

Routing brain traffic through the von Neumann bottleneck: Efficient cache usage in spiking neural network simulation code on general purpose computers

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Abstract

Simulation is a third pillar next to experiment and theory in the study of complex dynamic systems such as biological neural networks. Contemporary brain-scale networks correspond to directed graphs of a few million nodes, each with an in-degree and out-degree of several thousands of edges, where nodes and edges correspond to the fundamental biological units, neurons and synapses, respectively. When considering a random graph, each node's edges are distributed across thousands of parallel processes. The activity in neuronal networks is also sparse. Each neuron occasionally transmits a brief signal, called spike, via its outgoing synapses to the corresponding target neurons. This spatial and temporal sparsity represents an inherent bottleneck for simulations on conventional computers: Fundamentally irregular memory-access patterns cause poor cache utilization. Using an established neuronal network simulation code as a reference implementation, we investigate how common techniques to recover cache performance such as software-induced prefetching and software pipelining can benefit a real-world application. The algorithmic changes reduce simulation time by up to 50%. The study exemplifies that many-core systems assigned with an intrinsically parallel computational problem can overcome the von Neumann bottleneck of conventional computer architectures.

Keywords: spiking neural networks, large-scale simulation, cache performance, distributed computing, parallel computing, memory access bottleneck

1. Introduction

Irregular access to large amounts of memory challenges the von Neumann computer architecture. Distributed applications typically make use of systems with hybrid parallelization, using message passing libraries for the communication between compute nodes and multi-threading to employ the computational cores in each node. In this contribution, we investigate as an extreme real-world example application simulation code for biological neural networks.

Such networks correspond to graphs. The graph representing the neurons as nodes and their contact points, called synapses, as directed edges is sparse and complex. In the mammalian brain a neuron establishes several thousands or even more than ten-thousand of incoming synapses and the number of outgoing synapses is of the same order. This corresponds to a graph, where both in-degree and out-degree of each node are of the order of 1000 or even 10,000.

The brain mantle, called cerebral cortex, contains the neuronal cell bodies. Tangential to the cortical surface, the probability of two cortical neurons to establish a contact is approximately 0.1 within a distance of one millimeter, but it declines rapidly for longer distances. Half of a neuron's outgoing connections are not local but target neurons at distant locations forming a hierarchically organized architecture [for an example see 1]. Due to the sheer number of neurons in the brain, the probability of any pair sharing an edge is vanishingly small.

The interaction between neurons is mediated through synapses by point-like events, called spikes. Spike events are sparse in time given that neurons emit a single or few spikes per second while the time constants of single-neuron dynamics are in the range of milliseconds and also behavior is organized by the brain on a sub-second time scale. The time required to decide whether an image contains a living object is 180 ms for monkeys and 270 ms for humans, the interval between eye movements is about 250 ms, and humans utter about two words per second. Neuroscientists hypothesize that sparse, distributed brain activity supports this systems-level behavior.

Synapses outnumber the neurons by three to four orders of magnitude and thereby consume a significant part of com-

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puter memory in simulations of spiking neural networks. The strength of the interaction mediated by a synapse can change over time depending on the activities of the presynaptic and the postsynaptic neurons and third factors like neuromodulators and the membrane potential of the postsynaptic neuron. This dynamics, called synaptic plasticity, is a key mechanism of system-level learning. For the purpose of this investigation we assume that each synapse maintains a state variable representing the coupling strength, which we refer to as synaptic weight, but plastic processes are not considered. Moreover, synaptic transmission of spikes entails a delay, which is the time interval between the presynaptic neuron emitting a spike and the spike taking effect on the postsynaptic neuron. Depending on the spatial distance between presynaptic and postsynaptic neuron the delay can be shorter than 0.1 ms or longer than 10.0 ms. This study considers homogeneous delays of 1.5 ms.

The spikes emitted by model neurons represent the sharp voltage transients of biological neurons, called action potentials. Models describing neuronal networks at the level of resolution of neurons and synapses represent individual neurons by a small system of differential equations. Often the system is linear and all non-linearity is condensed in a threshold operation on the state vector generating the point-like event. In our case the subthreshold dynamics can be integrated exactly, limiting the workload in terms of floating point operations.

Over the past two decades simulation tools in computational neuroscience have increasingly embraced a conceptual separation of generic simulation engines and specific models of neuronal networks [2]. Many different models can thus be simulated with the same simulation engine. This enables the community to separate the life cycle of a simulation engine from those of specific models and to maintain and further develop simulation engines as an infrastructure. Furthermore, the separation facilitates the cross-validation of simulation engines.

Before the dynamical state of a model of a neuronal network can be propagated an instance of the model needs to be created in computer memory. Often network models are concisely defined by probabilistic construction rules rather than explicit adjacency lists. Therefore, in simulations of neuronal networks we distinguish between the phase of network construction and the actual simulation phase, where state propagation takes place. The former is a research topic on its own [3]. The present work concentrates on the simulation phase. While network construction may take relevant amounts of wall-clock time, the simulation time scales with the biological time span to be covered by the model, but network construction does not. Propagating the dynamical state of the network in time involves three repeating phases [4, 5]. The first, termed update, advances the state of the neurons by a time interval corresponding to the minimal synaptic delay in the system, where even smaller update steps at the level of individual neurons are possible. The second, communication, is concerned with distributing the spikes that have occurred in this time interval to the compute nodes and threads hosting the respective target neurons. The subsequent spike-delivery phase routes the spikes arriving at a compute node via the representations of the corresponding synapses to their target neurons. Our investigation concentrates

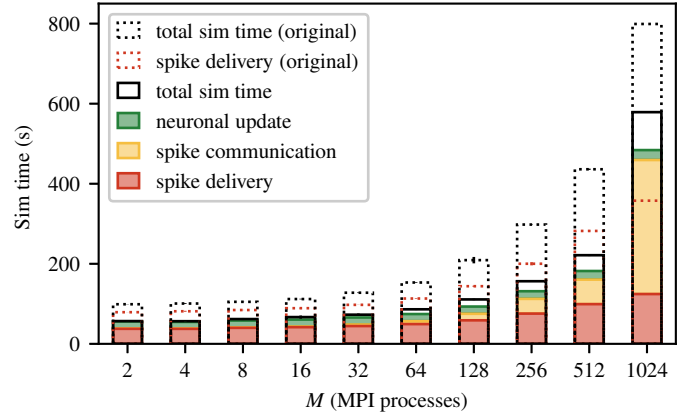


Figure 1: Contributions to simulation time (sim time) for spiking neural network simulations compared to the situation prior to optimization. Weak-scaling experiment running 2 MPI processes per compute node and 12 threads per MPI process, with a workload of 125,000 neurons per MPI process (network model see Section 2.2). The network dynamics is simulated for 1 s of biological time; spikes are communicated across MPI processes every 1.5 ms. Solid bars show total sim time (black outline), time spent on spike delivery (red), communication of spike data (yellow), and neuronal update (green) after refactoring [9] and using a combination of the optimizations of spike delivery discussed in this study (**bwRB***, **bwTS**). Dashed bars show total sim time (black dashed) and time spent on spike delivery (red dashed) of the original code (**Algorithm ORI** in [9]); original communication and update times (Figure 3 of [9]) omitted for clarity. Error bars (visible only at 128 MPI processes for the original sim time) indicate standard deviation over three repetitions. Timings obtained via manual instrumentation of the source code; measured on JURECA CM (Section 2.3).

on this final phase of the cycle.

The combination of irregular spiking activity and sparse connectivity, leads to a practically random memory-access pattern during spike delivery. Seemingly this is a worst case situation for the von Neumann architecture where for any computation the content of a respective memory unit has to be transported to the central processing unit and the result needs to be transported back. Other disciplines, such as graph processing [6] and main memory database systems [7, 8] suffer from frequent and unpredictable main memory access as well.

