

Research Article Convolutional Recurrent Neural Networks for Observation-Centered Plant Identification

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Traditional image-centered methods of plant identification could be confused due to various views, uneven illuminations, and growth cycles. To tolerate the significant intraclass variances, the convolutional recurrent neural networks (C-RNNs) are proposed for observation-centered plant identification to mimic human behaviors. The C-RNN model is composed of two components: the convolutional neural network (CNN) backbone is used as a feature extractor for images, and the recurrent neural network (RNN) units are built to synthesize multiview features from each image for final prediction. Extensive experiments are conducted to explore the best combination of CNN and RNN. All models are trained end-to-end with 1 to 3 plant images of the same observation by truncated back propagation through time. The experiments demonstrate that the combination of MobileNet and Gated Recurrent Unit (GRU) is the best trade-off of classification accuracy and computational overhead on the Flavia dataset. On the holdout test set, the mean 10-fold accuracy with 1, 2, and 3 input leaves reached 99.53%, 100.00%, and 100.00%, respectively. On the BJFU100 dataset, the C-RNN model achieves the classification rate of 99.65% by two-stage end-to-end training. The observation-centered method based on the C-RNNs shows potential to further improve plant identification accuracy.

1. Introduction

With the development of computer technology, there is increasing interest in the image-based identification of plant species. It can get rid of the dependence on the botany professionals and help the public identify plants easily. With the growth of demand, a growing number of researchers began to focus on plant identification and developed various sophisticated models.

Traditional automated plant identification builds classifiers on top of hand-engineered features whose objects are flowers, leaves, and other organs. For example, Lee et al. [1] designed a mobile system of flower recognition using color, texture, and shape features and the accuracy was 91.26%. Furthermore, more studies considered the features of leaves. Novotný and Suk [2], Zhao et al. [3], Liu et al. [4], and Li et al. [5] extracted shape features of leaves to recognize the plant species using the nearest neighbor (NN) classifier, the back propagation (BP) neural network classifier, and the support vector machine (SVM) classifier, respectively. En and Hu [6], Liu and Kan [7], and Zheng et al. [8] used shape



and texture features of leaves to identify the plants with the base classifier, deep belief network (DBN) classifier, and SVM classifier, respectively. Wang et al. [9] extracted a series of color, shape, and texture features from 50 foliage samples, and the accuracy identified by SVM classifier was 91.41%. Gao et al. [10] calculated the comprehensive similarity of geometric features, texture features, and corner distance matrix to recognize plant leaves and the accuracy on the Flavia dataset was 97.50%. Chen et al. [11] extracted the distance matrix and corner matrix of leaves and identified plant leaves in the top 20 of the highest similarity result set chosen by the kNN classifier. The accuracy on Flavia dataset reaches 99.61%. The handcrafted feature extractor is time-consuming and relies on the expertise to a certain extent; it also has poor generalization ability to big data. Therefore, conventional methods of plant identification are not suitable for images with huge species and complex background. Developments in deep learning have made significant contribution to in-field plant identification.

Deep learning is a machine learning method that extracts features automatically from raw data supervised by an end-to-end training algorithm. It has been applied to the field of plant identification in recent years. Dyrmann et al. [12] used the CNN to identify weeds and crops at the early growth stages and achieved 86.2% average accuracy rate. Grinblat et al. [13] used deep learning for leaf identification of white bean, red bean, and soybean and the accuracy was $92.6 \pm 0.2\%$. Zhang and Huai [14] used the CNN to identify the leaves of the self-expanding dataset based on the PlantNet dataset. The accuracy using the SVM classifier and the softmax classifier was 91.11% and 90.90% in simple background and 31.78% and 34.38% in complex background, respectively. Ghazi et al. [15] used three deep learning networks, that is, GoogLeNet, AlexNet, and VGGNet, to identify species on the LifeCLEF 2015 dataset and the overall accuracy of the best model was 80%. Sun et al. [16] designed a 26-layer deep learning model consisting of 8 residual building blocks for large-scale plant classification in natural environment and the recognition rate reached 91.78% on the BJFU100 dataset. The LifeCLEF plant identification challenge plays an important role in the field of plant automation recognition. The 2017th LifeCLEF plant identification challenge provides 10.000 plant species illustrated by a total of 1.1 M images. The winner of the challenge combined 12 trained models with four GoogLeNet, ResNet-152, and ResNetXT CNN models on three kinds of datasets. The Mean Reciprocal Rank (MRP) is 0.92 and the top 1 accuracy is 0.885 [17].

