

# Performance analysis of a parallel PDEVS simulator handling both conservative and optimistic protocols

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## ABSTRACT

With the ever increasing complexity of simulation models, parallel simulation becomes necessary to perform the simulation within reasonable time bounds. The built-in parallelism of Parallel DEVS is often insufficient to tackle this problem on its own. Several synchronization algorithms have been proposed, each with a specific kind of simulation model in mind. Due to the significant differences between these algorithms, current Parallel DEVS simulation tools restrict themselves to only one such algorithm. In this paper, we present a Parallel DEVS simulator, grafted on C++11, which offers both conservative and optimistic simulation. We evaluate the performance gain that can be obtained by choosing the most appropriate synchronization protocol. Our implementation is compared to adevs using hardware-level profiling on a spectrum of benchmarks.

## 1. INTRODUCTION

### 1.1 DEVS

The family of DEVS [28] formalisms serve as a common basis for most other discrete event formalisms. In this paper, we are in particular interested in Dynamic Structured DEVS [1] and Parallel DEVS [5]. We propose a DEVS implementation (called DEVS-Ex-Machina and referred to as "dxdev" throughout this paper) that is based on the the Direct Connection [4] algorithm. This algorithm flattens all hierarchically nested models which leads to, from a simulation kernel's perspective, a collection of Atomic Models linked to each other by ports.

### 1.2 Parallel computing

One way to speed up the PDEVS simulation is by using parallel execution. By its very nature, this can lower the runtime considerably and increase the bound on the state space, thereby enabling simulation of more complex systems in the same time-frame. While the shared memory parallelism offered by most modern hardware does not raise the state space

bounds, it can reduce the runtime and offers more direct communication and control between entities involved in synchronization compared to distributed simulation.

### 1.3 Motivation

One of the drawbacks of adevs [21], a well-known PDEVS simulator which is performing well with conservative shared memory synchronization, is that it lacks support for optimistic synchronization. Using optimistic synchronization can be significantly faster, especially in simulations where the runtime behaviour of the simulation is hard to predict. The matured parallelism features of C++11, on the other hand, created new opportunities as well. They allowed to write very efficient code giving access to powerful low-level threading primitives and yet guarantee standard-compliance and portability.

### 1.4 Solution

The usage of the Direct connection algorithm makes reusing existing kernels (such as adevs or PythonPDEVS [26], also known as PyPDEVS, hard. Dxex is based on PythonPDEVS [26](PyPDEVS) but differs in design and implementation. The core aim of the project is to offer a deterministic simulation kernel where the simulation modeller is shielded as much as possible from the kernel implementation, without sacrificing performance. As in PyPDEVS, a model has to be written only once for use in the different simulation kernels.

In order to monitor and investigate simulations, dxex uses a tracing framework which is derived from the one used in PyPDEVS.

### 1.5 Time

The theoretical DEVS formalisms all have  $\mathbb{R}$  as time base, but any implementation has to decide on an enumerable representation of time. In principle, any type with well-defined operators could be used, but, from a performance point of view, a type fitting in a machine word offers obvious advantages. In practice, dxex defaults to word sized unsigned integer. Although this significantly reduces the possible range of virtual time, it avoids approximation errors. The notion of  $\epsilon$ , moreover, is trivial for an integer representation, unlike floating point.

In addition, our time representation is enriched with a causality field the range of which is equal or greater than the maximum number of models in a single kernel ( $2^{48}$  by default). If  $A$  and  $B$  are imminent at time  $t$ , then  $t[1]_a < t[1]_b \oplus t[1]_b < t[1]_a$ , while  $t[0]_a == t[0]_b == t[0]$ . This allows a deterministic ordering of concurrent events.

## 2. BACKGROUND

In this section, we provide a brief introduction to two different synchronization protocols for parallel simulation, and the features offered by C++11 that aid in our implementation.

### 2.1 Conservative Synchronization

Conservative synchronization is defined by the invariant that no model will advance in time before it has received all input from any influencing model.

This requires the concepts of earliest output time (eot) and earliest input time (eit), which define the timespan within which a model can safely advance.

The eit of any model is the minimum of all eot values of (directly) influencing models. A model can simulate up to (but not including) eit and then waits until that value is increased. An important disadvantage here is that the relation that calculates the dependencies, is always defined at model (link) creation time, and not at runtime. A model that can influence another model theoretically, but never does in real life, can severely slow down the protocol.

Deadlocks between models that influence each other and end up waiting on each other, can be broken/avoided by a variety of means. In dxex, the CMB [3] null-message protocol is used.

In our implementation, null-time is the time stamp a model is guaranteed to have passed in simulation. More precisely, a null message of time  $t$  guarantees that any output with time stamp  $t - \epsilon$  has already been sent.

Conservative synchronization relies explicitly on information provided by the model creator in the form of a lookahead, a relative timespan during which the model is insensitive to outside events. This lookahead can be non-trivial to calculate; therefore, a simulation writer will in general not be able to determine the exact lookahead of the models involved in an experiment without having run the experiment.

