Larger GPU-accelerated brain simulations with procedural connectivity

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Large-scale simulations of spiking neural network models are an important tool for improving our understanding of the dynamics and ultimately the function of brains. However, even small mammals such as mice have on the order of 1×10^{12} synaptic connections which, in simulations, are each typically charaterized by at least one floating-point value. This amounts to several terabytes of data – an unrealistic memory requirement for a single desktop machine. Large models are therefore typically simulated on distributed supercomputers which is costly and limits large-scale modelling to a few privileged research groups. In this work, we describe extensions to GeNN – our Graphical Processing Unit (GPU) accelerated spiking neural network simulator – that enable it to 'procedurally' generate connectivity and synaptic weights 'on the go' as spikes are triggered, instead of storing and retrieving them from memory. We find that GPUs are well-suited to this approach because of their raw computational power which, due to memory bandwidth limitations, is often under-utilised when simulating spiking neural networks. We demonstrate the value of our approach with a recent model of the Macaque visual cortex consisting of 4.13×10^6 neurons and 24.2×10^9 synapses. Using our new method, it can be simulated on a single GPU – a significant step forward in making large-scale brain modelling accessible to many more researchers. Our results match those obtained on a supercomputer and the simulation runs 35% faster on a single high-end GPU than previously on over 1000 supercomputer nodes.

spiking neural networks | GPU | high-performance computing | brain simulation

The brain of a mouse has around 70×10^6 neurons, but this number is dwarfed by the 1×10^{12} synapses which connect them (1). In computer simulations of spiking neural networks, propagating spikes involves adding the 'weight' of synapses from each spiking presynaptic neuron to the input currents of connected postsynaptic neurons. Typically, the information describing which neurons are connected by a synapse and with what weight is generated before a simulation is run and stored in large arrays in random access memory (RAM). This creates high memory requirements for large-scale brain models, so that they can typically only be simulated on large distributed computer systems using software such as NEST (2) or NEURON (3). By careful design, these simulators can keep the memory requirements for each node constant, even when a simulation is distributed across thousands of nodes (4). However, high performance computer systems are bulky, expensive and consume a lot of power and are hence typically shared resources only accessible to a limited number of researchers for time-limited investigations.

Neuromorphic systems (5–10) take inspiration from the brain and have been developed specifically for simulating large spiking neural networks more efficiently. One particular relevant feature of the brain is that its memory elements – the synapses – are co-located with the computing elements – the neurons – throughout the entire system. In neuromorphic systems, this often translates to dedicating a large proportion of each chip to memory. However, while such on-chip memory is fast, it can only be fabricated at relatively low density so that many of these systems economize – either by reducing the maximum number of synapses per neuron to as few as 256 or by reducing the precision of the synaptic weights to 6 (10), 4 (5) or even 1 bit (6, 8). This allows some classes of spiking neural networks to be simulated very efficiently, but reducing the degree of connectivity to fit within the constraints of current neuromorphic systems inevitably changes the dynamics of brain simulations (11). Unlike most other neuromorphic systems, the SpiNNaker (7) neuromorphic supercomputer is fully programmable and combines large on-chip memory with external memories, distributed across the system, which enables real-time simulation of large-scale models (12). This is promising for the future but, due to its prototype nature, the availability of SpiNNaker hardware is limited and a physically large system is still needed even for moderately-sized simulations (9 boards for a model with around 10×10^3 neurons and 300×10^6 synapses (12)).

Modern GPUs have relatively little on-chip memory and, instead, dedicate the majority of their silicon area to arithmetic logic units (ALUs). GPUs use dedicated hardware to rapidly switch between tasks so that the latency of accessing external memory can be 'hidden' behind computation, as long as there is sufficient computation to be performed. For example, the memory latency of a typical modern GPU can be com-

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Significance Statement

Simulations are an important tool for investigating how brains work. However, in order to faithfully reproduce some of the features found in biological brains, large models are required. Simulating such models has, until now, required so much memory that it could only be done on large, expensive supercomputers. In this work, we present a new method for simulating large models that significantly reduces memory requirements. This method is particularly well-suited for use on Graphical Processing Units (GPUs), which are a common fixture in many workstations. We demonstrate that using our new method we can not only simulate a very large brain model on a single GPU, but also do so 35% faster than in previous supercomputer simulations.

