A Bio-inspired Convolutional Layer

Code at https://github.com/Einlar/biopytorch

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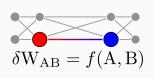
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1. Local Learning Rule

The change of a weight depends only on the activations of the neurons it connects.

- Only a forward pass is needed for training.
- Every layer except the very last is trained in an unsupervised way. The last layer can use the delta rule for a supervised task.

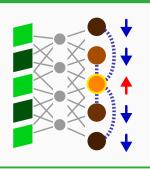


"Bio-inspired" algorithm by Krotov, Hopfield, *Unsupervised learning by competing hidden units*, 2019 [1].

2. Competition of Neurons

A neuron which is strongly activated for a pattern reduces the activation of all the other neurons in the same layer (lateral inhibition).

- Neurons "specialize" on different patterns.
- Only few neurons are active at a given time (sparse activation)

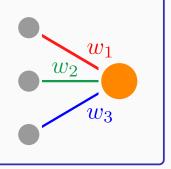


3. Normalization of Weights

Weights cannot grow indefinitely, and tend to the surface of a *p*-norm sphere.

$$|\mathbf{w}_1|^p + |\mathbf{w}_2|^p + |\mathbf{w}_3|^p \to R^p$$

 A lower p forces most weights to be small/zero (sparse weights).

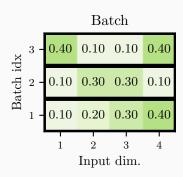


Bio-Learning Algorithm - Input

Let's inspect how the algorithm works through an example:

Consider the following batch of 3 samples $\{v_b\}_{i=1,2,3}$, each a vector with 4 entries.

Each sample is a ${\bf row}$ in the batch matrix V.



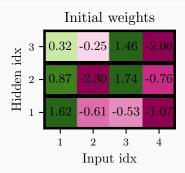
Bio-Learning Algorithm - Weights

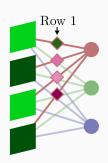
Consider 3 hidden neurons. Each of them has 4 weights connecting it to the 4 available inputs.

The weights of a neuron are a **row** in the weights matrix W.

Entries are normally distributed:

$$W_{\mu i} \sim \mathcal{N}(0,1)$$





Bio-Learning Algorithm - Lateral inhibition

For any sample v_b , compute the **current** received by the μ -th neuron as:

$$I_{\mu b} = \langle \boldsymbol{W}_{\mu}, \boldsymbol{v}_{b} \rangle_{p} \equiv \sum_{\nu=1}^{n} |W_{\mu \nu}|^{p-1} sgn(W_{\mu \nu}) V_{\nu b}$$

Then neurons **compete** with each other for their *post*- activations $g(I_{\mu\nu})$. For each sample:

- The neuron with the **highest** current "wins", and gets a positive post-activation, which will push it *towards* that sample.
- The neuron with the k-th highest current "loses", and gets a negative post-activation. In this way, at most k-1 neurons can be "close" to a pattern.

$$g(\mathrm{I}_{\mu b}) = egin{cases} 1 & \mu = r_{1b} \\ -\Delta & \mu = r_{k\mu} \\ 0 & \mathrm{otherwise} \end{cases}$$
 r_{sb} = index of the neuron with the s-th highest current for the b-th sample

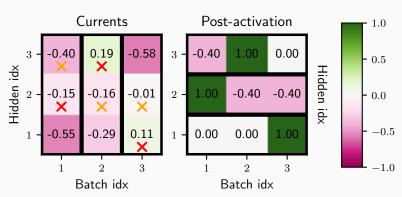


Figure 2 – Currents $I_{\mu b}$ and post-activations $g(I_{\mu b})$ for p=2, k=2, $\Delta=0.4$. For each sample in the batch (column in the left figure), the highest current is marked with a red \times , and the k-highest with an orange \times .