Conservative kernels can even operate if there are cyclic dependencies involved, but this can come at a severe performance penalty, as seen in section 4.

### 2.2 Optimistic Synchronization

In the case of optimistic synchronization, causality errors can occur but these will be reverted via a roll-back mechanism, the most common of which is Timewarp [15].

Whenever a kernel receives an event with a time stamp in the kernel's past, the state of the kernel (and all models) is reverted to that time.

Because an optimistic kernel does not wait for other kernels, it can run much faster than a conservative kernel. The cost of this approach, however, is higher complexity of the code and a larger memory usage for the state and message saving. In contrast to conservative synchronization, optimistic synchronization does not rely on any domain specific information. Therefore, it is only sensitive to the runtime use of

connections, not the probability that they might occur. If the (runtime) dependency graph contains a cycle, optimistic synchronization might lead to a series of cascading reverts. The cost of reverts can be reduced by lazy cancellation and/or lazy re-evaluation [8].

### 2.3 Global Virtual Time

To avoid exhausting memory in state/event saving, optimistic synchronization relies on the concept of global virtual time[15] or GVT. In optimistic simulations, GVT is defined as the lowest time stamp of any unprocessed event.

Intuitively, this is the simulation time point that is certain to be preserved, corresponding exactly with a sequential simulation. In conservative, the minimum simulation time of all kernels is the lower bound of time stamp or LBTS. In terms of null messages this corresponds to the least time stamp of any null message in transit. The GVT calculation is vital to safely commit unrecoverable transactions such as IO (e.g. tracing), releasing memory, etc.

### 2.4 C++11 Parallelism Features

C++11 offers a wide range of portable synchronization primitives in the Standard Library, whereas in earlier versions one had to resort to non-portable (C) implementations. More importantly, C++11 is the first version of the standard that actually defines a multi-threaded abstract machine memory model in the language. Our kernels use a wide range of threading primitives and atomic operations. As an example, eot/eit/nulltime are not exchanged as messages, but as reads/writes to atomic fields shared by all kernels. This avoids the otherwise unavoidable latency penalty by mixing simulation messages with synchronization messages. For an in-depth study, we refer to [6]. Most modern compilers support the full standard, allowing the kernels to be portable by default on any standard compliant platform.

## 3. FEATURES

### 3.1 Based on PyPDEVS

Dxex is based on PyPDEVS, and provides the following features:

1. Direct Connection
2. Dynamic Structured DEVS
3. Termination function: if specified, a termination function is applied to each model every simulation round to test whether the simulation can terminate. This feature is only available in single-threaded simulations.
4. The State/Message objects can have any payload type. Different allows different message types to be used within the same simulation.
5. Tracing: An asynchronous, thread safe and versatile tracing mechanism allows exact verification of the simulation.
6. Optimistic and Conservative synchronization

Furthermore, the implementation tries to adhere to the C++ principle that you don't pay performance-wise for what you

don't use. For this reason, the support for a termination function for the multi-threaded kernel was abandoned, as it is non-trivial to implement and had an adverse impact on the runtime, even when not in use. Another example is the state saving mechanism, which is only used for optimistic parallel simulation and has no performance impact in a conservative parallel simulation.

Our tracing implementation is not comparable to adevs's listener interface. To be usable in an optimistic simulation, the tracing of the simulation has to be reversible and only be committed at GVT points. Furthermore, the framework itself has to be thread safe and deterministic so that a simulation will always produce the exact same output.

The following features from PyPDEVS are not present

1. Activity tracking and relocation
2. Serialization: in this context this is the ability to save/load a complete simulation to disk, not the state saving mechanisms required for TimeWarp.
3. Interactive control
4. Distributed simulation

In dxex, the model allocation is realized by a derivable allocator object which the user can implement to arrange a more ideal (domain-specific) allocation. If this is omitted, a default (non-activity-aware) allocation stripes the models over the simulation kernels.

In addition, debugging tools such as a logger and a graph visualizer are included that can track activity with respect to the allocation for later inspection (but not are not online and as dynamic as is in PyPDEVS).

### 3.2 Different Synchronization protocols

For parallel executions, synchronization is required for PDEVS to prevent causality violations from happening or to recover from causality violations. Preventing causality violations (conservative) typically requires domain-specific information to compensate for the performance loss, while recovering from causality errors (optimistic) requires the calculation of a state to revert to.

#### *Conservative*

In case of conservative synchronization, any kernel will determine which kernels it is influenced by. This information is constructed from the incoming connections on all hosted models. The process is only 1 link deep, since an influenced kernel will in turn be blocked by others deeper in the graph.

A model should provide a lookahead function which returns, relative to the current time, the timespan during which the model cannot change state due to an external event. This information is collected for all models hosted on the kernel and the minimum time is set as the lookahead of that kernel. The kernel will calculate its earliest output time and write this value in shared memory. The eit of the kernel is then set as the minimal eit of all influencing kernels.

For the garbage collection (of sent messages), the LBTS/GVT is calculated as  $\min_{i \in \text{influencers}} (\text{nulltime}[i]) - \epsilon$ .

