

# Regularized Evolution for Image Classifier Architecture Search

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## Abstract

The effort devoted to hand-crafting neural network image classifiers has motivated the use of architecture search to discover them automatically. Although evolutionary algorithms have been repeatedly applied to neural network topologies, the image classifiers thus discovered have remained inferior to human-crafted ones. Here, we evolve an image classifier—*AmoebaNet-A*—that surpasses hand-designs for the first time. To do this, we modify the tournament selection evolutionary algorithm by introducing an age property to favor the younger genotypes. Matching size, AmoebaNet-A has comparable accuracy to current state-of-the-art ImageNet models discovered with more complex architecture-search methods. Scaled to larger size, AmoebaNet-A sets a new state-of-the-art 83.9% top-1 / 96.6% top-5 ImageNet accuracy. In a controlled comparison against a well known reinforcement learning algorithm, we give evidence that evolution can obtain results faster with the same hardware, especially at the earlier stages of the search. This is relevant when fewer compute resources are available. Evolution is, thus, a simple method to effectively discover high-quality architectures.

## Introduction

Until recently, most state-of-the-art image classifier architectures have been manually designed by human experts (Krizhevsky, Sutskever, and Hinton 2012; Szegedy et al. 2015; He et al. 2016; Huang et al. 2017; Hu, Shen, and Sun 2018). To speed up the process, researchers have looked into automated methods (Baker et al. 2017a; Zoph and Le 2016; Miikkulainen et al. 2017; Real et al. 2017; Xie and Yuille 2017; Suganuma, Shirakawa, and Nagao 2017; Liu et al. 2018a; Pham et al. 2018). These methods are now collectively known as *architecture-search algorithms*. A traditional approach is *neuro-evolution of topologies* (Miller, Todd, and Hegde 1989; Angeline, Saunders, and Pollack 1994; Stanley and Miikkulainen 2002). Improved hardware now allows scaling up evolution to produce high-quality image classifiers (Real et al. 2017; Xie and Yuille 2017; Liu et al. 2018b). Yet, the architectures produced by evolutionary algorithms / genetic programming have not reached the accuracy of those directly designed by human experts. Here we evolve image classifiers that surpass hand-designs.

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To do this, we make two additions to the standard evolutionary process. First, we propose a change to the well-established *tournament selection* evolutionary algorithm (Goldberg and Deb 1991) that we refer to as *aging* evolution or *regularized* evolution. Whereas in tournament selection, the best genotypes (architectures) are kept, we propose to associate each genotype with an age, and bias the tournament selection to choose the younger genotypes. We will show that this change turns out to make a difference. The connection to regularization will be clarified in the Discussion section. Second, we implement the simplest set of mutations that would allow evolving in the NASNet search space (Zoph et al. 2018). This search space associates convolutional neural network architectures with small directed graphs in which vertices represent hidden states and labeled edges represent common network operations (such as convolutions or pooling layers). Our mutation rules only alter architectures by randomly reconnecting the origin of edges to different vertices and by randomly relabeling the edges, covering the full search space.

Searching in the NASNet space allows a controlled comparison between evolution and the original method for which the space was designed, reinforcement learning (RL). Thus, this paper presents the first comparative case study of architecture-search algorithms for the image classification task. Within this case study, we will demonstrate that evolution can attain similar results with a simpler method, as will be shown in the Discussion section. In particular, we will highlight that in all our experiments evolution searched faster than RL and random search, especially at the earlier stages, which is important when experiments cannot be run for long times due to compute resource limitations.

Despite its simplicity, our approach works well in our benchmark against RL. It also evolved a high-quality model, which we name *AmoebaNet-A*. This model is competitive with the best image classifiers obtained by any other algorithm today at similar sizes (82.8% top-1 / 96.1% top-5 ImageNet accuracy). When scaled up, it sets a new state-of-the-art accuracy (83.9% top-1 / 96.6% top-5 ImageNet accuracy).

## Related Work

Review papers provide informative surveys of earlier (Yao 1999; Floreano, Dürr, and Mattiussi 2008) and more recent

(Elsken, Metzen, and Hutter 2018) literature on image classifier architecture search, including successful RL studies (Zoph and Le 2016; Baker et al. 2017a; Zoph et al. 2018; Liu et al. 2018a; Zhong, Yan, and Liu 2018; Cai et al. 2018) and evolutionary studies like those mentioned in the Introduction. Other methods have also been applied: cascade-correlation (Fahlman and Lebiere 1990), boosting (Cortes et al. 2017), hill-climbing (Elsken, Metzen, and Hutter 2017), MCTS (Negrinho and Gordon 2017), SMBO (Mendoza et al. 2016; Liu et al. 2018a), and random search (Bergstra and Bengio 2012), and grid search (Zagoruyko and Komodakis 2016). Some methods even forewent the idea of independent architectures (Saxena and Verbeek 2016). There is much architecture-search work beyond image classification too, but that is outside our scope.

Even though some methods stand out due to their efficiency (Suganuma, Shirakawa, and Nagao 2017; Pham et al. 2018), many approaches use large amounts of resources. Several recent papers reduced the compute cost through progressive-complexity search stages (Liu et al. 2018a), hypernets (Brock et al. 2018), accuracy prediction (Baker et al. 2017b; Klein et al. 2017; Domhan, Springenberg, and Hutter 2017), warm-starting and ensembling (Feurer et al. 2015), parallelization, reward shaping and early stopping (Zhong, Yan, and Liu 2018) or Net2Net transformations (Cai et al. 2018). Most of these methods could in principle be applied to evolution too, but this is beyond the scope of this paper.