Bio-Learning Algorithm - Hebbian term

The **hebbian** term for the change of weights is given by:

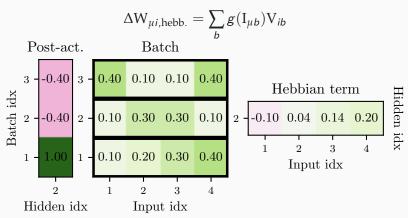


Figure 3 – To get the hebbian term, multiply every activation (left) by the corresponding sample (rows of center matrix). Then sum the rows. The figure shows the computation for $\mu=2$.

Bio-Learning Algorithm - Normalization term

The normalization term is given by:

$$\Delta W_{\mu i, \text{norm.}} = -\Big(\sum_{b} g(I_{\mu b})I_{\mu b}\Big)W_{\mu i}$$

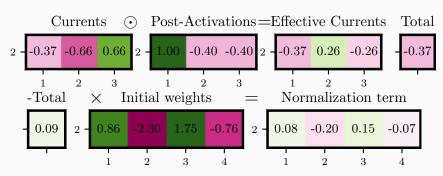


Figure 4 – Multiply the currents by the post-activations to get the effective currents, and sum the results to get the total. This is then used to multiply the initial weights to get the normalization term (for $\mu=2$ here).

Bio-Learning Algorithm - Weights update

The change in weights $\Delta W_{\mu i}$ consists of two terms: **hebbian** and **normalization**:

$$\Delta W_{\mu i} = \Delta W_{\mu i, \text{hebb.}} + \Delta W_{\mu i, \text{norm.}}$$

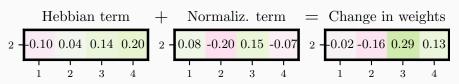


Figure 5 – Proposed change of weights for the second hidden unit ($\mu=2$) in the above example.

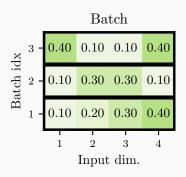
In practice, weights are updated iteratively, choosing a different minibatch at each iteration:

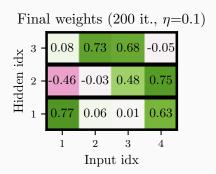
$$W_{\mu i}^{(t+1)} = W_{\mu i}^{(t)} + \eta_t \Delta W_{\mu i}^{(t)} \qquad W_{\mu i}^{(0)} \sim \mathcal{N}(\textbf{0}, \textbf{1})$$

where η_t is the (time-dependent) learning rate.

Bio-Learning Algorithm - Iterations

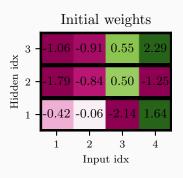
After 200 iterations with $\eta = 0.1$, the final weights converge:

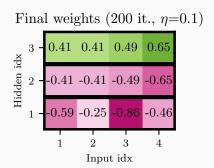




Bio-Learning Algorithm - Issues

However, a different weight initialization does **not** converge:





Note: the weights of the third neuron ($\mu=3$) are closer to the batch mean (positive), and so this neuron "dominates" over all the others, preventing them from adjusting with Hebbian learning.

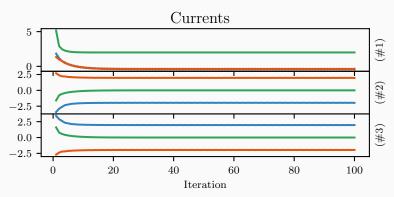


Figure 8 – Evolution of the currents during training. Each subplot refers to a different input sample, with the **blue**, **orange** and **green** lines referring respectively to the **first**, **second** and third neuron. Note how the current for $\mu = 3$ (the green one) is always the highest.

Bio-Learning Algorithm - Issues

To solve this issue, we would like the columns of ${\rm I}_{\mu b}$ to have (approximately) independent rankings.