#### *Optimistic*

The optimistic kernel requires from the hosted model that copying the state is done carefully (avoid unnecessary copies and make a copies whenever necessary).

The kernels use Mattern's [17] GVT algorithm with a maximum of 2 rounds per iteration to determine a GVT. This process runs asynchronously from the simulation itself. Once found, the controlling thread informs all kernels of the new value, which they can use to execute garbage collection of old states and (anti)messages.

The user must only provide one implementation of a model that can be used for both synchronization protocols. A lookahead function is desired to accelerate the conservative protocol, but is not required. In the absence of a user supplied lookahead, the kernel assumes it cannot predict beyond its current time  $t + \epsilon$ , creating a lockstep simulation. The user is shielded from the multi-threaded aspect of the kernel.

From the user's perspective, the multi-threaded aspect of the kernel is not exposed.

### 3.3 Performance Improvements

We now discuss a number of bottlenecks that were discovered during the profiling of several benchmarks.

#### *Heap*

In dxex, events are always passed with pointers and thus avoid a possibly expensive copy of the payload caused by allocation and copying overhead. In highly connected models, this allocation cost can become prohibitively expensive, so to reduce that overhead we use thread local memory pools for states and events, and, optionally, replace the system malloc with calls to tcmalloc [10]. In this way, allocating threads do not block each other. If desired, arena-pools are available for single-threaded simulation. A disadvantage is increased complexity in ownership semantics. The creating kernel is responsible for destruction, but this can only be guaranteed up to the GVT. Experiments with synchronized pools proved to be slower than implementations with the standard malloc/free.

Initially, dxex was using strings as identifiers as was also done in PyPDEVS. Profiling quickly indicated that this caused a real performance bottleneck. In C++, strings are heap allocated variable sized objects with an atomic reference count and not not immutable objects as in Python. Access of that reference count across threads turned out to be quite expensive, as are the calls to malloc/free the string implementation makes to create/destroy new objects or copy existing objects. Strings, however, are more intuitive to work with from a user's perspective, so, as a compromise, we allowed the user to reference models/ports by string name. Once the simulation starts however, all objects use integral identifiers. This also increased the usage of the constexpr feature of C++11 in, amongst others, time stamps and message headers.

#### *Raw pointers*

While an important C++11 feature in general, our initial usage of smart pointers for some types of objects was misplaced. Used across threads, the reference counting became very expensive, and the (de)allocation of memory caused significant contention between threads. So for the models and

the kernels, dxex is still using smart pointers whereas for the messages, a raw pointer to compacted memory is being used.

#### Locking

Locking between kernels uses mostly atomic operations and, occasionally, we can leverage memory orderings to only pay for synchronization when we need it. Messages, on the other hand, are exchanged via a shared set of queues each with a dedicated lock.

On a higher level, we avoid the sending of synchronization messages entirely by writing the time stamp directly into shared memory.

Sending of antimessages is fairly cheap in our implementation, since only the modified pointer to the original message is sent to the receiving kernel.

#### Schedulers

PyPDEVs has a wide range of schedulers to choose from with varying performance depending on the simulation type. Profiling showed that, the heap implementation used in adevs was faster than any of the schedulers we had tested before (in a C++ environment). Unlike most node based heaps, this scheduler uses a fixed size array where its heap is rebuilt or modified in place depending on the amount of items to update. Items are only updated, never removed.

## 4. PERFORMANCE

### 4.1 Sequential Simulation

In our benchmarks, we will consistently compare the performance of dxex with adevs. The implemented models for the each benchmark are functionally identical.

Dxex uses integer timestamps whereas adevs uses floating point time representation. Tests show there is little difference performance-wise (both types fit in a machine word).

The benchmarks have been compiled using the `-O3` option and link-time optimization, with all IO disabled. To compare the models in details, we have run the different benchmarks with different parameters.

The benchmarks have been run on a i7-2670QM, 16GiB RAM, Arch Linux (4.2.5 kernel) and G++ 5.2. Interpretation of the benchmark results is done using call-graphs generated by perf. A selection of those call-graphs is available in the code repository. For the runtime plots, perf is used in combination with R.

#### Devstone

The Devstone [11] benchmark is highly hierarchical. As can be seen in Figure 1, dxex is outperforming adevs considerably. This can be explained by the fact that, using Direct connection, the hierarchical structure is being flattened whereas, in the adevs case, considerable overhead is caused by traversing the coupled model's structure to pass events.

#### PHold

PHold [7] is a benchmark that focuses in particular on parallel execution; the sequential runtime is therefore only added as a baseline. A model selects the destination of a message at runtime using a state-saved RNG, which takes up a non-trivial amount of runtime. The results of the benchmark are visualized in Figure 2.

