Mixed Shared-Distributed Hash Tables
Approaches for Parallel State Space Construction *

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Abstract

We propose an algorithm for parallel state space construction based on an original concurrent data structure, called a localization table, that aims at better spatial and temporal balance. Our proposal is close in spirit to algorithms based on distributed hash tables, with the distinction that states are dynamically assigned to processors; i.e. we do not rely on an a-priori static partition of the state space.

In our solution, every process keeps a share of the global state space. Data distribution and coordination between processes is made through the localization table, that is a lockless, thread-safe data structure that approximates the set of states being processed. The localization table is used to dynamically assign newly discovered states and can be queried to return the identity of the processor that own a given state. With this approach, we are able to consolidate a network of local hash tables into an (abstract) distributed one without sacrificing memory affinity – data that are “logically connected” and physically close to each others – and without incurring performance costs associated to the use of locks to ensure data consistency.

We evaluate the performance of our algorithm on different benchmarks and compare these results with other solutions proposed in the literature and with existing verification tools.

1 Introduction

Model-checking is a very demanding activity in terms of computational resources. As a result, the extensive need for memory and computation power has resulted in the design of model checking algorithms that target parallel and distributed machines. Variations between these algorithms are often explained by differences between the targeted architectures – shared-memory versus distributed memory, clusters, ... – or differences on the criteria to optimize – achieving better spatial balance between processes, lowering synchronization costs, etc.

We propose an algorithm for parallel state space construction intended for shared memory, multiprocessor machines. (We focus here on the exhaustive generation of the state space of finite-state transition systems, often a preliminary step for model-checking.) The basic idea behind a state space construction algorithm is pretty simple: take a state that has not been explored (a fresh state); compute its successors and check if they have already been found before; iterate until there is no more new state to explore. Hence, a key point for performance is to use an efficient data structure for storing the set of generated states and for testing membership in this set.

In our approach, the goal is to build the state space of a system concurrently in such a way that: (1) the share of state space built by each processor is as uniform as possible; and (2) the processor occupancy is maximal. Our algorithm builds on previous work [19] and is based on the same simple design: the global state space is stored in a set of local containers (e.g. hash tables), each controlled by a different processor, while only a small part of the shared-memory is used for coordinating the state space exploration. This is close in spirit to algorithms based on distributed hash tables, with the distinction that we choose to dynamically assign states to processors, that is, we do not rely on an a-priori static partition of the state space.

Data distribution and coordination between processes is made through a localization table (LT), that is a lockless, thread-safe data structure that approximates the set of states being processed. The localization table is used to dynamically assign newly discovered states and behaves as an associative array that returns the identity of the processor that own a given state. With this approach, we are able to con-
solidate a network of local hash tables into an (abstract) distributed one without sacrificing memory affinity – data that are "logically connected" and physically close to each others – and without incurring performance costs associated to the use of locks to ensure data integrity.

The paper is organized as follows. In Section 2 we review the related work. Section 3 details our algorithm and defines the data structure for localization tables. Before concluding, we report on experimental results obtained on a set of typical benchmarks and compare our approach with solutions already proposed in the literature.

2 Related Work

Model Checking [7] is an automated verification method used to check whether the model of a system meets a given specification. This technique relies on the exploration of the state space of the model. State space construction can be classified as an irregular parallel problem because state graphs may be highly irregular, see [8] for a discussion on this topic. As a consequence, when parallelizing this problem, special attention should be taken to ensure a good load balancing among processors.

Several approaches have been proposed, since the early 1990s, for parallel and distributed state space exploration. These solutions adopt, in their vast majority, a common paradigm that could be labeled as “homogeneous” parallelism – a Single Program Multiple Data (SPMD) programming style – such that each processor performs the same steps concurrently. Most of these solutions, to mention a few [1, 17, 6, 9, 20], were intended for distributed computers and rely on slicing functions – that is functions that statically assign a state to a processor – and basically only differ by the nature of these functions. The choice of slicing functions has a major influence on the load balance and data locality of the algorithms.