A popular approach to evolution has been through *generational* algorithms, *e.g.* NEAT (Stanley and Miikkulainen 2002). In these, all models in the population must finish training before the next generation is computed. Generational evolution becomes inefficient in a distributed environment where a different machine is used to train each model: machines that train faster models finish earlier and must wait idle until all machines are ready. Real-time algorithms address this issue, *e.g.* rtNEAT (Stanley, Bryant, and Miikkulainen 2005) and tournament selection (Goldberg and Deb 1991). Unlike the generational algorithms, however, these discard models according to their performance or do not discard them at all, resulting in models that remain alive in the population for a long time—even for the whole experiment. We will present evidence that the finite lifetimes of aging evolution can give better results than direct tournament selection, while retaining its efficiency.

An existing paper (Hornby 2006) uses a concept of *age* but in a very different way than we do. In that paper, age is assigned to genes to divide a constant-size population into groups called *age-layers*. Each layer contains individuals with genes of similar ages. Only after the genes have survived a certain *age-gap*, they can make it to the next layer. The goal is to restrict competition (the newly introduced genes cannot be immediately out-competed by highly selected older ones). Their algorithm requires the introduction of two additional meta-parameters (size of the age-gap and number of age-layers). In contrast, in our algorithm, an *age* is assigned to the individuals (not the genes) and is only used to track which is the oldest individual in the population. This permits removing such oldest individual at each cycle (keeping a constant population size). Our approach,

therefore, is in line with our goal of keeping the method as simple as possible. In particular, our method remains similar to nature (where the young are less likely to die than the very old) and it requires no additional meta-parameters.

## Methods

This section contains a readable description of the methods. The Methods Details section gives additional information.

### Search Space

All experiments use the *NASNet search space* (Zoph et al. 2018). This is a space of image classifiers, all of which have the fixed outer structure indicated in Figure 1 (left): a feed-forward stack of Inception-like modules called *cells*. Each cell receives a *direct input* from the previous cell (as depicted) and a *skip input* from the cell before it (Figure 1, right). The cells in the stack are of two types: the *normal cell* and the *reduction cell*. All normal cells are constrained to have the same architecture, as are reduction cells, but the architecture of the normal cells is independent of that of the reduction cells. Other than this, the only difference between them is that every application of the reduction cell is followed by a stride of 2 that reduces the image size, whereas normal cells preserve the image size. As can be seen in the figure, normal cells are arranged in three stacks of  $N$  cells. The goal of the architecture-search process is to discover the architectures of the normal and reduction cells.

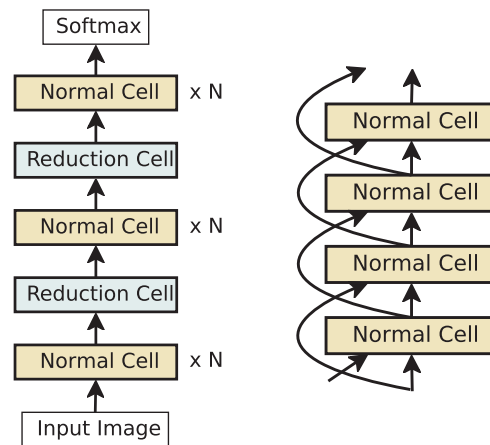


Figure 1: NASNet Search Space outer structure (Zoph et al. 2018). LEFT: the full outer structure, omitting skip inputs for clarity. RIGHT: detailed view with the skip inputs.

As depicted in Figure 1 (right) and Figure 2, each cell has two input activation tensors and one output. The very first cell takes two copies of the input image. After that, the inputs are the outputs of the previous two cells.

Both normal and reduction cells must conform to the following construction. The two cell input tensors are considered hidden states “0” and “1”. More hidden states are then constructed through *pairwise combinations*. A pairwise combination is depicted in Figure 2 (inside dashed circle). It consists in applying an operation (or *op*) to an existing

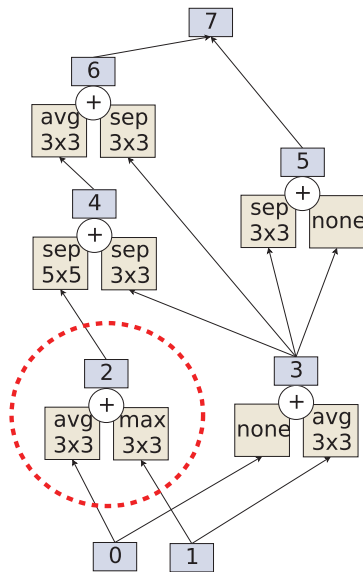


Figure 2: NASNet Search Space cell structure (Zoph et al. 2018). Example of a cell. Dotted line demarcates a pairwise combination.

hidden state, applying another op to another existing hidden state, and adding the results to produce a new hidden state. Ops belong to a fixed set of common convnet operations such as convolutions and pooling layers. Repeating hidden states or operations within a combination is permitted. In the cell example of Figure 2, the first pairwise combination applies a 3x3 average pool op to hidden state 0 and a 3x3 max pool op to hidden state 1, in order to produce hidden state 2. The next pairwise combination can now choose from hidden states 0, 1, and 2 to produce hidden state 3 (chose 0 and 1 in Figure 2), and so on. After exactly five pairwise combinations, any hidden states that remain unused (hidden states 5 and 6 in Figure 2) are concatenated to form the output of the cell (hidden state 7).