This is done by normalizing the rows of V_{bi} :

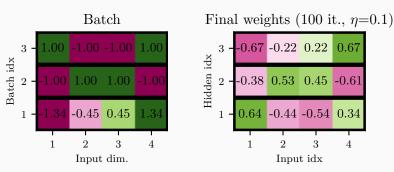
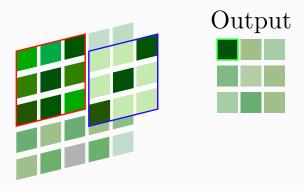


Figure 9 – The same weight initialization from before can converge if the batch samples are properly normalized to 0 mean and 1 std.

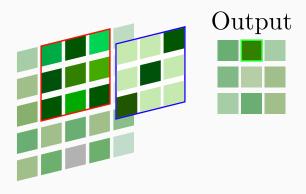
The main properties of a **convolutional layer**:

- Local Receptive Fields: each neuron receives inputs only from a patch of the image.
- Weight sharing: the same set of weights (kernel) is used for all the receptive fields.



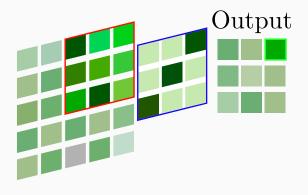
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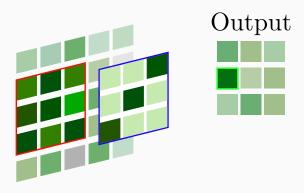
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Convolutions Why?

Convolutional Layers > Fully-Connected Layers

(for image classification)

Local Receptive Fields

- · Patterns in images are mostly local.
- Different scales can be captured by tweaking the kernel size, or by rescaling the inputs (e.g. through maxpooling).
- Inspired by retinal ganglion cells and V1 neurons in the primary visual cortex.

Weight sharing

- Ability to detect the same pattern in different position in the image.
- Images can be described by few "universal" patterns that can appear in any position.

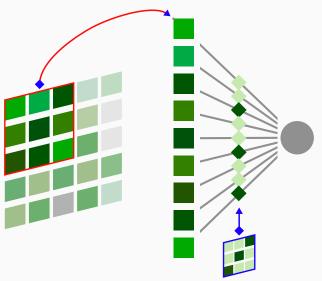
• Fewer weights to learn improve performance and help against overfitting.

How to *learn* Convolutional Layers with the bio-inspired rule?

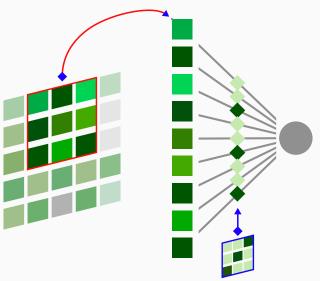
How to *learn* Convolutional Layers with the bio-inspired rule?

We just need the right perspective!

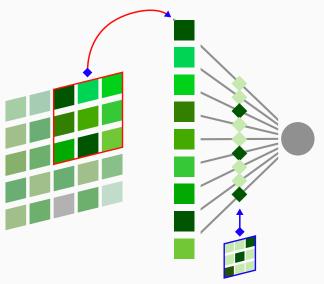
A **convolutional kernel** is just a **perceptron** working on patches:



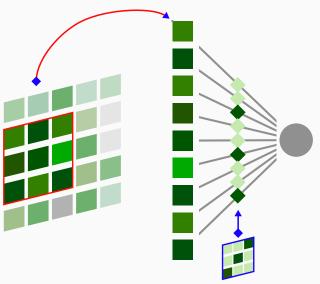
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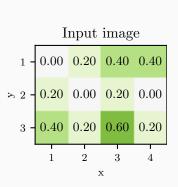
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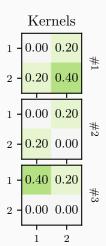


A **convolutional kernel** is just a **perceptron** working on patches:



Consider more kernels, acting on a one-channel image:





Convolutions Algorithm - Unfolding

By "unfolding" patches and kernels, the setup is the same as before:

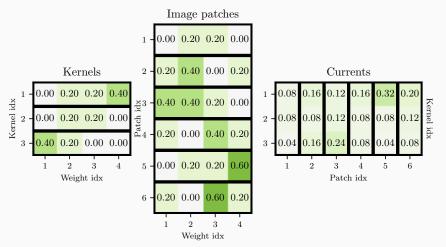


Figure 11 – Currents are the scalar product of each row of the "unfolded" kernels and each patch. This is the (reshaped) output of a 2d convolution.

