = 0.97, MAE = 0.77 and RMSE = 1.03), which may be due to its inherent disadvantages such as overfitting. The reason behind this abrupt performance is due to the observed training results of ANN18, which shows R² = 0.95, MAE = 0.93, and RMSE = 1.36 denoting an overfit. The problems of local minima as well as selection of suitable hyper-parameters suitable for our study may be the other causes of such results of ANN18 [102].

From Table 9 It can be seen that SVR occupied ranks between twenty-first to thirty-first among the 36 models. So, it can be inferred that SVR was unable to perform better than three other ML models tested in our dataset.

For evaluating the prediction quality, a scatter plot of predicted LA and measured LA values along with the reference line ($y = x$) of *D. nobile* for the four ML models viz., MLR, SVR, GBR, and ANN in the testing period was used (Figure 6). A deviation of the regression line of predicted LA and measured LA values from the reference line is suggestive of a bias.

In our study, the model showing the least bias was gradient boosting regression (GBR). Thus, it can be depicted that GBR models have a better capacity to predict LA in *D. nobile* than other ML models used. ANN and SVR models showed maximum deviations from the measured leaf area for all the input combinations studied.

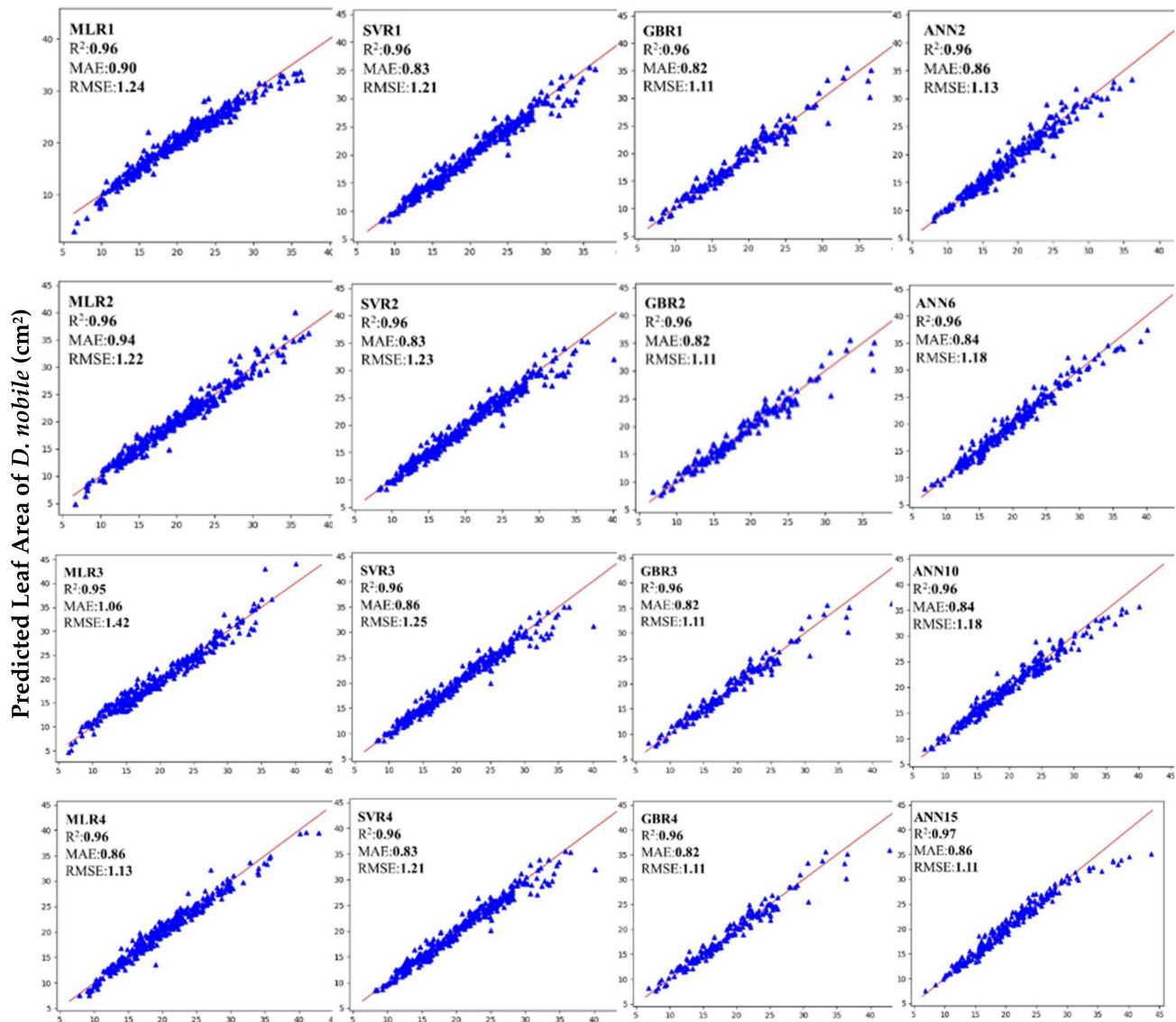


Figure 6. Cont.