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pletely hidden if each CUDA core performs approximately 10 arithmetic operations per byte of data accessed from memory. Unfortunately, propagating a spike in a spiking neural network simulation is likely to require accessing around 8 B of memory and performing many fewer than the required 80 instructions. This makes spike propagation highly memory bound. Nonetheless, we have shown in previous work (13) that, as GPUs have significantly higher total memory bandwidth than even the fastest CPU, moderately sized models of around 10×10^3 neurons and 1×10^9 synapses can be simulated on a single GPU with competitive speed and energy consumption. However, individual GPUs do not have enough memory to simulate truly large-scale brain models and, although small numbers of GPUs can be connected using the high-speed NVLink (14) interconnect, beyond such small GPU clusters, scaling will be dictated by the same communication overheads as for other CPU-based distributed systems.

In this work, we present a novel approach that converts large-scale brain simulation from a problem which is memory-bound on a GPU to one where the large amount of available computational power can be used to reduce both memory and memory bandwidth requirements and enable large-scale brain simulations on a single GPU workstation.

Results

In the following subsections, we first present two recent innovations in our GeNN simulator (15) which enable simulations of very large models on a GPU. We then demonstrate the power of the new features by simulating a recent model of the Macaque visual cortex (16) with 4.13×10^6 neurons and 24.2×10^9 synapses.

Procedural connectivity. The first crucial innovation that enables large-scale simulations on a GPU is what we call 'procedural connectivity'. In a brain simulation, neurons and synapses can be described by a variety of mathematical models but these are eventually all translated into time or event-driven update algorithms (17) which simulate their behaviour over time. Our GeNN simulator (15) uses code generation to convert neuron and synapse update algorithms – described using 'snippets' of C-like code – into CUDA code for efficient GPU simulation. Before a simulation can be run, its parameters, in particular the state variables and the synaptic connectivity, need to be initialised. Traditionally, this is done by running initialisation algorithms on the main CPU prior to the simulation. The results are stored in CPU and GPU memories and then used during the simulation. We have recently extended GeNN to use code generation from code snippets to also generate efficient, parallel code for model initialisation (13). Offloading initialisation to the GPU in this way made it around $20\times$ faster on a desktop PC (13), demonstrating that initialisation algorithms are well-suited for GPU acceleration. Here, we are going one step further. We realised that, if each synaptic connection can be re-initialised in less than the 80 operations required to hide the latency incurred by fetching its 8 B of parameter values from memory, it could be faster and vastly more memory efficient to regenerate synaptic connections on demand rather than storing them in memory. This is the concept of procedural connectivity and is applicable whenever synapses are static – plastic synapses which change their weights during a simulation will have to be simulated in the

traditional way. Although a similar approach was used by Eugene Izhikevich for simulating an extremely large thalamo-cortical model with 1×10^{11} neurons and 1×10^{15} synapses on a modest PC cluster in 2005 (18) – an incredible achievement – it has not been subsequently applied to modern hardware.

We implemented procedural connectivity in GeNN by repurposing our previously developed parallel initialisation methods. Instead of running them once for all synapses at the beginning of the simulation, we rerun the methods during the simulation for all outgoing synapses of each neuron that fires a spike and immediately use the identified connections and weights to run the post-synaptic code which calculates the effect of the spike onto other neurons. This is possible because the outgoing synaptic connections from each neuron are typically largely independent from those of other neurons as we shall see from typical examples below.