A given architecture is fully specified by the five pairwise combinations that make up the normal cell and the five that make up the reduction cell. Once the architecture is specified, the model still has two free parameters that can be used to alter its size (and its accuracy): the number of normal cells per stack ( $N$ ) and the number of output filters of the convolution ops ( $F$ ).  $N$  and  $F$  are determined manually.

## Evolutionary Algorithm

The evolutionary method we used is summarized in Algorithm 1. It keeps a population of  $P$  trained models throughout the experiment. The population is initialized with models with random architectures (“while  $|population|$ ” in Algorithm 1). All architectures that conform to the search space described are possible and equally likely.

After this, evolution improves the initial population in cycles (“while  $|history|$ ” in Algorithm 1). At each cycle, it samples  $S$  random models from the population, each drawn uniformly at random with replacement. The model with the highest validation fitness within this sample is selected as the

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### Algorithm 1 Aging Evolution (i.e. Regularized Evolution)

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```

population ← empty queue           ▷ The population.
history ← ∅                         ▷ Will contain all models.
while |population| < P do           ▷ Initialize population.
    model.arch ← RANDOMARCHITECTURE()
    model.accuracy ← TRAINANDEVAL(model.arch)
    add model to right of population
    add model to history
end while
while |history| < C do             ▷ Evolve for C cycles.
    sample ← ∅                       ▷ Parent candidates.
    while |sample| < S do
        candidate ← random element from population
                                ▷ The element stays in the population.
        add candidate to sample
    end while
    parent ← highest-accuracy model in sample
    child.arch ← MUTATE(parent.arch)
    child.accuracy ← TRAINANDEVAL(child.arch)
    add child to right of population
    add child to history
    remove dead from left of population   ▷ Oldest.
    discard dead
end while
return highest-accuracy model in history

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parent. A new architecture, called the *child*, is constructed from the parent by the application of a transformation called a *mutation*. A mutation causes a simple and random modification of the architecture and is described in detail below. Once the child architecture is constructed, it is then trained, evaluated, and added to the population. This process is called tournament selection (Goldberg and Deb 1991).

It is common in tournament selection to keep the population size fixed at the initial value  $P$ . This is often accomplished with an additional step within each cycle: discarding (or *killing*) the worst model in the random  $S$ -sample. We will refer to this approach as *non-aging evolution*. In contrast, in this paper we prefer a novel approach: killing the oldest model in the population—that is, removing from the population the model that was trained the earliest (“remove *dead* from left of *pop*” in Algorithm 1). This favors the newer models in the population. We will refer to this approach as *aging evolution*. In the context of architecture search, aging evolution allows us to explore the search space more, instead of zooming in on good models too early, as non-aging evolution would (see Discussion section for details).

In practice, this algorithm is parallelized by distributing the “while  $|history|$ ” loop in Algorithm 1 over multiple workers. Intuitively, the mutations can be thought of as providing exploration, while the parent selection provides exploitation. The parameter  $S$  controls the aggressiveness of the exploitation:  $S = 1$  reduces to a type of random search and  $2 \leq S \leq P$  leads to evolution of varying greediness.

New models are constructed by applying a mutation to existing models, transforming their architectures in random ways. To navigate the NASNet search space described

above, we use two main mutations that we call the *hidden state mutation* and the *op mutation*. A third mutation, the identity, is also possible. Only one of these mutations is applied in each cycle, choosing between them at random.

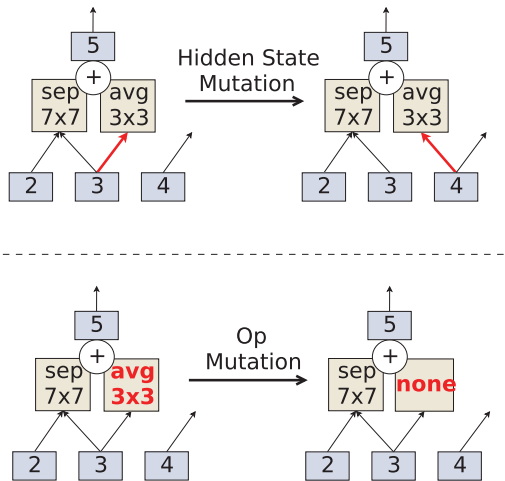


Figure 3: Illustration of the two mutation types.

The hidden state mutation consists of first making a random choice of whether to modify the normal cell or the reduction cell. Once a cell is chosen, the mutation picks one of the five pairwise combinations uniformly at random. Once the pairwise combination is picked, one of the two elements of the pair is chosen uniformly at random. The chosen element has one hidden state. This hidden state is now replaced with another hidden state from within the cell, subject to the constraint that no loops are formed (to keep the feed-forward nature of the convnet). Figure 3 (top) shows an example.

The op mutation behaves like the hidden state mutation as far as choosing one of the two cells, one of the five pairwise combinations, and one of the two elements of the pair. Then it differs in that it modifies the op instead of the hidden state. It does this by replacing the existing op with a random choice from a fixed list of ops (see Methods Details). Figure 3 (bottom) shows an example.

