

# Learning and Evolution in Artificial Neural Networks: A Comparison Study

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**Abstract.** This paper aims at learning and evolution in artificial neural networks. Here is presented a system evolving populations of neural nets that are fully connected multilayer feedforward networks with fixed architecture solving given tasks. The system is compared with gradient descent weight training (like backpropagation) and with hybrid neural network adaptation. All neural networks have the same architecture and solve the same problems to be able to be compared mutually. In order to test the efficiency of described algorithms, we applied them to the *Fisher's Iris data set* [1] that is the bench test database from the area of machine learning.

## 1 Learning in Artificial Neural Networks

Learning in artificial neural networks is typically accomplished using examples. This is also called training in artificial neural networks because the learning is achieved by adjusting the connection weights in artificial neural networks iteratively so that trained (or learned) artificial neural networks can perform certain tasks. The essence of a learning algorithm is the learning rule, i.e. a weight-updating rule, which determines how connection weights are changed. Examples of popular learning rules include the delta rule, the Hebbian rule, the anti-Hebbian rule, and the competitive learning rule, etc. More detailed discussion of artificial neural networks can be found in [2]. Learning in artificial neural networks can roughly be divided into supervised, unsupervised, and reinforcement learning. Without commonness, we are going to target multilayer feedforward neural networks that are adapted with backpropagation algorithm.

Supervised learning is based on direct comparison between the actual output of an artificial neural network and the desired correct output, also known as the target output. It is often formulated as the minimization of an error function such as the total mean square error between the actual output and the desired output summed over all available data. A gradient descent-based optimization algorithm such as backpropagation [2] can then be used to adjust connection weights in the artificial neural network iteratively in order to minimize the error. There have been some successful applications of backpropagation in various areas [3]–[5], but backpropagation has drawbacks due to its use of gradient descent. It often gets trapped in a local minimum of the error function and is incapable of finding a global minimum if the error function is multimodal and/or nondifferentiable.

## 2 NeuroEvolutionary Learning in Artificial Neural Networks

Evolutionary algorithms refer to a class of population-based stochastic search algorithms that are developed from ideas and principles of natural evolution. They include evolution strategies, evolutionary programming, genetic algorithms etc. [6]. One important feature of all these algorithms is their population based search strategy. Individuals in a population compete and exchange information with each other in order to perform certain tasks. Evolutionary algorithms are particularly useful for dealing with large complex problems, which generate many local optima. They are less likely to be trapped in local minima than traditional gradient-based search algorithms. They do not depend on gradient information and thus are quite suitable for problems where such information is unavailable or very costly to obtain or estimate. They can even deal with problems where no explicit and/or exact objective function is available. These features make them much more robust than many other search algorithms. There is a good introduction to various evolutionary algorithms for optimization in [6].

Evolution has been introduced into artificial neural networks at roughly three different levels [7]: connection weights; architectures; and learning rules. The evolution of connection weights introduces an adaptive and global approach to training, especially in the reinforcement learning and recurrent network-learning paradigm where gradient-based training algorithms often experience great difficulties. The evolution of architectures enables artificial neural networks to adapt their topologies to different tasks without human intervention and thus provides an approach to automatic artificial neural network design as both connection weights and structures can be evolved. The evolution of learning rules can be regarded as a process of learning to learn in artificial neural networks where the adaptation of learning rules is achieved through evolution. It can also be regarded as an adaptive process of automatic discovery of novel learning rules.

The evolutionary approach to weight training in artificial neural networks consists of two major phases. The first phase is to decide the representation of connection weights, i.e., whether in the form of binary strings or real strings. The second one is the evolutionary process simulated by an evolutionary algorithm, in which search operators such as crossover and mutation have to be decided in conjunction with the representation scheme. Different representations and search operators can lead to quite different training performance. In a binary representation scheme, each connection weight is represented by a number of bits with certain length. An artificial neural network is encoded by concatenation of all the connection weights of the network in the chromosome. The advantages of the binary representation lie in its simplicity and generality. It is straightforward to apply classical crossover (such as one-point or uniform crossover) and mutation to binary strings [6]. Real numbers have been proposed to represent connection weights directly, i.e. one real number per connection weight [6]. As connection weights are represented by real numbers, each individual in an evolving population is then a real vector. Traditional binary crossover and mutation can no longer be used directly. Special search operators have to be designed. Montana and Davis [8] defined a large number of tailored genetic operators, which incorporated many heuristics about training artificial neural networks. The idea was to retain useful feature detectors formed around hidden nodes during evolution.