Let us consider a concrete example for illustration. In weak scaling of simulations with the same number of neurons per MPI process, spike delivery dominates simulation time independent of the number of MPI processes employed (Figure 1). In the regime from 2 to 512 MPI processes the time required for spike delivery almost quadruples (factor of 3.9). The refactoring efforts described in a technical companion paper [9] and the optimizations discussed in this article reduce the dependence of spike delivery on the number of MPI processes. Not affected by these changes to the original spike-delivery algorithm are neuronal update and communication (cf. Figure 3 of [9]). Additionally, the absolute time for neuronal update remains unchanged throughout as the number of neurons per MPI process is fixed. Beyond 512 MPI processes the relative contribution of spike delivery to simulation time drops below 50% because the time required for communication increases.

All simulation code analyzed in this study already provides optimizations reducing the number of spikes that MPI processes need to exchange in small to medium scale simulations (Sec-

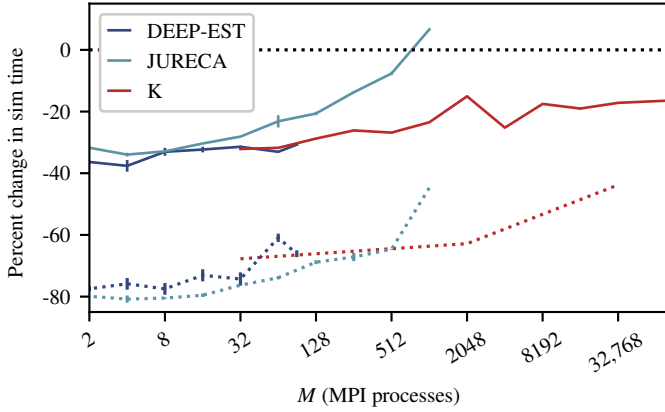


Figure 2: Relative change in simulation time after preparatory refactoring of spike-delivery algorithm (REF) as a function of the number of MPI processes M . Weak scaling of benchmark network model (Section 2.2) in linear-log representation. Same configuration as in Figure 1 for DEEP-EST CM (blue) and JURECA CM (turquoise) systems; error bars show standard deviation based on 3 repetitions. Data of K computer (red) for 1 MPI process per compute node, 8 threads per MPI process, and 18,000 neurons per MPI process (systems identified in Section 2.3). Black dotted line at zero indicates performance of original code (ORI in [9]). Dotted curves with corresponding colors indicate hypothetical limit to the decrease in sim time defined by the contribution of spike-delivery time to sim time.

tion 3.2; see Section 3.3 in 10). The code exploits the lesser degree of distribution across processes in this regime where the average number of outgoing synapses per neuron outnumbers the total number of MPI processes. This reduces both communication times and spike-delivery times, which in turn results in shorter overall simulation times (Figure 1). The present work addresses this practically relevant intermediate regime, where the effect of the optimizations gradually diminishes with the degree of distribution across processes, which impacts scalability. Nevertheless, the code scales well on modern supercomputers (see Figure 7C in 10; 5g-sort) as the optimizations reach their limit in the large-scale regime as the outgoing synapses of each neuron are almost fully distributed across processes. A technical companion paper [9] derives an analytical expression for this transition to a fully distributed network under the constraints of weak scaling.

The success of an optimization needs to be evaluated in the light of potential gain. The maximal gain by an optimization to a specific part of code is limited by the contribution of this part of code to the overall runtime. Conversely, if a different part of the code dominates the runtime, the optimization may go unnoticed. The spike-delivery algorithm that serves as a reference in this study (Section 3.2) already benefits from the refactoring described in a technical companion paper [9]. The refactoring significantly reduces simulation time (Figure 2) showing different effectiveness on the tested systems (Section 2.3). However, over the full range there is room for further improvement to spike delivery by at least the same amount.

Recently, Cremonesi et al. made major progress by deriving a detailed analytic performance model [11, 12]. The model reveals specific algorithmic and hardware bottlenecks for three classes of neuronal network models mapping out the field of

computational neuroscience. Latency of the communication network and memory bandwidth are identified as the most severe bottlenecks in large-scale simulations. The authors find that in models of the type discussed in the present study the spike-delivery phase is the most expensive one once the data do not fit into the cache hierarchy anymore. The most severe hardware constraint is the saturation of memory bandwidth which is driven by the memory-latency effect of the irregular access patterns. While the CPU is capable of issuing many memory accesses in advance to partially hide latency of processing spikes, the authors expect that two established strategies, pipelining and prefetching, cannot effectively hide the latency introduced by the non-contiguous data accesses.

Prefetching attempts to hide cache misses and memory stalls [13] by overlapping memory access with computation. In general, two types can be distinguished: hardware-induced and software-induced prefetching. The former relies on the underlying hardware to detect patterns in memory access such that the hardware can take care of getting the data into the cache just when it is needed. The latter relies on hints in the source code via prefetch instructions indicating what data should be loaded into cache. This is a promising technique when hardware predictions fail but the access pattern can still be known by the developer a priori.

In the past, several techniques for prefetching have been investigated. Especially pointer-based data structures which suffer from "pointer chasing", as for example large graphs [14, 15] and databases [16, 17, 18], profit from prefetching. The main challenge is to issue the prefetch instructions early enough such that the data is loaded timely into cache, but also late enough such that it does not clog the cache.

Recent work has focused on introducing code stages to deal with dynamic memory access and uncertainty in the number of lookups. This is achieved either by manually implementing state machines [19] or usage of coroutines [16, 17, 18]. Coroutines are resumable functions that can suspend their current execution. In both cases a prefetching instruction is inserted just before the memory stalls. After the prefetch instruction the program saves its current state and continues with other work. When entering the function the next time the required piece of memory is available and operation can resume where it left off. These techniques are promising for cases where irregularity such as variable-length pointer chains and potential early loop exits prevent usage of simpler prefetching techniques.

If the number of pointer dereferences is known ahead of time and is constant across lookups, a promising technique to employ is group prefetching [20]. Group prefetching is a loop-transformation method which breaks a single for loop into an outer and several inner loops allowing batchwise processing of code stages and critical data to be prefetched.

Another technique capable of hiding cache misses is software pipelining [21, 22, 23]. Here loops are transformed such that the instructions inside of the loop are carried out with an offset and overlapped with each other (details in Section 4.2). Thus, memory accesses and arithmetic operations inside the loop are no longer mutually dependent. As recent CPUs are superscalar, independent memory accesses and arithmetic op-

erations can be executed in parallel. This increases the number of instructions completed per cycle. As the modification of the for loop increases the number of instructions, the increase in instructions per cycle has to be greater than the increase in number of instructions to improve the overall performance.

The present study begins with a description of our setting in terms of hardware and software as well as the profiling framework (Section 2). We use the open-source community simulation code NEST (Section 2.1) to obtain performance data from a real-world application and as a framework for reference implementations of the optimization techniques discussed. Absolute performance data require a concrete neuronal network model close to the ones used in production. The network model introduced in Section 2.2 is prototypical for a wide class of models in neuroscience, it is scalable, and it has been used in a number of previous studies. A scientific community code needs to be developed and maintained over decades. Therefore it is important that new algorithms do not improve the performance on one architecture while making it impossible to adapt to the next generation of systems. Therefore, we assess the performance of a recent mainstream architecture, a common but older high-end cluster, and a dedicated supercomputer (Section 2.3). Section 3 describes the data structures representing neurons and synapses and how spikes travel through these data structures from arrival at the compute node to their ultimate delivery at the target neurons. Based on this, the subsequent two sections form the core of the investigation. Section 4 proposes new algorithms based on latency-hiding techniques indicated above for different phases of spike delivery. Section 5 presents a quantitative analysis of the effects of the new algorithms and their combination on different hardware architectures. Finally, Section 6 derives a combination of algorithms delivering a robust overall performance gain across problem sizes and hardware architectures. The study concludes by setting the results into the context of the generic problem of applications with essentially random memory access patterns and their implications for the interpretation of the von Neumann bottleneck and the design of neuromorphic computing systems.

Source code, simulation and analysis scripts are openly available [24] at Zenodo¹. The presented conceptual and algorithmic work is part of our long-term collaborative project to provide the technology for neural systems simulations [25]. Preliminary results have been presented in abstract form [26].