A smaller number of solutions target shared memory machines [10, 11, 12, 13, 1]. Like in our case, some of these solutions are based on the use of hash tables to store the states. An example is the version of DiVinE [2] for multicore machines, that is based on a static partitioning scheme where each process owns a private hash table. Another example is the parallel version of Spin [11, 10] which uses the stack-slicing strategy to share work in combination with a shared hash table protected by a fine-grained locking scheme. This work has been recently extended with a lockless shared hash table based on atomic primitives [13] (CAS – Compare & Swap). Additionally, we mention the work of Inggs et al. [12], which proposes a parallel algorithm based on a work stealing scheduling paradigm to provide dynamic load balancing. In their case, the data structure used to store the states is an “unsafe” shared hash table.

In the context of this work, we propose an extension of an algorithm that we defined in [19], which is based on two data structures: a lock-free, shared Bloom filter [4] to coordinate the data distribution; and local data containers to explicitly store the data (for instance, our initial implementation was based on AVL trees). The Bloom filter is used to represent, in a very compact way, the set of states that have already been found and to efficiently test whether a given state has already been found. Due to the probabilistic nature of the Bloom Filter, the algorithm is based on multiple iterations in order to perform an exhaustive exploration. In a first phase, the algorithm is guided by the Bloom filter until no new states can be found. During this phase, states found by a processor are stored locally in two dictionaries: one for states that, according to the Bloom filter, have also been found by another processor; the other for fresh states. Since the Bloom filter may, at times, falsely report that a state has already been visited, we need to handle these collision states in a second phase of the algorithm. The computation stops when there are no more states to explore and no more collisions.

Our Contributions: The algorithm proposed in this paper improves on our previous design and replaces the Bloom Filter by a dedicated data structure, the localization table. Unlike Bloom Filters, this data structure can be used to find the processor that owns a given data item – a state in our case – and not only if the object was already found. This simple addition significantly enhance the performance of our previous algorithm and also simplifies its logic. Indeed, it is now possible to solve possible collisions on-the-fly and to get rid of the collision resolution phase.

Our contributions are twofold. First, in the formal verification domain, we define a new algorithm for parallel state space construction. Our algorithm is able to exploit parallelism in all possible cases and, unlike algorithms based on slicing functions or heuristic rules, is compatible with dynamic load-balancing techniques. Second, in the parallel computing domain, the combination of local hash tables with a localization table provides an interesting implementation for concurrent hash maps that may be useful in other situations.

Our preliminary results are very promising. We observe performances close to those obtained using an algorithm based on lockless hash tables (that may be unsafe) and that outperforms an implementation based on the concurrent, unordered map provided in the Intel Threading Building Blocks [18], an industrial strength lockless hash table.

3 Description of the Algorithm

Our algorithm follows a “homogeneous” parallelization approach, where all processors execute the same program simultaneously and each processor handles its own local
view of the state space. Coordination between the processors is based on a Localization Table (LT for short), that is used to allocate newly discovered states to processors.

The LT is used to test whether a state has already been found and, if so, to keep track of the location – the processor id – where the state is held. The work performed by each processor is pretty simple: generate a state using the model of the system, say \( s \), and check in the LT where it could have potentially been assigned. If \( s \) is a newly discovered state, it will be assigned to the processor who generated it. Otherwise, the LT will return the location where the state \( s \) is assigned, say \( LT(s) \). This approach has the advantage to isolate the local hash tables; each processor has exclusive write access to its local table, whereas concurrent read accesses are unrestricted. As a result, we consolidate a network of distributed hash tables into a single, concurrent data structure. Another advantage is that we can easily resize local hash tables, as needed, without blocking the entire exploration.

We describe more precisely how the LT is implemented in Section 3.1. An advantage of our design is to be thread-safe: operations on a LT are simple and can be implemented using atomic actions. Another advantage is the small footprint of the LT. To summarize, our goal is to combine the advantages of distributed and shared hash tables for parallel state space construction in a single algorithm.