Kernels compete as neurons:

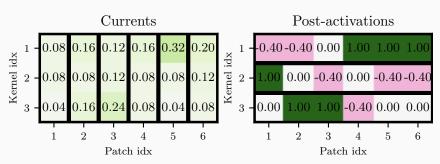
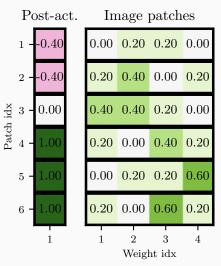


Figure 12 – The highest value in each column receives a (post)activation of 1. The k-th highest gets instead a $-\Delta$, while all the others are set to 0. Here $\Delta=0.4$ and k=2.

Convolutions Algorithm - Hebbian term

The **Hebbian term** is just the sum of the patches, weighted by the (post)activations:



Hebbian term



Unfolding patches is highly inefficient!

 Many patches: the number of patches scales as the square of the image size. Memory is quickly filled with a lot of redundant information.

For a 32 \times 32 image (CIFAR-10), with a 5 \times 5 kernel, there are 28² = 784 patches!

 Not optimized: modern CUDA kernels use many optimizations for 2d convolutions (e.g. Fourier transforms).

As a result, the above algorithm is *slow* and usually *crashes* due to the limited memory of a GPU.

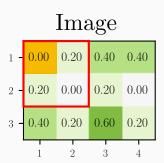
- Unfolding patches is highly inefficient!
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 - Not optimized: modern CUDA kernels use many optimizations for 2d convolutions (e.g. Fourier transforms).

As a result, the above algorithm is *slow* and usually *crashes* due to the limited memory of a GPU.

But it can be rewritten using a 2d convolution!

Consider the first entry of the Hebbian term.

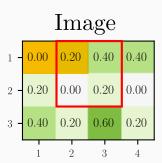
This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:

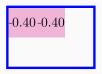




Consider the first entry of the Hebbian term.

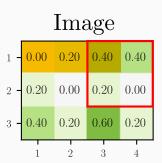
This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:

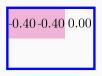




Consider the first entry of the Hebbian term.

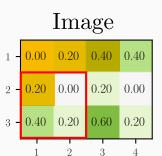
This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:





Consider the first entry of the Hebbian term.

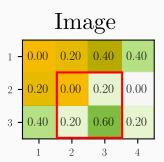
This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:





Consider the first entry of the Hebbian term.

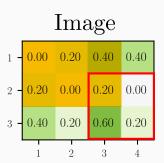
This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:





Consider the first entry of the Hebbian term.

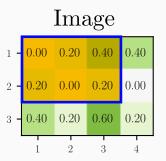
This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:



-0.40-0.40 0.00 1.00 1.00 1.00

Consider the first entry of the Hebbian term.

This is the scalar product between the **post-activations** vector and the **first column** of the image **patches**, which contains the **first** value of each receptive field:

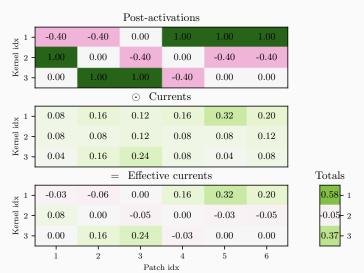


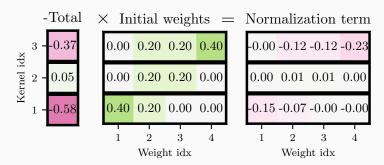


This is a 2d convolution with a 2×3 kernel, obtained by *reshaping* the post-activations!

