

# Easy and efficient spike-based Machine Learning with mlGeNN

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## ABSTRACT

Intuitive and easy to use application programming interfaces such as Keras have played a large part in the rapid acceleration of machine learning with artificial neural networks. Building on our recent works translating ANNs to SNNs and directly training classifiers with e-prop, we here present the mlGeNN interface as an easy way to define, train and test spiking neural networks on our efficient GPU based GeNN framework. We illustrate the use of mlGeNN by investigating the performance of a number of one and two layer recurrent spiking neural networks trained to recognise hand gestures from the DVS gesture dataset with the e-prop learning rule. We find that not only is mlGeNN vastly more convenient to use than the lower level PyGeNN interface, the new freedom to effortlessly and rapidly prototype different network architectures also gave us an unprecedented overview over how e-prop compares to other recently published results on the DVS gesture dataset across architectural details.

## CCS CONCEPTS

• **Computing methodologies** → *Bio-inspired approaches; Supervised learning; Vector / streaming algorithms.*

## KEYWORDS

spiking neural networks, efficient simulation, GPU, software

### ACM Reference Format:

James C Knight and Thomas Nowotny. 2022. Easy and efficient spike-based Machine Learning with mlGeNN. In *Neuro-Inspired Computational Elements Conference (NICE 2023)*, April 11–April 14, 2023, San Antonio, USA. ACM, New York, NY, USA, 6 pages. <https://doi.org/XXXXXXX.XXXXXXX>

## 1 INTRODUCTION

The development of efficient spiking neural network (SNN) simulators has been a key area of computational neuroscience research for several decades [1, 5, 9, 10, 25]. However, the prevalent SNN simulators are not well-suited to the types of models and the workflows required for spike-based machine learning (ML). Consequently, many ML researchers have chosen to build libraries [7, 8, 11, 21, 27] on top of frameworks such as PyTorch [20] and TensorFlow [6] which allow the definition of SNNs in an environment familiar to ML researchers. In these libraries, the activity of a population

of neurons is typically represented as a vector of activations and, for an SNN, this vector is populated with ones for neurons that are spiking and zeros for neurons that are quiescent. This representation allows the existing infrastructure of the underlying ML library to be used for SNNs but, as real neurons often spike at comparatively low rates, propagating the activity of inactive neurons through the network leads to many unnecessary computations. Additionally, the connections between populations of neurons can be sparse, in particular in biologically inspired SNNs. In standard ML libraries, sparse connections are typically implemented as a weight matrix containing many zeros which leads to further computational inefficiencies.

To avoid these inefficiencies we have developed mlGeNN – a new library for spike-based ML built on the GPU-optimised sparse data structures and algorithms provided by our GeNN simulator [13, 14, 25]. We previously presented an initial version of mlGeNN [23] which provided workflows for converting ANNs trained using TensorFlow [6] into SNNs that could be simulated using GeNN. However, while we found that performing inference using our converted SNNs was faster than with competing libraries, SNNs are inherently at a disadvantage for static (image) classification tasks as, by their nature, they turn a single inference step in an ANN library into SNN dynamics across several (often numerous) timesteps.

In this paper, we introduce a new Keras-inspired model description API for mlGeNN as well as functionality to train SNNs from scratch. This will allow users to tackle ML problems that are inherently more suitable for SNNs such as classification problems involving a temporal dimension, for instance speech or gesture recognition. Furthermore, it will enable them to explore the performance of recently published SNN learning rules such as e-prop [4] and EventProp [24].

To demonstrate the value of mlGeNN, we also present the results of an architecture exploration in which we use e-prop [4] to train one and two layer SNN classifier models on the DVS gesture dataset [2]. The best architectures found by our exploration achieve classification accuracies which are competitive with recent results for similar shallow architectures [22, 26]. Finally we explore the addition of sparse connectivity to our models and show that this can reduce training time by around 10× while still achieving competitive classification performance.

## 2 MLGENN

While fully describing the functionality of mlGeNN is beyond the scope of this paper, in the following sections we present an overview of some of our design decisions and the current feature set of mlGeNN. For more information about mlGeNN, we invite readers

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NICE 2023, April 11–April 14, 2023, San Antonio, USA

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ACM ISBN 978-1-4503-XXXX-X/18/06.