2. Benchmarking framework

2.1. Simulation engine

In the present work we evaluate the concepts and new algorithms in the framework of the simulation code NEST² [25], a widely used engine for spiking neuronal networks of natural density at the resolution of individual nerve cells and synapses.

NEST is an open source project governed by the public society NEST Initiative³ and a component of the ICT infrastructure⁴ created by the European Human Brain Project (HBP). The development is managed via GitHub where contributions undergo a formal code review and consistency is ensured by continuous integration using automated style checks and testing. The kernel of the simulator is written in C++ and uses *MT* coequal OpenMP [27] threads for parallelization which are arranged into *M* MPI [28] processes harboring *T* threads each. The differential equations and state transitions defining neuron and synapse models are expressed in the domain specific language NESTML [29, 30] which generates the required C++ code for dynamic loading into the simulation engine. The configuration of simulation experiments, including neuron and synapse models, network structure and recorded data, are specified interactively via Python using PyNEST [31, 32]. Neurons are distributed across parallel resources together with their incoming synapses in a round-robin fashion. This distribution implements a simple load balancing scheme as it minimizes the number of neurons from the same population, which may exhibit similar activity patterns, on the same thread. Already Morrison et al. [4] find that not only the propagation of the dynamical state of the neuronal network but also network construction needs to be parallelized to achieve sufficient performance for practical applications. More than a decade later Ippen et al. [3] improve the algorithms for multi-threading on many-core systems and point out that non-blocking memory allocation is essential for performance. However, as network creation is not in the focus, the present work stays with the system `malloc()`. The work is based on commit 059fe89 of release 2.18.

NEST uses a globally time-driven simulation scheme [4], where neurons are typically updated every 0.1 ms and spike times are constrained to this time grid. There is a biophysical delay between the emission of a spike by the source neuron and the arrival at the target neuron. Therefore, it suffices to exchange spike data between threads in intervals of the minimal delay in the neuronal network [5]. Consequently, the simulation cycle propagating the dynamical state of the network divides into three phases: update neurons, communicate spikes between threads, and deliver spikes to target neurons including the propagation of synaptic dynamics (see Figure 1). Different neuron models require solvers of different computational load ranging from precalculated exact propagator matrices for linear neuron models [33] to generic solvers for non-linear differential equations with adaptive time-stepping. Similarly the workload of synapses depends on the chosen model. Static synapses are stateless whereas for plastic synapses the state may depend on the activities of the pre- and the postsynaptic neuron [34]. Over the years the scalability of NEST has been demonstrated on a range of supercomputers [35, 36, 37, 10]. Recent revisions of the code employ `MPI_Alltoall` to send spikes only to MPI ranks where they have targets [10].

¹<https://www.zenodo.org>

²<https://www.nest-simulator.org>

³<https://www.nest-initiative.org>

⁴<https://www.ebrains.eu>

2.2. Network model

To measure and compare the suggested algorithmic improvements we use a balanced random network model [38] as a benchmark similar to the one used in previous studies on neuronal network simulation technology [39, 35, 36, 37, 3, 40, 10]. The parameters for this benchmark model are specified in the parameter tables 1, 2 and 3 in Jordan et al. [10]. Note that in contrast to previous studies, where excitatory-excitatory connections exhibited spike-timing dependent plasticity, the model considered here uses static synapses, which do not exhibit any dynamics and consequently have a fixed weight.

We consider the model a scalable version of a typical neuronal network simulation as the neuronal activity exhibits an asynchronous irregular spike pattern and does not depend significantly on the model’s size. Furthermore, the random connectivity of the network model represents a worst-case scenario in terms of network structure: Local connection patterns cannot, not even in principle, be exploited by representing subnetworks on a subset of available nodes as a single neuron connects with equal probability with any other neuron in the network.

All measurements of runtime in this study refer to the actual simulation time, or in short “sim time”, where the network state is propagated; measurements of network-construction time and initialization time are not part of this study.

2.3. Systems

The detailed specifications of the computer systems are given in [9]. JURECA CM [41] consists of 1872 compute nodes (dual Intel Xeon E5-2680 v3 Haswell 12-core CPUs at 2.5 GHz), DEEP-EST⁵ has 50 nodes (dual Intel Xeon Gold 6146 Skylake 12-core CPUs at 3.2 GHz), and the K computer [42] houses 82,944 nodes (8-core Fujitsu SPARC64 VIIIfx CPU at 2 GHz).

2.4. Measurements of runtime and profiling

For systematic benchmarking we rely on the Jülich Benchmarking Environment (JUBE)⁶[43], which is a software suite actively developed by the Jülich Supercomputing Centre. For measuring the time consumption of different parts of the code we rely on the Stopwatch class distributed with NEST. This class acts as a wrapper around `gettimeofday()` which is part of the header `<sys/time.h>` of the C POSIX library⁷.

We use the Microarchitecture Exploration analysis mode of the Intel VTune Profiler⁸, which provides detailed information on hardware usage. By specifying the option `uarch-exploration`, the running process is periodically interrupted enabling the sampling of hardware events from the processor. These events are used for calculating predefined ratios, which are reported once the program has finished. The study collects data only for the first 64 MPI processes and restricts measurements to the spike-delivery phase. Specifically,

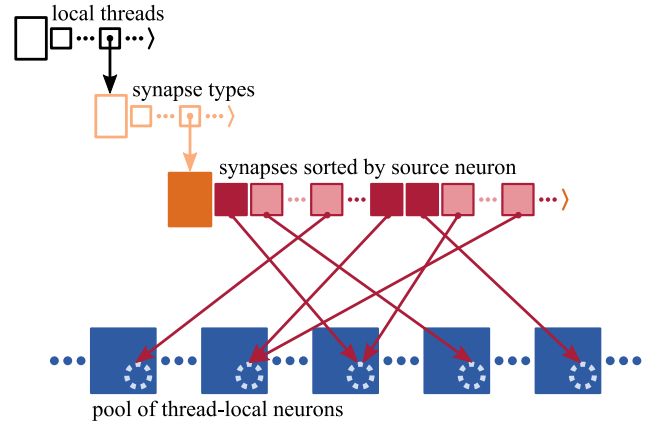


Figure 3: Memory layout of synapses and neurons on each MPI process. Each process stores the local synapses (pink filled squares) in a three-dimensional resizable array sorted by hosting thread and synapse type. At the innermost level, synapses are arranged in source-specific target segments (dark pink: first synapse; light pink: subsequent targets); only one innermost array shown for simplicity. Target neurons (blue filled squares) are stored in neuron-type and thread-specific memory pools; only one pool shown for simplicity. Each neuron maintains a spike ring buffer (dotted light blue circles). Synapses have access to their target neurons’ spike ring buffers through pointers (dark pink arrows). Adapted from Figure 1 of Pronold et al. [9].

we focus on the Clockticks per Instructions Retired (CPI) event ratio. The CPI measure is calculated via dividing the number of unhalted processor cycles (clockticks) by the number of instructions retired. CPI indicates to what extent latency affects the application’s execution, with smaller values corresponding to smaller latencies.

3. Memory access during spike delivery

In spiking neuronal network simulation code the temporally sparse event-based communication between neurons presents a challenging performance bottleneck for modern architectures optimized for dense data but hence also an optimization opportunity. Due to the connection data structures and the spike-delivery algorithm of our reference implementation NEST (Section 2.1), delivery of spikes to their targets involves frequent access to essentially random memory locations. Therefore, memory access is difficult to predict automatically, leading to long data access times due to ineffective use of caches. The following two sections describe the connection data structures and the spike-delivery algorithm that serves as a starting point for the latency-hiding techniques investigated in this study (Section 4) and as a reference in the quantitative analysis (Section 5). This reference algorithm is a result of the preparatory refactoring of the original data structures and algorithm presented in Pronold et al. [9], which already achieves a significant reduction in simulation time (Figure 2).