## Baseline Algorithms

Our main baseline is the application of RL to the same search space. RL was implemented using the algorithm and code in the baseline study (Zoph et al. 2018). An LSTM controller outputs the architectures, constructing the pairwise combinations one at a time, and then gets a reward for each architecture by training and evaluating it. More detail can be found in the baseline study. We also compared against random search (RS). In our RS implementation, each model is constructed randomly so that all models in the search space are equally likely, as in the initial population in the evolutionary algorithm. In other words, the models in RS experiments are *not* constructed by mutating existing models, so as to make new models independent from previous ones.

## Experimental Setup

We ran controlled comparisons at scale, ensuring identical conditions for evolution, RL and random search (RS). In particular, all methods used *the same* computer code for network construction, training and evaluation. Experiments always searched on the CIFAR-10 dataset (Krizhevsky and Hinton 2009).

As in the baseline study, we first performed architecture search over small models (*i.e.* small N and F) until 20k models were evaluated. After that, we used the *model augmentation* trick (Zoph et al. 2018): we took architectures discovered by the search (*e.g.* the output of an evolutionary experiment) and turned them into a full-size, accurate models. To accomplish this, we enlarged the models by increasing N and F so the resulting model sizes would match the baselines, and we trained the enlarged models for a longer time on the CIFAR-10 or the ImageNet classification datasets (Krizhevsky and Hinton 2009; Deng et al. 2009). For ImageNet, a stem was added at the input of the model to reduce the image size, as shown in Figure 6 (left). This is the same procedure as in the baseline study. To produce the largest model (see last paragraph of Results section), we increased N and F until we ran out of memory. Actual values of N and F for all models are listed in the Methods Details section.

## Methods Details

This section complements the Methods section with the details necessary to reproduce our experiments. Possible ops: none (identity); 3x3, 5x5 and 7x7 separable (sep.) convolutions (convs.); 3x3 average (avg.) pool; 3x3 max pool; 3x3 dilated (dil.) sep. conv.; 1x7 then 7x1 conv. Evolved with  $P=100$ ,  $S=25$ . CIFAR-10 dataset (Krizhevsky and Hinton 2009) with 5k withheld examples for validation. Standard ImageNet dataset (Deng et al. 2009), 1.2M 331x331 images and 1k classes; 50k examples withheld for validation; standard validation set used for testing. During the search phase, each model trained for 25 epochs;  $N=3/F=24$ , 1 GPU. Each experiment ran on 450 K40 GPUs for 20k models (approx. 7 days). To optimize evolution, we tried 5 configurations with P/S of: 100/2, 100/50, 20/20, 100/25, 64/16, best was 100/25. The probability of the identity mutation was fixed at the small, arbitrary value of 0.05 and was not tuned. Other mutation probabilities were uniform, as described in the Methods. To optimize RL, started with parameters already tuned in the baseline study and further optimized learning rate in 8 configurations: 0.00003, 0.00006, 0.00012, 0.0002, 0.0004, 0.0008, 0.0016, 0.0032; best was 0.0008. To avoid selection bias, plots do not include optimization runs, as was decided a priori. Best few (20) models were selected from each experiment and augmented to  $N=6/F=32$ , as in baseline study; batch 128, SGD with momentum rate 0.9, L2 weight decay  $5 \times 10^{-4}$ , initial lr 0.024 with cosine decay, 600 epochs, ScheduledDropPath to 0.7 prob; auxiliary softmax with half-weight of main softmax. For Table 1, we used N/F of 6/32 and 6/36. For ImageNet table, N/F were 6/190 and 6/448 and standard training methods (Szegedy et al. 2017): distributed sync SGD with 100 P100 GPUs; RMSProp opti-



mizer with 0.9 decay and  $\epsilon=0.1$ ,  $4 \times 10^{-5}$  weight decay, 0.1 label smoothing, auxiliary softmax weighted by 0.4; dropout probability 0.5; ScheduledDropPath to 0.7 probability (as in baseline—note that this trick only contributes 0.3% top-1 ImageNet acc.); 0.001 initial lr, decaying every 2 epochs by 0.97. Largest model used  $N=6/F=448$ . F always refers to the number of filters of convolutions in the first stack; after each reduction cell, this number is doubled. Wherever applicable, we used the same conditions as the baseline study.

## Results

### Comparison With RL and RS Baselines

Currently, reinforcement learning (RL) is the predominant method for architecture search. In fact, today’s state-of-the-art image classifiers have been obtained by architecture search with RL (Zoph et al. 2018; Liu et al. 2018a). Here we seek to compare our evolutionary approach against their RL algorithm. We performed large-scale side-by-side architecture-search experiments on CIFAR-10. We first optimized the hyper-parameters of the two approaches independently (details in Methods Details section). Then we ran 5 repeats of each of the two algorithms—and also of random search (RS).