<https://doi.org/XXXXXXX.XXXXXXX>

to visit [https://github.com/genn-team/ml\\_genn/](https://github.com/genn-team/ml_genn/) and explore our online documentation [16].

## 2.1 Describing models

Machine learning frameworks such as PyTorch [20] and TensorFlow [6] are designed to efficiently process directed and acyclical computational graphs. In order to express recurrent connectivity within an otherwise directed acyclical graph, special recurrent layers are required to ‘hide’ the cyclical nature of the recurrent connectivity. However, when beginning to consider more brain-like models with highly recurrent connectivity, the whole model is likely to become one big recurrent layer which is not likely to be an efficient representation.

By contrast, mlGeNN allows Spiking Neural Networks (SNNs) with arbitrary topologies to be defined in terms of homogeneous groups of neurons described by **Population** objects connected together with **Connection** objects. While this type of model description is common amongst SNN simulators used for computational neuroscience, the key difference in mlGeNN is that – like in ML frameworks – the model description is agnostic to how it eventually is going to be trained. **Connection** objects don’t need to provide a model of how their weights will be trained and additional learning-rule specific structures such as feedback connections do not need to be added by hand.

All populations and connections are owned by a **Network** object which acts as a context manager. A network with two populations of neurons could simply be created like:

```
network = Network()
with network:
    a = Population("poisson_input", 100)
    b = Population("integrate_fire", 100,
                  readout="spike_count")
```

```
Connection(a, b, Dense(1.0))
```

For simplicity, in this example, built-in neuron models with default parameters are specified using strings. However, if some of the default model parameters need to be altered or a model is required that does not have default parameters or is not built into mlGeNN, bespoke neuron parametrisations and models can be specified by instantiating a class derived from **Neuron**. For example, if the user wanted the Poisson population in the example to emit positive and negative spikes for positive and negative input values and the integrate-and-fire neuron to have a higher firing threshold, **PoissonInput** and **IntegrateFire** objects could be instantiated by the user like:

```
network = Network()
with network:
    a = Population(
        PoissonInput(signed_spikes=True), 100)
    b = Population(
        IntegrateFire(v_thresh=2.0), 100,
        readout="spike_count")

    Connection(a, b, Dense(1.0))
```

By default, **Connection** objects use a ‘delta’ synapse model where the accumulated weight of incoming spikes is directly injected into neurons. However, if a somewhat more realistic model is required, e.g. where inputs are *shaped* to mimic the dynamics of real ion channels, this can be swapped out, for example with an exponential synapse with a time constant of 2 ms:

```
Connection(a, b, Dense(1.0), Exponential(2.0))
```

While the flexibility to create networks with any topology is very useful, mlGeNN also provides a **SequentialNetwork** wrapper class – inspired by **Sequential** models in Keras – for specifying common feedforward models more tersely:

```
network = SequentialNetwork()
with network:
    a = InputLayer("poisson_input", 100)
    b = Layer(Dense(1.0), "integrate_fire", 100,
              readout="spike_count")
```