One of the problems faced by evolutionary training of artificial neural networks is the permutation problem [7] also known as the competing convention problem. It is caused by the many-to-one mapping from the representation (genotype) to the actual artificial neural network (phenotype) since two artificial neural networks that order their hidden nodes differently in their chromosomes will still be equivalent functionally. In general, any permutation of the hidden nodes will produce functionally equivalent artificial neural networks with different chromosome representations. The permutation problem makes crossover operator very inefficient and ineffective in producing good offspring. The role of *crossover* has been controversial in neuroevolution as well as among the evolutionary computation community in general. However, there have been successful applications using crossover operations to evolve neural networks [9]. Compared with the *mutation* only system, the performance of the system using crossover operations is in general better and that it also helps to compress the overall size of search space faster.

### 3 Comparison between Evolutionary Training and Gradient-based Training

The evolutionary training approach is attractive because it can handle the global search problem better in a vast, complex, multimodal, and nondifferentiable surface. It does not depend on gradient information of the error (or fitness) function and thus is particularly appealing when this information is unavailable or very costly to obtain or estimate. For example, the same evolutionary algorithms can be used to train many different networks: recurrent artificial neural networks [10], higher order artificial neural networks [11], and fuzzy artificial neural networks [12]. The general applicability of the evolutionary approach saves a lot of human efforts in developing different training algorithms for different types of artificial neural networks. The evolutionary approach also makes it easier to generate artificial neural networks with some special characteristics. For example, the artificial neural networks complexity can be decreased and its generalization increased by including a complexity (regularization) term in the fitness function. Unlike the case in gradient-based training, this term does not need to be differentiable or even continuous. Weight sharing and weight decay can also be incorporated into the fitness function easily.

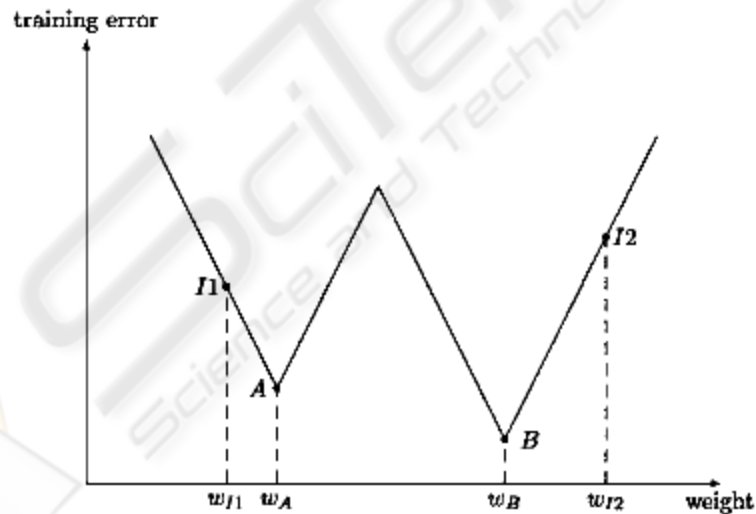
However, evolutionary algorithms are generally much less sensitive to initial conditions of training. They always search for a globally optimal solution, while a gradient descent algorithm can only find a local optimum in a neighbourhood of the initial solution.

### 4 Hybrid Training

Most evolutionary algorithms are rather inefficient in fine-tuned local search although they are good at global search. This is especially true for genetic algorithms. The efficiency of evolutionary training can be improved significantly by incorporating a local search procedure into the evolution, i.e. combining evolutionary algorithm's

global search ability with local search's ability to fine tune. Evolutionary algorithms can be used to locate a good region in the space and then a local search procedure is used to find a near-optimal solution in this region. The local search algorithm could be backpropagation [2] or other random search algorithms. Hybrid training has been used successfully in many application areas: Lee [13] and many others used genetic algorithms to search for a near-optimal set of initial connection weights and then used backpropagation to perform local search from these initial weights. Their results showed that the *hybrid algorithm* approach was more efficient than either the genetic algorithm or backpropagation algorithm used alone. If we consider that backpropagation often has to run several times in practice in order to find good connection weights due to its sensitivity to initial conditions, the hybrid training algorithm will be quite competitive. Similar work on the evolution of initial weights has also been done on competitive learning neural networks [14] and Kohonen networks [15].

