Multi-core On-the-fly Saturation Tom van Dijk Jeroen Meijer Jaco van de Pol TACAS 2019

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- TACAS 2001: Saturation original paper
- TACAS 2003: Saturation Unbound with on the fly transition learning
- QEST 2004: Saturation NOW network of workstations
- CAV 2007: Parallelising symbolic state-space generators multi-core, using work-stealing

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"Parallel symbolic state-space exploration is difficult, but what is the alternative?"

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"Parallel symbolic state-space exploration is difficult, but what is the alternative?"

• TACAS 2019: Multi-core On-the-fly Saturation (this work) almost 8× faster with 16 cores

Background

- On-the-fly model checking
- Using decision diagrams
- Saturation algorithm for transition relations
- The parallel decision diagram package Sylvan

Contribution

- How we parallelize saturation
- The tools that are available
- Our reproducible experiments

So what is on-the-fly model checking?

 $LTS{\ensuremath{\mathsf{SMIN}}}$ a backend talks to a model via a $N{\ensuremath{\mathsf{NTT-STATE}}}(s)$ interface

- Start with some initial state and no transitions
- Ask the model for successors
- Update transition relation with every new transition
- Ask the model for more successors
- Allows e.g. checking safety properties, LTL, etc, "on-the-fly"

On-the-fly model checking



Partitioned transition relation

- Often transitions are local and independent
- Example: Petri nets, composed systems, etc
- Disjunctive partition into multiple transition relations
- "Short" transitions only affect some variables

 $\rm LTSMIN$ implements a "Partitioned Next-State Interface" and models provide $\rm LTSMIN$ with a dependency matrix relating state variables and transitions

Advantage: few NEXT-STATE calls to learn exponentially larger model

On-the-fly model checking

Example Petri net



- Each transition is local and independent
- Few NEXT-STATE calls to learn all reachable transitions

LTSMIN

- On-the-fly compute product with automata of specifications
- On-the-fly check safety and reachability
- $\mathrm{LTS}_{\mathrm{MIN}}$ support various languages and backends via the interface
 - Promela, Petri nets (PNML), DVE, mCRL2, ProB languages, etc

TACAS 2015: Gijs Kant, Alfons Laarman, Jeroen Meijer, Jaco van de Pol, Stefan Blom and Tom van Dijk, LTSMIN: High-Performance Language-Independent Model Checking

- States and transitions are sets of Boolean vectors \mathbb{B}^k and \mathbb{B}^{2k}
- State $x \in \mathbb{B}^k$ is in the set if it is reachable
- Transition $(x,y)\in \mathbb{B}^{2k}$ is in the set if $x\rightarrow y$ is a transition

- Set of reachable states S(x)
- Transition relations: $\mathcal{T}_1(r_1, w_1') = \mathcal{T}_2(r_2, w_2') = \dots = \mathcal{T}_M(r_M, w_M')$

With x all state variables, and r_i the "read" variables of transition i, w_i' the "write" variables

Binary decision diagrams

- A BDD is a directed acyclic graph encoding a $\mathbb{B}^k\to\mathbb{B}$ function or a $S\subseteq\mathbb{B}^k$ set
- Every node has label x and two children then and else
- Semantics: if x then follow then else follow else
- Paths represent valuations of \mathbb{B}^k
- For sets: path to $\begin{bmatrix} 1 \\ -1 \end{bmatrix}$, then valuation is in the set



Motivation

- Transitions are local, only affect a part of the BDD
- Changing a node affects all ancestors
- Possibly waste of effort to recompute these ancestors!



State space exploration strategies

- Breadth-first: apply all transition relations once to current states, add all results to set of states
- Chaining: apply transition 1 and add to set of states, apply transition 2 and add to set of states, etc
- (and learn new transitions before applying each transition relation)

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Saturation strategy [Gianfranco Ciardo et al, since 2001]

- Always apply "deepest" transition until "saturated"
- Whenever we find new states, saturate deeper transitions first
- Saturate BDD nodes in-situ
- Considered an "optimal strategy"

How we parallelize saturation

- SIMPLIFY the saturation idea:
 - no in-situ updates of BDD nodes
 - no mutual recursion of "saturate" and "fire event"
 - implement a single BDD operation "saturate"
- Leverage the multi-core BDD framework Sylvan
 - Already parallelizes many BDD operations
 - relational product
 - other popular BDD operations required for model checking
 - Already offers all the infrastructure for new BDD operations
- Implement for standard BDDs but also for LDDs (multi-way decision diagrams implemented as linked list structures)

Sylvan

- https://www.github.com/trolando/sylvan (google "github sylvan", mirror at utwente-fmt repository)
- BDD but also framework for MTBDD with any type of terminal For example in probabilistic model checkers Storm, IscasMC, ePMC, with floating points, rational numbers, parameterised functions, etc
- Can also extend for other types of decision diagrams like LDDs
- Implements all common MTBDD operations internally parallelized Your program is sequential, automatically parallelized
- Relies on the work-stealing framework Lace to do scalable fine-grained load balancing (every suboperation is a task)
- Main datastructures (nodes table, operation caches) lock-free scalable
- With Sylvan, parallel saturation is easy!

