Original papers

# Analysis of transfer learning for deep neural network based plant classification models

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ABSTRACT

Plant species classification is crucial for biodiversity protection and conservation. Manual classification is time-consuming, expensive, and requires experienced experts who are often limited available. To cope with these issues, various machine learning algorithms have been proposed to support the automated classification of plant species. Among these machine learning algorithms, Deep Neural Networks (DNNs) have been applied to different data sets. DNNs have been however often applied in isolation and no effort has been made to reuse and transfer the knowledge of different applications of DNNs. Transfer learning in the context of machine learning implies the usage of the results of multiple applications of DNNs. In this article, the results of the effect of four different transfer learning models for deep neural network-based plant classification is investigated on four public datasets. Our experimental study demonstrates that transfer learning can provide important benefits for automated plant identification and can improve low-performance plant classification models.

## 1. Introduction

Recent studies estimate that there are currently between 220,000 and 420,000 flowering plant species on Earth (Scotland and Wortley, 2003; Mora et al., 2011; Govaerts, 2001). Although, many species are under threat of extinction, or have already been extinct due to pollution and natural disasters, still many species are waiting to be discovered. For supporting the research on biodiversity conservation and ecological monitoring, accurate identification of plants is necessary in, for example, endangered species monitoring, assessment of weed control actions, and analysis of species distribution under climate change (Wäldchen et al., 2018). Besides of their importance for nature and ecological balance, many plant species are raw material for the medicine and chemical industries. Hence, there is a continuing need for proper plant identification.

The objective of plant classification systems is to help non-expert and non-botanist users to identify the plants automatically. Otherwise, this identification process is too long, time consuming, and expensive. Speeding up this process provides several benefits such as low cost, less effort, and more time allocation for the other tasks. Plant recognition systems can be also used in making intelligent field guides, educational

tools, and agricultural practices and forestry automation. Besides of the agricultural domain, the proposed technology can be applied to other domains including healthcare, retail, and automotive.

The identification of plants is considered as a process which assigns a plant under study to a taxon based on two kinds of features, namely quantitative and qualitative characters. The flower length and plant height are examples of quantitative features, while the flower color and leaf shape are examples of qualitative features (Wäldchen et al., 2018). Although there are many similarities between individuals of the same species, the exact identification of plants is certainly not trivial and requires sophisticated knowledge. Thus, experts in plant species identification need to be consulted, but they are however costly and not always available. Moreover, the number of plant classification experts seem be also decreasing (Hopkins and Freckleton, 2002).

Gaston and O'Neill (2004) indicated 15 years ago that automated species identification systems was not the norm at that time (Wäldchen et al., 2018) due to the difficulty of the implementation, the high labor intensiveness, and the high cost. Yet, with the recent dramatic advancements in Machine Learning (ML), Cloud Computing, Internet of Things (IoT), and Distributed Computing this situation has definitely changed, and automated identification systems have now become

feasible and are also more and more applied.

The automated identification of plants in several domains can now replace manual processing of huge plant species catalogs by subject matter experts (Grinblat et al., 2016) and thus provides several benefits including reduction of cost and time, but also more accurate identification. The success of automated plant identification has led to its broad applications in biological taxonomy studies, definition of industrial raw materials, implementation of automated systems for the precision agriculture, detection of diseased plants, and educational purposes. With the help of plant identification systems in precision agriculture, for example, it is possible to spray herbicides only on weeds (Pahikkala et al., 2015), which make the agriculture more cost-effective and eco-friendly. Food safety, remote-sensing for farming, and deforestation control are some of the other application areas which can get benefit from the automated plant classification approaches.

To realize automated classification of plant species, various machine learning algorithms have been proposed. Among these machine learning algorithms, Deep Neural Networks (DNNs) have been applied to different data sets. DNNs have been however often applied in isolation and no systematic effort has been made to reuse and transfer the knowledge of different applications of DNNs. Transfer learning in the context of machine learning implies the reuse of the earlier acquired knowledge in similar tasks. In principle, transfer learning can be applied for any machine learning algorithm in any application domain. Yet, although several studies have focused on the use of deep learning approaches for the plant classification problem (Grinblat et al., 2016; Lee et al., 2017; Dyrmann et al., 2016; Yalcin and Razavi, 2016; Strothmann et al., 2017), there is not a comprehensive study yet, which evaluates several transfer learning scenarios for deep learning algorithms. Hence, we focus on the following research questions (RQs):

- RQ1: To what extent can plant classification benefit from transfer learning in deep learning?
- RQ2: How do the different transfer learning scenarios perform at improving the performance of plant classification models using deep learning?

To answer these questions, we designed and implemented five classification models including the baseline model (end-to-end CNN model) by applying four transfer learning strategies on deep learning-based models. The first model (baseline approach) is an end-to-end Convolutional Neural Networks (CNN) model which were tested on the four datasets separately. As the second model (cross-dataset fine tuning), we trained the same algorithm with three datasets and fine-tuned the model on the remaining one. For the third model (fine tuning), we fine-tuned the pre-trained CNN models which were previously trained with ImageNet database. As the fourth model, features were extracted from pre-trained networks and then, two traditional classification algorithms were applied based on these features. Finally, a combination of RNN (Recurrent Neural Network) and CNN algorithms is tested for the classification of plants. In this paper we analyze and compare these different approaches and provide the lessons learned.

Section 2 introduces the related work, Section 3 provides the transfer learning concept, Section 4 presents the adopted approach, Section 5 indicates the experimental results, Section 6 discusses the findings and the potential threats to validity. Section 7 explains the conclusion.

## 2. Related work

The use of deep learning algorithms in machine vision problems is becoming more and more popular. The plant and the leaf identification have been previously studied with different techniques by various researchers.