Finally, in the same way that Keras can easily be extended by subclassing built-in classes and implementing new functionality using TensorFlow constructs, mlGeNN can be extended by subclassing built-in classes and providing PyGeNN model descriptions. A full overview over the model description syntax is beyond the scope of this work but is described in more detail in our documentation [16] and previous work [13]. Nonetheless, the following example illustrates how a minimal Integrate-and-Fire neuron model could be defined for use in mlGeNN:

```
genn_model = {
    "var_name_types": [("V", "scalar")],
    "sim_code": "$ (V) += $(Isyn);",
    "threshold_condition_code": "$ (V) >= 1.0",
    "reset_code": "$ (V) = 0.0;",
    "is_auto_refractory_required": False}

class IntegrateFire(Neuron):
    v = ValueDescriptor("V")

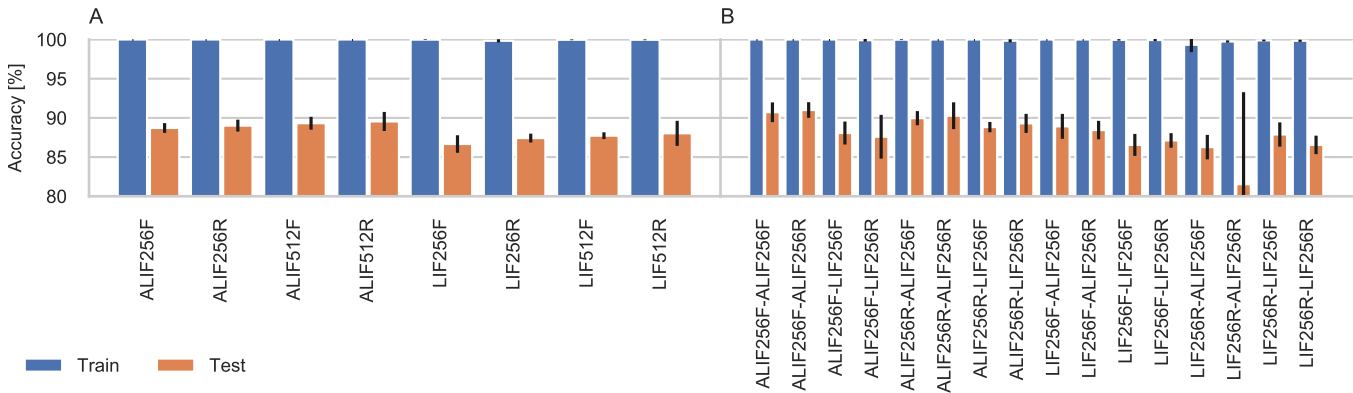
    def __init__(self, v = 0.0, softmax = False,
                 readout=None):
        super(IntegrateFire, self).__init__(
            softmax, readout)
        self.v = v

    def get_model(self, population, dt):
        return NeuronModel.from_val_descriptors(
            genn_model, "V", self, dt)
```

## 2.2 Using models for training and inference

Once a network has been defined, mlGeNN provides a range of *compilers* to produce GeNN models which can be used for training or inference. The simplest compiler is the **InferenceCompiler** which builds a GeNN model with static weights and parallel batching support for efficiently performing inference using an SNN model:

```
compiler = InferenceCompiler(
    evaluate_timesteps=1500, batch_size=512)
```



**Figure 1: Accuracy on the testing and training set of the DVS gesture dataset [2] on one (A) and two (B) layer classifiers trained with e-prop [4]. All models were trained for 100 epochs with a batch size of 512. X-axis labels describe the model architecture with the neuron model used for each layer, its size and whether it is connected in a purely feedforward (F) manner or whether it has recurrent connections (R). Bars signify the mean and error bars the standard deviation of 5 models trained with different seeds.**

```
compiled_net = compiler.compile(network)
```

The resulting `compiled_net` object can then be used to evaluate the network on a dataset:

```
metrics, _ = compiled_net.evaluate(
    {input: inputs}, {output: labels})
```

By default, the `evaluate` method calculates sparse categorical accuracy against the provided labels but mlGeNN also provides alternative metrics for regression tasks and custom metrics can easily be implemented.

Additionally, mlGeNN also provides an e-prop compiler which can be used to train models with the e-prop learning rule [4] using the implementation described in Knight and Nowotny [15]. This compiler checks the validity of the model and automatically adds the appropriate learning rules to its connections as well as feedback connections and error signal calculation logic:

```
compiler = EPropCompiler(
    example_timesteps=1500,
    losses="sparse_categorical_crossentropy",
    optimiser="adam", batch_size=512)
```

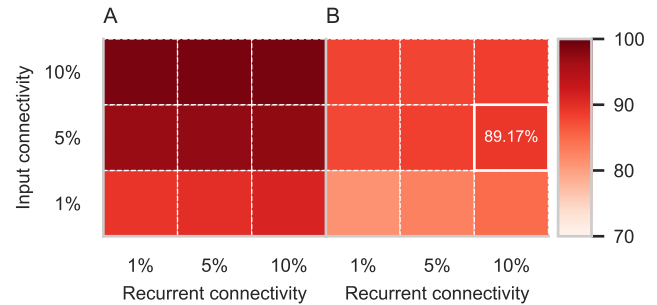
```
compiled_net = compiler.compile(network)
```

Note that here again, loss functions and optimisers with default parameters are specified using strings but can also be customised by instantiating appropriate classes. After compiling with the `EPropCompiler`, the `compiled_net` object can then be used to train the network on a dataset:

```
metrics, _ = compiled_net.train(
    {input: inputs}, {output: labels},
    num_epochs=args.num_epochs, shuffle=True)
```

