

Anytime system level verification via parallel random exhaustive hardware in the loop simulation[☆]



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ABSTRACT

System level verification of cyber-physical systems has the goal of verifying that the *whole* (i.e., software + hardware) system meets the given specifications. Model checkers for hybrid systems cannot handle system level verification of actual systems. Thus, Hardware In the Loop Simulation (HILS) is currently the main workhorse for system level verification. By using model checking driven exhaustive HILS, System Level Formal Verification (SLFV) can be effectively carried out for actual systems.

We present a *parallel random exhaustive* HILS based model checker for hybrid systems that, by simulating *all* operational scenarios *exactly once* in a *uniform random* order, is able to provide, at *any time* during the verification process, an *upper bound* to the probability that the System Under Verification exhibits an error in a yet-to-be-simulated scenario (Omission Probability).

We show effectiveness of the proposed approach by presenting experimental results on SLFV of the Inverted Pendulum on a Cart and the Fuel Control System examples in the Simulink distribution. To the best of our knowledge, no previously published model checker can *exhaustively* verify hybrid systems of such a size and provide at *any time* an upper bound to the Omission Probability.

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1. Introduction

The cost for fixing a design error in a system becomes larger and larger as the design proceeds from the requirement analysis to the implementation (see, e.g., [2, Chapter 1]) since the later in the design phase an error is detected the more reworking it may trigger. The above observation has motivated the development of methods and tools to verify correctness of a system already in the early phases of its design, namely during the requirement analysis or during the functional specification phases. The goal of all such approaches is to catch design errors well *before* the system implementation begins.

Of course, all such approaches are *model based*, that is they work on a model describing the system behaviour since no system implementation exists in the early design phases. Accordingly, *System Verification* is carried out by simulating a system model and analysing its behaviour under a *suitable set* of simulation scenarios.

For example, in a digital hardware setting, model based approaches have been used since a long time. In fact, even before

considering going to silicon, a heavy simulation activity is performed, aimed at verifying that the system model (defined, e.g., using Verilog, VHDL or SystemC¹) meets the system requirements for most (possibly all) *uncontrollable inputs* (that is, *primary inputs* and *faults* the system is expected to withstand).

Along the same line of reasoning, in a purely software setting, before generating low level code, model based approaches are used to verify that the software model (defined, e.g., using AADL [3,4]) meets the given requirements.

If all possible simulation scenarios are considered, then we can prove *correctness* of the system (i.e., absence of simulation scenarios violating the system requirements), otherwise we can only show that the system design is faulty (by exhibiting a simulation scenario violating the system requirements). In other words, a *simulation campaign* that does not consider all possible simulation scenarios can only show that the system design has a bug. To show correctness of the system design we need an *exhaustive* simulation campaign, that is one considering all possible simulation scenarios. A verification approach able to show system correctness is usually referred to as *formal verification*. One of the most successful techniques to carry out formal verification is *Model Checking* (see, e.g., [5]).

[☆] This paper is an extended and revised version of [1].

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¹ http://www.mentor.com/products/fv/hdl_designer/

The need for model checking stems from the high cost that a bug may have for certain systems. This is the case for *mission critical* systems, that is, systems for which a system malfunctioning may entail loss of money, as well as for *safety critical* systems, that is, systems for which a system malfunctioning may entail loss of human lives. Examples of mission critical systems are: decision support systems, satellites, processors (e.g., the infamous P5 FDIV bug costed about \$475 million to INTEL). Examples of safety critical systems are: railway interlocking, avionics control software.

Many Cyber-Physical Systems (CPSs) are indeed mission or safety critical systems. Accordingly, in this paper we focus on formal verification techniques for CPSs.

A CPS consists of hardware (e.g., motors, electrical circuits, etc.) and software components. Thus, in order to verify a CPS design, we need methods and tools that can model and effectively support simulation of hardware as well as software components.

From a formal point of view, CPSs can be modelled as hybrid systems (see, e.g., [6] and citations thereof). Many *Model Based Design* software tools offer support for modelling and simulation of CPSs. Well known examples are Simulink², VisSim³ and Modelica⁴. All such tools take as input a (mathematical) model of the behaviour of the CPS along with a simulation scenario and provide as output the time evolution (*trace* or *simulation run*) of the system at hand.