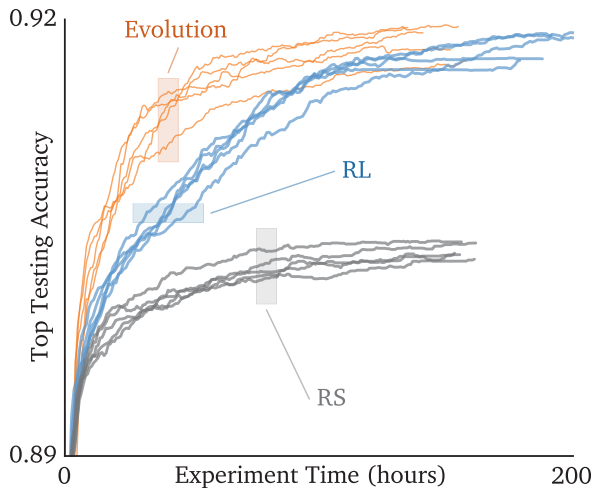


Figure 4: Time-course of 5 identical large-scale experiments for each algorithm (evolution, RL, and RS), showing accuracy before augmentation on CIFAR-10. All experiments were stopped when 20k models were evaluated, as done in the baseline study. Note this plot does not show the compute cost of models, which was higher for the RL ones.

Figure 4 shows the model accuracy as the experiments progress, highlighting that evolution yielded more accurate models at the earlier stages, which could become important in a resource-constrained regime where the experiments may have to be stopped early (for example, when 450 GPUs for 7 days is too much). At the later stages, if we allow to run for the full 20k models (as in the baseline study), evolution produced models with similar accuracy. Both evolution and

RL compared favorably against RS. It is important to note that the vertical axis of Figure 4 does not present the compute cost of the models, only their accuracy. Next, we will consider their compute cost as well.

As in the baseline study, the architecture-search experiments above were performed over small models, to be able to train them quicker. We then used the *model augmentation* trick (Zoph et al. 2018) by which we take an architecture discovered by the search (e.g. the output of an evolutionary experiment) and turn it into a full-size, accurate model, as described in the Methods.

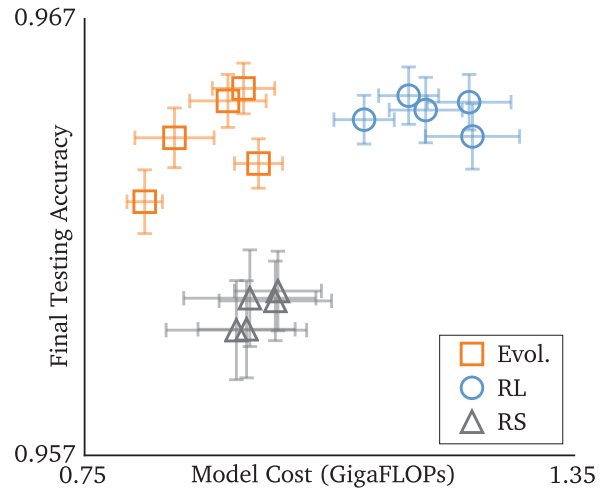


Figure 5: Final augmented models from 5 identical architecture-search experiments for each algorithm, on CIFAR-10. Each marker corresponds to the top models from one experiment.

Figure 5 compares the augmented top models from the three sets of experiments. It shows test accuracy and model compute cost. The latter is measured in FLOPs, by which we mean the total count of operations in the forward pass, so lower is better. Evolved architectures had higher accuracy (and similar FLOPs) than those obtained with RS, and lower FLOPs (and similar accuracy) than those obtained with RL. Number of parameters showed similar behavior to FLOPs. Therefore, evolution occupied the ideal relative position in this graph within the scope of our case study.

Table 1: CIFAR-10 testing set results for AmoebaNet-A, compared to top model reported in the baseline study.

Model	# Params	Test Error (%)
NASNet-A (baseline)	3.3 M	3.41
AmoebaNet-A (N=6, F=32)	2.6 M	3.40 ± 0.08
AmoebaNet-A (N=6, F=36)	3.2 M	3.34 ± 0.06

So far we have been comparing evolution with our reproduction of the experiments in the baseline study, but it is also informative to compare directly against the results reported