In the absence of knowledge of the exact microscopic connectivity in the brain, there are a number of typical connectivity schemes that are used in brain models. We will now discuss two typical examples and how they can be implemented efficiently on a GPU. One very common connectivity scheme is the 'fixed probability connector' which is described by a fixed probability P_{conn} that a neuron in the presynaptic population will be connected to a neuron in the postsynaptic population. In this case, the postsynaptic targets of any presynaptic neuron can be sampled from a Bernoulli process with success probability P_{conn}. One simple way of sampling from the Bernoulli process is to repeatedly draw samples from the uniform distribution $\mathrm{Unif}[0, 1]$ and generate a synapse if the sample is less than P_{conn}. However, for sparse connectivity, it is much more efficient to sample from the geometric distribution $\mathrm{Geom}[P_{\mathrm{conn}}]$ which is the distribution of the number of Bernoulli trials required to get the next success (i.e. a synapse). The geometric distribution can be sampled in constant time by inverting the cumulative density function (CDF) of the equivalent continuous distribution (the exponential distribution) to obtain $\frac{log(\mathrm{Unif}[0,1])}{log(1-P_{\mathrm{conn}})}$ (19, p499). Note, that when directly drawing from the uniform distribution, the sampling for each potential synapse is independent from any other potential synapse and all these operations could be performed in parallel. For the more efficient 'geometric sampling' employed here, the sampling for the post-synaptic targets of a presynaptic neuron must be done serially, but is still independent from the sampling for any other presynaptic neuron.

Another common scheme for defining connectivity is the 'fixed number total connector'. In this scheme the synaptic connections between two neuronal populations are characterised by a fixed total number N_{syn} of randomly placed synapses. In order to initialise this connectivity in parallel, the number of synapses that originate from each of the N_{pre} presynaptic neurons must first be calculated by sampling from the multinomial distribution $\mathrm{Mult}[N_{\mathrm{syn}}, \{P_n, P_n, \ldots, P_n\}]$, where $P_n = \frac{1}{N_{\mathrm{pre}}}$, on the host CPU up front because these numbers need to add to N_{syn} and are hence not independent. However, once the numbers of outgoing synapses are determined, the postsynaptic targets for a presynaptic neuron can be generated very efficiently in parallel by sampling from the discrete uniform distribution $\mathrm{Unif}[0, N_{\mathrm{post}}]$ where N_{post} is the size of the postsynaptic population. Note, that this can only be done because the targets of each presynaptic neuron are independent from those of any other presynaptic neuron. Where synaptic
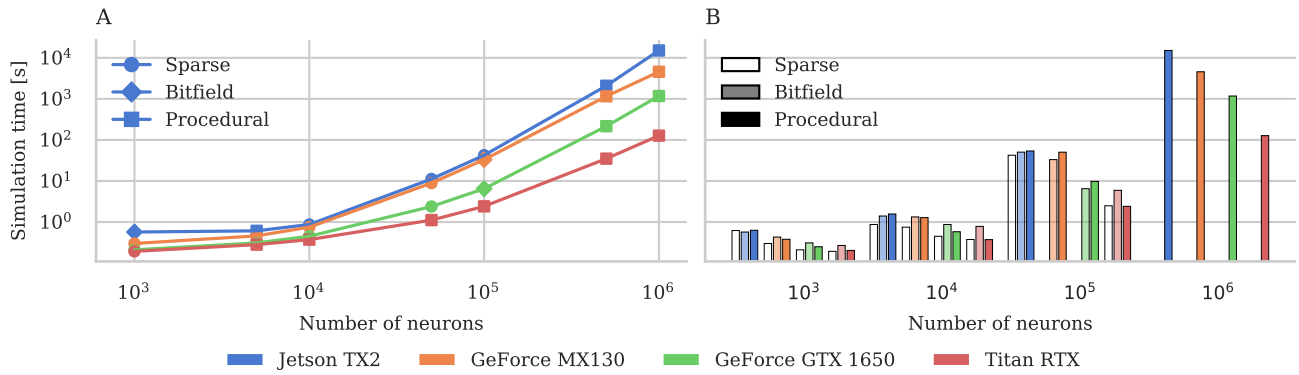
Fig. 1. Simulation time performance scaling on a range of modern GPUs (colors). **A** The best performing approach at each scale on each GPU (indicated by the symbols). For the largest models, the procedural method is always best. **B** Raw performance of each approach on each GPU. Missing bars indicate insufficient memory to simulate.

weights and delays are not constant across synapses, but are described by some statistical distribution, they can also be sampled independently from each other and hence in parallel.