Devstone single core

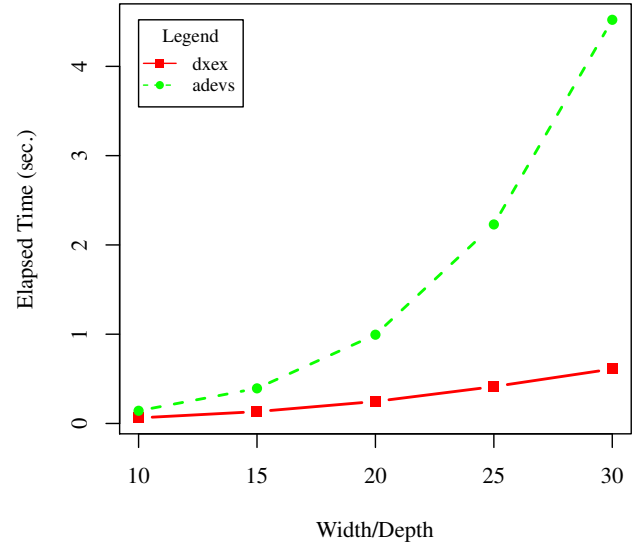


Figure 1. Devstone

#### Interconnect

Interconnect [24] is a benchmark where all models broadcast, creating a complete graph in dependencies between models. As the model count increases, we see (see Figure 3 the expected quadratic increase in runtime for both adevs and dxex, but an increasing penalty for dxex. Profiling shows this is entirely due to the heap allocation of messages, which even though minimized by using memory pools remains significant.

### 4.2 Parallel Simulation

By default, the benchmarks use 4 kernels for parallel simulation.

#### Devstone

The flattened models are allocated to kernels by giving each kernel a distinct section of the chain, resulting in a low ratio of inter-kernel to intra-kernel messages. For the optimistic case, this can cause more reverts since the kernels will start to drift faster as the model-count increases. Furthermore, optimistic synchronization is quite sensitive to an increase in kernels, since the delay before a revert propagates, increases. This benchmark requires a specific warm-up time; for  $n = d \times w$  models, it takes  $\text{timeadvance}() \times (n - 1)$  transitions to activate the last model in the chain. When using parallel processing, this can be reduced to  $\text{timeadvance}() \times (n^{\frac{\text{kernels}-1}{\text{kernels}}} - 1)$  before the last kernel becomes active.

Figure shows that dxex outperformance adevs (in both optimistic and conservative case).

#### PHold

In Phold, the allocation is specified in the benchmark itself. Each kernel manages a single node with a constant set of sub-nodes. The parameter  $R$  (which is here set to 10) determines

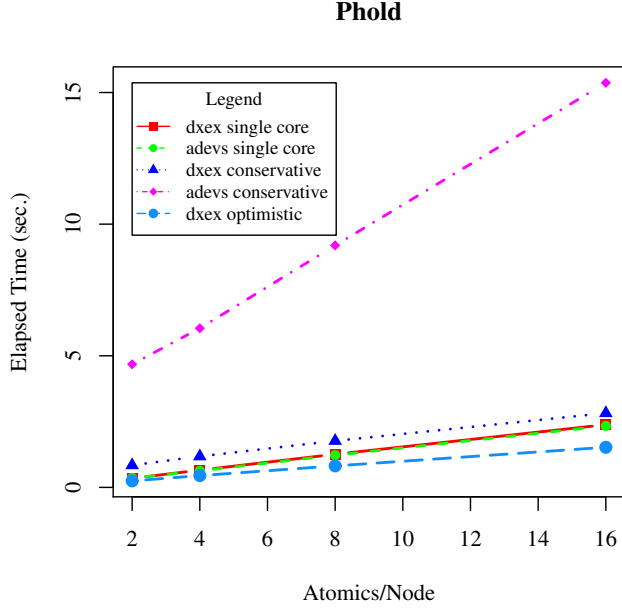


Figure 2. PHold

the percentage of remote destination models.

The dynamic dependency graph is a very sparse version of the static dependency graph, penalizing the conservative case. The lookahead is  $\epsilon$ , so the conservative case spends most of its time crawling in steps of  $\epsilon$ . Since the dependency graph between kernels is a complete graph, this is not a simulation that scales in our implementation. For  $N$  kernels, each kernel has to query the null-time of  $N - 1$  kernels, resulting in  $O(N^2)$  polling behaviour. In a non-cyclic simulation with a non-trivial lookahead (like in, e.g., Devstone), that choice does pay off (see Figure ). Adevs's lesser performance is due in part to their lookahead management, which after profiling shows to spend a non-trivial amount of time in exception handling code.

The optimistic case suffers little from the above problems; due to the high interconnectivity, however, a cascading revert is still possible. With the percentage of remotes equal to 100, PHold reflects interconnect in behaviour, which is why the  $R$  value is not dimensioned here. A revert is very expensive in PHold due to our usage of C++11's random nr generators. The cost of a revert is dominated by the recalculation of destination models, not in allocating/deallocating states/messages. Once a revert happens the drift between kernels increases fast, increasing the likelihood of more reverts. Despite all this, optimistic can for low  $R$  values quickly exploit the uncertainty that slows down conservative in this benchmark.

#### Interconnect

In Interconnect, the set of atomic models form a complete graph (w.r.t. connections); each model broadcasts messages to the entire set.