Convolutions Algorithm - Normalization term

For the **normalization term**, the procedure is the same as before:





This is just a multiplication (with broadcasting), so it does not need any optimization.

Convolutions Algorithm - Technical details

What about the **implementation**?

- Implemented with PyTorch 1.9 [2].
- Supports training on GPU, and all the parameters of a PyTorch nn.Conv2d (stride, dilation, groups...).

What about **performance**?

- Each training pass requires two 2d-convolutions on the input tensor, with the first one being the forward pass (plus some faster operations).
 - $2-3\times$ slower than a Conv2d forward.
 - Up to 5× faster than a "unfold" implementation (e.g. dcasbol/biolearn_torch), with significantly less memory required!
- Few epochs (< 5) are needed to converge if the hyperparameters (p, k, Δ , η) are chosen well.

Convolutions Algorithm - CIFAR-10

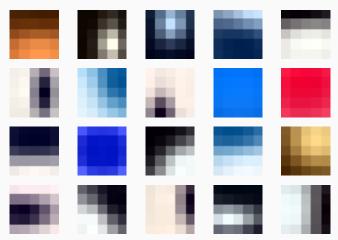
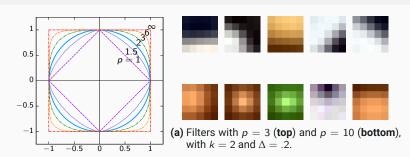


Figure 20 – First 20 5 \times 5 filters (out of 96) after training for 2 epochs on 40k RGB samples from CIFAR-10, with k=2, p=5 and $\Delta=0.2$. Learning rate starts at .007 and decays exponentially to 10^{-4} with $\lambda=.8$.

Convolutions Hyperparameters

1. Lebesgue norm p

- Lower p means sparser weights (fewer weights in a kernel can be significantly ≠ 0).
- Low p leads generally to more localized filters (e.g. edges/corners).
 High p leads to diffuse filters (e.g. on-center, gradients).
- Convergence (p-norm \rightarrow 1) is best for *middle* values: p = 5, 6.



Convolutions Hyperparameters

2. Ranking param k

- The neurons receiving the *k*-th highest currents are *pushed away* from those patterns.
- Thus, the number of neurons *encoding* a similar pattern is $\leq k 1$.
- Lower *k* leads to sparser activations (i.e. fewer kernels are active for a given sample).
- Convergence (p-norm → 1) is best for k = 4, 5. A lower k = 2, 3 works with higher Δ.

3. Anti-Hebbian strength Δ

- Specifies how much losing neurons are pushed away from a pattern.
- Higher values of ∆ can generate more diverse patterns, but if ∆ is too high, the training does not converge.
- Convergence (p-norm → 1) is very difficult for Δ > .2, and is generally easier for Δ ≤ .1.

BioConv2d - Hyperparameters Optimization

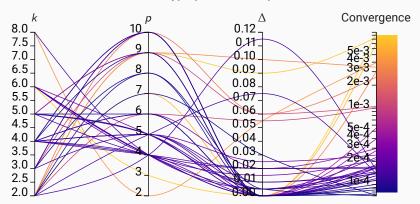


Figure 22 – Convergence is measured as $\max_{\mu} \left| \sum_{i} |W_{\mu i}|^{p} - 1 \right|$ (since $\sum_{i} |W_{\mu i}|^{p} \to R^{p} = 1$).

 \sim 500 trials are done on CIFAR-10 through Optuna [3], and the results are plotted with Wandb. Only "converged" runs (convergence $\leq .01$) are shown.

CIFAR-10 Dataset [4]

- 60k 32 × 32 RGB images of 10 classes, divided in 40k images for training, 10k for validation and 10k for testing.
- Initial **normalization** to [0, 1] range, by dividing pixel values by 255.
- Data augmentation: "Reflect" padding of 4 + Random Crop, Random horizontal flips (p = .5).



Figure 23 - CIFAR-10 classes.