It is interesting to consider finding good initial weights as locating a good region in the weight space. Defining that basin of attraction of a local minimum as being composed of all the points, sets of weights in this case, which can converge to the local minimum through a local search algorithm, then a global minimum can easily be found by the local search algorithm if an evolutionary algorithm can locate a point, i.e. a set of initial weights, in the basin of attraction of the global minimum. Fig. 1 illustrates a simple case where there is only one connection weight in the artificial neural networks. If an evolutionary algorithm can find an initial weight such as  $w_{I2}$ , it would be easy for a local search algorithm to arrive at the globally optimal weight  $w_B$  even though  $w_{I2}$  itself is not as good as  $w_{I1}$ .



**Fig. 1.** An illustration of using an evolutionary algorithm to find good initial weights such that a local search algorithm can find the globally optimal weights easily.  $w_{I2}$  in this figure is an optimal initial weight because it can lead to the global optimum  $w_B$  using a local search algorithm.

## 5 Experiments

In order to test the efficiency of described algorithms, we applied it to the Iris flower data set or Fisher's Iris data set is a multivariate data set introduced by Sir Ronald Aylmer Fisher as an example of discriminated analysis [1]. It is sometimes called *Anderson's Iris data* set because Edgar Anderson collected the data to quantify the geographic variation of Iris flowers in the Gaspé Peninsula. The dataset consists of 50 samples from each of three species of Iris flowers (*Iris setosa*, *Iris virginica* and *Iris versicolor*). Four features were measured from each sample, they are the length and the width of sepal and petal. Based on the combination of the four features, Fisher developed a linear discriminated model to determine which species they are (see Table 1). There are three-layer feedforward neural networks with architecture is 4 - 4 - 3 (e.g. four units in the input layer, four units in the hidden layer, and three units in the output layer) in all experimental works, because the *Fisher's Iris data set* [1] is not linearly separable and therefore we cannot use neural network without hidden units. All nets are fully connected. The input values of the training set from the Table 1 are transformed into interval  $<0; 1>$  to be use backpropagation algorithms for adaptation.

**Weight Evolution in Artificial Neural Networks:** the initial population contains 30 individuals (weight representations of three-layer neural networks). There are connection weights represented by real numbers in each chromosome. We use the genetic algorithm with the following parameters: probability of mutation is 0,01 and probability of crossover is 0,5. Adaptation of each neural network starts with randomly generated weight values in the initial population.

**The Gradient Descent** adaptation deals through backpropagation with the following parameters: learning rate is 0.4, momentum is 0.

**The Hybrid Training** combines parameters from genetic algorithms and backpropagation. It makes one backpropagation epoch with probability 0,5 in each generation.

## 6 Conclusions

History of error functions is shown in the Figure 2. There are shown average values of error functions in the given population. The "*gradient descent adaptation*" represents an adaptation with the backpropagation. There are shown average values of error functions, because the adaptation with backpropagation algorithm was applied 10 times for each calculation.

All networks solve *Fisher's Iris data set* [1] in our experiment. Now we can compare results from all experiments, e.g. weight evolution, gradient descent adaptation, and hybrid training. Other numerical simulations give very similar results. If we can see from Figure 2, the hybrid training shows the best results from all of them. In general, hybrid algorithms tend to perform better than others for a large number of problems, because they combine evolutionary algorithm's global search ability with local search's ability to fine tune

**Table 1.** The set of patterns (the Fisher's Iris Data training set), where **Se** means *setosa*, **Vi** means *virginica*, and **Ve** means *versicolor*.