System Level Verification of CPSs aims at verifying that the *whole* (i.e., software + hardware) system meets the given specifications. *System Level Formal Verification (SLFV)* has the goal of *exhaustively* verifying that the above holds for *all* possible operational scenarios.

For digital circuits, formal verification is usually carried out using model checking techniques (e.g., see [7]). Unfortunately, model checkers for hybrid systems cannot handle SLFV of real world CPSs. Thus, HILS is currently the main workhorse for system level verification of CPSs, and is supported by model based design tools (e.g., the previously mentioned Simulink, VisSim and Modelica).

In HILS, the *actual software* reads/sends values from/to mathematical models (*simulation*) of the physical systems (e.g., engines, analog circuits, etc.) it will be interacting with. Notwithstanding the word *hardware*, in HILS the only hardware present is the one devoted to support the system simulation, that is: computational and communication devices. This is because HILS is used in a model based design setting to validate the system design *before* any hardware is built (the whole goal of model based design). For example, Simulink, VisSim, Modelica, ESA Satellite Simulation Infrastructure SIMULUS⁵ all provide simulation software supporting HILS, where the only hardware involved is just the computer on which the simulator is actually running.

Simulation can be very time consuming. Accordingly, in order to reduce system design time, Opal-RT⁶ and dSpace⁷ among others provide modelling and simulation software along with FPGA-based hardware to support real-time simulation. We note that in all cases the only hardware present in HILS is the one supporting the simulation itself.

1.1. Motivations

SLFV is an exhaustive HILS where *all* relevant simulation scenarios are considered. Using a parallel model checking driven

approach, exhaustive HILS enables formal verification of actual systems. Examples of such systems are the Inverted Pendulum on a Cart (IPC) and the Fuel Control System (FCS) in the Simulink distribution (see Section 6.1.1).

Considering that parallel exhaustive HILS based SLFV may take days of computation (e.g., see [8]), from a practical point of view it would be very useful to have available at *any time* during the verification process, *quantitative* information about the degree of assurance attained. Such an information would enable us to evaluate if it is worth to continue the verification activity, or instead stop it since the degree of assurance attained can be considered adequate for the application at hand (*graceful degradation*).

The above considerations suggest looking for a HILS based model checking approach satisfying the following requirements: (i) it is *parallel*, in order to make exhaustive HILS computationally feasible; (ii) it is *exhaustive*, since our focus is SLFV; (iii) it is *any time*, to support *graceful degradation*.

The work in [9] presents a Propositional Satisfiability (SAT) based model checker for finite state systems which returns, at *any time* during the verification process, the *coverage* (i.e., the fraction of operational scenarios verified so far). Unfortunately, while coverage measures the *amount* of verification work done, it does not provide any information about the *degree of assurance* attained by the verification process. This is because formal verification aims at finding *hard to find* errors, i.e., errors that were not detected while verifying operational scenarios designed by experts. As a result, formal verification addresses the search of errors that we are *unlikely* to consider. For this reason, we can model the problem as an adversary system, that is a system where, knowing our verification strategy, the adversary places the error in operational scenarios we are less likely to visit. In such a framework, *any* deterministic ordering of the operational scenarios would not increase the degree of assurance until the end of the verification. In fact, the adversary would choose to place the single error in the *last* scenario picked by the verification procedure.

To provide a formally sound information about the degree of assurance attained by the verification process, approaches have been proposed which verify the operational scenarios in a *random* order. In particular, the work in [10] presents a Monte-Carlo based model checker for finite state systems that provides, at *any time* during the verification process, an upper bound to the probability that the System Under Verification (SUV) exhibits an error in a yet-to-be-simulated scenario (Omission Probability). The Omission Probability (OP) provides indeed the information we are looking for. Unfortunately, while Monte-Carlo based approaches guarantee randomness (thereby enabling OP computation) they are not exhaustive (within a finite time).

To the best of our knowledge, no model checker is available, neither for finite state systems nor for hybrid systems, which, at the same time, is both *random* and *exhaustive*, thereby enabling effective *anytime* SLFV. In this paper we advance the state of the art by presenting a *parallel random exhaustive* HILS based model checker along with experimental results showing its effectiveness.