In the remainder of the text, we use \( N \) to denote the number of processors and \( 1..N \) for the natural interval between 1 and \( N \). We define the localization table in Sect. 3.1. The work-sharing techniques used in our algorithm is discussed in Sect. 3.2 while Sect. 3.3 gives some pseudo-code and further explanations about the algorithm.

### 3.1 Localization Table

Storing the relation \((s, LT(s))\) – associating each state, \( s \), with the processor that owns it – in a single table would require a very large amount of memory. Actually, it would defeat the need to store the state itselfs. Instead, the idea is to use a notion of a key associated to a state and to store the association between keys and processors. In our implementation, keys are computed using hash-functions and we will use a scheme based on multiple keys.

A localization table is essentially a “table” that associates a processor id – a value in \( 1..N \) – to every key in the table. A straightforward implementation is to use an integer vector for the underlying table.

We can implement the table using a vector \( V \) of size \( n \) and, for computing the key of a state, an independent hash function, \( h \), with image in \( 1..n \). In this case, we can check if a state \( s \) has already been found by looking into the local table of processor \( V[h(s)] \). While this implementation is simple, its disadvantage is that it is not possible to ensure a fine dynamic distribution of the states if \( h \) is not uniform. Indeed, if processor \( id_1 \) finds a new state, \( s \), such that \( V[h(s)] = id_2 \), then we need to transfer \( s \) between the two processors. A solution is to increase the size of the vector – but this has a direct impact on the memory consumption – or to use better hashing functions – but this has an impact on the performances.

We propose another implementation of the localization table that improves upon the choice of a vector. Inspired from our previous experience with a Bloom Filter (BF), the idea is to use a finite family of hashing functions \( h_1, \ldots, h_k \). To test if a state \( s \) is in the LT, we search if the key \( h_1(s) \) is in \( LT \). If it not, we know that the state is fresh. If \( LT(h_1(s)) = id_1 \) then we check if \( s \) is in the local table of processor \( id_1 \). If \( LT(h_1(s)) \neq id_1 \), we continue searching with the key \( h_2(s) \) and so forth.

This is only a rough description of how the LT works. Next, we define more formally the operation of our data structure. In particular, we explain how to deal with states that are not in the processors \( LT(h_1(s)), \ldots, LT(h_k(s)) \), that we call collision states. By convention, our algorithm will route a collision state to the last processor found, that is \( LT(h_k(s)) \).

A Localization Table, \( L \), is defined by two parameters: its size \( n \); and a family of \( k \) independent hashing functions \( h_1, \ldots, h_k \) with image in \( 1..m \), where we choose \( m \) such that \( n \ll m \) (see the discussion about the ratio \( n/m \) in Sect. 4.2).

A localization table \( L \) of size \( n \) is an array of size \( n \) containing pairs of the form \((p, d)\), where \( p \) is a processor id \((p \in 1..N)\) and \( d \) is a key \((d \in 1..m)\). To look for values inside of \( L \), we use a fixed surjective function \( map \), from \( 1..m \) to \( 1..n \). Hence, to check the value associated to the key \( h_i(s) \), we look in the array \( L \) at index \( map(h_i(s)) \).

Initially, an empty \( LT \) is an array initialized with the value \((0, 0)\). Assume that the processor \( id \) attempts to insert a state \( s \) into \( L \). The function takes as input a state and a processor id and returns a pair made of: a status, to determine if the element is new (or old); and the identifier of the processor who owns the state. The insertion operation is performed by looking successively at the elements with index \( map(h_i(s)) \) in \( L \) for all \( i \in 1..k \). There are four possible cases at each step:

- if \( L[map(h_i(s))] = (0, 0) \) then we know for sure that the state is fresh (it has never been added before). We can stop our iteration and write the pair \((id, h_i(s))\) in \( L \). This can be done using an atomic compare and swap operation;