## 2.3 Recording and debugging

One of the challenges of working with SNNs compared to standard ANNs is that spiking neurons are stateful, meaning that the ability to efficiently record and visualise neuronal states over time is vital

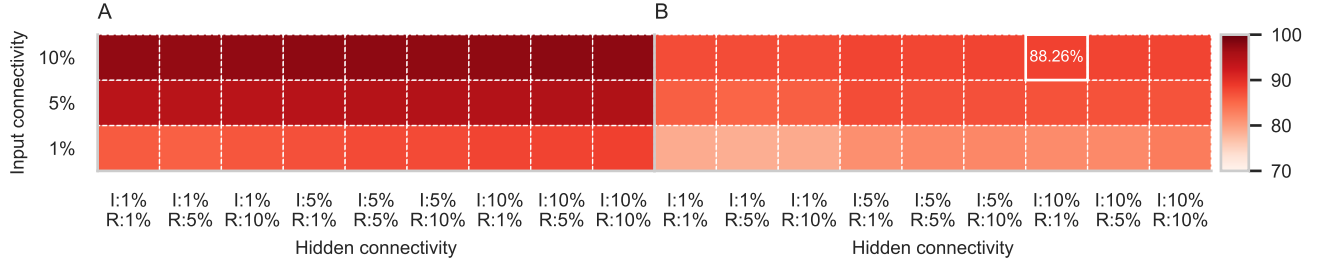


**Figure 2: Accuracy on the testing and training set of the DVS gesture dataset [2] using sparse models with ALIF512R architectures (one recurrent hidden layer of 512 adaptive LIF neurons) and varying levels of sparsity trained with e-prop [4]. All models were trained for 100 epochs with a batch size of 512. Connectivity percentages describe the fixed probability of there being a synapse between any two neurons (with no self connections in the recurrent connectivity). ‘Input connectivity’ describes connections from the input to the hidden layer. ‘Recurrent connectivity’ describes recurrent connections in the hidden layer.**

for understanding model behaviour and for debugging. Inspired by Keras, mlGeNN has a callback system which – as well as being used internally to implement progress bars, triggers processes such as weight updates during model training etc – can be used by the user to configure the recording of state variables and spikes.

The `VarRecorder` callback can be used to record a population state variable over time and the `SpikeRecorder` callback can be used to record spikes (using the efficient system described in Knight et al. [13]). For example, here we record the spikes emitted by the first layer `a` and the membrane voltages of the second layer `b` of the previous example:

```
callbacks = [SpikeRecorder(a, key="a_spikes"),
```



**Figure 3: Accuracy on the training (A) and testing (B) set of the DVS gesture dataset [2] on sparse models with ALIF256F-ALIF256R architectures (two hidden layers, one feedforward with 256 adaptive LIF neurons, one recurrent with 256 adaptive LIF neurons) and varying levels of sparsity trained with e-prop [4]. All models were trained for 100 epochs with a batch size of 512. Connectivity percentages describe the fixed probability of there being a synapse between any two neurons (with no self connections in the recurrent connectivity). ‘Input connectivity’ describes connections from the input to the first hidden layer. ‘Hidden connectivity’ describes *internal* connection from the first to the second hidden layer (‘I’) and recurrent connections in the second hidden layer (‘R’).**

```
VarRecorder(b, "v", key="b_v")
metrics, cb_data = compiled_net.evaluate(
    {input: inputs}, {output: labels},
    callbacks=callbacks)
```

The `key` strings are used to uniquely identify recorded data produced by callbacks and, after the simulation has completed, allow the user to retrieve the recorded variables from the `cb_data` dictionary returned by the `evaluate` and `train` methods of compiled models. For example, the membrane voltages emitted by all neurons during the presentation of the first input pattern could be accessed using `cb_data["b_v"][0]`.

In machine learning workflows, where models may be trained for millions of timesteps, it is also important to be able to *filter* what data is recorded to particular areas/stages of training to save memory and reduce overheads. Both `SpikeRecorder` and `VarRecorder` support filtering by neuron or example, e.g. the syntax `SpikeRecorder(a, neuron=np.s_[0::2])` could be used to record spikes from every other neuron and, in a similar vein, the command `SpikeRecorder(a, example_filter=1000)` could be used to only record during the 1000th example.