In order to use these parallel initialisation schemes for procedural connectivity, we require reproducible pseudorandom numbers that can be generated independently for each presynaptic neuron. In principle this could be done with 'convential' pseudorandom number generators (PRNGs), but each presynaptic neuron would need to maintain its own PRNG state which would lead to a significant memory overhead. Instead, we use the 'counter-based' Philox4×32-10 PRNG [20]. Counter-based PRNGs are designed for parallel applications and essentially consist of a pseudo-random bijective function which takes a counter as an input (for Philox4×32-10 a 128 bit number) and outputs random numbers. In contrast to convential PRNGs, this means that generating the nth random number in a stream has exactly the same cost as generating the 'next' random number, allowing us to trivially divide up the random number stream between multiple parallel processes (in this case presynaptic neurons).

For an initial demonstration of the performance and scalability of procedural connectivity, we used a network that was initially designed to investigate signal propagation through cortical networks [21], but subsequently has been widely used as a scalable benchmark [17]. The network consists of N integrate-and-fire neurons, partitioned into $\frac{4N}{5}$ excitatory and $\frac{N}{5}$ inhibitory neurons. The two populations of neurons are connected to each other and with themselves with a fixed $P_{\mathrm{conn}} = 10\%$ connection probability.

We ran simulations of this network at scales ranging from 1×10^3 to 1×10^6 neurons (100×10^3 and 100×10^9 synapses respectively) on a representative selection of modern NVIDIA GPU hardware: Jetson TX2, a low-power embedded system with 8 GB (shared memory); Geforce MX130, a laptop GPU with 2 GB; Geforce GTX 1650, a low-end desktop GPU with 4 GB; and Titan RTX, a high-end workstation GPU with 24 GB. In Fig. 1 we compare the duration of these simulations using our new procedural approach against the standard approach of storing synaptic connections in memory using two different data structures. Both data structures are described in more detail in our previous work [13] but briefly, in the 'sparse' data structure, a presynaptic neuron's postsynaptic targets are represented as an array of indices whereas, in the 'bitfield' data structure, they are represented as a N_{post} ar-

ray of bits where a '1' at position i indicates that there is a connection to postsynaptic neuron i and a '0' that there is not. None of the devices have enough memory to store the 100×10^9 synapses required for the largest scale using either data structure but, at the 100×10^3 neuron scale, the bitfield data structure allows the model to fit into the memory of several devices it otherwise would not. However, not only is the new procedural approach the *only* way of simulating models at the largest scales but, as Fig. 1 illustrates, the performance of the precedural approach is competitive with and sometimes better than the standard approach even at smaller scales. All of the synapses in this model have the same synaptic weight meaning that they can be hard-coded into the procedural connectivity kernels. However, if weights vary across synapses, the 'bitfield' representation cannot be used and the memory constraints for 'sparse' representations become even more severe.