3.1. Memory layout of synapses and neurons

Each synapse is represented on the same MPI process and thread as its target neuron [4], where a model synapse has a

⁵<https://www.deep-projects.eu>

⁶<https://www.fz-juelich.de/jsc/jube>

⁷<https://pubs.opengroup.org/onlinepubs/9699919799/idx/head.html>

⁸<https://software.intel.com/vtune>

memory footprint of few tens of Bytes, while a model neuron easily consumes a Kilobyte or more. Each MPI process makes use of a three-dimensional resizable array to store the process-local synapses sorted by hosting thread and synapse type (Figure 3). The data structure takes into account that each neuron typically connects to many target neurons (out-degree): In the innermost arrays, synapses are sorted by source neuron and thereby arranged in target segments, each consisting of at least one target synapse potentially followed by subsequent targets. The number of unique source neurons and hence the average length of target segments depends on the distribution of synapses across MPI processes and threads as well as the number of synapse types. In the limit of sparsity where the number of neurons in the network exceeds the number of thread-local synapses, the length of target segments approaches one.

In order to account for synaptic transmission delays each neuron maintains a ring buffer accommodating the incoming spikes until they are due to take effect on the neuronal dynamics [4]. In the reference algorithm, a synapse has access to its target neuron’s spike ring buffer through a pointer [cf. Section 4.2, 9].

Neurons of the same type that are also hosted by the same thread belong to the same memory pool consisting of multiple chunks that allow for contiguous storage of many objects. As synapses from many different source neurons converge on a target neuron (in-degree), the memory locations of target neurons cannot be ordered according to source neurons and are hence independent of the order of synapses in the target segments.

3.2. Spike-delivery algorithm

The distributed time-driven simulation of spiking neuronal networks proceeds in a cycle of updates to all neurons, communication of all recent spikes across MPI processes, and delivery of the spikes to their process-local synaptic and neuronal targets. After every spike communication, each MPI process holds a receive buffer filled with spike data that need to be dispatched to the process-local targets. Each spike entry addresses an entire target segment of synapses (Section 3.1). To this end, the spike entry needs to be able to locate the beginning of the target segment within the three-dimensional data structure storing the process-local synapses (Figure 3), i.e. the first synapse of the target segment. Therefore, each spike entry conveys identifiers for the hosting thread and the type of the synapse, as well as the synapse’s index within the innermost resizable array.

In the reference algorithm, the delivery of spikes from the MPI receive buffer to the local targets hosted by different threads is a two-step process allowing for an entirely thread-parallel delivery with a single synchronization point [cf. Section 4.1, 9]. First, the threads sort the spike entries by hosting thread and synapse type in parallel using a dedicated intermediate data structure, called spike-receive register, only then the threads dispatch the spikes, now exclusively reading relevant entries.

Starting with the first synaptic target, the hosting thread subsequently processes all synapses of a spike’s target segment. The number of spike entries in the receive buffer depends on the degree of distribution of synapses across MPI processes and threads as well as the number of synapse types. Arranging

bwRB*: Delivery of spikes in batches of size B_{RB} , including group prefetching of spike ring buffers. **TS** marks iteration over a synaptic target segment. **SYN** marks access to an individual target synapse; **RB** marks access to the spike ring buffer of the corresponding target neuron; **RB*** marks group prefetching prior to access. Based on [REF](#).

Data: *spike_reg, synapses*

create arrays *target_rb, delay, weight* of size B_{RB}

$i \leftarrow 0$

foreach *spike* **in** *spike_reg* **do**

$l_{cid} \leftarrow \text{spike}.l_{cid}$

$subsq \leftarrow \text{true}$

while $subsq$ **do**

TS

SYN

$(subsq, \text{target_rb}[i], \text{delay}[i], \text{weight}[i]) \leftarrow \text{synapses}[l_{cid}].\text{Send}()$

$i \leftarrow i + 1$

$l_{cid} \leftarrow l_{cid} + 1$

if $i == B_{RB}$ **then**

RB*

for $i \leftarrow 0$ **to** $B_{RB} - 1$ **do**

 prefetch *target_rb*[i]

RB

for $i \leftarrow 0$ **to** $B_{RB} - 1$ **do**

$\text{target_rb}[i].\text{AddValue}(\text{delay}[i], \text{weight}[i])$

$i \leftarrow 0$

process remaining entries in *target_rb*

synapses in source-specific target segments such that only one spike needs to be communicated to address an entire segment is an effective optimization for small to medium-scale simulations (see Section 3.3 in 10). However, when increasing the network size in a weak-scaling experiment, ever more source neurons have ever fewer thread-local targets, or conversely, ever more thread-local synapses originate from different source neurons such that they can no longer be combined into long target segments addressed by a single source-specific spike entry. This is the limit of sparsity discussed in Section 3.1.

The synapse object stores all information relevant for the subsequent delivery to the target neuron, foremost a pointer to the target neuron’s spike ring buffer. During delivery of the spike to the target neuron the algorithm retrieves the pointer from the synapse as well as synaptic properties such as delay and weight, which define time and amplitude of the spike’s impact on the neuron, respectively. Taking into account the delay, the algorithm adds the weight of the incoming spike to the correct position in the neuronal spike ring buffer.

For each relevant spike entry the hosting thread accesses target synapses and neurons in an alternating fashion. All target synapses of a spike entry are in contiguous locations in memory as they are part of the same target segment, but the corresponding target neurons are in nonadjacent memory locations. Furthermore, with every spike entry, the thread proceeds to a different synaptic target segment, most likely not in

REF: Reference algorithm delivering spikes to local targets. **TS** marks iteration over a synaptic target segment. **SYN** marks access to an individual target synapse; **RB** marks access to the spike ring buffer of the corresponding target neuron. Variables typeset in italics, functions in typewriter. See Pronold et al. [9] for a more detailed presentation.

Data: *spike_reg, synapses*

```

foreach spike in spike_reg do
  lcid  $\leftarrow$  spike.lcid
  subsqs  $\leftarrow$  true
  while subsqs do
    (subsqs, target_rb, delay, weight)  $\leftarrow$ 
      synapses[lcid].Send()
    lcid  $\leftarrow$  lcid + 1
    target_rb.AddValue(delay, weight)

```

a proximate memory location. With increasing sparsity of the network in a weak-scaling experiment, such switches between target segments become more frequent as ever more spike entries need to be delivered to ever shorter target segments. In the sparse limit, both accessing target synapses and the corresponding target neurons requires the hosting thread to jump to random memory locations. The pseudocode of the reference algorithm (REF) presents this memory bottleneck of the spike-delivery algorithm in an abstract way omitting intricacies caused by support for multi-threading and different synapse types.

4. Latency-hiding techniques

Based on the reference algorithm (REF), we investigate three techniques to hide memory fetch latency and reduce the number of cache misses during spike delivery. To assist the reader in following the algorithmic changes, we provide pseudocode for the reference algorithm (REF), each of the potential adaptations (**bwRB***, **lagRB**, and **bwTS**), and a combination of two adaptations (**bwTSRB***). The algorithms **bwRB*** and **lagRB** encode competing adaptations of the reference algorithm REF, whereas **bwTS** encodes an adaptation that can be combined with either of the former. The pseudocode is reduced to the essential elements of the spike delivery; in particular, we apply the following simplifications: the code does not take into account multiple threads or synapse types and the initial thread-parallel transfer of spike entries from the MPI receive buffer to the spike-receive register is omitted (Section 3.2). The performance of the algorithms is analyzed in Section 5.

The reference algorithm has access to a resizable array of thread-local *synapses* and to a spike register (*spike_reg*), which contains all spike entries that need to be delivered [see Section 4.1 of 9, for details]. For each spike entry, the location of the first target synapse is extracted and assigned to the variable *lcid*, which is then used in the enclosed while loop to iterate over the spike's entire synaptic target segment within the *synapses* array (**TS**). Each synapse stores an indicator (*subsqs*)

lagRB: Delivery of spikes introducing a lagged access to ring buffers with respect to the corresponding target synapses; lag is B_{RB} . **TS** marks iteration over a synaptic target segment. **SYN** marks access to an individual target synapse; **RB** marks access to the ring buffer of the target neuron. Based on REF.