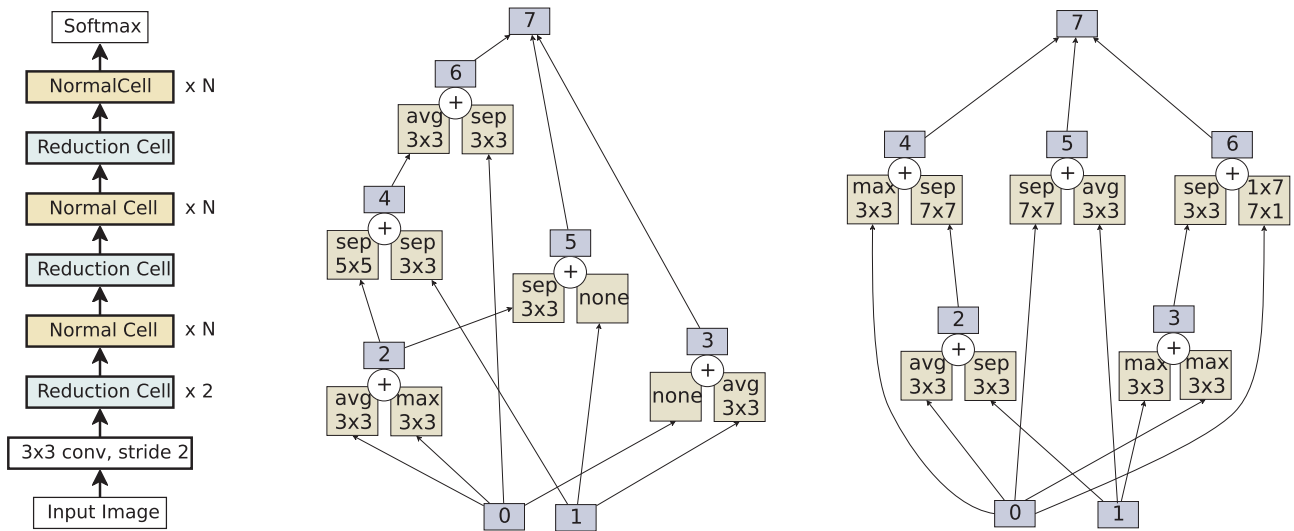


Figure 6: AmoebaNet-A architecture. The overall model (Zoph et al. 2018) (LEFT) and the AmoebaNet-A normal cell (MID-DLE) and reduction cell (RIGHT).

by the baseline study. We select our evolved architecture with highest validation accuracy and call it *AmoebaNet-A* (Figure 6). Table 1 compares its test accuracy with the top model of the baseline study, NASNet-A. Such a comparison is not entirely controlled, as we have no way of ensuring the network training code was identical and that the same number of experiments were done to obtain the final model. The table summarizes the results of training AmoebaNet-A at sizes comparable to a NASNet-A version, showing that AmoebaNet-A is slightly more accurate (when matching model size) or considerably smaller (when matching accuracy). We did not train our model at larger sizes on CIFAR-10. Instead, we moved to ImageNet to do further comparisons in the next section.

## ImageNet Results

Following the accepted standard, we compare our top model’s classification accuracy on the popular ImageNet dataset against other top models from the literature. Again, we use AmoebaNet-A, the model with the highest validation accuracy on CIFAR-10 among our evolution experiments. We highlight that the model was evolved on CIFAR-10 and then transferred to ImageNet, so the evolved *architecture* cannot have overfit the ImageNet dataset. When re-trained on ImageNet, AmoebaNet-A performs comparably to the baseline for the same number of parameters (Table 2, model with  $F=190$ ).

Finally, we focused on AmoebaNet-A exclusively and enlarged it, setting a new state-of-the-art accuracy on ImageNet of 83.9%/96.6% top-1/5 accuracy with 469M parameters (Table 2, model with  $F=448$ ). Such high parameter counts may be beneficial in training other models too but we have not managed to do this yet.

## Discussion

This section will suggest directions for future work, which we will motivate by speculating about the evolutionary process and by summarizing additional minor results. The details of these minor results have been relegated to the supplements, as they are not necessary to understand or reproduce our main results above.

**Scope of results.** Some of our findings may be restricted to the search spaces and datasets we used. A natural direction for future work is to extend the controlled comparison to more search spaces, datasets, and tasks, to verify generality, or to more algorithms. Supplement A presents preliminary results, performing evolutionary and RL searches over three search spaces (SP-I: same as in the Results section; SP-II: like SP-I but with more possible ops; SP-III: like SP-II but with more pairwise combinations) and three datasets (gray-scale CIFAR-10, MNIST, and gray-scale ImageNet), at a small-compute scale (on CPU,  $F=8$ ,  $N=1$ ). Evolution reached equal or better accuracy in all cases (Figure 7, top).

**Algorithm speed.** In our comparison study, Figure 4 suggested that both RL and evolution are approaching a common accuracy asymptote. That raises the question of which algorithm gets there faster. The plots indicate that evolution reaches half-maximum accuracy in roughly half the time. We abstain, nevertheless, from further quantifying this effect since it depends strongly on how speed is measured (the number of models necessary to reach accuracy  $a$  depends on  $a$ ; the natural choice of  $a = a_{max}/2$  may be too low to be informative; *etc.*). Algorithm speed may be more important when exploring larger spaces, where reaching the optimum can require more compute than is available. We saw an example of this in the SP-III space, where evolution stood out (Figure 7, bottom-right). Therefore, future work could explore evolving on even larger spaces, where the initial relative speed of evolution may be even more significant.

Table 2: ImageNet classification results for AmoebaNet-A compared to hand-designs (top rows) and other automated methods (middle rows). The evolved AmoebaNet-A architecture (bottom rows) reaches the current state of the art (SOTA) at similar model sizes and sets a new SOTA at a larger size. All evolution-based approaches are marked with a \*. We omitted Squeeze-and-Excite-Net because it was not benchmarked on the same ImageNet dataset version.

Model	# Parameters	# Multiply-Adds	Top-1 / Top-5 Accuracy (%)
Incep-ResNet V2 (Szegedy et al. 2017)	55.8M	13.2B	80.4 / 95.3
ResNeXt-101 (Xie et al. 2017)	83.6M	31.5B	80.9 / 95.6
PolyNet (Zhang et al. 2017)	92.0M	34.7B	81.3 / 95.8
Dual-Path-Net-131 (Chen et al. 2017)	79.5M	32.0B	81.5 / 95.8
GeNet-2 (Xie and Yuille 2017)*	156M	-	72.1 / 90.4
Block-QNN-B (Zhong, Yan, and Liu 2018)*	-	-	75.7 / 92.6
Hierarchical (Liu et al. 2018b)*	64M	-	79.7 / 94.8
NASNet-A (Zoph et al. 2018)	88.9M	23.8B	82.7 / 96.2
PNASNet-5 (Liu et al. 2018a)	86.1M	25.0B	82.9 / 96.2
AmoebaNet-A (N=6, F=190)*	86.7M	23.1B	82.8 / 96.1
AmoebaNet-A (N=6, F=448)*	469M	104B	83.9 / 96.6