- if \( L[map(h_i(s))] = (id', d) \) and \( d \neq h_i(s) \) then we cannot decide if the state \( s \) has been found and we continue to the next iteration, with the key function \( h_{i+1} \);
• if $$L[map(h_i(s))] = (id', h_i(s))$$ then we answer that
  $$s$$ is in the local table of processor $$id'$$ with high
  confidence. With this approach, states with the same hash
  value are not handled at the LT level. In our algo-
  rithm, these collisions will be spotted when the pro-
  cessor tries to recover the state from the local table of
  $$id'$$. In order to keep the consistency of the LT and to
  prevent states from being stored more than once, we
  choose to assign $$s$$ to the processor $$id'$$ and this is han-
  dled like a collision state;

• if we cannot decide after checking the values of
  $$L[map(h_k(s))]$$ for $$i \in 1..k$$, we also say that $$s$$ is a
  collision state and we choose to assign $$s$$ to the proces-
  sor $$id'$$ such that $$L[map(h_k(s))] = (id', d_i)$$. The operation for checking whether a state $$s$$ is already
  in the LT is very similar to the insertion function. We
  test successively if there is an index $$i \in 1..k$$ such that
  $$L[map(h_i(s))] = (id_i, h_i(s))$$, stopping if one of the
  position in $$L$$ is empty. If this is the case, we know that $$s$$ is
  not in the LT. If we find no match after $$k$$ attempts, then we
  consider that $$s$$ is a collision state that belongs to
  $$id_i$$.

In Figure 1 we illustrate the insertion and test operation
for three data items: $$x$$, $$y$$, and $$z$$; performed in this order
by the processors $$P_1$$, $$P_2$$ and $$P_3$$. The figure display a LT
of size $$n = 4$$ with two independent hash functions $$h_1$$ and
$$h_2$$. We assume that $$k = 2$$ and that $$m = 31$$. The insertion
of $$x$$ requires only one operation since the slot at position
$$map(h_1(x))$$ is initially empty. As a result the slot is asso-
ociated to processor $$P_1$$ for elements with key 17. Element $$y$$
  is inserted at the second attempt, since the slot in position
  $$map(h_1(y))$$ is already filled and that $$h_1(y) \neq d_1$$. Finally,
  element $$z$$ cannot be properly inserted – it is a collision –
  and it is assigned to processor $$P_2$$. The state space exploration proceeds until no new states
  are stored in the local table of
  processor $$id_i$$.

  • the **private minimum workload**, which defines the minimal
    charge of work that should be kept private. The processor
    will share work only if the charge in its private
    stack is larger than this value;

• the **share workload**, which defines the ratio of work
  that should be added in the shared stack if the load in
  the private stack is larger than the private minimum
  workload.

Our implementation of the work-stealing paradigm is inter-
esting in its own right since it differs from [12] by its use
of unbounded shared stacks and the use of a “share work-
load” parameter. The description of this extension is also
interesting since it shows that common optimizations and
load-balancing techniques are not precluded by our algo-

3.3 Algorithm

To conclude this section, we give a high-level view of
our algorithm that can be described by the pseudo-code of
Listing 1. The diagram of Fig. 2 describes the shared and
local data structures used in the algorithm. Each processor
manages a “private work” stack of unexplored states and a
local hash table to store the states assigned to him. The
shared values are: the Localization Table; one bitvector
of size $$N$$ to store the state of the processors (idle or busy),
used to detect termination; $$N$$ stacks – one for each proces-
sor – for the work sharing technique described in Sect. 3.2;
and finally $$N$$ collision stacks used to route collisions states
to their correct processors.

The state space exploration proceeds until no new states
  can be added to the LT and all stacks are empty. Given
  a state $$s$$, a processor, say my_id, will check the LT to
test whether $$s$$ is new and, otherwise, what is the owner of
  $$s$$. This information is returned by a call to the function
test_or_insert($$s$$, my_id). During the exploration, states that
  are labeled as new by the LT are stored in the local table of
  depending on the processor occupancy. First, we use an ac-
  tive technique very similar to the work-stealing paradigm
  of [12]. This mechanism uses two stacks: a private stack
that holds all states that should be worked upon; and a
shared stack for states that can be borrowed by idle proces-
sors. This shared stack is protected by a lock to take care of
  concurrent access. The second technique can be described
  as passive and has the benefit to avoid useless synchroniza-
  tion and contention caused by the active technique. In the
  passive mode, an idle processor waits for a wake-up signal
from another processor willing to give away some work in
stead of polling other shared stacks. The shift between the
  passive and active modes is governed by two parameters:

• the **private minimum workload**, which defines the minimal
  charge of work that should be kept private. The processor
  will share work only if the charge in its private
  stack is larger than this value;

• the **share workload**, which defines the ratio of work
  that should be added in the shared stack if the load in
  the private stack is larger than the private minimum
  workload.