The initial approaches on these problems used the color information to distinguish the plant from the soil (Woebbecke et al., 1995; Gerhards and Christensen, 2003). Some studies utilized the vein morphology (Sack et al., 2008; Scoffoni et al., 2011). The leaf veins include many textural and shape properties which can be useful for the vision-based plant identification. Some leaf-based studies also used the shape information (Agarwal et al., 2006; Neto et al., 2006). Several studies combined the shape and texture information extracted from the leaf (Husin et al., 2012) while other studies utilized the color and texture information (Pydipati et al., 2006). A method is proposed by Larese et al. (2014) in which computer vision techniques are used to extract the morphological leaf vein features. Extracted features are later used with machine learning algorithms to predict the three different plant species. These researchers improved this method later on in a new research project (Larese et al., 2014).

Spectroscopic methods are also applied in plant genre classification. The reflectances of the infrared, multispectral, and visible bands are used for the feature extraction in these methods (Mattila et al., 2013; Wang et al., 2007; Tyystjärvi et al., 2011). Muthevi and Uppu (2017) utilized Local Binary Patterns (LBP) in their study. Different types of LBP extraction methods, Signed component of CLBP (SCLBP) were applied for the feature extraction. Murat et al. (2017) combined different shape descriptors for the leaf classification of the tropical shrub species. The proposed method was tested with Flavia and Swedish Leaf datasets. Yousefi et al. (2017) used view rotation invariant features which are extracted from the Fast Fourier Transform and Discrete Wavelet Transform. Yu et al. (2016) developed a method utilizing the leaf contour and venation. Horaisová and Kukal (2016) proposed a method for the invariant leaf detection.

The use of deep learning in plant and leaf identification is a relatively new approach. Grinblat et al. (2016) proposed a CNN model to work with leaf veins. Leaf veins are extracted as a binary mask and an end-to-end CNN model is trained with these masks. dos Santos Ferreira et al. (2017) combined the superpixel segmentation algorithm and CNN to build a weed segmentation system. CNNs are used in weed detection on images which are segmented with the superpixel algorithm. Barré et al. (2017) proposed a CNN architecture for the leaf identification. They applied the Foliage, LeafSnap, and Flavia datasets and built a CNN model for the classification. An architecture similar to well-known CNN models like AlexNet and CifarNet was proposed in their study. Jeon and Rhee (2017) used the transfer learning from pre-trained CNN models. Pre-trained GoogleNet was used for the feature extraction. Leaf images in different scales were given to GoogleNet as input and the activation values of different layers were stored as features.

According to the literature survey outlined in this section, there is no study which evaluates the effect of several transfer learning scenarios for plant classification models in detail. Hence the work that we present in this article can be considered complementary to the earlier studies.

The classification accuracy comparison of some related studies with the best result in our experiments are shown in Table 1. In Table 1, datasets are represented in columns and methods are represented in rows. Although different experimentation setups are used in these studies like k-fold cross validation and test-train data splitting, the best results obtained are shown in Table 1. The results obtained from our experiments are promising and comparable with the recent deep based studies. The deep learning based studies (Lee et al., 2017; Yalcin and Razavi, 2016; Jeon and Rhee, 2017) have a superiority in classification accuracy compared to traditional ones (Murat et al., 2017; Yousefi et al., 2017; Caglayan et al., 2013). Usage of transfer learning and CNN in visual recognition tasks gives better results and we can observe this from our results. If a comparison is made between deep approaches, transfer learning with a pre-trained model methods are more successful than end-to-end CNN approaches. The reason of this is need of large data to obtain an accurate CNN model. Most of the related studies are tested with one or two datasets. We applied different datasets with different sizes to observe effectiveness of different transfer and deep learning approaches. In traditional methods color, shape and textural

**Table 1**
Related studies summary.

|  | Datasets | Best results | Method |
|---|---|---|---|
| Our Study | Flavia | 99.00 | DF - VGG16/LDA |
|  | Swedish | 98.80 | CNN - RNN |
|  | UCI Leaf | 96.20 | DF - Alexnet/LDA |
|  | Plantvillage | 99.80 | FT - VGG16 |
| Lee et al. (2017) | Flavia | 99.40 | CNN, Fine-Tuning |
|  | MK | 97.47 |  |
| Jeon and Rhee (2017) | Flavia | 99.60 | CNN |
| Caglayan et al. (2013) | Flavia | 96.30 | Hand Crafted Shape & Color Features |
| Yalcin and Razavi (2016) | TARBIL | 97.47 | CNN, SVM |
| Murat et al. (2017) | Flavia | 95.25 | HOG, Moments, ANN, RF, SVM |
|  | Swedish | 99.89 |  |
| Yousefi et al. (2017) | Flavia | 97.50 | Fourier & Wavelet Descriptors, MLP |

features are extracted manually and combined afterwards. By using a CNN all these features are extracted and weighted. A more compact and descriptive feature set is obtained. All of these datasets consist of cropped leaf images. So all of these methods and ours are unable to work with wild (unedited and segmented) leaf, plant images. This problem can be solved with a two layer approach. Detection can be made in first layer with an R-CNN model. Then a CNN or other kind of classifier can be applied for specie recognition.

## 3. Transfer learning

Machine learning models are mostly built to work in isolation, which need to be rebuilt when the features and data change. However, very often the previously acquired knowledge in machine learning can be applied for similar tasks. Instead of rebuilding the models which usually requires lots of effort, transfer learning aims to reuse the model and acquired knowledge, and likewise to decrease the model development time dramatically, and improve the model performance of the isolated learning model. Transfer learning has been used in many applications such as software defect prediction (Nam et al., 2017), sentiment classification (Wang and Mahadevan, 2011), and activity recognition (Cook et al., 2013).

### 3.1. Transfer learning approaches based on domains

Pan and Yang (2010) published a review paper on transfer learning, introducing four transfer learning approaches, including:

1. Instance-transfer: Re-weighting the labeled data for the target domain
2. Feature-representation-transfer: Selecting a good feature set to reduce the difference between two domains
3. Parameter transfer: Discovering parameters in one domain and re-using these parameters in the target domain
4. Relational-knowledge-transfer: Mapping of knowledge between two domains

With respect to the nature of the target task Pan and Yang (2010) distinguish the following three settings:

1. Inductive transfer learning: target task is different than the source one.
2. Transductive transfer learning: target task and source task are the same, but the domains are different.
3. Unsupervised transfer learning: target task is different than the source task, but they are related to each other.