1.2. Main contribution

Our System Under Verification (SUV) is a *Hybrid System* (e.g., see [6] and citations thereof) whose inputs belong to a finite set of uncontrollable events (*disturbances*) modelling failures in sensors or actuators, variations in the system parameters, etc. We focus on *deterministic systems* (the typical case for control systems) and model nondeterministic behaviours (such as faults) with disturbances. Accordingly, in our framework, a *simulation scenario* is just a finite sequence of disturbances and a *simulation campaign* is a finite sequence of simulation instructions (namely: *save* a simulation state, *restore* a saved simulation state, *remove* a saved

² <http://www.mathworks.com>.

³ <http://www.vissim.com>.

⁴ <http://www.modelica.org>.

⁵ http://www.esa.int/Our_Activities/Operations/gse/ESA_operations_software_licensable_products_-_overview.

⁶ <http://www.opal-rt.com/about-hardware-loop-simulation>.

⁷ <https://www.dspace.com/en/inc/home.cfm>.

simulation state, *inject* a disturbance, *advance* the simulation of a given time length).

A system is expected to *withstand* all disturbance sequences that may arise in its operational environment. Correctness of a system is thus defined with respect to such *admissible* disturbance sequences. In our setting, the set of admissible disturbance sequences (*disturbance model*) can be defined as the language accepted by a suitable Finite State Automaton, which in turn can be defined using the modelling language of any finite state model checker.

In such a framework we address *Bounded SLFV of safety* properties. That is, given a time step τ (time quantum between disturbances) and a time horizon $T = \tau h$ we return *PASS* if there is no *admissible* disturbance sequence of length h and time step τ that violates the given safety property. We return *FAIL*, along with a counterexample, otherwise. Therefore, SLFV is an *exhaustive* (with respect to admissible disturbance sequences) HILS. In other words, we are aiming at (*black box*) *bounded model checking* where the SUV behaviour is defined by a simulator (Simulink in our examples).

In such a setting, our main contribution can be summarised as follows. We present an *anytime parallel random exhaustive* HILS based model checker that effectively conjugates exhaustiveness with randomness, thereby enabling the computation of the Omission Probability.

While a simulation run for digital hardware takes order of milliseconds on a normal desktop computer, in our setting a simulation run takes order of seconds since it entails heavy numerical computations aimed at solving a system of many *Ordinary Differential Equations* (modelling the hardware components of the CPS). Indeed (see [Section 6](#)) simulation of operational scenarios takes almost 100% of the overall verification time. Resting on such observation we build on the SLFV approach discussed in [\[8\]](#). In particular:

1. From the disturbance model we generate all admissible simulation scenarios and evenly split them into disjoint sets (*slices*).
2. For each slice, we compute a highly optimised *simulation campaign* that exploits simulator save/restore/remove commands in order to save on the simulation time while scheduling execution of *all* simulation scenarios *exactly once* and in a *uniform random order*. This guarantees exhaustiveness and allows us to compute, at *any time* during the verification process, an upper bound to the OP.
3. We run simulation campaigns in parallel. This guarantees a very efficient parallelism, since no communication among processes is needed. This step is supported by simulation tools (Simulink in our case study).

We also show that, thanks to the fact that we have first generated all admissible simulation scenarios, attaining point 2 above (i.e., anytime OP computation) can be done in a *not-too-complicated* (from both technical and computational points of view) and *non-invasive way*, by simply introducing a new module into the workflow of [\[8\]](#) that seamlessly interoperates with the others.

1.3. Summary of experimental results

We implemented our approach and present ([Section 6](#)) experimental results on two case studies, namely the Inverted Pendulum on a Cart (IPC) and the Fuel Control System (FCS) examples in the Simulink distribution. Overall, we compute optimised simulation campaigns under four operational environments (disturbance models), which entail from 3 208 276 to 35 641 501 simulation scenarios.