3.2 Work-Sharing Techniques

Our algorithm relies on two different work-sharing tech-
niques to balance the work load between processors. We use
these mechanisms alternately during the exploration phase
depending on the processor occupancy. First, we use an ac-

tive technique very similar to the work-stealing paradigm
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from another processor willing to give away some work in
stead of polling other shared stacks.
The occupancy rate measures the utilization of the machine CPUs whereas the system occupancy rate may vary between 90% (Hanoi model) and 51% (Kanban model), indicating the model. For instance, our efficiency varies between 90% (Hanoi model) and 51% (Kanban model), whereas the system occupancy rate is consistently over 95%. Clearly, the algorithm depends on the “degree of concurrency” of the model – it is not necessary to use lots of processors for a model with few concurrent actions – but this is an inherent limitation with parallel state space construction [8], which is an irregular problem.

### 4 Experiments

We implemented our algorithm using the C language with Pthreads [5] for concurrency and we developed a library for the Localization Table with support for concurrent insertions. The Experimental results presented here are obtained using a Sun Fire x4600 M2 Server, configured with 8 dual core opteron processors and 208GB of RAM memory, running the Solaris 10 operating system. This machine is classified as Non-Uniform Memory Access (NUMA) because the shared memory is physically divided among the processors.

For the benchmark, we used models taken from two sources. We have three classical examples: Dining Philosophers (PH); Flexible Manufacturing System (FMS); and Kanban – taken from [14] – together with 5 Puzzles models: Peg-Solitaire (Peg); Sokoban; Hanoi; Sam Lloyd’s puzzle (Fifteen); and 2D Toads and Frogs puzzle (Frog) – taken from the BEEM database [16]. All these examples are based on finite state systems modelled using Petri Nets [15]. This means that, in these cases, a state is a marking, that is a tuple of integers. Our algorithm may be adapted to other formalisms, for instance including data, time, etc.

With our computer setup, we are able to tackle examples with approximately 10 billions of states, but we selected models with less than 500 millions of states in order to complete our experiments in reasonable time. (A complete run of our benchmark takes more than a week to finish.)

We study the performance of our implementation on different aspects. While speedup is the obvious criteria, we also study the memory footprint of our approach and the physical distribution of states among processors.

#### 4.1 Speedup and Physical Distribution

In Fig. 3 we give the observed speedup of our algorithm on a set of examples. We give the absolute speedup, measured as the ratio between the execution time using $N$ processors ($T_N$) and the time of an optimized, sequential version. Our implementation delivers some promising speedups. The results also show different behaviors according to the model. For instance, our efficiency may vary between 90% (Hanoi model) and 51% (Kanban model), whereas the system occupancy rate is consistently over 95%. Clearly, the algorithm depends on the “degree of concurrency” of the model – it is not necessary to use lots of processors for a model with few concurrent actions – but this is an inherent limitation with parallel state space construction [8], which is an irregular problem.

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1. Efficiency is computed as the ratio between speedup, $T_N$, and the number $N$ of processors.
2. The occupancy rate measures the utilization of the machine CPUs.
Concerning the use of memory, we can measure the quality of the distribution of the state space using the mean standard deviation (σ) of the number of states among the processors. In our experiments, we observe that the value of σ is quite small and that it stays stable when we change the number of processors (see Fig. 8). For instance, we have σ ≈ 1.5% for the Hanoi model and σ ≈ 7% for Kanban. The difference between values of σ can be explained by the difference in the “degree of concurrency”. It may also be affected by the processor’s performance, that is, a processor that handles “simpler states”, or smaller work units, may dynamically assign more states than others. Finally, our experiments are also affected by the Non-Uniform Memory Access (NUMA) architecture of our machine, where the latency and bandwidth characteristics of memory actions depend on the processor or memory region being accessed.