Sepal Length	Sepal Width	Petal Length	Petal Width	Species	Sepal Length	Sepal Width	Petal Length	Petal Width	Species	Sepal Length	Sepal Width	Petal Length	Petal Width	Species
5,1	3,5	1,4	0,2	Se	6,3	3,3	6	2,5	Vi	7	3,2	4,7	1,4	Ve
4,9	3	1,4	0,2	Se	5,8	2,7	5,1	1,9	Vi	6,4	3,2	4,5	1,5	Ve
4,7	3,2	1,3	0,2	Se	7,1	3	5,9	2,1	Vi	6,9	3,1	4,9	1,5	Ve
4,6	3,1	1,5	0,2	Se	6,3	2,9	5,6	1,8	Vi	5,5	2,3	4	1,3	Ve
5	3,6	1,4	0,2	Se	6,5	3	5,8	2,2	Vi	6,5	2,8	4,6	1,5	Ve
5,4	3,9	1,7	0,4	Se	7,6	3	6,6	2,1	Vi	5,7	2,8	4,5	1,3	Ve
4,6	3,4	1,4	0,3	Se	4,9	2,5	4,5	1,7	Vi	6,3	3,3	4,7	1,6	Ve
5	3,4	1,5	0,2	Se	7,3	2,9	6,3	1,8	Vi	4,9	2,4	3,3	1	Ve
4,4	2,9	1,4	0,2	Se	6,7	2,5	5,8	1,8	Vi	6,6	2,9	4,6	1,3	Ve
4,9	3,1	1,5	0,1	Se	7,2	3,6	6,1	2,5	Vi	5,2	2,7	3,9	1,4	Ve
5,4	3,7	1,5	0,2	Se	6,5	3,2	5,1	2	Vi	5	2	3,5	1	Ve
4,8	3,4	1,6	0,2	Se	6,4	2,7	5,3	1,9	Vi	5,9	3	4,2	1,5	Ve
4,8	3	1,4	0,1	Se	6,8	3	5,5	2,1	Vi	6	2,2	4	1	Ve
4,3	3	1,1	0,1	Se	5,7	2,5	5	2	Vi	6,1	2,9	4,7	1,4	Ve
5,8	4	1,2	0,2	Se	5,8	2,8	5,1	2,4	Vi	6,7	3,1	4,4	1,4	Ve
5,7	4,4	1,5	0,4	Se	6,4	3,2	5,3	2,3	Vi	5,6	2,9	3,6	1,3	Ve
5,4	3,9	1,3	0,4	Se	6,5	3	5,5	1,8	Vi	5,6	3	4,5	1,5	Ve
5,1	3,5	1,4	0,3	Se	7,7	3,8	6,7	2,2	Vi	5,8	2,7	4,1	1	Ve
5,7	3,8	1,7	0,3	Se	7,7	2,6	6,9	2,3	Vi	5,6	2,5	3,9	1,1	Ve
5,1	3,8	1,5	0,3	Se	6	2,2	5	1,5	Vi	6,2	2,2	4,5	1,5	Ve
5,4	3,4	1,7	0,2	Se	6,9	3,2	5,7	2,3	Vi	5,9	3,2	4,8	1,8	Ve
5,1	3,7	1,5	0,4	Se	5,6	2,8	4,9	2	Vi	6,1	2,8	4	1,3	Ve
4,6	3,6	1	0,2	Se	7,7	2,8	6,7	2	Vi	6,3	2,5	4,9	1,5	Ve
5,1	3,3	1,7	0,5	Se	6,3	2,7	4,9	1,8	Vi	6,1	2,8	4,7	1,2	Ve
4,8	3,4	1,9	0,2	Se	6,7	3,3	5,7	2,1	Vi	6,4	2,9	4,3	1,3	Ve
5	3	1,6	0,2	Se	7,2	3,2	6	1,8	Vi	6,6	3	4,4	1,4	Ve
5	3,4	1,6	0,4	Se	6,2	2,8	4,8	1,8	Vi	6,8	2,8	4,8	1,4	Ve
5,2	3,5	1,5	0,2	Se	6,1	3	4,9	1,8	Vi	6,7	3	5	1,7	Ve
5,2	3,4	1,4	0,2	Se	6,4	2,8	5,6	2,1	Vi	6	2,9	4,5	1,5	Ve
4,7	3,2	1,6	0,2	Se	7,2	3	5,8	1,6	Vi	5,7	2,6	3,5	1	Ve
4,8	3,1	1,6	0,2	Se	7,4	2,8	6,1	1,9	Vi	5,5	2,4	3,8	1,1	Ve
5,4	3,4	1,5	0,4	Se	7,9	3,8	6,4	2	Vi	5,5	2,4	3,7	1	Ve
5,2	4,1	1,5	0,1	Se	6,4	2,8	5,6	2,2	Vi	5,8	2,7	3,9	1,2	Ve
5,5	4,2	1,4	0,2	Se	6,3	2,8	5,1	1,5	Vi	6	2,7	5,1	1,6	Ve
4,9	3,1	1,5	0,2	Se	6,1	2,6	5,6	1,4	Vi	5,4	3	4,5	1,5	Ve
5	3,2	1,2	0,2	Se	7,7	3	6,1	2,3	Vi	6	3,4	4,5	1,6	Ve
5,5	3,5	1,3	0,2	Se	6,3	3,4	5,6	2,4	Vi	6,7	3,1	4,7	1,5	Ve
4,9	3,6	1,4	0,1	Se	6,4	3,1	5,5	1,8	Vi	6,3	2,3	4,4	1,3	Ve
4,4	3	1,3	0,2	Se	6	3	4,8	1,8	Vi	5,6	3	4,1	1,3	Ve
5,1	3,4	1,5	0,2	Se	6,9	3,1	5,4	2,1	Vi	5,5	2,5	4	1,3	Ve
5	3,5	1,3	0,3	Se	6,7	3,1	5,6	2,4	Vi	5,5	2,6	4,4	1,2	Ve
4,5	2,3	1,3	0,3	Se	6,9	3,1	5,1	2,3	Vi	6,1	3	4,6	1,4	Ve
4,4	3,2	1,3	0,2	Se	5,8	2,7	5,1	1,9	Vi	5,8	2,6	4	1,2	Ve
5	3,5	1,6	0,6	Se	6,8	3,2	5,9	2,3	Vi	5	2,3	3,3	1	Ve
5,1	3,8	1,9	0,4	Se	6,7	3,3	5,7	2,5	Vi	5,6	2,7	4,2	1,3	Ve
4,8	3	1,4	0,3	Se	6,7	3	5,2	2,3	Vi	5,7	3	4,2	1,2	Ve
5,1	3,8	1,6	0,2	Se	6,3	2,5	5	1,9	Vi	5,7	2,9	4,2	1,3	Ve
4,6	3,2	1,4	0,2	Se	6,5	3	5,2	2	Vi	6,2	2,9	4,3	1,3	Ve
5,3	3,7	1,5	0,2	Se	6,2	3,4	5,4	2,3	Vi	5,1	2,5	3	1,1	Ve
5	3,3	1,4	0,2	Se	5,9	3	5,1	1,8	Vi	5,7	2,8	4,1	1,3	Ve