Each processor (actually, a core of a 8-core machine) runs an instance of our (random) simulation campaign computation algorithm and takes as input a slice of our set of simulation scenarios. We present experimental results with 16, 32, 64 machines totalling 128, 256, 512 parallel processes. Our approach takes negligible time to generate an optimised simulation campaign for a given slice, with respect to the time needed to actually execute the simulation campaign (e.g., minutes vs. hours, see [Section 6](#)).

Our experimental results show that, by exploiting parallelism, our random exhaustive simulation campaigns effectively counteract the simulation time overhead due to randomisation. The above ensures feasibility of our parallel random exhaustive approach for actual systems, such as the Inverted Pendulum on a Cart (IPC) and Fuel Control System (FCS) examples in the Simulink distribution.

As for the Omission Probability (OP), the worst case scenario is when just one error trace is present. Our experimental results ([Section 6.6](#)) show that, even in such a case, our upper bound to the OP decreases about *linearly* with respect to the coverage, that is the fraction of scenarios simulated. This is the *best* one can hope to obtain in our setting.

Finally, simulation of scenarios in random order allows us to use the coverage as a reliable estimator for the *completion time* of the whole verification task. Our experimental results show that the error in the completion time estimation decreases quickly with respect to coverage.

1.4. Paper outline

[Section 2](#) gives background notions to make our paper self-contained. [Section 3](#) presents our formal framework, by formalising the notion of OP and by providing an upper bound for it, computable from the number of the yet-to-be-simulated traces in each slice. [Section 4](#) outlines our algorithm which enables the computation, from a sequence (slice) of disturbance traces, a highly optimised simulation campaign which simulates the input traces in uniform random order and exploits the save/restore/remove capabilities of the simulator. [Section 5](#) is devoted to the formal proofs of our results. Finally, [Section 6](#) presents experimental results assessing effectiveness of our approach.

2. Background and preliminaries

In this section we give some background notions. Unless otherwise stated, all definitions are based on [\[8,11,12\]](#). Throughout the paper, we use $\mathbb{R}^{\geq 0}$ for the set of non-negative reals, \mathbb{R}^+ for the set of strictly positive reals, and $\text{Bool} = \{0, 1\}$ for the set of Boolean values (0 for *false* and 1 for *true*). \mathbb{N}^+ denotes the set of positive natural numbers.

2.1. Modelling the operational environment

Our System Under Verification (SUV) is a Discrete Event System (DES), namely a continuous time Input-State-Output deterministic dynamical system [\[11\]](#) whose inputs are *discrete event sequences*. A discrete event sequence ([Definition 1](#)) is a function $u(t)$ associating to each (continuous) time instant $t \in \mathbb{R}^{\geq 0}$ a *disturbance event* (or, simply, *disturbance*). Disturbances, encoded by natural numbers in the interval $[0, d]$ (for a given $d \in \mathbb{N}^+$), represent uncontrollable events (e.g., faults). We use event 0 to represent the event carrying no disturbance. As no system can withstand an infinite number of disturbances within a finite time, we require that, in any time interval of finite length, a discrete event sequence $u(t)$ differs from 0 only in a finite number of time points ([Fig. 1a](#)).

Definition 1 (Discrete event sequence). Let $d \in \mathbb{N}^+$. A *discrete event sequence* over the natural interval $[0, d]$ is a

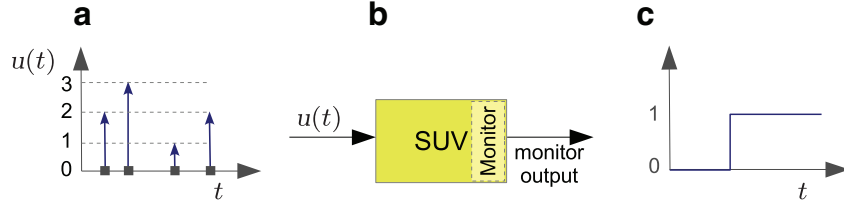


Fig. 1. (a) A discrete event sequence ($d = 3$); (b) Our SUV embedding a monitor; (c) The SUV monitor output.