4.2 Localization Table Configuration and Memory Footprint

The LT data structure is configured using two parameters: its dimension (n) and the number of hash-functions keys (k). The values of these parameters have an impact on the performance. If the dimension is to small, the LT will get quickly saturated and the number of collisions will increase. Ideally, a LT of size n is sufficient for a space of n states. However, hash functions are not perfect (uniform), which affects our structure just like with standard hash table. In our experiments, we observe that LT behaves well for load factors (ratio between the number of states in the LT and its dimension) lower than 0.7.

In Fig. 4, we display the ratio (in percentage) between the number of collisions and the size of the state space, on the Kanban model (∼ 4.10^8), for three different values of the load factor and for different values of k. This Fig. also depicts the absolute number of collisions for each experiment. For instance, we have approximately 58.10^5 collisions (ratio of 1.5%) when LT is set with n = 2^30 ≈ 10^9 (load factor of 0.36) and k ≤ 8.

Likewise, in Fig. 5 we show the impact of different load factors (choice of the size of the LT) on the execution time of the algorithm for a fixed model. Once again, we observed that our LT gives better results when the load factor is in average smaller than 0.7.

For the speedup results given in this Section, we have adjusted the dimension of the LT to obtain load factors smaller than 0.5 for every models and we have chosen to limit ourselves to at most four hash-functions keys (k ≤ 4). We decided to fix these settings beforehand in order to not artificially improve our results and also to show the memory efficiency of our solution. To illustrate this point, we may observe that in the experiments of Fig. 5, the size of the LT is of 1GB (that is approximately one billions data items) for a load factor of 0.36, which is the only significant memory overhead used by our solution.

4.3 Comparison With Other Tools and Other Algorithms

We conclude this section with a comparison with other algorithms. We developed our own implementation of some classical parallel algorithms based on the use of hash tables. In Fig. 6, we briefly describe the different implementations used for this comparison. We decided to skip a comparison with our previous work [19] because our current results are at least twice as good.
# Table of Algorithms

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<th>Description</th>
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<tr>
<td>Vector</td>
<td>Vector of integers like structure: Localization Table with only one key</td>
</tr>
<tr>
<td>Static</td>
<td>States are distributed using a static slicing function</td>
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<tr>
<td>Lockless</td>
<td>Lockless shared hash table as the shared space</td>
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<tr>
<td>TBB</td>
<td>Unordered hash map as the shared space, from Intel-Threading Building Blocks library</td>
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Figure 6. Algorithms selected.

Figure 7 shows the average (absolute) speedups, over all models, for the different implementations. The Lockless implementation has the best performance but it is an unsafe solution, since states may be skipped [3]. All the other implementations are safe. We include the results for Lockless since it provides a good reference for performance. Our algorithm (LT) performs better than all the other implementations for all models. As we mentioned earlier, the difference in performance between Vector and LT is mainly due to the non-uniformity of hash functions. This difference is significant especially for Sokoban and Kanban models (see Kanban analysis at Fig 5). Concerning Static, an explanation for the better performances is that we exchange less states between processors: in some experiments with Static, we can observe that up to 96% of the states have not been found by their owner. The gain in performance compared to TBB (based on an lockless, non-lossy hash table found in the Intel-TBB [18]) may be explained by the adequacy of our data structure to our targeted application (state space generation). Indeed, in this application, we have many more reads than writes (state spaces have more transitions than states). The LT has several benefits in this case: (1) it delivers a low complexity mechanism to grant exclusive write access for the local hash-table; (2) the structure is cache-friendly since data are stored in-place (avoiding pointers); and (3) the use of local hash tables improves memory affinity, which is important for NUMA machines.