The previous studies identify plants from a single image (i.e., perform the image-centered plant identification). Although these methods are straightforward, they are not consistent with human cognitive habits. A series of factors like seasonal color changes, shape distortion, leaf damage, and uneven illumination cause significant intraclass variances. In fact, botanists often observe plants from multiple views or varied specimens, because the combination of multiple views can effectively improve the recognition accuracy. To circumvent the limitations of image-centered identification, observation-centered identification is proposed to mimic human behaviors. Each observation includes several images of the same species with different views. The convolutional recurrent neural networks (C-RNNs) are trained end-toend to automatically extract and synthesize features in one observation. The recognition accuracy on the Flavia and BJFU100 datasets is further improved by the C-RNN models.

2. The Proposed Method

Almost all public image datasets are image-centered, while the observation-centered data are constructed based on the image-centered dataset. So, the plant images of the same species taken in the nearby time and location are selected as one observation. The observation-centered training and testing sets are built using different sets of samples from an image-centered dataset. To improve the generalization of the model, a sequence of images in the training set are permutated randomly.

As shown in Figure 1, the C-RNN model inputs all images in one observation and outputs one prediction, which works in a many-to-one approach. The C-RNN model consists of CNN backbones and RNN units. The CNN backbones





FIGURE 1: The C-RNN model for observation-centered plant identification. The green arrows are the forward inference and the dashed red arrows are the truncated back propagation through time.

extract features from each image belonging to one observation, which include the residual network (ResNet) [18], InceptionV3 [19], Xception [20], and MobileNet [21]. Then, the RNN units including the simple RNN [22], Long Short Term Memory (LSTM) [23, 24], and Gated Recurrent Unit (GRU) [25] synthesize all features to implement observation-centered plant identification through the softmax layer. And the fully connected layer with the rectified linear unit (ReLU) connects the CNN backbones and the RNN units. Table 1 shows the MobileNet with GRU as the example to illustrate the network architecture.

As the RNNs with fixed architecture are differentiable end-to-end, the derivative of the loss function can be calculated with respect to each of the parameters in the model [26]. The categorical cross-entropy loss is used in the C-RNN model for the multiclassification, and the loss function is defined to be

$$L = \sum_{t} L_t \left(y_{nt}, \, \widehat{y}_{nt} \right) = -\sum_{t} y_{nt} \log \, \widehat{y}_{nt}, \tag{1}$$

where \hat{y}_{nt} is the *t*th element in the classification score vector \hat{y}_n , y_{nt} is the *t*th element in the classification label vector y_n , and *n* is the step size of RNN unit.

The process of truncated BPTT is shown as the red dashed arrows in Figure 1, and the gradient calculation formulas for the weights *V*, *W*, and *U* are expressed as follows [27]:

$$\frac{\partial L}{\partial V} = \frac{\partial L}{\partial \hat{y}_3} \frac{\partial \hat{y}_3}{\partial V} = \frac{\partial L}{\partial \hat{y}_3} \frac{\partial \hat{y}_3}{\partial z_3} \frac{\partial z_3}{\partial V} = (\hat{y}_3 - y_3) \otimes s_3, \quad (2)$$

$$\frac{\partial L}{\partial W} = \frac{\partial L}{\partial \hat{y}_3} \frac{\partial \hat{y}_3}{\partial s_3} \frac{\partial s_3}{\partial W},\tag{3}$$

$$\frac{\partial L}{\partial U} = \frac{\partial L}{\partial \hat{y}_3} \frac{\partial \hat{y}_3}{\partial s_3} \frac{\partial s_3}{\partial U},\tag{4}$$

$$z_3 = V s_3, \tag{5}$$

$$s_3 = \tanh\left(Ux_3 + Ws_2\right),\tag{6}$$

where *L* is the loss function, *s* is the simplification of the RNN unit calculation process, and \otimes is the outer product of two

2

Journal of Electrical and Computer Engineering

Layer name	Input size	Output size	Filter shape
cnn_l	$224 \times 224 \times 3$	1 × 1024	MobileNet Alpha = beta = 1
cnn_2	$224 \times 224 \times 3$	1×1024	MobileNet Alpha = beta = 1
cnn_3	$224 \times 224 \times 3$	1×1024	MobileNet Alpha = beta = 1
time_ distributed (fc)	3×1024	3×1024	1024×1024
rnn (gru)	3×1024	1×128	1024×128
fc (softmax)	1×128	1×32	128×32

TABLE 1: C-RNN architecture.

vectors. Considering (6), s is related to W and U, and it can be deduced that

$$\frac{\partial s_t}{\partial W} = \frac{\partial s_t}{\partial W} + \frac{\partial s_t}{\partial s_{t-1}} \frac{\partial s_{t-1}}{\partial W}$$

$$= \left(1 - s_t\right)^2 \left(s_{t-1} + W \frac{\partial s_{t-1}}{\partial W}\right), \quad (7)$$

$$\frac{\partial s_t}{\partial U} = \frac{\partial s_t}{\partial U} + \frac{\partial s_t}{\partial s_{t-1}} \frac{\partial s_{t-1}}{\partial U} = \left(1 - s_t\right)^2 \left(x_t + W \frac{\partial s_{t-1}}{\partial U}\right).$$