Allocation cannot avoid cycles and the resulting dependency graph between kernels remains a complete graph. The run-time dependency graph is almost immediately identical to the

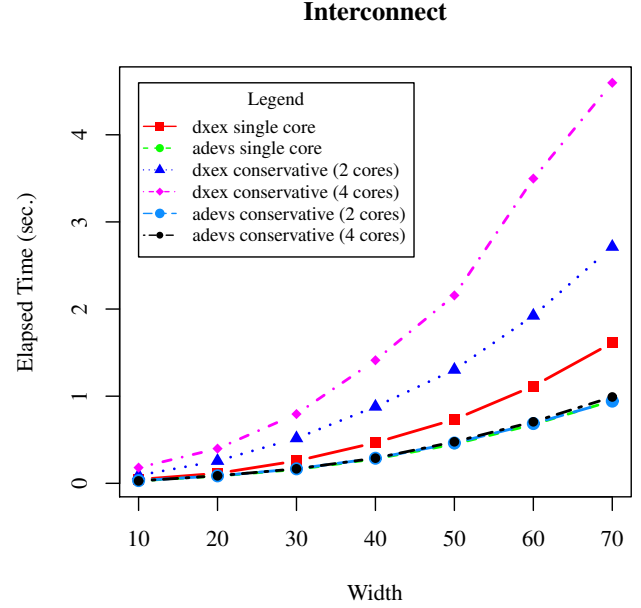


Figure 3. Interconnect

static graph.

The conservative case still shows the same issues as in PHold, with the key difference, for a fixed time advance, the lookahead is equal to the timespan between transitions. The scaling issue is identical as with PHold.

In this benchmark, the optimistic case runs very quickly out of memory. With  $c$  kernels and  $N$  atomic models, a single revert undoing  $k$  transitions will lead to  $(N - 1) \times k$  messages that need to be recreated, plus  $(N - 1) \frac{c-1}{c} \times k$  anti-messages that need to be sent.

#### Priority network model

The priority benchmark is composed of a single server generating a stream of  $0 \leq m \leq n$  messages at fixed time intervals, interleaved with a probability  $p$  for a priority message to  $n$  receivers (see Figure 5).

This defaults the lookahead for the receivers to  $\epsilon$ , but this time there is no scaling effect, nor are there cycles in the dependency graph. This model therefore highlights the basic strengths/weaknesses of both synchronization protocols. Receiving models are allocated on another kernel than the server and have an internal transition so that they will not wait for the incoming messages.

A key difference here with the other benchmarks is that a state (in the Receiver instances) is very cheap to copy/create. The kernel holding the server will never revert since it is a source in the dependency graph. The optimistic case will therefore not suffer the same performance hit in recreating the states as it does in PHold. The overhead in the optimistic case is entirely due to the factor  $m$ , which will quickly dominate in increasing buffers of received messages. Interestingly, the

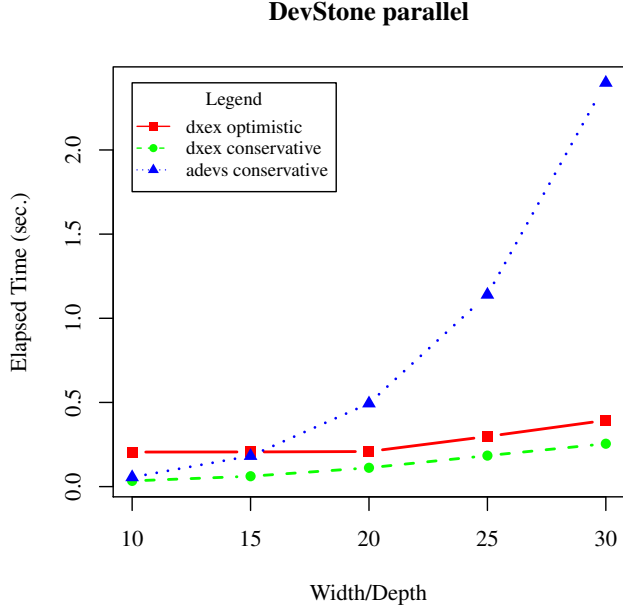


Figure 4. DevStone parallel

parameter  $p$  does not clearly favour either synchronization protocol (see Figure ). While this removes any possibility for a lookahead, the conservative case can quickly bridge the timespan between fixed messages since there is no cyclic dependency. Lightweight states prevent performance loss due to reverts in the optimistic case; only the overhead in event handling in the optimistic case eventually becomes the deciding factor.