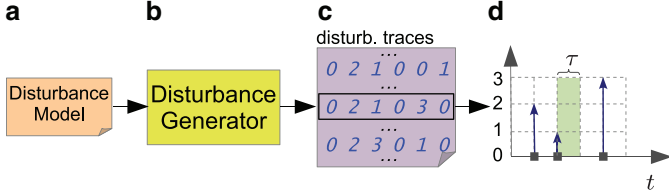


Fig. 2. (a) Disturbance model; (b) CMurphi-based disturbance generator; (c) Generated sequence of disturbance traces ($d = 3, h = 6$); (d) The discrete event sequence associated to the trace in the black rectangle in part (c), given time quantum τ .

function $u : \mathbb{R}^{\geq 0} \rightarrow [0, d]$ such that, for all $t \in \mathbb{R}^{\geq 0}$, the set $\{\tilde{t} \in \mathbb{R}^{\geq 0} \mid 0 \leq \tilde{t} \leq t \text{ and } u(\tilde{t}) \neq 0\}$ has finite cardinality. We denote with \mathcal{U}_d the set of discrete event sequences over $[0, d]$ (following control engineering notation for input functions to dynamical systems, see e.g. [11]).

In Definitions 2 and 3 we specify the concepts of *restriction* and *concatenation*, respectively, for functions belonging to \mathcal{U}_d .

Definition 2. Let \mathcal{U}_d be the set of discrete event sequences over the interval $[0, d]$. Given a function $u \in \mathcal{U}_d$ and two real numbers $0 \leq t_1 < t_2$, we denote with $u|_{[t_1, t_2]}$ the function $u|_{[t_1, t_2]} : [t_1, t_2] \rightarrow [0, d]$, such that $u|_{[t_1, t_2]}(t) = u(t)$ for all $t \in [t_1, t_2]$. We denote $\mathcal{U}_d^{[t_1, t_2]}$ the restriction of \mathcal{U}_d to the domain $[t_1, t_2]$.

Definition 3. Assume that $t_1, t_2, t_3 \in \mathbb{R}^{\geq 0}$ such that $t_1 < t_2 < t_3$. If $\omega \in \mathcal{U}_d^{[t_1, t_2]}$ and $\omega' \in \mathcal{U}_d^{[t_2, t_3]}$, their *concatenation*, denoted as $\omega\omega'$, is the function $\tilde{\omega} \in \mathcal{U}_d^{[t_1, t_3]}$ defined as:

$$\tilde{\omega}(t) = \begin{cases} \omega(t) & \text{if } t \in [t_1, t_2] \\ \omega'(t) & \text{if } t \in [t_2, t_3] \end{cases}$$

System level verification follows an *Assume-Guarantee* approach aimed at showing that the SUV meets its specification (*Guarantee*) as long as the SUV operational environment behaves as expected (*Assume*). In this work we focus on *bounded* system level verification. As a consequence, we model (Definition 4) the SUV oper-

ational environment as the sequence of disturbances our SUV is expected to withstand within a *finite* time horizon, and we bound the *time quantum* between two consecutive disturbances.

Definition 4 (Disturbance trace). Let $h, d \in \mathbb{N}^+$. An (h, d) disturbance trace δ is a finite sequence $\delta : [0, h-1] \rightarrow [0, d]$. Given $\tau \in \mathbb{R}^+$ (*time quantum*), an (h, d) disturbance trace δ is univocally associated to a discrete event sequence u_δ^τ , defined as follows: for all $t \in \mathbb{R}^{\geq 0}$, if there exists $j \in [0, h-1]$ such that $t = \tau j$, then $u_\delta^\tau(t) = \delta(j)$, else $u_\delta^\tau(t) = 0$ (no disturbance).

Thus, a disturbance trace δ defines an operational scenario (namely, u_δ^τ) for our SUV. Fig. 2d shows the discrete event sequence associated to a disturbance trace. We represent our SUV *operational environment* as a finite set of (h, d) disturbance traces $\Delta = \{\delta_0, \dots, \delta_{n-1}\}$, since $U_\Delta^\tau = \{u_{\delta_0}^\tau, \dots, u_{\delta_{n-1}}^\tau\}$ (for a given $\tau \in \mathbb{R}^+$) defines the operational scenarios our SUV should withstand. Note that, by taking h large enough (as in Bounded Model Checking BMC) and τ small enough (to faithfully model our SUV operational scenarios), we can achieve any desired precision. On such considerations rests the effectiveness of the approach.