Kernel merging. NVIDIA GPUs are typically programmed in CUDA using a Single Instruction Multiple Thread (SIMT) paradigm where programmers write 'kernel' functions containing serial C-like code which is then executed in parallel across many virtual threads. We call our second innovation "kernel merging" and it relates to the way these kernels are implemented. While the procedural connectivity approach presented in the previous section allows us to simulate models which would otherwise not fit into the memory of a single GPU, there are additional problems when using code generation for models with a large number of neuron and synapse populations. GeNN and all other SNN simulators which use code generation to generate all of their simulation code [22] (as opposed to, for example NESTML [23], which uses code generation only to generate neuron simulation code) generate seperate pieces of code for each population of neurons and synapses. This allows optimizations such as hard-coding constant parameters and, although generating code for models with many populations will result in large code size, C++ CPU code can easily be divided between multiple modules and compiled in parallel, minimizing the effects on build time. However, GPUs can only run a small number of kernels – which are equivalent to modules in this context – simultaneously (128 on the latest NVIDIA GPUs [24, p278]). Therefore, in GeNN, multiple neuron populations are simulated within each kernel, resulting in code of the form shown in the following pseudocode which illustrates how 3 populations of 100 neurons each can
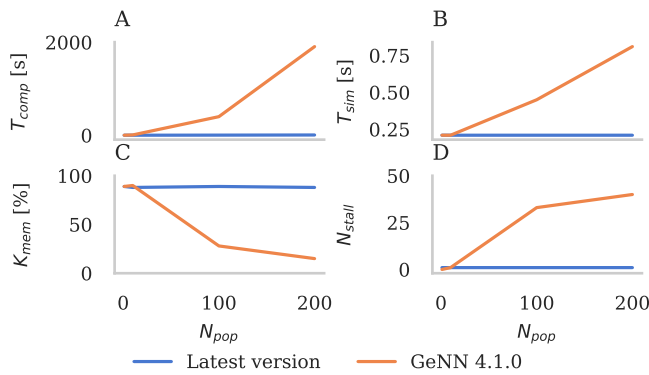
Fig. 2. Performance of a simulation of $1\,000\,000$ LIF neurons driven by a gaussian input current, partitioned into varying numbers (N_{pop}) of populations and running on a workstation equipped with a Titan RTX GPU. **A** Compilation time (T_{comp}) using GCC 7.5.0. **B** Simulation time (T_{sim}) for an $1\,\text{s}$ simulation. **C** Memory throughput (K_{mem}) reported by NVIDIA Nsight compute profiler 'Speed of light' metric. **D** Number of 'No instruction' stalls reported by NVIDIA Nsight compute profiler (N_{stall}).

be simulated in a single kernel:

```
void updateNeurons() {
  if(thread < 100) {
    // Update neuron population A
  } else if(thread >= 100 && thread < 200) {
    // Update neuron population B
  } else if(thread >= 200 && thread < 300) {
    // Update neuron population C
  }
}
```

This works well for a small number of populations but, as Fig. 2A illustrates, when we partition a model consisting of $1\,000\,000$ LIF neurons into an increasingly large number of (smaller and smaller) populations, compilation time increases super-linearly and quickly becomes impractical. Furthermore, as Fig. 2B shows, the simulation also runs much more slowly when the model is partitioned into a large number of populations. Normally, we would expect this model to be memory bound as each thread in the model reads $32\,\text{B}$ of data and, as discussed above, hiding the latency of these memory accesses would require approximately 320 arithmetic operations – many more than required to sample an input current from the normal distribution and update a LIF neuron. Fig. 2C – obtained using data from the NVIDIA Nsight compute profiler (25) – shows that this is true for small numbers of populations. In this case, the memory system is around $90\,\%$ utilised. However, when the model is partitioned into a larger number of smaller populations, the memory is used less efficiently and the kernel becomes latency bound, i.e. neither memory *nor* compute are used efficiently. Investigating further, we found that this drop in performance was accompanied by an increasing number of "No instruction" stalls (Fig. 2D) which are events that prevent the GPU from doing any work during a clock cycle. The profiler documentation suggests that these particular events are likely to be caused by "Excessively jumping across large blocks of assembly code" (25, p47), which makes sense when we are generating kernels with hundreds of thousands of lines of code. Several neural modelling tools including Brian2 (26) provide modellers with tools to work with 'slices' of neuron populations, allowing models to be defined with fewer popula-

tions. However, if a model is defined by connecting these slices together, the resulting connectivity is the result of *multiple* simple connection rules of the type discussed in the previous section, making it much more difficult to apply our procedural connectivity approach. Furthermore, such an approach places the responsibility for structuring a model in such a way that it can be simulated efficiently onto the modellers, who often prefer to concentrate on the science and organise populations according to anatomy or physiology.