Data: *spike_reg, synapses*

```

create arrays target_rb, delay, weight of size  $B_{RB} + 1$ 
i  $\leftarrow$  0
j  $\leftarrow$  0
foreach spike in spike_reg do
  lcid  $\leftarrow$  spike.lcid
  is_init  $\leftarrow$  true
  subsqs  $\leftarrow$  true
  while subsqs do
    (subsqs, target_rb[i], delay[i], weight[i])  $\leftarrow$ 
      synapses[lcid].Send()
    i  $\leftarrow$  i + 1
    if i ==  $B_{RB} + 1$  then
      i  $\leftarrow$  0
    lcid  $\leftarrow$  lcid + 1
    if is_init and i ==  $B_{RB}$  then
      is_init  $\leftarrow$  false
    else
      target_rb[j].AddValue(delay[j], weight[j])
      j  $\leftarrow$  j + 1
      if j ==  $B_{RB} + 1$  then
        j  $\leftarrow$  0
  process remaining entries in target_rb

```

of whether the target segment continues or not. To deliver a spike to the target synapse at position *lcid*, the synapse member function Send() is called on *synapses*[*lcid*] returning the indicator *subsqs*, the pointer to the spike ring buffer, and the synaptic delay and weight (**SYN**). The pointer is then used to call AddValue(), a member function of the neuronal spike ring buffer requiring the delay and the weight (**RB**). Taking into account the delay, the spike ring buffer adds the weight of the incoming spike to the correct position. This implements the delivery of the spike to the target neuron. Note the dependency of the algorithmic step **RB** on step **SYN**, which is readily visible as a result of the prior refactoring [Section 4.2 of 9].

4.1. Batchwise access to spike ring buffers

We next consider a loop-transformation method that allows for group prefetching [20]. The method breaks a single for loop of L iterations containing several code stages into an outer and several inner loops of size L_B and B , respectively, where $L = L_B B$. For example, an original loop $\{X_i Y_i\}^L$ over two operations X_i and Y_i depending on running index i is transformed into $\{X_{(j-1)B+1} \dots X_{jB} Y_{(j-1)B+1} \dots Y_{jB}\}^{L_B}$. The code stages of the original for loop are hence processed in a batchwise manner. All code stages are handled B times before moving to the next

step of the outer loop. By introducing code stages of size B , a prior loop prefetching all critical data can be inserted, making use of memory-level parallelism.

We adapt the reference algorithm such that access to synapses and access to spike ring buffers is carried out in batches of size B_{RB} (**bwRB***). First, B_{RB} synapses are accessed to retrieve the pointers to the corresponding target neurons' spike ring buffers as well as the synaptic weights and delays (**SYN**). The retrieved synaptic information is temporarily buffered in three auxiliary arrays (*target_rb*, *delay*, *weight*) of size B_{RB} . Each time B_{RB} synapses have been accessed such that the auxiliary buffers are filled, batchwise access to the B_{RB} collected target spike ring buffers is triggered to add the corresponding weights to the correct buffer positions (**RB**). An optional prior step is the software-induced prefetching of the B_{RB} spike ring buffers into the cache (**RB***). To this end, in the NEST implementation (Section 2.1) we make use of the GCC built-in function `__builtin_prefetch(const void *addr, ...)`⁹; the prefetching can be enabled at compile time. The method is referred to as group prefetching as it consecutively prefetches data from an entire batch of B_{RB} individual memory addresses before operating on the group of data instead of using a per-memory-address approach alternating between prefetching and processing. In the quantitative analysis (Section 5) we refer to this set of optimizations as either **bwRB** or **bwRB***, where the asterisk indicates prefetching.

The batchwise progression is agnostic with regard to the boundaries of synaptic target segments (Section 3.1), which means that in case of short target segments, it can take several iterations of the for loop over spike entries to process B_{RB} synaptic targets. Leaving and re-entering of the enclosed while loop traversing the synaptic target segment of a specific spike entry (**TS**) does not affect the progression. Eventually, all spike entries in the register are processed and either delivered to their synaptic targets or added to the auxiliary array. After the loop over the spike register exits the algorithm delivers the remaining entries in the auxiliary arrays to their targets.

4.2. Lagging access to spike ring buffers

This optimization exploits the idea of software pipelining for spike delivery. In software pipelining [21, 22, 23] loops are reformed in such a way that the instructions inside of the loop are carried out with an offset of $B \geq 1$ and overlapped with each other. For example, an original loop of L iterations $\{X_i Y_i\}^L$ over two operations X_i and Y_i depending on index i is transformed into $\{X_i\}^B \{X_i Y_{i-B}\}_{B+1}^L \{Y_i\}_{L-B+1}^L$. By doing so, the operation X inside of the central loop is from a different iteration than Y .

We adapt the reference algorithm such that access to synapses and access to spike ring buffers is still carried out in an alternating fashion but algorithmic progression is always B_{RB} synapses ahead of spike ring buffers (**lagRB**). This means when a ring buffer is accessed the synaptic information and the respective pointers for the next B_{RB} spike ring buffers are already available in local arrays. To this end, B_{RB} synapses

bwTS: Delivery of spikes with processing of spike entries and corresponding target segments in batches of size B_{TS} . **TS** marks iteration over a synaptic target segment using a fixed count loop according to the target-segment size ts_size . **SYN** marks access to a target synapse; **RB** marks access to the spike ring buffer of the corresponding target neuron. Based on [REF](#).

Data: *spike_reg*, *synapses*

create arrays *lcid*, *ts_size* of size B_{TS}

$l \leftarrow 0$

repeat *spike_reg*.Size() / B_{TS} **times**

for $k \leftarrow 0$ to $B_{TS} - 1$ **do**

$lcid[k] \leftarrow spike_reg[l + k].lcid$

$l \leftarrow l + B_{TS}$

for $k \leftarrow 0$ to $B_{TS} - 1$ **do**

$ts_size[k] \leftarrow synapses[lcid[k]].GetTSSize()$

for $k \leftarrow 0$ to $B_{TS} - 1$ **do**

repeat $ts_size[k]$ **times**

 (*target_rb*, *delay*, *weight*) $\leftarrow$

$synapses[lcid[k]].Send()$

$lcid[k] \leftarrow lcid[k] + 1$

$target_rb.AddValue(delay, weight)$

TS
SYN

RB

process remaining entries in *spike_reg*

are initially accessed to retrieve the pointers to the corresponding target neurons' spike ring buffers as well as the synaptic weights and delays (**SYN**). The retrieved synaptic information is temporarily buffered in auxiliary arrays (*target_rb*, *delay*, *weight*) of size $B_{RB} + 1$. Once B_{RB} synapses have been accessed such that the auxiliary buffers are almost full (if-condition is met), the initialization phase ends and alternating access to synaptic targets (**SYN**) and neuronal spike ring buffers (**RB**, else-clause) starts, using the auxiliary arrays as ring buffers and the indices i and j for writing to the arrays and reading from the arrays, respectively. In the quantitative analysis (Section 5) we refer to this optimization as **lagRB**.

As for **bwRB***, leaving and re-entering of the enclosed while loop traversing the synaptic target segment of a specific spike entry (**TS**) does not interrupt the cyclic processing of the auxiliary arrays. Eventually, all spike entries in the register are processed and delivered to their synaptic targets, but the auxiliary arrays still contain synaptic information that needs to be delivered to the spike ring buffers.

4.3. Batchwise access to target segments

The previous two algorithms, **bwRB*** and **lagRB**, are concerned with disentangling access to synapses and neuronal spike ring buffers. They implicitly make hopping from one synaptic target segment to the next more seamless as they progress writing and reading the auxiliary arrays (*target_rb*, *delay*, *weight*) irrespective of the boundaries of single target segments. The optimizations of the present algorithm, **bwTS**, directly target the iteration over spike entries in the spike

⁹<https://gcc.gnu.org/onlinedocs/gcc/Other-Builtins.html>

bwTSRB*: Delivery of spikes combining processing of spike entries and corresponding target segments in batches of size B_{TS} (**bwTS**) with access to target synapses and neuronal ring buffers in batches of size B_{RB} , including group prefetching of spike ring buffers (**bwRB***). **TS** marks iteration over a synaptic target segment using a fixed count loop according to the target-segment size ts_size . **SYN** marks access to an individual target synapse; **RB** marks access to the ring buffer of the corresponding target neuron; **RB*** marks group prefetching prior to access. Based on [REF](#).