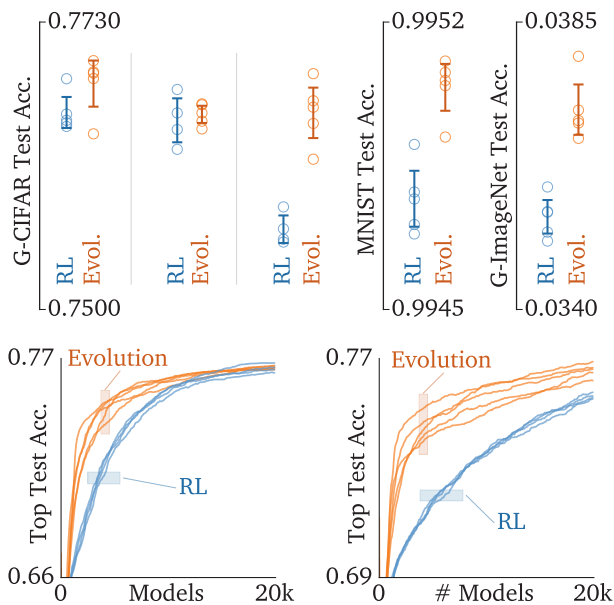


Figure 7: TOP: Comparison of the final model accuracy in five different contexts, from left to right: G-CIFAR/SP-I, G-CIFAR/SP-II, G-CIFAR/SP-III, MNIST/SP-I and G-ImageNet/SP-I. Each circle marks the top test accuracy at the end of one experiment. BOTTOM: Search progress of the experiments in the case of G-CIFAR/SP-II (LEFT, best for RL) and G-CIFAR/SP-III (RIGHT, best for evolution).

**Model speed.** The speed of individual models produced is also relevant. Figure 5 demonstrated that evolved models are faster (lower FLOPs). We speculate that asynchronous evolution may be reducing the FLOPs because it is indirectly optimizing for speed even when training for a fixed number of epochs: fast models may do well because they “reproduce” quickly even if they initially lack the higher accuracy

of their slower peers. Verifying this speculation could be the subject of future work. As mentioned in the Related Work section, in this work we only considered asynchronous algorithms (as opposed to generational evolutionary methods) to ensure high resource utilization. Future work may explore how asynchronous and generational algorithms compare with regard to model accuracy.

**Benefits of aging evolution.** Aging evolution also seemed advantageous in small-compute-scale experiments, shown in Figure 8 and presented in more detail in Supplement B.

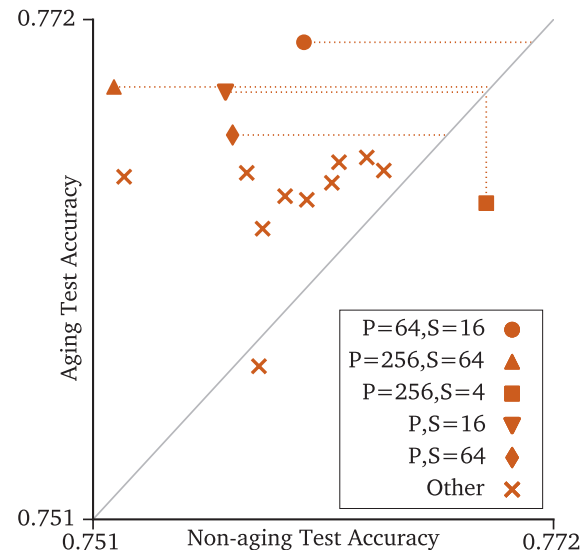


Figure 8: Small-compute-scale comparison between our aging tournament selection variant and the non-aging variant, for different population sizes (P) and sample sizes (S). Aging tends to be beneficial (most markers above the  $y=x$  line).

In Supplement B, we also show that the benefits of aging

evolution tend to hold when varying the dataset or the search space. In order to reduce compute requirements, all these additional experiments were carried out on CPU instead of GPU and used a gray-scale version of CIFAR-10.

**Understanding aging evolution and regularization.** We can speculate that aging may help navigate the training noise in evolutionary experiments, as follows. Noisy training means that models may sometimes reach high accuracy just by luck. In non-aging evolution (NAE, *i.e.* standard tournament selection), such lucky models may remain in the population for a long time—even for the whole experiment. One lucky model, therefore, can produce many children, causing the algorithm to focus on it, reducing exploration. Under aging evolution (AE), on the other hand, all models have a short lifespan, so the population is wholly renewed frequently, leading to more diversity and more exploration. In addition, another effect may be in play, which we describe next. In AE, because models die quickly, the only way an architecture can remain in the population for a long time is by being passed down from parent to child through the generations. Each time an architecture is inherited it must be re-trained. If it produces an inaccurate model when re-trained, that model is not selected by evolution and the architecture disappears from the population. The only way for an architecture to remain in the population for a long time is to re-train well repeatedly. In other words, AE can only improve a population through the inheritance of architectures that re-train well. (In contrast, NAE can improve a population by accumulating architectures/models that were lucky when they trained the first time). That is, AE is forced to pay attention to *architectures* rather than *models*. In other words, the addition of aging involves introducing additional information to the evolutionary process: architectures should re-train well. This additional information prevents overfitting to the training noise, which makes it a form of *regularization* in the broader mathematical sense<sup>1</sup>. Regardless of the exact mechanism, in Supplement C we perform experiments to verify the plausibility of the conjecture that aging helps navigate noise. There we construct a toy search space where the only difficulty is a noisy evaluation. If our conjecture is true, AE should be better in that toy space too. We found this to be the case. We leave further verification of the conjecture to future work, noting that theoretical results may prove useful here.