### 3.2. Transfer learning approaches based on feature spaces

Weiss et al. (2016) published a more recent survey paper on transfer learning in which they indicate that over 700 papers were published after 2010 on transfer learning. They categorize the recent existing approaches into three categories:

1. Homogeneous transfer learning: Transferring knowledge across similar feature spaces. For instance, in software defect classification problem, when the source and target datasets consist of the same set of software metrics, homogeneous transfer learning can be applied in this case because features of the two domains are the same.
   (a) Instance-based: In this category of techniques, instances are re-weighted to reduce the effect of misleading source data.
   (b) Feature-based (symmetric or asymmetric): In symmetric feature-based approach, features of source and target data are transformed into a common space. In asymmetric feature-based approach, features are transformed into the target domain.
   (c) Parameter-based: Target model parameters are learned based on the source model.
   (d) Relational-based: Source and target domains are investigated through a relational pattern.
   (e) Hybrid-based: The above-mentioned approaches might be integrated to build a hybrid approach such as the integration of instance and parameter-based techniques.
2. Heterogeneous transfer learning: Transferring knowledge across different feature spaces. For instance, in software defect classification problem, when software metrics of the source domain are different than the software metrics of the target domain, heterogeneous transfer learning approaches are needed. For this purpose, Nam et al. (2017) used Kolmogorov-Smirnov test to match the features from the source to the features of the target domain based on the closeness of the distribution between two domains. A recent survey paper on heterogeneous transfer learning investigates 38 methods and discusses those approaches in two categories as follows Day and Khoshgoftaar (2017):
   (a) Symmetric transformation: Source and target are transformed into a common feature set called domain-invariant feature subspace.
   (b) Asymmetric transformation: Features of the source are mapped into the target feature space.
3. Negative transfer: If the information from the source domain impacts the target learner adversely, that portion of the dataset should not be used for the target domain. Negative transfer techniques help to select the best information.

### 3.3. Transfer learning from pre-trained models

For deep neural networks, in some cases there may not be enough data to train the network or creating the labeled data might be expensive. Hence, transfer learning can be applied to adopt the knowledge that has been learned in earlier settings. For example, there are various CNN models such as AlexNet (Krizhevsky et al., 2012), GoogleNet (Szegedy et al., 2015), and VGG (Simonyan and Zisserman, 2014), which can be used later on for similar tasks.

The two common transfer learning strategies in deep learning are deep feature extraction and fine-tuning. In the deep feature extraction, the input data is provided to the pre-trained network and activation values of various layers are stored and used as features. In fine-tuning, deep neural network is trained for a similar task, in which labeling is relatively easier. While the first layers of the pre-trained network can be fixed, fine-tuning can be done on the final layers of the model to learn the properties of the new dataset. The pre-trained model is re-trained with the new small dataset and weight values of the model are updated

according to a new task. Fine-tuning process occurs on the network using back-propagation with labels.

Learning to transmit is often faster than training a new neural network because all the parameters in the new network are not estimated from scratch. In the lower layers of the network, more general features exist such as color blobs and Gabor filters and they can be transferred to other tasks as well. However, in higher layers, features are more task-specific. Deep learning systems provide high performance for several problems, but they require huge amount of data and time for their training. In this case, reusing these pre-trained models for similar tasks is quite helpful.

## 4. Methodology

In this article, we address the two common strategies (deep feature extraction and fine-tuning) of transfer learning for deep learning based plant classification models. In our scenarios, the visual recognition is the domain for the source and the target. Therefore, our experiments apply homogeneous transfer learning methods based on the categorization of Weiss et al. (2016). Based on the Pan and Yang (2010) survey paper, our setting is inductive transfer learning because source and target tasks are different. While the target task is the identification of plants, the source task is the identification of different objects.

We present five Deep Neural Networks for the plant classification problem. These models are shown in Fig. 1. First we develop a CNN from scratch. Subsequently, we show the experimentation for the four different transfer learning models which are based on fine tuning and deep feature extraction. These include fine tuning, the cross-dataset fine-tuning which is applied over the trained CNN and the third approach is the fine-tuning of CNN models on leaf datasets. In the fourth approach, pre-trained CNN models are used for feature extraction and classical machine learning algorithms with the extracted deep features are applied for the classification. The fifth approach is the combination of RNN and CNN algorithms, which is explained in later sub sections.

With respect to the classification in the previous section the Fine Tuning, and Cross-Dataset Fine Tuning models are parameter-based transfer learning approaches. The Deep Feature Learning, and CNN-RNN Classification are feature-based transfer learning approaches.
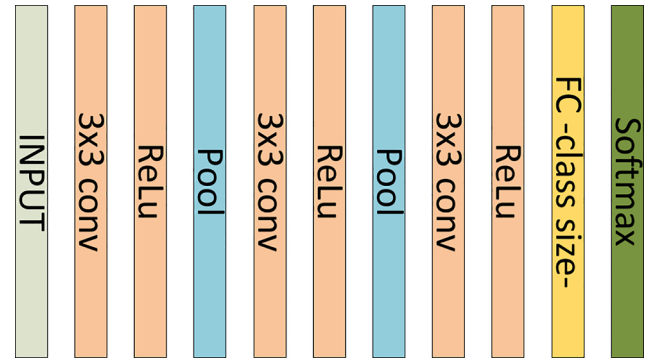