Data: *spike_reg, synapses*

```

create arrays lcid, ts_size of size  $B_{TS}$ 
create arrays target_rb, delay, weight of size  $B_{RB}$ 
l ← 0
i ← 0
repeat spike_reg.Size() /  $B_{TS}$  times
  for k ← 0 to  $B_{TS} - 1$  do
    lcid[k] ← spike_reg[l + k].lcid
    l ← l +  $B_{TS}$ 
    for k ← 0 to  $B_{TS} - 1$  do
      ts_size[k] ← synapses[lcid[k]].GetTSSize()
      for k ← 0 to  $B_{TS} - 1$  do
        TS
        SYN
        repeat ts_size[k] times
          (target_rb[i], delay[i], weight[i]) ←
            synapses[lcid[k]].Send()
          i ← i + 1
          lcid[k] ← lcid[k] + 1
          if i ==  $B_{RB}$  then
            RB*
            for i ← 0 to  $B_{RB} - 1$  do
              prefetch target_rb[i]
            for i ← 0 to  $B_{RB} - 1$  do
              RB
              target_rb[i].AddValue(delay[i],
                weight[i])
            i ← 0
    process remaining entries in spike_reg and target_rb

```

receive register (*spike_reg*) in the outermost loop and the concomitant processing of synaptic target segments of different lengths in different memory locations. The quantitative analysis ([Section 5](#)) refers to this set of optimizations as **bwTS**.

We adapt the reference algorithm to process spike entries in the spike receive register in batches of size B_{TS} , which in turn entails batchwise access to B_{TS} distinct target segments (**bwTS**). Like in **bwRB***, this optimization is based on a loop-transformation method ([Section 4.1](#)) but targets a different part of the algorithm. Moreover, we replace the while loop iterating over every target synapse of a specific target segment with a fixed count loop (**TS**). This requires however that the length of the target segment is available when entering the fixed count loop. Here, we decide on a straightforward solution: the in-

formation is provided by the first target synapse of each target segment. To this end, we extend synapse objects with a member variable to store the target-segment size, which needs to be determined just once when all synapses have been created (not shown in the algorithm); the public member function `GetTSSize()` returns the size. Note that while the algorithm only requires this capability for the first synapse of each target segment, all synapse objects are equipped with the extra member variable as they are stored in a container for homogeneous objects ([Section 3.1](#)). In our reference implementation ([Section 2.1](#)), we ensure that the extra member variable does not increase the per-synapse memory usage by reducing the storage size of another synaptic member variable, namely the synaptic delay. This does not affect the precision of results for our benchmark network model as the reduced storage size is sufficient to fully represent the homogeneous and relatively short delays of the model ([Section 2.2](#)). This solution does, however, not generalize to all types of network models as, for example, in case of longer delays, an increase in per-synapse memory usage due to the additional member variable might be inevitable.

For each batch of B_{TS} spike entries, the algorithm carries out three consecutive for loops each with a fixed number of B_{TS} iterations. In the first for loop iterating over B_{TS} spike entries, the locations of the first target synapses of the corresponding target segments are extracted from the spike entries and buffered in the auxiliary array *lcid*. In the subsequent for loop the B_{TS} locations are used to access the first target synapses to retrieve the corresponding target-segment sizes, which are buffered in the auxiliary array *ts_size*. In the final for loop and the enclosed fixed count loop, alternating access to synaptic targets (**SYN**) and neuronal spike ring buffers (**RB**) is carried out like in the reference algorithm. If the number of spike entries in the spike receive register is not divisible by B_{TS} without remainder, the remaining spike entries are processed in a similar fashion using three consecutive for loops, where the fixed number of iterations is given by the number of remaining entries.

4.4. Combined batchwise access to target segments and spike ring buffers

The combined algorithm **bwTSRB*** adopts the loop structure of **bwTS**, but the instructions inside the fixed count loop iterating over a specific synaptic target segment (**TS**) are adapted according to **bwRB***. The combined algorithm requires all auxiliary arrays of the two individual algorithms, where indices *l* and *k* and index *i* are used as in **bwTS** and **bwRB***, respectively.

5. Results

We quantitatively evaluate the effect of the three different techniques of latency hiding **bwRB***, **lagRB**, and **bwTS** on simulation time relative to the refactored code ([Section 3.2](#)), where the optimization **bwTS** can be combined with either **bwRB*** or **lagRB** because they modify different parts of the code. As the combined optimizations may either support each other or in the worst case reduce the impact of the best individual optimization, we also present performance data for a combined scenario

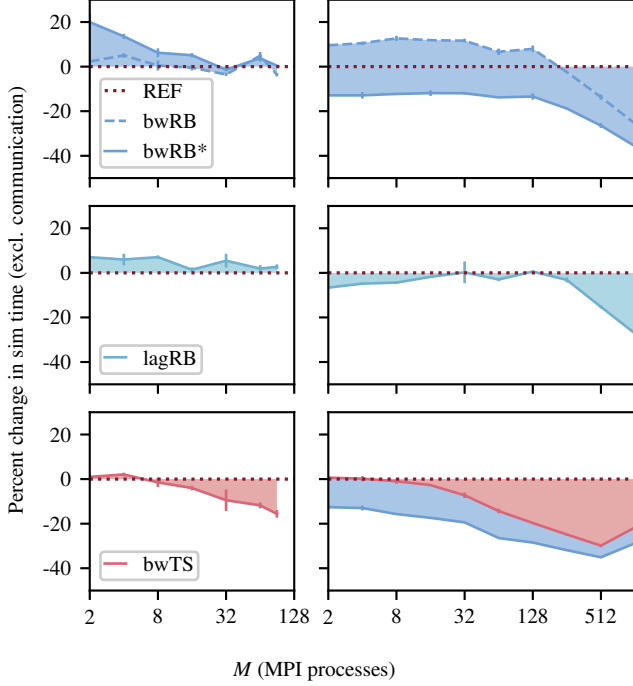


Figure 4: Cumulative change in simulation time relative to refactored code disregarding communication as a function of number of MPI processes. Left column DEEP-EST CM and right column JURECA CM: linear-log representation for number of MPI processes $M \in \{2; 4; 8; 16; 32; 64; 90\}$ and $M \in \{2; 4; 8; 16; 32; 64; 128; 256; 512; 1,024\}$, respectively. Dark carmine red dotted line at zero percent (REF, Section 3.2) indicates performance of reference code. First row, batchwise access to spike ring buffers (Section 4.1) with batches of size 16 without group prefetching (bwRB, dashed light blue curve) and with prefetching (bwRB*, solid light blue curve). Second row, software pipelining with a lag of 16 (lagRB, turquoise curve). Third row, batchwise access to target segments with batches of size 16 (bwTS, solid coral red curve), for JURECA CM (right) further combined with batchwise access to spike ring buffers using prefetching (bwRB*, same coloring as in top row: solid light blue curve; see bwTSRB* for combined pseudocode). Shadings fill area to respective reference. Weak scaling of benchmark network model as in Figure 1.

of bwTS and bwRB*. Figure 4 summarizes the quantitative results. The section concludes by demonstrating that the improved performance is indeed a result of a reduction of clock ticks per instruction retired (Section 5.1).

Plain batchwise access to spike ring buffers without group prefetching (bwRB, Section 4.1) has no effect on the performance on the DEEP-EST CM and enabling prefetching (bwRB*) actually worsens the situation for small numbers of MPI processes. The situation is entirely different on JURECA CM which uses an older generation of processors. Here, plain batchwise access has a negative effect on the performance but group prefetching leads to an overall performance improvement. At larger numbers of MPI processes already the batchwise processing increasingly improves performance and the additional gain by group prefetching remains constant. For tested batch sizes between 1 and 64, we observe the least decline in performance for batch sizes of 8 or larger on DEEP-EST CM and the best performance for batch sizes of 16 or larger on JURECA CM (data not shown). However, a comprehensive analysis is outside the scope of this study; we do not

claim that this observation generalizes to other architectures.