### 4.3 Memory Usage

#### Platform and tools

Both dxex and adevs use tcmalloc as memory allocator. Additionally, dxex uses memory pools to further

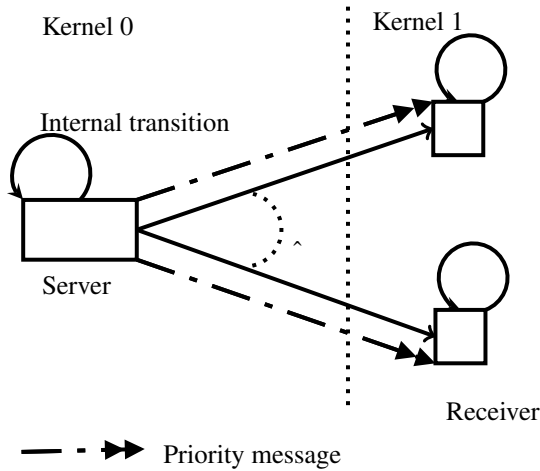


Figure 5. The priority network model

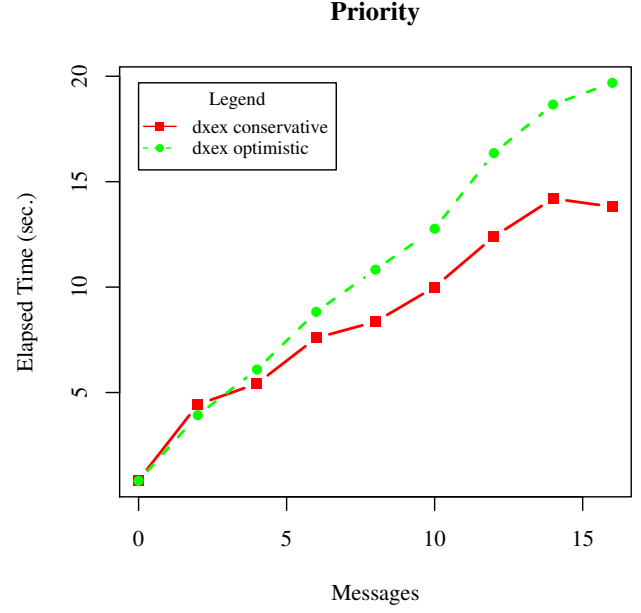


Figure 6. Priority

reduce the frequency of expensive system calls (malloc/free/sbrk/mmap/...). Tcmalloc will only gradually release memory back to the OS, whereas our pools will not do so at all. If memory has been allocated once, it is from a performance point of view better to keep that memory in the pool. This is one reason why memory utilization is best measured by peak allocation. Profiling is done using Valgrind’s massif tool [20]. The platform used for memory profiling has an i5-3317U Intel CPU and 8GiB RAM with a page size of 4,096KiB, running Fedora 22 (kernel 4.2.6).

#### Measure

Adevs passes messages by value, dxex passes a pointer. The runtime effects of this choice have already been demonstrated in the Interconnect benchmark, so, in this section, we measure memory usage in number of allocated pages combining text, stack and heap memory for the program profiled. For the OS and/or user, this is the actual memory footprint of the application. It is important to note that, especially in the optimistic case, not all this memory is always in use by the kernels. During simulation, the pools will generally not return deallocated memory to the OS, but keep it for later reuse.

#### Results

##### Devstone

adevs	adevs con	dxex	dxex con	dxex opt
44	70	42	75	363

Table 1. Devstone 40x40 t5e5, unit MiB, 4 kernels (if parallel)

Since, in the conservative case, messages are passed by pointer, a GVT/LBTS implementation is required to organise the garbage collection. This inevitable delay explains the higher memory usage compared to adevs.

Optimistic’s TimeWarp requires state/event saving, and its GVT algorithm is more complex (with a resulting higher latency) than the LBTS calculation in the conservative case. Moreover, the differences in LP virtual times are far larger compared to conservative time synchronization. All these factors explain the heavier memory usage. Devstone (flattened) is allocated in a chain. Leafs in the dependency graph will therefore do a lot of unnecessary simulation before having a revert, leading to an increased memory pressure. Unlike conservative and sequential execution, memory usage in the optimistic case varies greatly depending on scheduling of kernel threads and drifting between kernels.

#### PHold

adevs	adevs con	dxex	dxex con	dxex opt
40	x	37	61	682

**Table 2.** Phold n 4 s 16 t1e6 r 10, unit MiB, 4 kernels (if parallel)

With only 10% of all messages being inter-kernel, we expect conservative to have memory consumption near that of the single threaded implementation, since intra-kernel messages are reclaimable after each round. The counterintuitive high memory usage can be explained by conservative’s stalled round behaviour which occurs whenever a kernel cannot advance (eit == time). In such a round messages are sent out but the kernel does not execute any transitions until it has received all input from all influencing kernels.

With lookahead  $\epsilon$  this then leads to a high frequency of polling on shared null-times, which are used to determine lbtS and thus garbage collection. The lbtS calculation will not wait until a new value is found, since this can create unwanted contention with the simulation. The resulting longer time intervals within which no new lbtS is found delay memory deallocation. Adevs’ conservative fails to complete the benchmark under valgrind even with a significantly reduced load. Optimistic exhibits the expected high memory usage.

#### Interconnect

In section 4.2 the parallel performance of this benchmark is further explained. Interconnect highlights dxex’s lower performance due to message allocation overhead.

adevs	adevs con	dxex	dxex con	dxex opt
39	39	35	43	259

**Table 3.** Interconnect w 20 t5e5, unit MiB, 2 kernels (if parallel)

#### Priority Network

The priority network model is detailed in section 4.1. The high memory usage of the conservative case is largely due to the garbage collection implementation. Kernel 0 is the initiator of the LBTS calculation but since it holds only models independent of any other (server), it will finish simulation very fast leaving the other kernel bereft of an updated LBTS. The kernel holding the server will after simulation wait on the receiving kernel until it can prove all sent messages can be deallocated.