So, the gradients of *W* and *U* are expressed as follows [27]:

$$\frac{\partial L}{\partial W} = \sum_{k=0}^{3} \frac{\partial L}{\partial \hat{y}_{3}} \frac{\partial \hat{y}_{3}}{\partial s_{3}} \left(\prod_{j=k+1}^{3} \frac{\partial s_{j}}{\partial s_{j-1}}\right) \frac{\partial s_{k}}{\partial W},$$

$$\frac{\partial L}{\partial U} = \sum_{k=0}^{3} \frac{\partial L}{\partial \hat{y}_{3}} \frac{\partial \hat{y}_{3}}{\partial s_{3}} \left(\prod_{j=k+1}^{3} \frac{\partial s_{j}}{\partial s_{j-1}}\right) \frac{\partial s_{k}}{\partial U}.$$
(8)

So, the C-RNN model can be trained end-to-end and the identification error is minimized by stochastic gradient descent with truncated back propagation through time (BPTT).

3. Experiment and Results

In this section, the method of constructing the observationcentered dataset is described in detail. Then, extensive experiments are conducted on the leaves observation-centered dataset with the combinations of different CNN backbones and RNN units to show a significant performance improvement compared with traditional transfer learning.

3.1. The Dataset. The leaves observation-centered dataset is based on the Flavia dataset [28]. There are 1,907 samples of 32 species in Flavia and each sample is an RGB leaf scan image with white background. All of the images for the same class were taken in the same experimental environment, so they could be thought of as the observation-centered data. Figure 2 shows the observation-centered example for leaves with different input numbers.

To reduce variability and overfitting, 10-fold crossvalidation is adopted. The dataset is divided into 10 complementary subsets randomly. In one fold of the experiment, one





FIGURE 2: Observation-centered sample with (a) one, (b) two, and (c) three leaves. Samples containing less than three leaves are padded with zero.

subset is randomly chosen as the test set and the rest belong to the training set. Considering the size of the dataset, there are 1717 images for training and 190 images for test in the first 9 folds and there are 1710 images for training and 197 images for test in the 10th fold.

To simulate collection of multiple leaves from one tree, each image is combined with 0 to 2 randomly chosen leaves of the same species to construct one observation-centered sample. The observation-centered samples with less than 3 leaves are padded with 0 (as shown in Figure 2). All leaves in the test set are held out from the training set for each fold.

3.2. Results. The models are implemented by the deep learning framework Keras [29] with TensorFlow [30] backend. All the experiments were conducted on an Ubuntu 16.04 Linux



FIGURE 3: Effects of (a) ResNet50, (b) InceptionV3, (c) Xception, and (d) MobileNet CNN with three different RNN units on the 10-fold test accuracy. The pentagrams in (b), (c), and (d) represent the outlier of the results.

server with an AMD Ryzen 7 1700x CPU (16 GB memory) and an NVIDIA GTX 1070 GPU (8 GB memory). The batch size for each network is 32, the learning rate is 0.0001, and the optimizer is RMSprop in training phase. The test accuracy is compared after 50 epochs.

3.2.1. Effects of RNN Units and CNN Backbones. To explore the best RNN units, the models are implemented with different RNN units: simple RNN, LTSM, and GRU. Besides, the effects of different CNN backbones are also considered. Four state-of-the-art CNN backbones—ResNet50, InceptionV3, Xception, and MobileNet models—are pretrained on the ImageNet dataset with 1.2 million images of 1000 classes. Then, the last fully connected layers are replaced by the aforementioned RNN units.

Figure 3 shows the 10-fold test accuracy of each CNN with 3 different RNN units at the last epoch. All of the RNN units work well with InceptionV3, Xception, and MobileNet, and the test accuracy is close to 100%; only several outliers exist as Figures 3(b), 3(c), and 3(d) indicate. The effect of ResNet50 was marginal due to the unstable convergence and the test accuracy is about 80% as shown in Figure 3(a).

TABLE 2: 10-fold test mean accuracy with different CNN and RNN units.

	Simple RNN	LTSM	GRU
ResNet50	0.7987	0.8276	0.7973
InceptionV3	1.0000	0.9995	0.9995
Xception	0.9989	0.9995	1.0000
MobileNet	0.9995	1.0000	1.0000

To explore the best combination of CNN and RNN, the means of 10-fold test accuracy at the last epoch are listed in Table 2. The combinations of InceptionV3 with simple RNN, Xception with GRU, MobileNet with LTSM, and MobileNet with GRU outperform the others. All of the four models achieve 100% mean accuracy. Considering that the InceptionV3 with simple RNN has 2,249,880 trainable parameters, the Xception with GRU has 2,545,056 trainable parameters, the MobileNet with LTSM has 1,644,064 trainable parameters, and the MobileNet with GRU has 1,496,480 parameters, the MobileNet with GRU model with fewer parameters is considered as the best model.