### 3 RESULTS

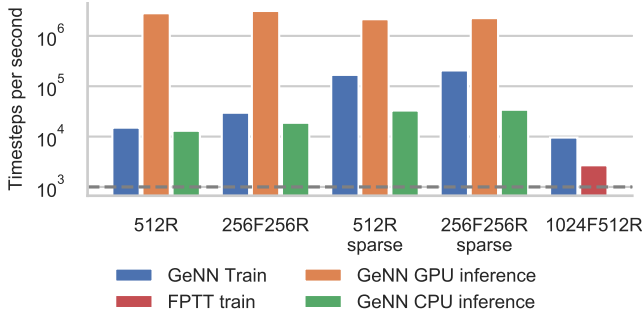
In order to demonstrate the value of mlGeNN for spike-based ML research, in the following section we present the results of an exploration of SNN architectures, trained using e-prop [4] on the DVS gesture dataset [2] – which has recently been used to evaluate EGRU [22] models and the FPTT [26] learning rule. This dataset consists of 1342 recordings of 11 different hand and arm gestures, collected from 29 subjects under three different lighting condition using a DVS 128 event-based camera [18]. We used the Tonic library [17] to access and spatially downsample the dataset to  $32 \times 32$  resolution. While mlGeNN does not depend on Tonic, it includes a `preprocess_tonic_spikes` helper function which converts spike trains from Tonic datasets into mlGeNN’s internal format. All code used to generate the results in this section is publicly available at [https://github.com/neworderofjamie/ml\\_genn\\_eprop\\_paper](https://github.com/neworderofjamie/ml_genn_eprop_paper).

#### 3.1 Accuracy

Figure 1 shows the accuracy of a wide selection of one and two layer classifier models trained with e-prop. These include configurations with and without recurrent connections and using both adaptive and standard Leaky Integrate-and-Fire (ALIF and LIF) neuron models. Of the single layer classifiers, the variant with a single layer of 512 recurrently connected ALIF neurons achieves the highest accuracy of  $(89.55 \pm 1.22) \%$  on the test set. Although this model only has a single hidden layer and around  $1.3 \times 10^6$  parameters, it out-performs a two layer EGRU model [22] with over  $4\times$  as many parameters which achieved 88.02 % accuracy.

Models using ALIF neurons also achieve higher accuracies than those using simpler LIF neurons in all of the two layer configurations. Interestingly, like the architecture employed by Yin et al. [26], the best performing model – with a test accuracy of  $(91.00 \pm 1.00) \%$  – has a feedforward first layer followed by a recurrent second layer rather than two recurrent layers. While this performance is slightly lower than the  $(91.89 \pm 0.16) \%$  reported by Yin et al. [26], our model has approximately a third the number of parameters ( $0.66 \times 10^6$  vs  $1.8 \times 10^6$ ).

One of the significant advantages of mlGeNN is that it can exploit sparse connectivity to reduce training and inference time so, in the final part of our architectural exploration, we explore the effect of sparsity on the performance of our two best performing architectures (ALIF512R and ALIF256F-ALIF256R). The results of this exploration are shown in figures 2 and 3 and, while increasing sparsity clearly impacts performance, sparse configurations of both architectures are still highly competitive. In fact, the best performing sparse versions of both architectures out-perform the two layer EGRU model [22], even though they have a small fraction of the number of parameters. The best-performing ALIF512 model – with 5 % connectivity in the input connections and 10 % connectivity in the recurrent connections – achieves an accuracy of  $(89.17 \pm 0.57) \%$  with  $60\times$  fewer parameters and the best-performing ALIF256F-ALIF256R model – with 10 % connectivity in both feed-forward connections



**Figure 4: Throughput of inference and training of the best four models on the DVS gesture dataset [2]. All GeNN models used ALIF neurons. The horizontal dashed line indicates the performance required to operate in real-time. Aside from for the 1024F512R architecture, GPU training was performed with a batch size of 512 and inference with a batch size of 264 (the full size of the test set). 1024F512R was trained with a batch size of 64 to match Yin et al. [26]. GPU training and inference were performed on an NVIDIA RTX A5000 GPU with 24 GB of memory. CPU inference was performed on an Intel Xeon Gold 6134 CPU. FPTT training throughput was calculated based on the reported [26] training time of 0.4 s epoch divided by the number of frames (timesteps) in each example.**

and 1 % in the recurrent connectivity – achieves  $(88.17 \pm 0.57) \%$  with 88× fewer parameters.