## 5. RELATED WORK

### 5.1 PythonPDEVS

dxex	dxex con	dxex opt
35	450	651

**Table 4.** Priority model n 128, m 16, p 10, t2e8, unit MiB, 2 kernels

Dxex is closely related to PyPDEVS in design and philosophy. PyPDEVS allows anyone who grasps the PDEVS formalisms to immediately simulate his/her model without having to consider the kernel implementation. A Python based implementation offers the advantage of very fast prototype/run/evaluate cycles, this can’t be matched by a C++ simulator. It should be noted that, once the kernels have been compiled (in shared libraries), the actual compilation time of the model is small enough to make prototyping possible. Advanced features such as activity based relocation and the performance gains this results in, are still unique to PyPDEVS.

### 5.2 Adevs

Adevs is still under active development, allowing for an exact comparison in performance and features. It remains a very fast simulation engine for the PDEVS formalism, but it lacks an optimistic synchronization implementation. Due to the fact that adev does not flatten the Coupled Models, it’s performance degrades significantly if the hierarchy becomes deeper. Dxex, on the other hand, manages lookahead with more overhead than adevs, leading to a performance difference as the model count increases. Finally, adevs employs the full dependency graph this in contrast with dxex kernels, which only observe 1-edge removed nodes.

### 5.3 CD++

Different projects on CD++ offer conservative (CCD++) as well as optimistic (PCD++) parallel simulation. In contrast to dxex, neither projects offer both synchronization protocols. CD++ relies on the WARPED kernel. It is a middleware that provides memory, event, file, time and communication scheduling. We did not use the WARPED kernel (nor the underlying MPI) because dxex is designed specifically for a shared memory architecture and, as such, any middleware, however feature-rich, would have lead to an unacceptable overhead.

## 6. FUTURE WORK

### 6.1 Memory

Our optimistic implementation can benefit from a faster GVT algorithm such as used in [9] or, more recently in [2]. With access to the memory subsystem already in place (pools), the optimistic case could be constrained using per kernel quota, preventing the simulation from exhausting memory. While less sensitive to memory pressure, the same approach could improve the conservative case.

### 6.2 Activity

As shown in [25], activity and allocation of models across kernels is a key aspect in achieving high performance in any parallel implementation. Allocating models so that there are no dependency cycles between their containing kernels is a first step, but not always possible. For the optimistic case, one can use re-allocation to break (runtime dependency) cycles and/or perform load-balancing. If kernels are unevenly

balanced, they will begin to drift fast, causing increasingly more reverts. In dxex, we implement a limited framework to track model activity for debugging purposes; this could be extended to enable the above strategies.

### 6.3 Hybrid

The optimistic implementation could use (null/eot/eit) from the conservative case to detect and/or reduce the cost of reverts without completely stalling on influencing kernels. The eot calculation in the conservative case can be enhanced by using the entire dependency graph between kernels (not models). The implementation could detect dependency cycles between kernels and, instead of waiting on dependent kernels, advance as a synced group with better scaling. LBTS calculation could leverage this as well to reduce memory usage, though this is not that stringent in the conservative case. Ultimately, the simulation could switch at runtime between protocols based on the information provided by activity tracking. This requires the above mentioned framework, and can only be done at timepoints where a GVT/LBTS time is agreed upon between the kernels. From the model's perspective, no changes are needed, although a non-trivial lookahead is obviously desired.

## 7. CONCLUSIONS

In this paper, we presented dxex, a new C++-based PDEVs implementation supporting both sequential and parallel processing models.

We showed that, in the sequential case, dxex outperforms adevs in hierarchical simulations and that adevs produces better results in broadcast-like simulations.

In the parallel case, the results are more subtle. The sensitivity to broadcast simulations is carried over in the parallel implementation, resulting in performance loss for both synchronization protocols. The optimistic case can still achieve good performance if the volume of inter-kernel messages is low. Both synchronization protocols are fast in non-cyclic simulations, whereas the conservative case is better suited whenever state-saving becomes expensive; the optimistic case, on the other hand, is better suited when a lookahead is not available. Finally, the current optimistic implementation is ill-suited for memory constrained systems, especially in simulations where the frequency of reverts is high.