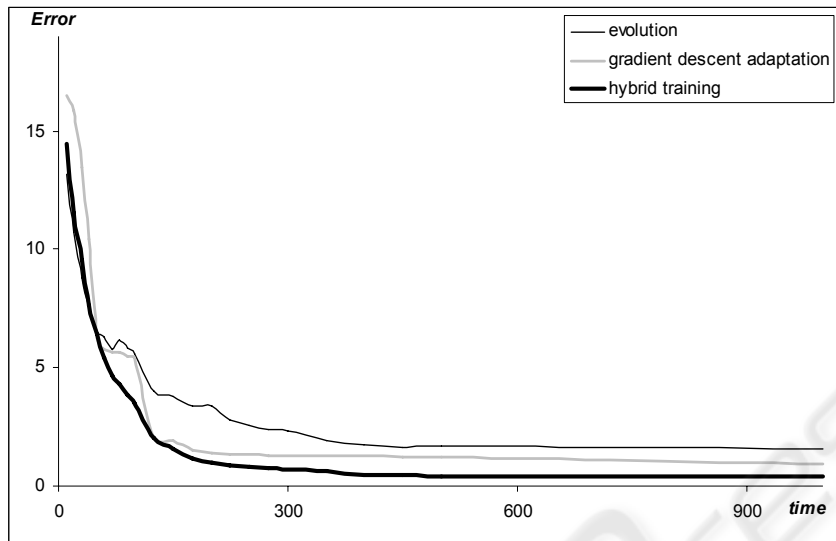


Fig. 2. The error function history.

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