To address the issue of too many populations, we developed a new code generator for GeNN which first 'merges' the model description, grouping together populations which can be simulated using the same generated code. From this merged description, structures are generated to store the pointers to state variables and parameter values which are still allowed to differ between merged populations:

```
struct NeuronUpdateGroup {
  unsigned int numNeurons;
  float* V;
};
```

An array of these structures is then declared for each merged population and each element is initialised with pointers to state variables and parameter values:

```
NeuronUpdateGroup neuronUpdateGroup[3];
neuronUpdateGroup[0] = {100, VA};
neuronUpdateGroup[1] = {100, VB};
neuronUpdateGroup[2] = {100, VC};
```

where VA is a pointer to the array containing the state variable 'V' of populations 'A' and so on. In order for a thread to determine which neuron in which population it should simulate, we generate an additional data structure – an array containing a cumulative sum of threads used for each population:

```
unsigned int startThread[3] = {0, 100, 200};
```

Each thread performs a simple binary search within this array to find the index of the neuron and population it should simulate. As Fig. 2 shows, this approach solves the observed issues with compilation time and simulation performance.

The multi-area model. Due to lack of computing power and sufficiently detailed connectivity data, previous models of the cortex have either focussed on modelling individual local microcircuits at the level of individual cells (28, 29) or modelling multiple connected areas at a higher level of abstraction (30). However, recent data (31) has shown that cortical activity has distinct features at both the global and local levels which can only be captured by modelling interconnected microcircuits at the level of individual cells. The recent multi-area model (16, 32) is an example of such multiscale modeling. It uses scaled versions of a previous, 4 layer microcircuit model (29) to implement $1\,\text{mm}^2$ 'patches' for 32 areas of the macaque visual cortex. The patches are connected together according to inter-area axon tracing data from the CoCoMac (33) database, further refined using additional anatomical data (34) and heuristics (35) to obtain estimates for the number of synapses between areas. The synapses are distributed between populations in the source and target area using layer-specific tracing data (36) and cell-type-specific dendritic densities (37). Individual populations are connected by the fixed number connectors described above. For a full
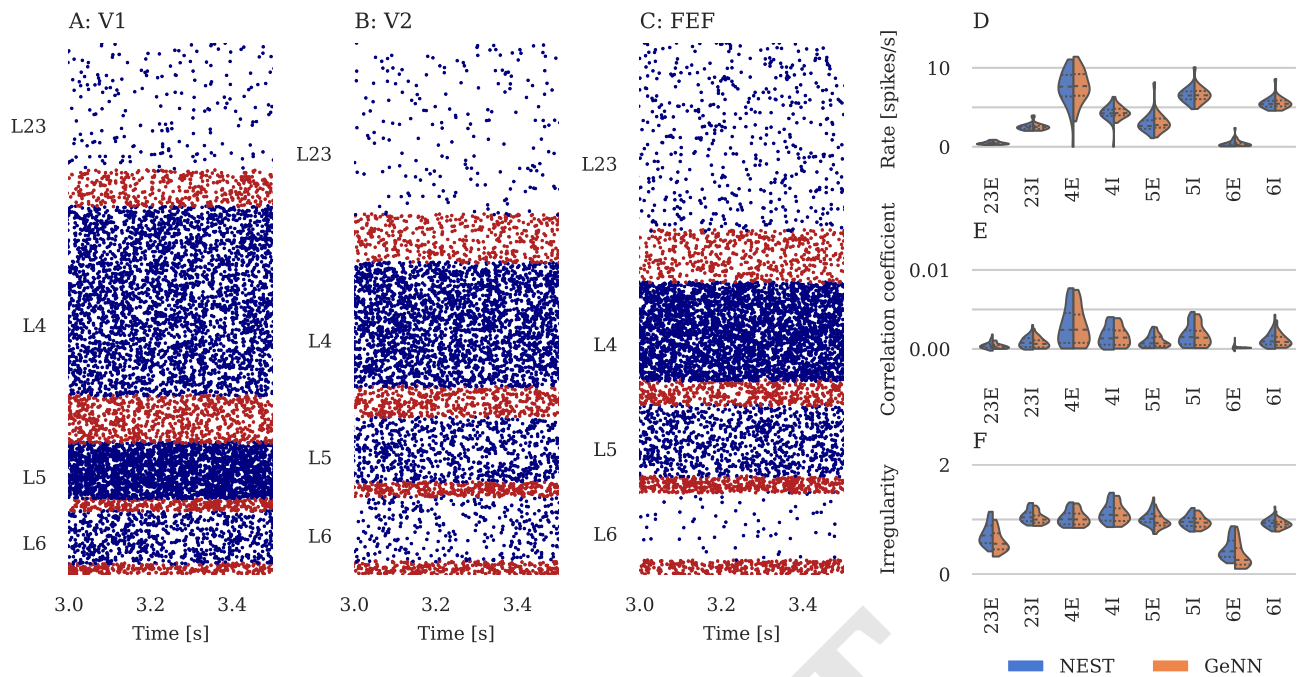
Fig. 3. Results of full-scale multi-area model simulation. **A-C** Raster plots of spiking activity of 3 % of the neurons in area V1 **A**, V2 **B**, and FEF **C**. Blue: excitatory neurons, red: inhibitory neurons. **D-F** Spiking statistics for each population across all 32 areas simulated using GeNN and NEST shown as split violin plots. Solid lines: medians, Dashed lines: Interquartile range (IQR). **D** Population-averaged firing rates. **E** Average pairwise correlation coefficients ofs piking activity. **F** Irregularity measured by revised local variation LvR (27) averaged across neurons.