As it is typically infeasible to define a SUV operational environment by explicitly listing all its disturbance traces, we define an operational environment with a *disturbance model* which is in turn defined as the language accepted by a suitable finite state automaton. The following example illustrates this point.

Example 1. Consider a disturbance model consisting of one disturbance (namely, a fault) which is always recovered within 4 s (i.e., 4 seconds). Let us assume that between two consecutive disturbances (faults) there must be at least 5 s and that disturbances can arise only at time steps multiple of $\tau = 1$ s (*time quantum*). We also assume that the verification time horizon is set to 6 s.

In Fig. 3a we show disturbance traces represented as strings of zeros (no disturbance) and ones (disturbance), with time flowing from left to right. The 8 strings terminated by \checkmark denote all the disturbance traces accepted by the disturbance model (*admissible*

000000	✓	010000	✓
000001	✓	010001	⊗
000010	✓	01001	⊗
000011	⊗	0101	⊗
000100	✓	011	⊗
000101	⊗	100000	✓
00011	⊗	100001	✓
001000	✓	10001	⊗
001001	⊗	1001	⊗
00101	⊗	101	⊗
0011	⊗	11	⊗

(a) Admissible disturbance traces (✓) and shortest disturbance sequences that cannot be extended to an admissible disturbance trace (⊗)

```
function disturbanceModel(h)
  cnt ← 0; /* counter */
  t ← 0; /* time */
  while t ≤ h do
    d ← read(); t ← t + 1;
    if cnt > 0 then cnt ← cnt - 1;
    if d = 1 then
      if cnt > 0 then return ⊗;
      else cnt ← 4;
    return ✓;
end
```

(b) Finite state automaton recognising the language of admissible disturbance traces (disturbance model)

Fig. 3. Example 1.

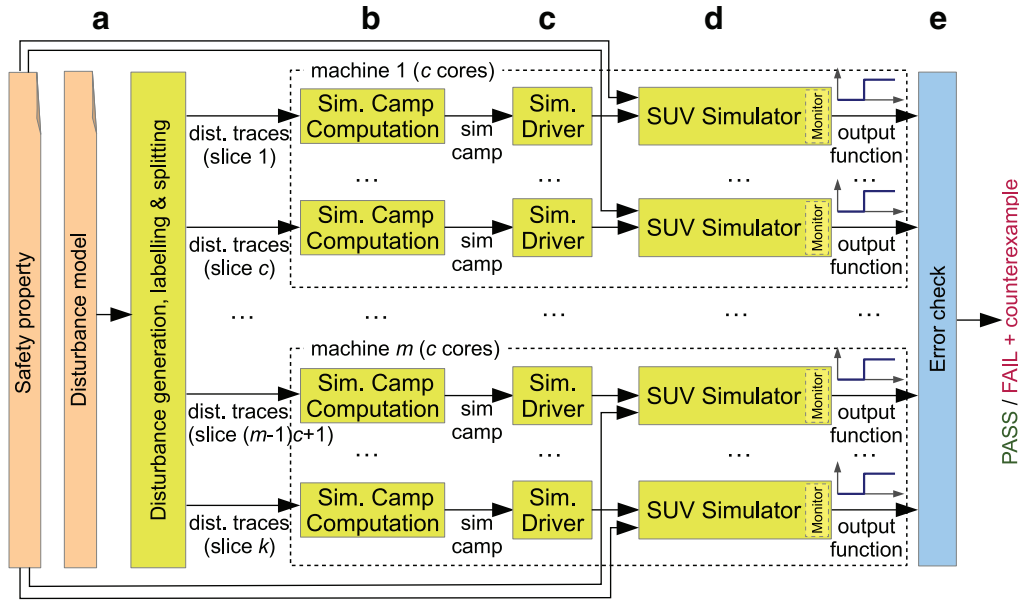


Fig. 4. Parallel HILS based dSLFV [8]: k parallel processes are run on m multi-core machines (we show a possible deployment with machines having c cores each, i.e., $k = mc$).

disturbance traces). The 14 strings terminated by \otimes are the shortest non-admissible sequences of disturbances, that is disturbance sequences that *cannot* be extended to admissible disturbance traces.