**Fig. 2.** The architecture of the proposed CNN model.

### 4.1. End-to-end CNN model

An end-to-end CNN model is trained on benchmark datasets. The architecture of the designed CNN model is given in Fig. 2. There are 15 layers in this CNN model. The first layer is the image input layer. Images with the size of $100 \times 100$ are provided as input. The leaf images with different width and height are re-sized before given to the CNN algorithm. There are three convolution layers in our network. Convolution layers are the main layers of a CNN. In these layers, there are filters to learn different feature types. Each filter is slid over the input images and the convolution is applied. The computed results of the convolution operations are mapped as the output. The following layers after a convolution layer are batch normalization and ReLu layers. Batch normalization layers are used for adjusting and normalizing the activation values computed by the previous layers. ReLu layers perform the threshold operation on the input to eliminate the effect of dark and noisy regions. The duty of max-pooling layers is to reduce the input dimensions to lower the computational complexity. This process is done by applying a mathematical MAX operation over the values, which corresponds to the filter. The fully connected layers are the final layers of a CNN model. These layers can be considered as the layers of standard neural networks. The class values for a given input are computed in these layers. Activation values of these layers
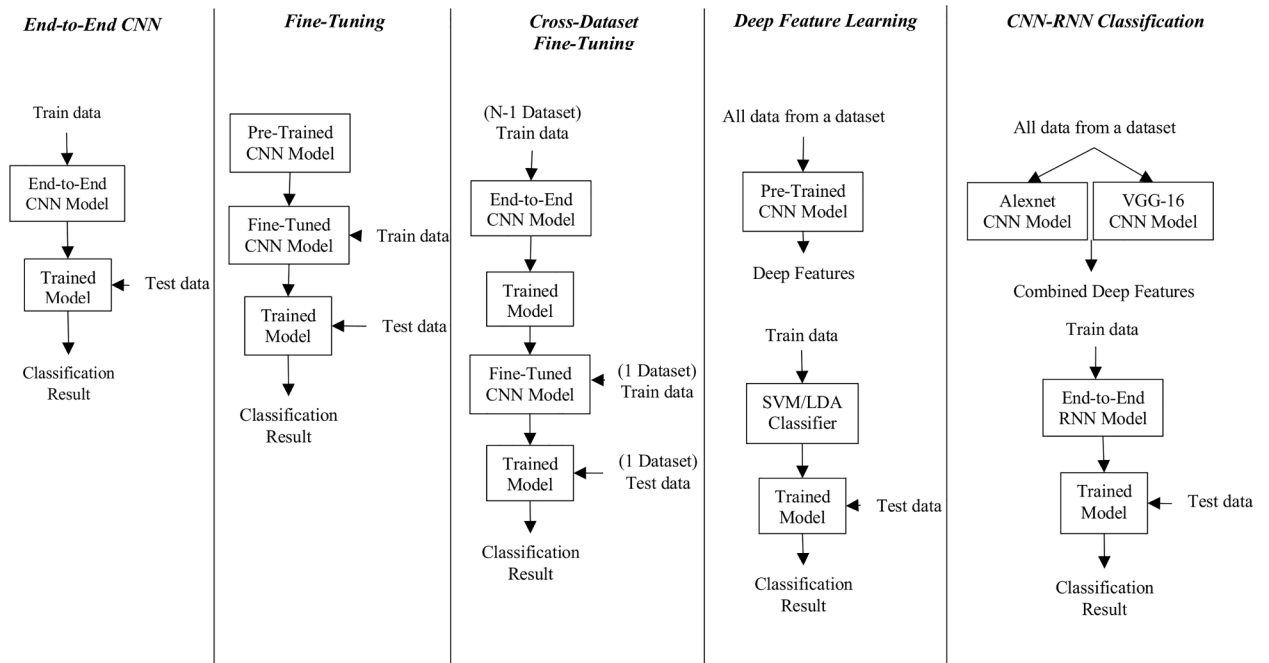


**Fig. 1.** General schema for experimental studies.

correspond to a different layer of abstractions. Top end layers are the softmax and classification layers. Softmax layer applies softmax function and the classification layer selects the label with the maximum possibility as its output.

### 4.2. Cross-dataset fine tuning

In this model, the CNN model built from scratch is trained with a combination of datasets. During this process, one of the datasets is excluded from the training sets and used for testing. All the datasets are used in testing once. Fine tuning is a widely used concept in transfer learning. In transfer learning, a model trained for a purposed task is used for a second task. In this approach, we train CNN models by using three of our four datasets and then, fine-tune the trained model with the excluded dataset. The fine-tuning operation is made by removing the last three layers and replacing with the new ones. These layers are fully connected, softmax, and classification layers. The number of outputs of the fully-connected layer should be equal to the number of the class count in the training dataset. The softmax and the classification layers are the same as the previous ones. The CNN model with new layers is trained and updated according to the new dataset.

### 4.3. Fine tuning

If there is not a huge amount of training data in a problem domain, training a Convolutional Neural Network from scratch is not efficient. For this kind of cases, the common method is to use a pre-trained model obtained from a large dataset for similar problems (Shin et al., 2016; Shie et al., 2015). This time we apply fine-tuning for VGG-16 and AlexNet models trained with a large scale dataset. These models are well-known CNN models trained with ImageNet image database. The architectures of these models are given in Figs. 3 and 4. With the use of these models, we transfer the knowledge obtained from a large dataset to our own problem domain. The fine-tuning approach is similar to the method applied in cross-dataset fine tuning. Both VGG16 and AlexNet have an output of 1000 classes. Last three layers of these models are removed and replaced with the layers suitable for leaf datasets. The last fully connected, softmax, and classification layers are removed. A new fully connected layer that has the same number of outputs with the number of classes in the new training dataset is placed. Options for the new fully connected layer are set according to the new data. Finally, the new network structure is trained with the new training datasets that contain leaf images. AlexNet and VGG-16 are fine-tuned with datasets separately.

### 4.4. Deep feature extraction

Another approach is extracting deep features from pre-trained models. In this approach, input images are given to VGG-16 and AlexNet directly. Then, activation values of the last fully connected layers of these pre-trained CNNs are obtained (fully connected layer 8 (Fc8)). Input layers of the AlexNet and VGG-16 are fixed and therefore, all the input data is re-sized to $227 \times 227$ for AlexNet and $224 \times 224$ for VGG-16. 1000 features are gathered from the Fc8 layer. After the
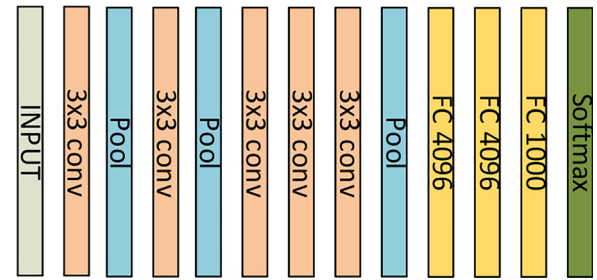


**Fig. 4.** Architecture of the AlexNet network.

deep feature extraction, classical machine learning algorithms are applied for the classification model development. Linear Discriminant Analysis (LDA) (Burks et al., 2000) and linear kernel SVM (Priya et al., 2012) are used for the training of models because they were previously applied in this problem successfully.