description of the multi-area model please see (16, 32). In 2018, this model was simulated using NEST (2) on one rack of an IBM Blue Gene/Q supercomputer (a 2 m high enclosure containing 1024 compute nodes, weighing over 2 t and requiring around 80 kW of power). On this system, initialization of the model took around 5 min and simulating 1 s of biological time took approximately 12 min (16).

The multi-area model consists of 4.13×10^6 neurons in 254 populations and 24.2×10^9 synapses in 64 516 populations. Without kernel merging, it would therefore be unlikely that the model would compile or simulate at a workable speed using GeNN. Additionally, unlike the model we benchmarked previously, each synapse in this model has an independant weight and synaptic delay sampled from a normal distribution so that the bitfield data structure cannot be used. Even if we assume that 16 bit floating-point would provide sufficient weight precision, that delays could be expressed as 8 bit integers and that neuron populations are all small enough to be indexed using 16 bit indices, our sparse data structure would still require 5 B per synapse, such that the synaptic data would need over 100 GB of GPU memory. While a cluster of GPUs connected using NVLink could be built with this much memory, it is more than any single GPU has available. However, using procedural connectivity, we are able to simulate this model on a single workstation with a Titan RTX GPU.

In order to validate our GeNN simulations of the multi-area model, we ran a 10.5 s simulation of the model. Initialization of our model took 6 min – 3 min of which was spent generating and compiling code – and simulation of each biological second took 7.7 min – 35 % less than in the supercomputer simulation. Fig. 3A-C shows some example spike rasters from three of the modelled areas, illustrating the asynchronous irregular nature of the model's ground state. Next, we calculated the per-layer distributions of rates, spike-train irregularity and cross-correlation coefficients across all areas (disregarding the first 500 ms of simulation) and compared them to the previously published values of the same measures obtained from the supercomputer simulations. We calculated irregularity using the revised local variation LvR (27), averaged over a subsample of 2000 neurons and cross-correlation from spike histograms with 1 ms bins, calculated from a subset of 2000 non-silent neurons. The violin plots in Fig. 3D-F show the comparison of the distributions of values obtained from the two simulations – which are essentially identical.