An alternative latency-hiding technique for the same part of the code is lagged access to ring buffers (lagRB, Section 4.2), where we use a lag of 16. The algorithm thus retrieves information from a target synapse 16 steps ahead from the position in the target segment where it currently accesses the corresponding ring buffer, thereby decoupling these operations. Again we observe almost no effect on the DEEP-EST CM. JURECA CM exhibits an increasing gain only for large numbers of MPI processes. For tested lags between 1 and 16, we observe little differences on DEEP-EST CM and the best performance for lags of 2 and larger on JURECA CM (data not shown).

The introduction of the spike-receive register [9] opens the opportunity for batchwise processing of the register and the synaptic target segments (bwTS, Section 4.3). The batchwise processing of target segments has an increasing gain reaching 20% at 90 MPI processes on the DEEP-EST CM, and the total gain reaches more than 40% on JURECA CM. For tested batch sizes between 1 and 64, we observe little differences on DEEP-EST CM and the best performance for batch sizes of up to 16 on JURECA CM (data not shown).

Finally, on JURECA CM we investigate a combined implementation of batchwise access to target segments and batchwise access to spike ring buffers including group prefetching (bwTSRB*) as the two techniques modify different parts of the code. We do not further consider lagged access to ring buffers (Section 4.2), as lagRB is only effective for large numbers of MPI processes, while its alternative bwRB* exhibits a sustained gain over the full range. The lower right panel of Figure 4 shows that bwRB* is effective already at small numbers of MPI processes but continues to improve the performance of bwTS, leading to a combined gain of 50% at larger numbers of MPI processes.

The preparatory refactoring described in Pronold et al. [9] and the latency-hiding techniques investigated in this study jointly improve simulation time with respect to the original code (Figure 1). While refactoring is most effective for small numbers of MPI processes where synaptic target segments are long (Figure 2), latency hiding is most effective for larger numbers of MPI processes where target segments significantly shorten. The differential behavior of the combined technologies leads to a sustained performance gain of 30% to 50% over the whole range of investigated MPI processes.

5.1. Dependence of effectiveness on microarchitecture

The latency-hiding techniques introduced in Section 4 promise to speed up the overall code by reducing the latency of memory access. Indeed our measurements show that our particular application becomes substantially faster, despite additional lines of code required to implement the techniques. Profiling tools (Section 2.4) provide direct access to the latency that processor instructions experience. This allows us to investigate whether the new algorithms are faster due to the reduction of latency. Specifically we are investigating clock ticks per instruction retired (CPI), a metric available in the tool VTune (Section 2.4). All algorithms exhibit a decrease in CPI compared to REF (Figure 5), except for less than 32 MPI processes bwRB* on DEEP-EST CM and bwRB for JURECA

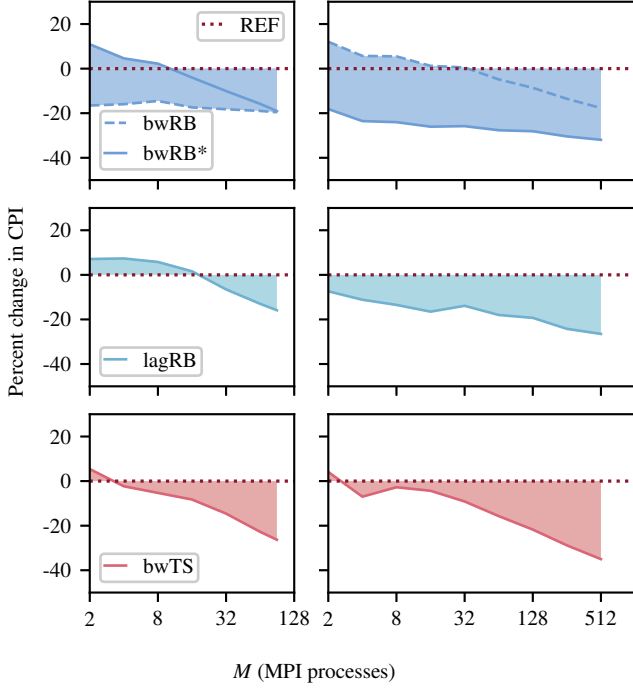


Figure 5: Relative change in clock ticks per instruction retired (CPI) during spike delivery as a function of the number of MPI processes. Same arrangement of panels and labeling as in Figure 4.

CM. Larger networks show larger gains and **lagRB** and **bwTS** show similar behavior on both machines. The dependence on network size is particularly pronounced for **bwTS**. This directly relates to the decreasing length of target segments, diminishing the algorithmic potential for ordering synaptic data for sequential processing. As **bwTS** addresses the bottleneck of hopping from one target segment to the next, its effect increases as target-segment lengths decrease. The characteristics of **bwRB** (Section 4.1) and **bwRB*** on DEEP-EST and JURECA appear inverted. This hints that the processor in DEEP-EST CM, in contrast to JURECA CM, automatically prefetches the necessary data on time. Adding additional explicit prefetching instructions may interfere with the automatic prefetching and hence degrade performance. While the changes in simulation time (Figure 4) follow the changes in CPI, the reduction in sim time does not fully reflect the success observed in CPI. All latency-hiding techniques come with additional lines of source code. This may lead to an increased number of processor instructions that the reduction in CPI has to make up for, but the compiler may now also find more opportunities for reducing the number of instructions.

6. Discussion

At the outset of this investigation stands the observation that for a typical neuronal network the spike-delivery phase dominates total simulation time in state-of-the-art simulation code and that the time required for spike delivery increases under weak scaling such that it still dominates at 1024 MPI processes

(Figure 1). This article explores techniques to rearrange the elementary algorithmic steps required to deliver the incoming essentially random spike data to the thread-local targets such that they can be more efficiently processed by conventional computer hardware. Each of the techniques under consideration is characterized individually, but some address different steps of spike delivery. Combining the most promising techniques achieves a significant performance boost.

The combined spike-delivery algorithm starts with the thread-parallel sorting of the spike data from the MPI receive buffer into a thread-specific data structure called spike receive register (SRR), where sorting is performed according to hosting thread and synapse type of each spike’s synaptic targets. This preparatory refactoring improves parallelization but is not related to the optimization of cache performance on the level of the individual threads. The technical companion paper [9] provides the detailed description of the SRR together with pseudo code. Sorting into the spike receive register is followed by the thread-parallel processing of the SRR in fixed-size batches cycling through three stages (**bwTS**, Section 4.3): for each batch of spike entries, collecting first the location and second the length of the spike’s synaptic target segment in separate arrays before continuing with the actual delivery as a third stage. In the third stage, also synapses and corresponding neuronal spike ring buffers are processed in batches (**bwRB***, Section 4.1) by, irrespective of target-segment boundaries, weight, delay, and pointer to ring buffer, first collecting separate arrays for each batch of synapses before adding the weights to the buffers.