### 4.5. CNN-RNN

The final approach is to extract deep features from the pre-trained models and combine them with an RNN (Recurrent Neural Network) model. Although RNNs perform well for the sequential data, there are many studies which show that they can also be used for the non-sequential data (Karpathy, 2015). In this approach, input images are given to VGG-16 and AlexNet directly. The images are re-sized to fit to the input layers of the pre-trained models. The activation values of the last fully connected layers of these pre-trained CNNs are obtained (fully connected layer 8 (Fc8)) as in the previous approach. 1000 features are gathered from these layers. After the deep feature extraction, deep features from different CNNs are concatenated and reshaped as a matrix with dimensions $40 \times 50$. These matrices are given to an RNN model as input. The RNN model used in this step has an input layer with the size 20, an LSTM layer that contain 1000 hidden units, a fully connected layer that has the number of outputs as the number of classes, and a softmax layer.

## 5. Experimental results

In this section, we first explain the datasets and then, provide our experimental results and the statistical analysis.

### 5.1. Datasets

Four publicly available plant datasets are used during our experiments. We first performed our experiments on Flavia and Swedish Leaf datasets because they mostly contain clean images and no variations of luminance. Also, they are widely used datasets in this domain and publicly available. Later, we added the UCI Leaf dataset, which is hosted in UCI Machine Learning Repository because it is mostly applied by researchers for the comparison of algorithms. We continued to our experiments with Plantvillage dataset because it includes 50 times more samples compared to the samples in Flavia and Swedish Leaf and we
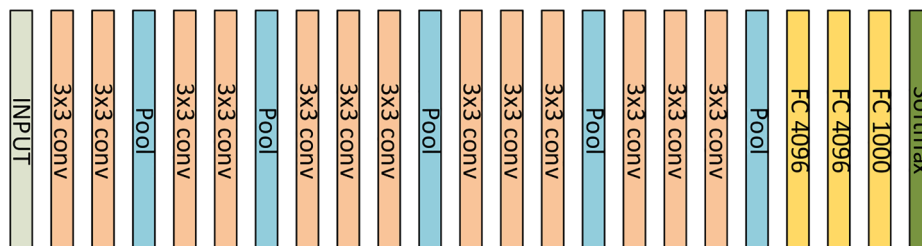


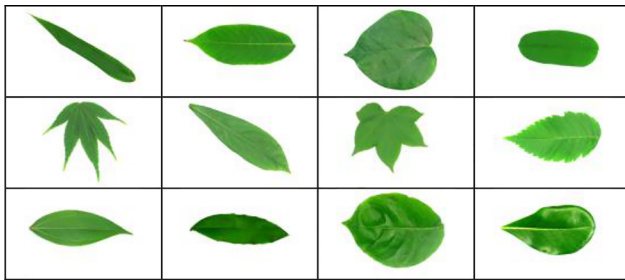**Fig. 3.** The architecture of VGG16 network.
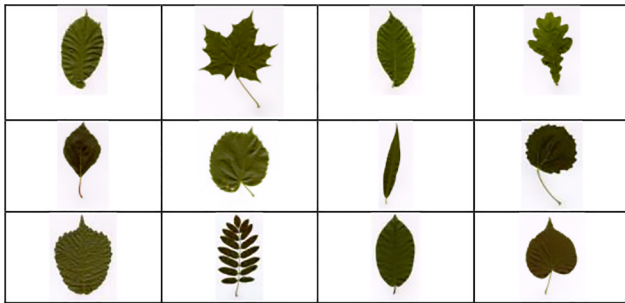
**Fig. 5.** Sample images from Flavia Dataset.



**Fig. 6.** Sample images from Swedish Leaf Dataset.



**Fig. 8.** Sample images from Plantvillage Dataset.

**Table 2**
Properties of the experimented datasets.

| Dataset | # of Classes | # of Samples | Image background | Color |
|---|---|---|---|---|
| Flavia | 32 | 1907 | White | RGB |
| Swedish Leaf | 15 | 1125 | White | RGB |
| UCI Leaf | 40 | 443 | Pink | RGB |
| Plantvillage | 14*(38**) | 54,306 | Transparent | RGB |

\* Healthy.
\*\* Healthy + diseased.

aimed to analyze our models on this large dataset. However, our models are not dependent on these datasets and can be applied on different plant datasets.

The first one is the Flavia Dataset created by Wu et al. (2007), which contains 32 species and 1900 images. Sample images are presented in Fig. 5. The second one is the Swedish Leaf Dataset generated by Söderkvist (2001). There are leaf images which belong to 15 tree classes. Sample images are presented in Fig. 6. The third one is the UCI Leaf Dataset built by Silva et al. (2013). There are 40 different species of plants included in this dataset with a total number of 443 images. Sample images are presented in Fig. 7. The fourth dataset is PlantVillage dataset created by Mohanty et al. (2016). There are 14 different species of plants and 38 classes (healthy-diseased), having a total number of 54,306 images. The healthy plant images are included during the experiments. Sample images are presented in Fig. 8. The properties of the datasets are presented in Table 2.