### 3.2 Performance

Efficiency is one of the main reasons why we believe dedicated tools such as mlGeNN are necessary for spike-based ML so it is important to investigate the computational speed of our models both for inference and training. Figure 4 compares the throughput – the rate that timesteps can be processed at – of the best performing models identified in the previous section for both training and inference. Because e-prop training time is dominated by the time-driven update of synaptic eligibility traces, the ALIF512R architecture trains slower than the ALIF256F-ALIF256R architecture as it has more parameters to update every timestep. Furthermore, the sparse ALIF512R and ALIF256F-ALIF256R architectures train 11× and 7× faster respectively than their dense counterparts – demonstrating how effectively GeNN can take advantage of sparse connectivity.

In our previous work [15], we showed that the latency and energy delay product of inference models simulated using GeNN were significantly lower than for an equivalent LSTM running on the same hardware. This was particularly stark when performing inference on CPUs – a likely scenario in edge applications – due to the significantly lower amount of computation required by SNNs compared to LSTMs. Figure 4 demonstrates that the DVS gesture classification models developed in the previous section also perform well on CPU with the dense models running more than 10× faster than real-time on a workstation CPU, suggesting they would still be capable of running in real-time on a lower-power edge CPU. Furthermore, both sparse models offered around double the throughput

on CPU compared to their dense counterparts. However, this performance improvement is not visible for GPU inference. Further analysis suggests that this is because, due to the very small size of the DVS gesture test set (264 images), the batch size that these models are simulated with is limited, which prevents the GPU from being fully occupied because sparse models have less parallelism to exploit compared to their dense counterparts.

Finally, we show the training throughput of an ALIF1024F512R model trained with a batch size of 64 to match the architecture described by Yin et al. [26]. Although e-prop requires time-driven updates of synaptic eligibility traces (rather than event-driven updates for which GeNN has a more significant algorithmic advantage) and these updates are more complex than those required for FPTT, mlGeNN provides over 3.5× the training throughput.

## 4 CONCLUSIONS

We hope that the mlGeNN library described in this paper will be as valuable to the community as it was in generating the results presented here. While it is by no means a precise measure, the simplicity and ease of using mlGeNN is illustrated by the fact that the code used for all simulations presented here was less than 300 lines whereas, when we used PyGeNN directly in our previous work [15] one 900 line model was required for training and a separate 500 line model for inference.

The e-prop learning rule is an approximation to gradient descent, both in terms of using a surrogate gradient and in neglecting some terms relating to recurrent connections. The results we have presented in this paper demonstrate that these approximations do not necessarily prevent e-prop from offering competitive performance on relatively complex datasets. However, e-prop has properties that could be seen as disadvantageous for fast and efficient training. Firstly, e-prop training requires time-driven updates of synaptic variables which dominates the time taken to train models. We have demonstrated that by using sparse connectivity, this problem can be reduced and, in initial unpublished experiments, we have found that by combining sparse connectivity with the Deep-R [3] rewiring rule, much of the performance lost due to the sparser connectivity can be recovered. Secondly and unsurprisingly for a visual task, both the original experiments on the DVS gesture dataset [2] and more recent work using the Decolle and FPTT learning rules [12, 26], showed that a convolutional SNN can achieve significantly higher performance than shallow recurrent networks. However, because the eligibility traces which drive e-prop weight updates are calculated from the product of pre and postsynaptic activity, each synapse needs individual state variables and they cannot be shared across the corresponding synapses in a convolutional layer. Neither the FPTT [26] learning rule – where the additional  $\Phi_t$  state variable only depends on past weights so can be shared – nor EventProp [24] – which does not require any additional synaptic state – have these issues so are much more suitable for training convolutional architectures. Based on the promising initial results presented by Nowotny et al. [19], we are working to develop an mlGeNN compiler for the EventProp [24] learning rule which will allow convolutional models to be trained and training times to be reduced by taking advantage of temporal sparsity through purely event-driven training.



## ACKNOWLEDGMENTS

This work was funded by the EPSRC (grant numbers EP/V052241/1 and EP/S030964/1) and the EU's Horizon 2020 program (grant agreement 945539). Compute time was provided through Gauss Centre for Supercomputing application number 21018 and EPSRC (grant number EP/T022205/1) and local GPU hardware was provided by an NVIDIA hardware grant award.

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