Discussion

In this work we have presented a novel approach for large-scale brain simulation on GPU devices which entirely removes the need to store connectivity data in memory. We have shown that this approach allows us to simulate a cortical model with 4.13×10^6 neurons and 24.2×10^9 synapses (16, 32) on a single modern GPU. While this represents a significant step forward in terms of making truly large-scale brain modelling tools accesible to a large community of brain researchers, this model still has around $20\times$ fewer neurons and $40\times$ fewer synapses than the brain of even a small mammal such as a mouse (1). Our implementation of the multi-area model requires a little over 12 GB of GPU memory, with the majority (8.5 GB) being used for the circular dendritic delay buffers (described in our previous work (13)). These are a per-neuron (rather than per-synapse) data structure but, because the inter-area connections in the model have delays of up to 500 simulation timesteps (0.1 ms), the delay buffers become quite large.

One important aspect of large-scale brain simulations not

Table 1. Model parameters.

Parameter	Procedural connectivity benchmark	Merging benchmark	Multi-area model
τ_m [ms]	20	20	2
V_rest [mV]	-60.0	-70.0	-65
V_th [mV]	-50.0	-51.0	-50
R_m [MΩ]	20	20	40
τ_syn [ms]	$5/10$[1]	—	0.5
τ_ref [ms]	5	2	2
I_{ext_j} [nA]	0.55	1.00 ± 0.25	Poisson[2]
w_{ij} [nA]	$\frac{3.2}{N} / \frac{40.8}{N}$[1]	—	Various[2]

[1]Excitatory/Inhibitory.
[2]Please refer to original works (16, Table 1,2)

addressed in this work is synaptic plasticity and its role in learning. As discussed in our previous work (13), GeNN supports a wide variety of synaptic plasticity rules. In order to modify synaptic weights, they need to be stored in memory rather than generated procedurally. However, connectivity could still be generated procedurally, potentially halving the memory requirements of models with synaptic plasticity. This would be sufficient for synaptic plasticity rules that only require access to presynaptic spikes and postsynaptic neuron states (38, 39) but, for many Spike-Timing-Dependent Plasticity (STDP) rules, access to *postsynaptic* spikes is also required. GeNN supports such rules by automatically generating a lookup table structure (see our previous work (13)). While this process could be adapted to generate a lookup table from procedural connectivity, this would further erode memory savings. However, typically not all synapses in a simulation are plastic and those that are not could be simulated fully procedurally.

In this work, we have discussed the idea of procedural connectivity in the context of GPU hardware but, we believe that there is also potential for developing new types of neuromorphic hardware built from the ground up for procedural connectivity. Key components such as the random number generator could be implemented directly in hardware leading to truly game-changing compute time improvements.

Materials and Methods

In all experiments presented in this work, neurons are modelled as leaky integrate-and-fire (LIF) units with the parameters listed in Table 1. The membrane voltage V_i of neuron i is modelled as

$$\tau_\mathrm{m}\frac{dV_i}{dt} = (V_i - V_\mathrm{rest}) + R_\mathrm{m}(I_{\mathrm{syn}_j} + I_{\mathrm{ext}_j}), \qquad [1]$$

where τ_m and R_m represent the time constant and resistance of the neuron's cell membrane, V_rest defines the resting potential, I_{syn_j} represents the synaptic input current and I_{ext_j} represents an external input current. When the membrane voltage crosses a threshold V_th a spike is emitted, the membrane voltage is reset to V_rest and updating of V is suspended for a refractory period τ_ref. In the models where there are synaptic connections, pre-synaptic spikes lead to exponentially-decaying input currents I_{syn_j}

$$\tau_\mathrm{syn}\frac{dI_{\mathrm{syn}_i}}{dt} = -I_{\mathrm{syn}_i} + \sum_{i=0}^{n} w_{ij} \sum_{t_j} \delta(t - t_j), \qquad [2]$$

where τ_syn represents the decay time constant and t_j are the arrival times of incoming spikes from n presynaptic neurons. The continuous terms of the Eq. 1 and 2 are seperately solved algebraically so that the synaptic input current I_{in_i} is treated as a constant throughout each simulation timestep.

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