*5.2. Results*

All the deep learning based models have 100 epochs during the training phase. Datasets are divided into test and training sets with a ratio of 30%, and 70% respectively. During deep feature learning experiments, linear kernel SVM and LDA classifiers are used with fivefold (k = 5) cross-validation. As explained in the previous sections,
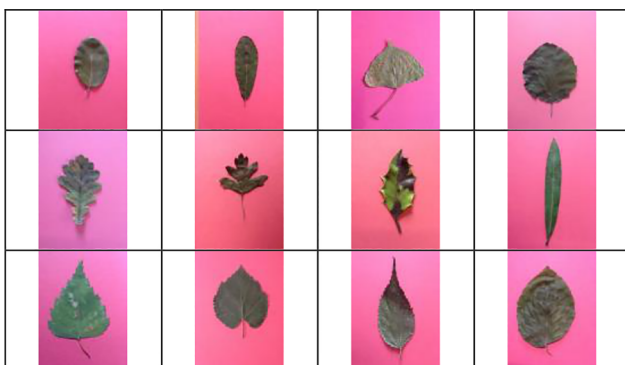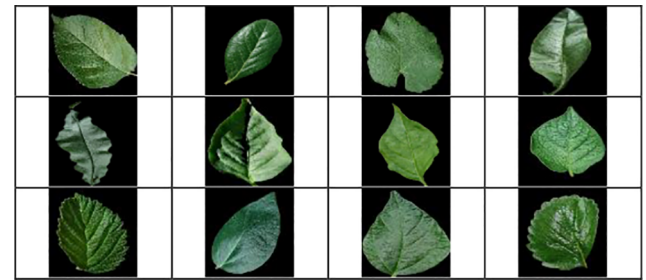
experiments were performed on several datasets with different sizes. Some of the datasets have a smaller number of samples. The use of a large k value would cause to have a limited number of samples in the test set for the classes with a small number of samples. Increasing the number of k would probably reduce the overfitting, but this time computational cost would increase. To cope with the trade-off between the accuracy and the performance, we set the k value to 5.

Since each of these approaches were discussed in the previous section, we provide only a figure to reflect the differences between different configurations.

Results of the end-to-end CNN model is provided in Table 3. According to this table, it is observed that the same CNN model performs better on the datasets with higher sample amount, and this observation is consistent with the general idea of deep learning models, which state that the models trained with huge amount of data provide better classification outcomes. UCI Leaf dataset has the smallest size and therefore, the classification accuracy on this dataset is not as good as the performance of the model on the other datasets.

The cross-dataset fine-tuning experiment results are presented in Table 4. UCI Leaf Dataset result is improved by 4%, however, the performance on other datasets is similar to the end-to-end CNN results. This experiment demonstrated that increasing the data size for training positively affects the performance of the model on datasets which have lower sample size.

In Table 5, results of the fine-tuning of the pre-trained models are presented. The classification performance of all the datasets improves compared to the previous approaches. Pre-trained models are trained with millions of images from the ImageNet dataset. Even if models are trained with images from different domains, fine-tuning of these models provides better results than the end-to-end models on especially datasets with limited elements. According to this experimental analysis, it is



**Fig. 7.** Sample images from UCI Leaf Dataset.

**Table 3**
Classification accuracy of the end-to-end CNN model.

| Dataset | CA |
|---|---|
| Flavia | 91.08 |
| SwedishLeaf | 96.06 |
| UCI Leaf | 76.15 |
| PlantVillage | 97.40 |

**Table 4**
Classification accuracy of the cross-dataset fine-tuning of the CNN model.

| Test dataset | Training datasets | CA |
|---|---|---|
| Flavia | Swedish Leaf UCI Leaf Plantvillage | 91.43 |
| Swedish Leaf | Flavia UCI Leaf Plantvillage | 96.06 |
| UCI Leaf | Flavia Swedish Leaf Plantvillage | 80.60 |
| Plantvillage | Flavia Swedish Leaf UCI Leaf | 96.93 |

**Table 5**
Classification accuracy of the fine-tuning of pre-trained models.

| Dataset | Pre-trained model | CA |
|---|---|---|
| Flavia | Alexnet | 97.89 |
|  | VGG16 | 98.16 |
| Swedish Leaf | Alexnet | 95.56 |
|  | VGG16 | 99.11 |
| UCI Leaf | Alexnet | 89.41 |
|  | VGG16 | 90.56 |
| Plantvillage | Alexnet | 98.60 |
|  | VGG16 | 99.80 |

observed that the VGG16 network provides relatively better performance than the Alexnet, but the difference between these two approaches is not very significant.

Experimental results of classification with deep features are provided in Table 6. LDA and SVM methods perform similarly on all the datasets except the UCI Leaf dataset. The best classification outcomes are obtained from these experiments, especially for the UCI dataset. In the first experiment, the performance of CNN was around 76% on the UCI leaf dataset, but the performance of the model with deep features is beyond this performance value.

**Table 6**
Classification accuracy of classifiers with deep features.

| Dataset | PT-Model/Classifier | CA |
|---|---|---|
| Flavia | Alexnet/LDA | 99.00 |
|  | Alexnet/SVM | 97.50 |
|  | VGG16/LDA | 99.10 |
|  | VGG16/SVM | 97.70 |
| Swedish Leaf | Alexnet/LDA | 95.80 |
|  | Alexnet/SVM | 97.80 |
|  | VGG16/LDA | 96.10 |
|  | VGG16/SVM | 98.80 |
| UCI Leaf | Alexnet/LDA | 96.20 |
|  | Alexnet/SVM | 88.90 |
|  | VGG16/LDA | 94.80 |
|  | VGG16/SVM | 89.60 |
| Plantvillage | Alexnet/LDA | 98.70 |
|  | Alexnet/SVM | 97.80 |
|  | VGG16/LDA | 98.70 |
|  | VGG16/SVM | 98.00 |

**Table 7**
Classification accuracy of CNN-RNN models on datasets.

| Dataset | CA |
|---|---|
| Flavia | 92.65 |
| Swedish Leaf | 99.11 |
| UCI Leaf | 70.79 |
| Plantvillage | 98.77 |

Results of the classification when the CNN-RNN combination is applied are shown in Table 7. According to this table, the same classification model performed better on datasets which have higher sample amount as in the results of end-to-end CNN experiments. The model trained with the UCI Leaf dataset provides the lowest classification accuracy (70.79%). The performance of the models developed here is comparable with the performance of our other experiments.