**Simplicity of aging evolution.** A desirable feature of evolutionary algorithms is their simplicity. By design, the application of a mutation causes a random change. The process of constructing new architectures, therefore, is entirely random. What makes evolution different from random search is that only the good models are selected to be mutated. This selection tends to improve the population over time. In this sense, evolution is simply “random search plus selection”. In outline, the process can be described briefly: “keep a population of  $N$  models and proceed in cycles: at each cycle, copy-mutate the best of  $S$  random models and kill the oldest in the population”. Implementation-wise, we believe the methods of this paper are sufficient for a reader to understand

evolution. The sophisticated nature of the RL alternative introduces complexity in its implementation: it requires back-propagation and poses challenges to parallelization (Salimans et al. 2017). Even different implementations of the same algorithm have been shown to produce different results (Henderson et al. 2018). Finally, evolution is also simple in that it has few meta-parameters, most of which do not need tuning (Real et al. 2017). In our study, we only adjusted 2 meta-parameters and only through a handful of attempts (see Methods Details section). In contrast, note that the RL baseline requires training an agent/controller which is often itself a neural network with many weights (such as an LSTM), and its optimization has more meta-parameters to adjust: learning rate schedule, greediness, batching, replay buffer parameters, *etc.* (These meta-parameters are all in addition to the weights and training parameters of the image classifiers being searched, which are present in both approaches.) It is possible that through careful tuning, RL could be made to produce even better models than evolution, but such tuning would likely involve running many experiments, making it more costly. Evolution did not require much tuning, as described. It is also possible that random search would produce equally good models if run for a very long time, which would be very costly. Finally, the evolutionary algorithm could be improved through additional complexity; for example, the mutation probabilities could be learned to improve speed.

**Interpreting architecture search.** Another important direction for future work is that of analyzing architecture-search experiments (regardless of the algorithm used) to try to discover new neural network design patterns. Anecdotally, for example, we found that architectures with high output vertex fan-in (number of edges into the output vertex) tend to be favored in all our experiments. In fact, the models in the final evolved populations have a mean fan-in value that is 3 standard deviations above what would be expected from randomly generated models. We verified this pattern by training various models with different fan-in values and the results confirm that accuracy increases with fan-in, as had been found in ResNeXt (Xie et al. 2017). Discovering broader patterns may require designing search spaces specifically for this purpose.

**Additional AmoebaNets.** Using variants of the evolutionary process described, we obtained three additional models, which we named *AmoebaNet-B*, *AmoebaNet-C*, and *AmoebaNet-D*. We describe these models and the process that led to them in detail in Supplement D, but we summarize here. *AmoebaNet-B* was obtained through platform-aware architecture search over a larger version of the NAS-Net space. *AmoebaNet-C* is simply a model that showed promise early on in the above experiments by reaching high accuracy with relatively few parameters; we mention it here for completeness, as it has been referenced in other work (Cubuk et al. 2018). *AmoebaNet-D* was obtained by manually extrapolating the evolutionary process and optimizing the resulting architecture for training speed. It is very efficient: *AmoebaNet-D* won the Stanford DAWN Bench competition for lowest training cost on ImageNet (Coleman et al. 2018).

<sup>1</sup>[https://en.wikipedia.org/wiki/Regularization\\_\(mathematics\)](https://en.wikipedia.org/wiki/Regularization_(mathematics))



## Supplements

The supplements can be found online at:  
<https://arxiv.org/abs/1802.01548>

## Conclusion

This paper used an evolutionary algorithm to discover image classifier architectures. Our contributions are the following:

- We proposed *aging evolution*, a variant of tournament selection by which genotypes die according to their age, favoring the young. This improved upon standard tournament selection while still allowing for efficiency at scale through asynchronous population updating. We open-sourced the code.<sup>2</sup> We also implemented simple mutations that permit the application of evolution to the popular NASNet search space.
- We presented the first controlled comparison of algorithms for image classifier architecture search in a case study of evolution, RL and random search. We showed that evolution had somewhat faster search speed and stood out in the regime of scarcer resources / early stopping. Evolution also matched RL in final model quality, employing a simpler method.
- We evolved AmoebaNet-A (Figure 6), a competitive image classifier. On ImageNet, it is the first evolved model to surpass hand-designs. Matching size, AmoebaNet-A has comparable accuracy to top image-classifiers discovered with other architecture-search methods. At large size, it sets a new state-of-the-art accuracy. We open-sourced code and checkpoint.<sup>3</sup>

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<sup>3</sup>[https://tfhub.dev/google/imagenet/amoebanet\\_a\\_n18\\_f448/classification/1](https://tfhub.dev/google/imagenet/amoebanet_a_n18_f448/classification/1)

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