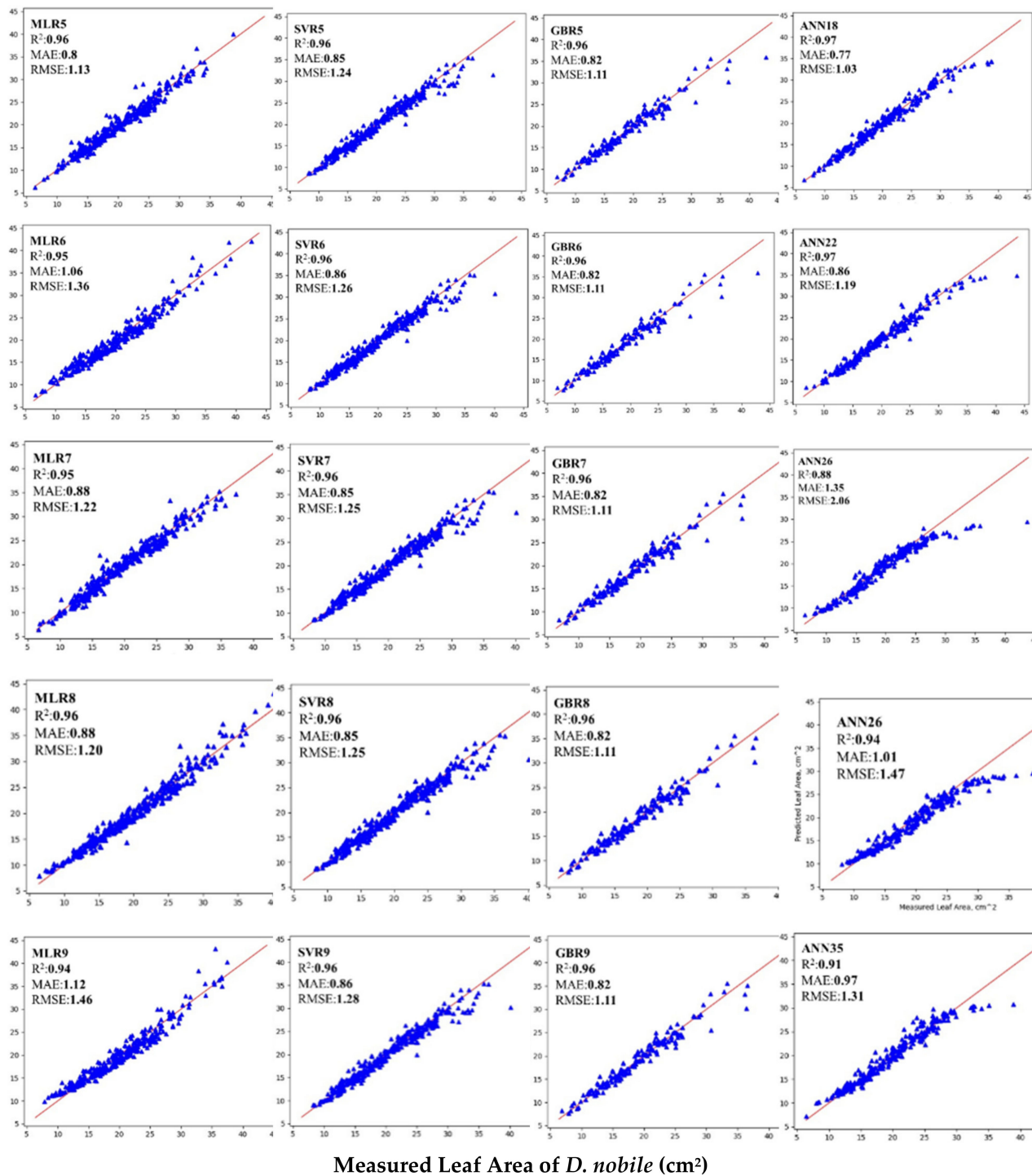


Figure 6. Comparison of predicted and measured values of leaf area of *D. nobile* using MLR, SVR, GBR, and ANN models in different input combinations. The x -axis represents measured values of leaf area by ImageJ software and y -axis denotes the predicted values by the two models above. The red line represents the reference line ($y = x$). Blue triangles represent the regression line of the predicted and measured leaf area. R^2 refers to the coefficient of determination, MAE refers to mean absolute errors and RMSE refers to root mean square errors (cm²).

The main limitations of performance metrics viz., R^2 , MAE, and RMSE metrics as seen in our study is that if algorithm A outperforms B based on only r^2 , then again it takes down A and brings up B on the basis of either MAE or RMSE. This is in accordance with the

theorem of No Free Lunch (NFL) and the solution depends on the domain of the problem and knowledge of algorithms by the analysts [103,104]. Thus, ranking of the algorithms by average rank (AR) was chosen, whereby for each 36 ML models which includes the four ML techniques viz., MLR, SVR, GBR, and ANN ranks were assigned based on R^2 , MAE, and RMSE. The average ranks were obtained by averaging the ranks for each model. The final rank was derived by arranging the average ranks in ascending order. Table 9 depicts a comparative statistical analysis of MLR, SVR, ANN, and GBR algorithms and assigned ranks on R^2 , MAE, and RMSE each based on the training results and testing results to calculate the average rank. The final rank depicted that GBR7 secured the first rank among the 36 models as well as GBR algorithm has a frequency of occurrence of nearly 80% among the top 10 ML models. This provided a conclusive decision to select GBR as the best model for leaf area predictive modeling in the case of the *Dendrobium* orchid. Subsequent to the GBR algorithm, MLR occupied nearly 20% frequency among the top 10 ML models. Figure 7 depicts the frequency of the occurrence of the ML models studied in our experiment among the top ten ranks. It can be seen from the graph that GBR occupies 80% and MLR occupies 20% among the top ten ranks. On the other hand, SVR and ANN occupy nil percentage in the proposed rank group.