In Fig. 9, all the results are presented with a bar graph. In Tables 8–11, the statistical significance analysis is performed for the analyzed methods which are represented as follows: M1: End-to-end CNN, M2: cross-dataset, M3: FT - Alexnet, M4: FT - VGG16, M5: DF - Alexnet/LDA, M6: DF - Alexnet/SVM, M7: DF - VGG16/LDA, M8: DF - VGG16/SVM, M9: CNN - RNN. McNemar's test is used to measure the statistical significance of the methods according to classification outcomes. All the classes are tested one-versus-all manner (one taken as positive, others are considered as negative class). Significant values (p < 0.05) are implied with boldface font.

From Tables 4–11, we demonstrated that the transfer learning approaches perform better than the end-to-end models for all the benchmarking datasets. This difference is very clear for the UCI Leaf dataset which has a smaller sample size compared to the other datasets. VGG16/LDA methods overall performance is the best one. For Flavia, Plantvillage, and UCI Leaf datasets, VGG16/LDA provide relatively higher performance. The performance is lower in Swedish Leaf dataset, but this method can be considered as the most successful method on these datasets. The end-to-end CNN model provides the worst performance for all the datasets. For Flavia, Plantvillage, and UCI Leaf, all the methods provide better results compared to the performance of end-to-end CNN and the cross-dataset methods.

The Plantvillage is the dataset which has the highest number of samples. The difference between deep learning based methods trained on this dataset is marginal. FT-VGG16 is the best performing one in this dataset, and all comparisons with this method are statistically significant, thus it is the most preferable method.

The similar situation is also valid for the Swedish Leaf dataset. The difference and the statistical significance between methods are lower in these datasets compared to the other ones. When compared to the End-to-End CNN models, nearly all of the methods provided statistically significant results in all benchmark datasets. Considering the classification performance of this approach, its preference is low compared to other methods.

When the all comparisons are considered, the least statistical significance is seen in the comparisons with the FT-Alexnet method. This method provided close but not the best results in many of the datasets, except UCI Leaf.

## 6. Discussion

The related work section showed that the effect of transfer learning on the performance of deep learning models has not been investigated. We comparatively assessed the performance of five transfer learning scenarios on the deep learning models. We presented our Research Questions (RQs) in Section 1 and here, we discuss our research findings. To repeat the research questions were the following:
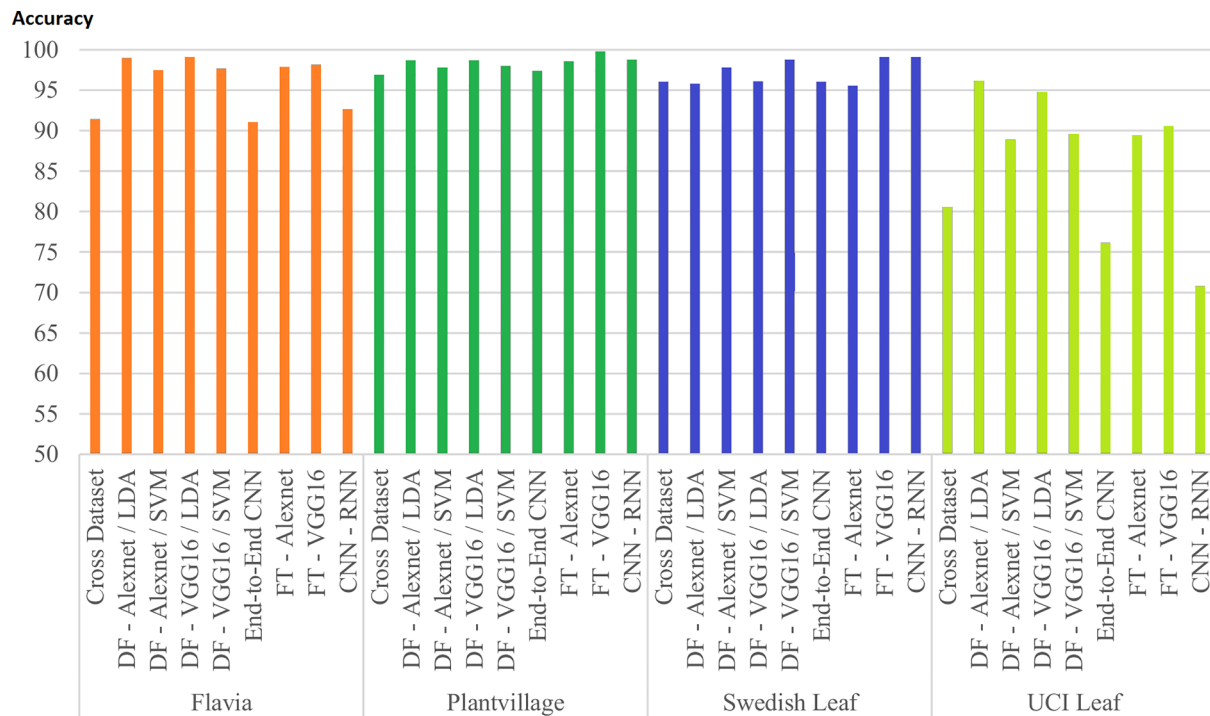
**Fig. 9.** Graph of experimental results.

**Table 8**
Statistical significance test results on Flavia dataset. The significant values are implied by boldface.

|     | M2   | M3       | M4       | M5       | M6       | M7       | M8       | M9       |
|-----|------|----------|----------|----------|----------|----------|----------|----------|
| M1  | 0.75 | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | 0.14     |
| M2  | –    | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | 0.25     |
| M3  | –    | –        | 0.62     | **0.02** | 0.51     | **0.01** | 0.74     | **0.00** |
| M4  | –    | –        | –        | 0.07     | 0.25     | **0.04** | 0.41     | **0.00** |
| M5  | –    | –        | –        | –        | **0.00** | 0.79     | **0.01** | **0.00** |
| M6  | –    | –        | –        | –        | –        | **0.00** | 0.74     | **0.00** |
| M7  | –    | –        | –        | –        | –        | –        | **0.00** | **0.00** |
| M8  | –    | –        | –        | –        | –        | –        | –        | **0.00** |

**Table 10**
Statistical significance test results on the Swedish Leaf dataset. The significant values are implied by boldface.