Concerning the memory distribution, we display the average mean-standard deviation for all implementations in Fig. 8. The results show that the best distribution, by far, is from the Static version. We can observe that all other implementations have similar distributions. (The anomalous values for $N = 16$ can be explained by the fact that, in this case, we use all the processors of our computer.) In the context of this work, we use no heuristics to ensure an uniform partition of states, so the quality of the distribution depends on the model “degree of concurrency” and the performance of processors.

We have also compared our implementation with “state of the art” verification tools that provide a parallel implementation. We looked both at the Spin and DiVinE tools. We give some performance results but it is difficult to make a fair comparison. For one thing, it has proved difficult to port available implementations on the configuration used for our experiment. For instance, our benchmark with DiVinE and Spin are based on Linux instead of Solaris, which means that we take advantage of more efficient libraries. On the other hand, a major discrepancy lies in the fact that we compare an algorithm with a tool. For instance, we do not make use of any “general optimizations” techniques, such as local caches, data-alignment optimizations, etc. Also, while Spin work with compiled models, we currently use interpreted models. DiVinE accepts both models but we use for this comparison their interpreted variant. After these words of caution on the significance of the comparison, we give some results obtained on the Sokoban model. Using the parallel version of Spin on our benchmarks, we observe a maximum speedup of $3.6$ using 8 cores (73s). Nonetheless, the sequential performance of Spin (264s) is about 3 times better than our prototype implementation of LT. In our experiments, LT is marginally faster than Parallel Spin when both are running on 12 cores and is faster using the 16 available cores. The computation for Spin is of 81.2s for 12 cores and 82.3 for 16 cores, while we generate the state space in 80s with 12 cores and 64s with 16 cores using LT. Concerning DiVinE – whose sequential performance is about 40% better than our prototype implementation – LT matches the performance of DiVinE when both are using 10 cores and outperforms it of about 20% using 16 cores. More
precisely, the running time for DiVinE is of 96.5s with 10 cores – while LT time is of 96s – and 84.9s with 16 cores. It is possible to connect this result with the comparison given in Fig. 7. Indeed, DiVinE is based on a static slicing function to distribute the states – as in the static implementation of Fig. 6 – and the difference of performance between LT and Static is of about 30% on 10 cores and of almost 70% for 16 cores. These preliminary results against two of the most popular parallel model-checker are very encouraging since we have a prototype implementation of LT.

5 Conclusion

We define a new parallel state space construction algorithm targeted at shared memory machines. The main innovation lies in a new data structure, named Localization Table, that is used to coordinate a network of local hash tables in order to obtain an efficient concurrent hash map. This new structure replaces the Bloom Filter that was used in one of our previous work [19]. The LT is used to dynamically assign newly generated states and behaves as an associative array that returns the identity of the processor that owns a given state.

A first implementation of our algorithm shows promising results as we observed speedups consistently better than with other parallel algorithms. For instance, our experimental results show that efficiency varies between 90% and 50%, depending on the “degree of concurrency” of the model. In addition, our memory footprint is almost negligible when compared to the total memory used for storing the state space. For example, in the worst case (the Kanban model, with 25GB) we consume less than 4% of the memory for the LT and the different stacks used by our algorithm. That is approximately 1GB of memory. The same benchmark also shows that our implementation fares well when compared with related tools. Indeed, our experiments show that our solution performed very well against an industrial strength lockless hash table, the concurrent hash map implementation provided in the Intel-TBB. This may be explained by the fact that we provide a concurrent data structure for encoding sets that is optimized for the case where deletions are very rare and the same item may be inserted several times, whereas the Intel-TBB provides a general implementation. This is a very encouraging since we obtained these results with minimal optimizations (i.e., without resorting to global caches, data-alignment optimizations, etc.), so there is still room for improvements. Altogether, our solution fulfilled our goal of having both the best temporal and spatial balance as possible.

For future work, we are investigating a probabilistic version of our current exhaustive algorithm. In this context, the term probabilistic denotes an algorithm that builds an underapproximation of the global state space, with a very high probability of building the exact state space (by very high, we mean a probability of failure less than $10^{-5}$).

References