FIGURE 4: The mean of training loss, training accuracy, and test accuracy of the MobileNet with GRU model in one fold.

The MobileNet backbone for the best model has 27 convolutional layers counting depth-wise and pointwise convolutions as separate layers, and the width multiplier and the resolution multiplier are 1. The number of the FC layer units after MobileNet backbone is 1024 and the number of the GRU unit hidden nodes is 128. The total number of the network parameters is 4,725,344 and there are 3,228,864 untrained parameters. Figure 4 shows the training process of the best model. The training loss declines rapidly in the first 10 epochs and decreases slowly in the next 10 epochs, and then it approaches 0 at the 50th epoch. The training accuracy and test accuracy rise rapidly in the first 5 epochs and increase slowly in the next 5 epochs, and then they approach 1 finally. The increments in both training and test accuracy indicate that the model is not overfitted.

3.2.2. Comparison with Transfer Learning. To make a fair comparison, the MobileNet retrained by traditional transfer learning is compared to the MobileNet with GRU model when inputting 1 to 3 leaves. When the number of inputting leaves is smaller than 3, the extra input channels of the RNN units were filled with 0.

Figure 5 shows the 50th epoch test accuracy. The 10-fold test mean accuracy of retrained MobileNet is 99.79%, and that of our model inputting 1 leaf, 2 leaves, and 3 leaves is 99.53%, 100.00%, and 100.00%, respectively. The result shows that the traditional transfer learning is better when inputting only 1 leaf, while the C-RNNs show advantage when observation-centered images are available.

3.2.3. Comparison with Non-End-to-End Method. In order to evaluate the effect of end-to-end training for the C-RNN in the observation-centered identification, a comparative experiment with a non-end-to-end method, that is, majority voting, was carried out. Majority voting got 0.667 for the average classification score in LifeCLEF 2015 plant task [31]. The experiments have been performed on the BJFU100 [16] dataset. The BJFU100 dataset consists of 10,000





FIGURE 5: The MobileNet with GRU model compared with traditional transfer learning.



FIGURE 6: Each column shows one observation-centered sample from the BJFU100 dataset.

images with 100 species of ornamental plants in Beijing Forestry University campus by a mobile phone. Two images of the same class taken in the nearby time and location are selected as one observation. Figure 6 shows the typical observation-centered samples selected from the BJFU100 dataset.

The CNN uses the softmax classifier to identify one image at one time. Then, majority voting gives the final prediction by counting votes from all CNN predictions. The CNN in the majority voting is trained end-to-end on a single image, while the error of voting cannot be corrected by back propagation and stochastic gradient descent.

The MobileNet is the CNN for both C-RNN and majority voting. Firstly, the MobileNet is trained on a single image of BJFU100 for 86 epochs with the test accuracy of 95.8%. The majority voting has no trainable parameters and directly inputs the prediction of MobileNet, while the C-RNN model can be further fine-tuned by second-stage endto-end training. The trained MobileNet without the final layer is transferred into the C-RNN model for further 20-epoch



FIGURE 7: The test accuracy of C-RNN evolves as training advances.

end-to-end training. The training parameters and networks parameters are the same as the above experiments except that the batch size is changed to 16.

The single MobileNet, majority voting, MobileNet with simple RNN, MobileNet with LSTM, and MobileNet with GRU achieve the test accuracy of 95.8%, 98.95%, 99.35%, and 99.65%, respectively. Figure 7 shows the function of the test accuracy of MobileNet with GRU and the majority voting during training. The C-RNN model trained by two-stage end-to-end training outperforms the majority voting method by 0.7%.

4. Conclusions

In this paper, the C-RNN models were proposed for observation-centered plant identification. The CNN backbones extract features and the RNN units integrate features and implement classification. The combination of MobileNet and GRU is the best trade-off of classification accuracy and computational overhead on the Flavia dataset. The test accuracy reaches 100%, while it has fewer parameters. Experiments on the BJFU100 dataset show that the C-RNN model trained by two-stage end-to-end training further improves the accuracy of majority voting method by 0.7%. The proposed C-RNN model mimics human behaviors and further improves the performance of plant identification, which has great potential in in-field plant identification.

In the future work, an observation oriented dataset with a large amount of plants in an unconstrained environment will be constructed for further experiments.

Conflicts of Interest

The authors declare that there are no conflicts of interest regarding the publication of this paper.

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