|     | M2   | M3   | M4       | M5   | M6       | M7   | M8       | M9       |
|-----|------|------|----------|------|----------|------|----------|----------|
| M1  | 1.00 | 0.52 | **0.00** | 0.74 | **0.01** | 0.96 | **0.00** | **0.00** |
| M2  | –    | 0.52 | **0.00** | 0.74 | **0.01** | 0.96 | **0.00** | **0.00** |
| M3  | –    | –    | **0.00** | 0.76 | **0.00** | 0.49 | **0.00** | **0.00** |
| M4  | –    | –    | –        | **0.00** | **0.01** | **0.00** | 0.44 | 1.00 |
| M5  | –    | –    | –        | –    | **0.00** | 0.70 | **0.00** | **0.00** |
| M6  | –    | –    | –        | –    | –        | **0.01** | **0.05** | **0.01** |
| M7  | –    | –    | –        | –    | –        | –    | **0.00** | **0.00** |
| M8  | –    | –    | –        | –    | –        | –    | –        | 0.44     |

**Table 9**
Statistical significance test results on Plantvillage dataset. The significant values are implied by boldface.

|     | M2   | M3       | M4       | M5       | M6       | M7       | M8       | M9       |
|-----|------|----------|----------|----------|----------|----------|----------|----------|
| M1  | 0.47 | **0.03** | **0.00** | **0.02** | 0.50     | **0.02** | 0.31     | **0.01** |
| M2  | –    | **0.00** | **0.00** | **0.00** | 0.17     | **0.00** | 0.08     | **0.00** |
| M3  | –    | –        | **0.00** | 0.82     | 0.12     | 0.82     | 0.24     | 0.70     |
| M4  | –    | –        | –        | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** |
| M5  | –    | –        | –        | –        | 0.08     | 1.00     | 0.16     | 0.87     |
| M6  | –    | –        | –        | –        | –        | 0.08     | 0.72     | 0.06     |
| M7  | –    | –        | –        | –        | –        | –        | 0.16     | 0.87     |
| M8  | –    | –        | –        | –        | –        | –        | –        | 0.12     |

**Table 11**
Statistical significance test results on the UCI Leaf dataset. The significant values are implied by boldface.

|     | M2       | M3       | M4       | M5       | M6       | M7       | M8       | M9       |
|-----|----------|----------|----------|----------|----------|----------|----------|----------|
| M1  | **0.01** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** |
| M2  | –        | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** | **0.00** |
| M3  | –        | –        | 0.33     | **0.00** | 0.68     | **0.00** | 0.87     | **0.00** |
| M4  | –        | –        | –        | **0.00** | 0.16     | **0.00** | 0.41     | **0.00** |
| M5  | –        | –        | –        | –        | **0.00** | 0.08     | **0.00** | **0.00** |
| M6  | –        | –        | –        | –        | –        | **0.00** | 0.56     | **0.00** |
| M7  | –        | –        | –        | –        | –        | –        | **0.00** | **0.00** |
| M8  | –        | –        | –        | –        | –        | –        | –        | **0.00** |

- RQ1: To what extent can plant classification benefit from transfer learning in deep learning?
- RQ-2: How do the different transfer learning scenarios perform at improving the performance of plant classification models using deep learning?

To answer these questions we have compared four different transfer learning approaches with the base, end-to-end CNN approach. For our comparison we mainly focused on classification accuracy (CA). Our study showed that transfer learning does have an impact on the classification accuracy. We have shown that transfer learning approaches perform better than the end-to-end models for the benchmarking datasets. In particular, a clear performance benefit can be observed for datasets which have less number of data points.

Experimental studies consist of potential threats to validity (Claes et al., 2000). Regarding the internal validity, we did not apply just one transfer learning approach to evaluate the impact of transfer learning on deep learning models. Instead, we designed and implemented four different transfer learning scenarios for comparative assessment of models. While we covered several transfer learning scenarios during

our experiments, new studies might apply other approaches with different settings and reach new results. We focused on deep learning models instead of traditional machine learning models because we aimed to benefit pre-trained deep learning-based plant classification models.

Regarding the external validity, our conclusions are valid for the datasets explained in Datasets subsection and observations might be different on a new set of datasets. We preferred the datasets which follow FAIR (Findable/Accessible/Interoperable/Reusable) principles (Wilkinson et al., 2016) and therefore, results might be different on datasets which do not adopt FAIR principles.

Regarding the construct validity, we did our experiments on widely-used datasets which are still applied by other researchers in this field. Conclusion validity addresses threats which impact the ability to conclude appropriately. We split the datasets into training and test sets by using the widely-used ratio (70–30%) and fivefold cross-validation was applied during deep feature learning experiments with linear kernel SVM and LDA classifiers. Also, 100 epochs were used for all the deep learning based models. These validation techniques were applied to avoid the randomness in data. Apart from this techniques, a statistical analysis was also performed to check the statistical significance of our experimental results. Results were reported based on accuracy parameter. Four public datasets were analyzed during our experiments and further experiments are required for new datasets.

## 7. Conclusion

Correct classification of plant species has many advantages not only in agriculture, but also in several other domains such as, for example biodiversity, health and forest studies. Instead of manually processing of plants by experts, automated plant identification systems enable stakeholders to quickly deal with the huge amount of plants and lessen the required time and cost of these operations.

While there are some studies on the use of deep learning algorithms for the plant classification, there has not been an in-depth study which applies several transfer learning scenarios for deep learning models. Hence our study can be considered as complementary to the existing studies paving the way for further research on transfer learning for plant classification.

This study demonstrated that the transfer learning improves the performance of deep learning models and especially, models which apply deep features and use fine-tuning provide better performance compared to the other transfer learning strategies. This result implies that instead of only applying an end-to-end CNN model for the plant classification, the other transfer learning approaches must also be considered in the low accuracy performance case.

In the near future, we are planning to investigate the performance of these models for different applications in agriculture such as plant disease detection and weeds detection. Another dimension is to perform new experiments when more public datasets become available.

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## Appendix A. Supplementary material

Supplementary data associated with this article can be found, in the online version, at https://doi.org/10.1016/j.compag.2019.01.041.

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