As a result of the redesign the time required for spike delivery is halved (reduced by factor of 2.1) for two MPI processes compared to the original algorithm and reduced to a third (by a factor of 2.9) for 1024 MPI processes. In the original algorithm, the increased network sparsity at larger numbers of MPI processes causes an increase in spike-delivery time by a factor of 4.5 between two and 1024 MPI processes. The new algorithm reduces this to a factor of 3.3. As a consequence, simulation time is reduced by 43% for low numbers of MPI processes, by 49% for 512 MPI processes, where communication is not yet dominating, and by 28% for 1024 MPI processes, where communication time starts to dominate overall simulation time. The measurements were obtained using ParaStationMPI v5.2.2-1, while initial tests using the more recent version v5.4 result in a significant reduction in communication time (data not shown). However, following up on this observation is not within the scope of this manuscript due to JURECA CM being decommissioned. In summary, with the new algorithm spike delivery still dominates the simulation time below 1024 MPI processes under weak scaling, but it increases less rapidly such that now even below 1024 MPI processes the major cause for the loss in performance is the increase in communication time. A second qualitative change sets in at 1024 MPI processes, where the absolute simulation time is no longer dominated by spike delivery but by communication time. Hence, the new algorithm overcomes the previously reported barrier of spike delivery limiting the performance on supercomputers (Figure 12 in 10). Now progress can be made by optimizing communication, for example, by exploiting the spatial organization of neuronal net-

works. Cortical neuronal networks are characterized by a local coupling with a space constant of a few hundred micrometers and delays in the range of a tenth of a millisecond combined with long distance coupling between brain areas and delays beyond a millisecond. If cortical areas were represented on one or a few compute nodes, the communication between nodes hosting different areas could be reduced to much larger intervals than required between nodes hosting neurons of the same area. Topology-aware distribution of neuronal networks across compute nodes has been exploited in other simulation codes [44, 45] but without taking into account different delay categories.

Performance profiling of the algorithms `bwRB*`, `lagRB`, and `bwTS` indicates that they indeed reduce latency (Section 5.1). In accordance with the measurements of sim times (Section 5), which show that the techniques are more successful on JURECA CM (Intel Haswell) than on DEEP-EST CM (Intel Skylake), the relative reduction in the performance metric clock ticks per instruction retired (CPI) is also more pronounced on JURECA CM than on DEEP-EST CM. We speculate that the improved cache utilization of the newer generation processor (Skylake) renders `bwRB*` and `lagRB` ineffective or even detrimental, at least in the regime up to 90 MPI processes investigated here. On JURECA CM the effect on sim time increases beyond 128 MPI processes, which is beyond the current capacity of DEEP-EST CM. Therefore, emulating simulations on more than 128 MPI processes on DEEP-EST CM using the NEST dry-run mode [40] may provide further insights. A comprehensive comparison between the two generations of processors based on microbenchmarks as presented in [46] for the related microarchitectures Intel Broadwell and Intel Cascade Lake is out of the scope of this study. It is our hope though that our efforts to present the changes to code in an abstract fashion make this study relevant for a broader computer science community and might even inspire the definition of future microbenchmarks. The present work focuses on the level of source code only. It may be illuminating to explore in future studies the effect of algorithmic changes on the level of the resulting assembler code.

The incoming spike events of a compute node specify the thread hosting the target neuron as well as the location of the synaptic targets, but are unsorted with respect to the hosting thread and synapse type. Nevertheless the present work shows that the processing of spikes can be completely parallelized requiring only a single synchronization between the threads at the point where the spikes have been sorted according to target threads and synapse types, which is when all spikes have been transferred from the MPI receive buffer into the novel spike receive register. This suggests that spike delivery can fully profit from a further increase in number of threads per compute node. Due to the large number of outgoing synapses per neuron, early work on distributed simulations of spiking neuronal networks employing tens of compute nodes was concerned with the problem of efficiently delivering each spike emitted by a specific source neuron on a specific MPI process to many targets on all MPI processes [4]. With increasing parallelization using more compute nodes and threads, however, each neuron has ever fewer thread-local targets resulting in ever shorter synaptic

target segments in the data structure storing the local synapses. Therefore, later work concentrated on reducing the memory overhead caused by the increasing number of shorter target segments [37]. However, shorter target segments are also a burden computationally because memory locations need to be switched often leaving little opportunity for vectorization. The present study overcomes the problem of short segments by reorganizing the spike delivery algorithm such that it operates across the boundaries of target segments and in this way becomes more independent of the degree of parallelization. Spikes are routed to their thread-local targets in a batchwise fashion using the technique of loop transformation. This explicit declaration of virtually independent code blocks apparently helps compilers to generate efficient machine code.

Due to the simplicity of the neuron model of the benchmark and the lack of synaptic plasticity, the application has little workload in terms of the propagation of neuron and synapse dynamics and thus exposes bottlenecks in the delivery of spikes to local targets and the communication between compute nodes. The synaptic delay of about a millisecond assumed in the present benchmark model is used in prominent neuronal network models of the balanced random class [38] and representative for the connectivity at the brain scale [1]. In models of the local network below a square millimeter of cortical surface [cf. 47] the minimal delay (Section 2.1) is an order of magnitude smaller than the delay considered in this study. Consequently, the simulation of such models requires a ten-fold increase in the number of communication calls and therefore the communication phase takes up a larger fraction of the total simulation time. Nevertheless, for a given spike rate the amount of spikes that need to be delivered in a given stretch of biological time is independent of the number of communication calls. Therefore, we expect the optimizations discussed in the present work to still be effective, while their impact on total simulation time will be lower. Neuroscientists have started to investigate models combining the local structure of the brain with its organization over long distances [48]. The minimal delay in such a brain-scale network is the minimal delay in the local structure. Unless the placement of neurons on compute nodes respects the architecture of the network, global communication in intervals as for the local network is required, limiting the success of optimizations of spike delivery.

Our analysis is restricted to static neuronal networks. In such networks all synaptic connections and their weights are determined at the time of network construction. In nature synaptic efficacies change over time: the phenomenon called synaptic plasticity. A further mechanism, removing and creating synapses, is called structural plasticity. A wide class of models of synaptic plasticity can be formulated as an update scheme driven by the presynaptic spike events in which the synaptic weight is only computed for those times where it becomes visible for the postsynaptic neuron [39]. In this way the computations become part of the spike delivery phase and contribute substantially to its duration. Depending on the particular plasticity model, alongside the location of the ring buffer, additional information on the target neuron needs to be accessed. This may be the time of the last spike, the full record

of spikes since the last presynaptic spike, or the membrane potential [49]. Our companion paper [9] discusses the resulting tradeoff between memory consumption and speed. A domain specific language like NESTML [29] could avoid cluttering the model pool with solutions for different optimization goals by providing directives influencing the balance between the generation of more compact or faster code depending on the needs. How the more complex data handling for plastic synapses and the floating point operations on the synaptic weights interact with the batchwise processing needs to be evaluated.

The communication scheme between compute nodes assumed in the present study sends for each spike a separate event to each thread containing at least one target neuron. This enables a straightforward readout of the MPI receive buffer containing the incoming spikes (Section 3.2). However, for typical cortical connection densities where each neuron has on the order of 10,000 postsynaptic targets for simulations with less than 10,000 compute nodes and more than one thread per node, compute nodes receive the same spike event multiple times. This overhead may challenge the bandwidth of the communication network and limit the speed of the communication phase. Future work needs to investigate whether duplicate spike events can be efficiently compressed such that the costs of decompression in the spike-delivery phase are smaller than the gain in the communication phase.

The present study analyses spike-delivery times and overall simulation times in a weak-scaling scenario. Neuroscientists and industry may, however, also be interested in reducing the simulation time of an application for a fixed size network studied over long stretches of biological time as required for system-level learning or for the capability to interact with the real world in robotics. Therefore strong-scaling scenarios are of interest. With an increasing number of threads, the number of target neurons and therefore synapses a thread needs to take care of decreases. Consequently the amount of memory a thread operates on decreases. This suggests that it becomes easier for the core running the thread to predict memory access and if the amount of cache memory per core decreases proportionally to the number of cores faster execution is expected. In this way, the combination of a highly parallel spike-delivery algorithm with a many-core architecture overcomes the von Neumann bottleneck of conventional applications. Thus our findings reconsidering the spike-delivery algorithm support a more positive view on the prospects of improving the performance of neuronal network simulations by prefetching and pipelining than the analytical results of [11, 12]. It remains to be seen whether the optimizations by batchwise processing explored in the present work become less relevant in situations where each core needs to manage a small amount of memory anyway.

The study removes several bottlenecks in routing spikes in a compute node to ever more distributed targets. The spike-delivery phase of a simulation code for spiking neuronal network remains a specific sorting problem in space and time, but our study exposes that its fully parallelizable nature will certainly benefit from future fine-grained processing hardware.

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