Common architecture, taken from [5]:

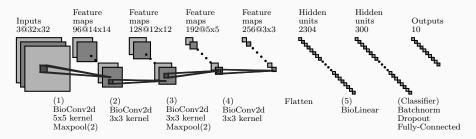


Figure 24 – The first 5 layers (4 convolutional, one linear) are trained using the bio-learning rule, while the last layer (classifier) is trained with SGD.

Each hebbian layer is preceded by a batch normalization (without affine transformation, so with no learnable parameters), and followed by a ReLU activation.

The **classifier** can be *adapted* to the output of each intermediate layer, to inspect the change in performance given by adding each layer.

Benchmark Common hyperparameters

The following hyperparameters are kept **fixed**:

- Batch size: 64.
- Hebbian learning rate
 - BioConv2d layers: start at .007, decay exponentially to 10^{-4} with $\lambda = 0.8$.
 - BioLinear layer: start at .1, decay exponentially to .005 with $\lambda = .1$.
- Hebbian layers are trained until they reach convergence ≤ .01 at the end of an epoch, or if convergence starts increasing for more than 10 batches.
- Training of Hebbian layers can be either sequential (wait for convergence of one layer to train the next), or all at once (train all layers on the same batches).
- **Optimizer** (classifier): SGD with momentum = 0.9, learningrate starting at 0.05, with Cosine Annealing with Warm Restarts ($T_0 = 10$, $T_{\text{mult}} = 2$) [6].

Performance and best hyperparameters

#layers	1	2	3	4	5
Accuracy (val)	66.97	65.44	63.87	58.79	45.91
Accuracy (test)	66.21	65.22	63.08	58.86	45.45
р	2	8	8	8	8
k	9	3	5	7	2
Δ	.08	.34	.25	.235	.335
Dropout	.2	.25	.05	.1	.1
Params	195k	302k	387k	804k	1.475M

Table 1 – Results of hyperparameter optimization (0ptuna, ~ 100 -200 trials for each case). Note how higher p and Δ are necessary for deeper architectures, but with lower dropout. However, hebbian learning quickly loses performance when adding layers, because the learning rule has "no way to know" which features should be kept for improving classification.

Test accuracy with different learning rules

#layers	1	2	3	4	5
Hebbian (Krotov)	66.21	65.22	63.08	58.86	45.45
Hebbian ([5])	63.92	63.81	58.28	52.99	41.78
Full SGD ([5])	60.71	66.30	72.39	82.69	84.95

Table 2 – Comparison with results from [5]. Krotov's learning rule seems slightly better. Moreover, contrary to [5], there is no "teaching signal" added to neurons in the unsupervised layer to develop class-specificity.

Nonetheless, a network fully trained with SGD can achieve higher performance, especially for deeper layers.

Benchmark Discussion

Convolutions work better!

- The network with 1 fully-connected unsupervised layer from [1] achieved only 55.26% accuracy, but with 6.166M parameters.
- The convolutional analogue achieves 66.21% accuracy, with only 195k parameters.
- Hebbian layers requires < 5 epochs to train (very fast).
- Performance is very good (better than SGD!) for shallow networks, but decreases for deeper ones.
 - Possible applications: transfer learning, fine-tuning, weight initialization
 - Supervised "top-down" signals are necessary to solve the performance lost to depth.

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 Unsupervised learning by competing hidden units.
 Proceedings of the National Academy of Sciences, 116(16):7723-7731, 2019.
- [2] Adam Paszke, Sam Gross, Francisco Massa, Adam Lerer, James Bradbury, Gregory Chanan, Trevor Killeen, Zeming Lin, Natalia Gimelshein, Luca Antiga, Alban Desmaison, Andreas Kopf, Edward Yang, Zachary DeVito, Martin Raison, Alykhan Tejani, Sasank Chilamkurthy, Benoit Steiner, Lu Fang, Junjie Bai, and Soumith Chintala.
 Pytorch: An imperative style, high-performance deep learning library.
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