Algorithm

Saturate set of states S given current transition i (of k sorted transitions)

- Leaf cases: return S if S = 0, S = 1 or no more relations left (i = k)
- Check if result is in the cache
- Root BDD node of S: $\langle x, \mathsf{Then}, \mathsf{Else} \rangle$
- Is x accessed by transition i
- Yes?
 - First saturate deeper: $S \leftarrow \mathtt{saturate}(S, i+1)$
 - Then apply relation i to S with multi-core operation relprod
 - Repeat until no change
- No?
 - Run in parallel: saturate(Then, i) and saturate(Else, i)
 - Compute new BDD node of the result
- Store final result in the cache

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Learning: update transition just before running relprod

Parallel is easy!

- Task parallelism in Sylvan: spawn 2 tasks, then wait until they are done!
- Implemented with the work-stealing framework Lace
- Scalable parallel datastructures nodes tables and operation cache

Evaluation based on Petri nets of the Model Checking Contest

- Compare with state-of-the-art MEDDLY (sequential) (Algorithms are fundamentally different, but are we at least close?)
- Measure parallel speedup on 16 cores and 48 cores
- Compare with/without learning (on-the-fly or offline)
- (and compare with BDDs + compare to chaining and bfs)

Tools

- LTSMIN: on-the-fly transition learning; using LDDs
- LDDMC: offline (pre-learned) using LDDs
- MEDMC+MEDDLY: offline non-parallel version of Babar, Miner
- (and BDDMC + implementations of LDD chaining and bfs)

Empirical evaluation

Procedure

- Take Petri nets of MCC 2016 (491 input files)
- Use two good variable orders Sloan and Force for BDD variable ordering
- Use LTSMIN with *generous timeout* to generate transition system 413 out of 982 potential transition systems
- All tools use precisely the same model and variable order
- Compare runtimes only for inputs that all solvers can handle 301 inputs
- Please reproduce our results: (Apache 2.0) https://www.github.com/trolando/ParallelSaturationExperiments Artifact Evaluation Accepted
- Please download the tools: (Apache 2.0) https://www.github.com/trolando/sylvan https://www.github.com/utwente-fmt/ltsmin

Method		Number of solved models with # workers							
		1	2	4	8	16	Any		
LTSMIN	otf, par	387	397	399	404	407	408		
LDD _{MC}	par	388	393	399	402	402	404		
Meddly	seq	375	-	-	-	-	375		

Table: Number of benchmarks (out of 413) solved within 20 minutes with each method with the given number of workers.

Method	Order	Total time (sec) with # workers				Total speedup				
		1	2	4	8	16	2	4	8	16
LTSmin	Sloan	1850	1546	698	398	313	1.2	2.7	4.6	5.9
LDDmc	Sloan	932	609	311	194	151	1.5	3.0	4.8	6.2
Meddly	Sloan	572	–	-	–	–	-	–	-	-
LTSmin	Force	2704	1162	712	401	343	2.3	3.8	6.8	7.9
LDDmc	Force	856	602	348	216	180	1.4	2.5	4.0	4.7
Meddly	Force	1738	-	-	-	–	-	–	-	–

Table: Cumulative time and parallel speedups for each method-#workers combination on the models where all methods solved the model in time. These are 301 models in total: 151 models with Force, 150 models with Sloan.

Results

Model (with ldd-sat)	Order		Time (sec)		Speedup	
		1	24	48	24	48
Dekker-PT-015	Sloan	77.3	4.7	2.4	16.3	32.5
PhilosophersDyn-PT-10	Force	273.8	16.8	12.4	16.3	22.1
Angiogenesis-PT-10	Sloan	333.2	28.5	16.5	11.7	20.2
SwimmingPool-PT-02	Force	25.0	2.1	1.4	11.6	17.8
BridgeAndVehicles-PT-V20P10N20	Force	1035.8	101.8	60.7	10.2	17.1
Model (with otf-ldd-sat)						
Dekker-PT-015	Sloan	174.5	7.4	3.3	23.6	52.2
SwimmingPool-PT-07	Sloan	1008.0	69.2	42.0	14.6	24.0
SmallOperatingSystem-PT-MT0256DC0064	Sloan	957.3	52.9	40.0	18.1	23.9
Kanban-PT-0050	Sloan	940.6	78.7	48.9	11.9	19.2
TCPcondis-PT-10	Force	68.4	5.7	3.8	11.9	17.8

Table: Parallel speedup for a selection of benchmarks on the 48-core machine (only top 5 shown)

Conclusions

- \bullet Saturation is parallelized in $\mathrm{LTS}_{\mathrm{MIN}}$ and SylvaN
 - Parallel saturation now available for many modeling languages
 - Promela, Petri nets (PNML), DVE, mCRL2, ProB languages, etc
- Online available under Apache 2.0 License
 - github.com/utwente-fmt/ltsmin
 - github.com/trolando/sylvan
 - github.com/trolando/ParallelSaturationExperiments
- Competitive: often as good as or better than Meddly, especially with 2 or more cores
- Scalable: up to 7.9× on-the-fly on a 16-core machine (with the FORCE variable ordering); even some superlinear speedups
- Similar scalability with BDDs, with on-the-fly learning
- Parallel saturation is easy with Sylvan!