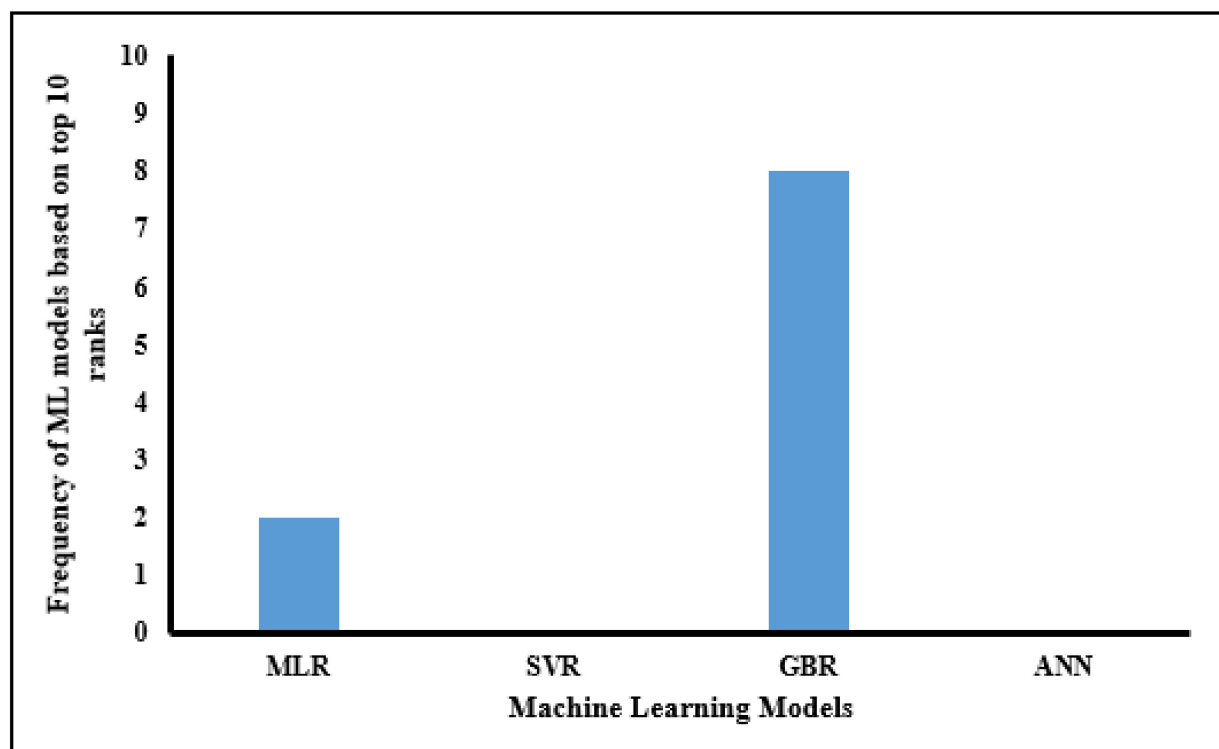


Figure 7. Frequency of ML models based on top 10 ranks by AR ranking methodology.

Light interception by the plant is totally dependent on the two-dimensional structure viz., length (L) and width (W) [105] and thus highlighting the importance of only those two proxy variables [13,40,62,87,106–109]. In this study, a non-destructive methodology of LA modeling in *D. nobile* through ML techniques was proposed that can be employed to develop a web portal or mobile app for providing an interactive user interface to obtain rapid, precise, and accurate leaf area estimation results from simple measurements such as leaf length (L) and leaf width (W). The proposed best model, i.e., GBR can play a crucial role in real time LA estimation of *D. nobile* avoiding the need for expensive instruments and destructive sampling procedures.

4. Conclusions

Leaf area (LA) is a biophysical variable of utmost importance which helps in maintaining various physiological processes ultimately helping in the adaptation of the plant to its environment. LA estimation is necessary for calculating various physiological indices. In summary, the main goal of our study is to develop a simple, non-destructive, low cost, reliable machine learning LA prediction model for *D. nobile*. This is the first attempt in the *Orchidaceae* family for leaf area estimation using ML techniques. To the best of our knowledge, the use of GBR and SVR models for predicting leaf area have not been reported in the earlier studies. Our experiment showed that GBR outperforms the other models viz., MLR, SVR, and ANN in terms of R^2 , MAE, and RMSE in the testing period. Thus, GBR provided the best predictive capacity for leaf area estimation in the case of *D. nobile* with higher R^2 (0.96) and lower MAE (0.82–0.91) and RMSE (1.10–1.11) cm^2 values. Model ranks based on average rank (AR) also statistically suggested the outperformance of GBR over the other state-of-the-art ML models such that it occupies first rank and a frequency of around 80% of the top ten models tested in our study. Limitations of our study include identification of the theoretical basis of the causal relationship between the rigidly stable behavior of GBR in the training phase irrespective of different input combinations. The leaf area models developed in our study were carried on healthy plants under control conditions only, ignoring the biotic and abiotic stress situation present in real situation. Sampling for images can be performed in the future in real conditions. In the future, a substantial increase in the training dataset has the possibility to be used for a deep learning set up for developing a nearly perfect model. Machine vision field can be considered for automated area calculation of *D. nobile* using leaf images directly.

Supplementary Materials: The following supporting information can be downloaded at: <https://www.mdpi.com/article/10.3390/app12094770/s1>, Supplementary S1: <http://doi.org/10.17632/8tk2sc4ytg.1>. (published on 28 June 2021) and KRISHI, ICAR research data repository for knowledge management [<http://krishi.icar.gov.in/jspui/handle/123456789/71908> (published on 6 May 2022)]

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