## ACKNOWLEDGMENTS

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## REFERENCES

1. Barros, F. J. Modeling formalisms for dynamic structure systems. *ACM Transactions on Modeling and Computer Simulation* 7 (1997), 501–515.
2. Bauer, D., Yaun, G., Carothers, C. D., Yuksel, M., and Kalyanaraman, S. Seven-o'clock: A new distributed gvt algorithm using network atomic operations. In *Proceedings of the 19th Workshop on Principles of Advanced and Distributed Simulation*, PADS '05, IEEE Computer Society (Washington, DC, USA, 2005), 39–48.
3. Chandy, K. M., and Misra, J. Asynchronous distributed simulation via a sequence of parallel computations. *Commun. ACM* 24, 4 (Apr. 1981), 198–206.
4. Chen, B., and Vangheluwe, H. Symbolic flattening of DEVS models. In *Summer Simulation Multiconference* (2010), 209–218.
5. Chow, A. C. H., and Zeigler, B. P. Parallel DEVS: a parallel, hierarchical, modular, modeling formalism. In *Proceedings of the 26th Winter Simulation Conference*, SCS (1994), 716–722.
6. De Munck, S., Vanmechelen, K., and Broeckhove, J. Revisiting conservative time synchronization protocols in parallel and distributed simulation. *Concurrency and Computation: Practice and Experience* 26, 2 (2014), 468–490.
7. Fujimoto, R. M. Performance of Time Warp under synthetic workloads. In *Proceedings of the SCS Multiconference on Distributed Simulation* (1990).
8. Fujimoto, R. M. *Parallel and Distributed Simulation Systems*, 1st ed. John Wiley & Sons, Inc., New York, NY, USA, 1999.
9. Fujimoto, R. M., and Hybinette, M. Computing global virtual time in shared-memory multiprocessors. *ACM Trans. Model. Comput. Simul.* 7, 4 (Oct. 1997), 425–446.
10. Ghemawat, S., and Menage, P. TCMalloc : Thread-Caching Malloc. <http://goog-perftools.sourceforge.net/doc/tcmalloc.html>, Nov. 2005.
11. Glinsky, E., and Wainer, G. DEVStone: a benchmarking technique for studying performance of DEVS modeling and simulation environments. In *Proceedings of the 2005 9th IEEE/ACM International Symposium on Distributed Simulation and Real-Time Applications* (2005), 265–272.
12. Glinsky, E., and Wainer, G. New parallel simulation techniques of DEVS and Cell-DEVS in CD++. In *Proceedings of the 39th annual Symposium on Simulation* (2006), 244–251.
13. Himmelspach, J., and Uhrmacher, A. M. Sequential processing of PDEVs models. In *Proceedings of the 3rd European Modeling & Simulation Symposium* (2006), 239–244.
14. Jafer, S., and Wainer, G. Conservative vs. optimistic parallel simulation of devs and cell-devs: A comparative study. In *Proceedings of the 2010 Summer Computer Simulation Conference*, SCSC '10 (2010), 342–349.
15. Jefferson, D. R. Virtual time. *ACM Trans. Program. Lang. Syst.* 7, 3 (July 1985), 404–425.
16. Kim, K. H., Seong, Y. R., Kim, T. G., and Park, K. H. Distributed simulation of hierarchical DEVS models: Hierarchical scheduling locally and time warp globally. *Transactions of the SCS* 13, 3 (1996), 135–154.



17. Mattern, F. Efficient algorithms for distributed snapshots and global virtual time approximation. *Journal of Parallel and Distributed Computing* 18, 4 (1993), 423–434.
18. Muzy, A., and Nutaro, J. J. Algorithms for efficient implementations of the DEVS & DSDEVS abstract simulators. In *1st Open International Conference on Modeling and Simulation (OICMS)* (2005), 273–279.
19. Muzy, A., Varenne, F., Zeigler, B. P., Caux, J., Coquillard, P., Touraille, L., Prunetti, D., Caillou, P., Michel, O., and Hill, D. R. C. Refounding of the activity concept? towards a federative paradigm for modeling and simulation. *Simulation* 89, 2 (2013), 156–177.
20. Nethercote, N., and Seward, J. Valgrind: A framework for heavyweight dynamic binary instrumentation. *SIGPLAN Not.* 42, 6 (jun 2007), 89–100.
21. Nutaro, J. J. ADEVs.  
<http://www.ornl.gov/~1qn/adevs/>, 2015.
22. OpenMP Architecture Review Board. OpenMP application program interface version 4.5, 2015.
23. Troccoli, A., and Wainer, G. Implementing Parallel Cell-DEVS. In *Annual Simulation Symposium* (2003), 273–280.
24. Van Tendeloo, Y. Research internship i: Efficient devs simulation.
25. Van Tendeloo, Y., and Vangheluwe, H. Activity in pythonpdevs. In *Activity-Based Modeling and Simulation* (2014).
26. Van Tendeloo, Y., and Vangheluwe, H. The Modular Architecture of the Python(P)DEVS Simulation Kernel. In *Spring Simulation Multi-Conference, SCS* (2014), 387 – 392.
27. Van Tendeloo, Y., and Vangheluwe, H. PythonPDEVs: a distributed Parallel DEVS simulator. In *Proceedings of the 2015 Spring Simulation Multiconference, SpringSim '15*, Society for Computer Simulation International (2015), 844–851.
28. Vangheluwe, H. DEVS as a common denominator for multi-formalism hybrid systems modelling. *CACSD. Conference Proceedings. IEEE International Symposium on Computer-Aided Control System Design* (2000), 129–134.
29. Zeigler, B. P., Praehofer, H., and Kim, T. G. *Theory of Modeling and Simulation*, second ed. Academic Press, 2000.