Fig. 3b shows the pseudo-code for a finite state automaton recognising such a language.

We define a finite state automaton for a disturbance model using the modelling language of a finite state model checker (namely, CMurphi [13]), along the lines of [8].

2.2. Modelling the property to be verified

Along the lines of [14], we model the property to be verified with a continuous-time *monitor* which observes the state of the system to be verified and checks whether the property under verification is satisfied (Fig. 1b). The output of the monitor is 0 as long as the property under verification is satisfied and becomes and stays 1 (*sustain*) as soon as the property fails, thus ensuring that we never miss a property failure report, even when sampling the monitor output only at discrete time points (Fig. 1c). The use of monitors gives us a flexible approach to model the property to be verified. In particular, it is easy to model bounded safety and bounded liveness properties as monitors. Figs. 8 and 9 show the Simulink/Stateflow representations of our two case studies (Inverted Pendulum on a Cart and Fuel Control System, respectively), along with their property monitors (see Section 6).

2.3. Modelling the SUV

Since the monitor output is all we need to carry out our verification task, we can model our SUV *along with* the property to be verified as a DES with an *embedded monitor* (Fig. 1b). We call *Monitored Discrete Event System (MDES)* such a DES.

According to our *black-box* approach to SUV modelling, given a time quantum $\tau \in \mathbb{R}^+$, Definition 5 formalises an (h, d) MDES as a function \mathcal{H} associating, to each (h, d) disturbance trace δ , a Boolean value $\mathcal{H}(\delta)$ representing the output of the SUV monitor at time $T = \tau h$ (the time horizon), when the system (starting from its initial state) is given as input the discrete event sequence $u_\delta^\tau(t)$ associated to δ . For any disturbance trace δ , $\mathcal{H}(\delta)$ is 1 (*error*) if and

only if $u_\delta^\tau(t)$ violates the property under verification within time horizon $T = \tau h$ (with the SUV starting from its initial state).

Definition 5. ((h, d) **Monitored DES**) Let $h, d \in \mathbb{N}^+$. A (h, d) *Monitored Discrete Event System (MDES)* is a function $\mathcal{H} : ([0, h - 1] \rightarrow [0, d]) \rightarrow \text{Bool}$ mapping all (h, d) disturbance traces to Boolean values.

2.4. System Level Formal Verification (SLFV)

Definition 6 formalises our bounded System Level Formal Verification problem.

Definition 6. A *System Level Formal Verification (SLFV) problem* is a tuple $P = (h, d, \Delta, \mathcal{H})$ where: $h, d \in \mathbb{N}^+$, $\Delta = \{\delta_0, \dots, \delta_{n-1}\}$ is an (h, d) set of disturbance traces, and \mathcal{H} is a (h, d) MDES.

The *answer* to SLFV problem P is *FAIL* if there exists a disturbance trace δ in Δ such that $\mathcal{H}(\delta) = 1$ (in such a case also the *counterexample* δ is returned), *PASS* otherwise.

Note that, notwithstanding the fact that the number of states of our SUV is infinite and we are in a continuous time setting, to answer a SLFV problem we only need to check a *finite* number of disturbance traces. This is because we are bounding: (a) our time horizon to $T = \tau h$, and (b) the set of time points at which disturbances can take place, by taking τ as the time quantum among disturbance events.

2.5. Parallel HILS based deterministic SLFV

In the black-box parallel approach shown in [8], the MDES \mathcal{H} defining our SUV (plus the property to be verified) is defined using the modelling language of a suitable *simulator* (e.g., MatLab and Stateflow for Simulink). The answer to a SLFV problem $(h, d, \Delta, \mathcal{H})$ is computed by simulating *each* operational scenario δ in the operational environment Δ , thus by performing an exhaustive (with respect to Δ) Hardware In the Loop Simulation (HILS). The overall workflow is shown in Fig. 4 and described in the remainder of this section. We will refer to this approach as *Deterministic SLFV (dSLFV)*, where the word “deterministic” stems from the fact that the disturbance traces are verified in a deterministic order.

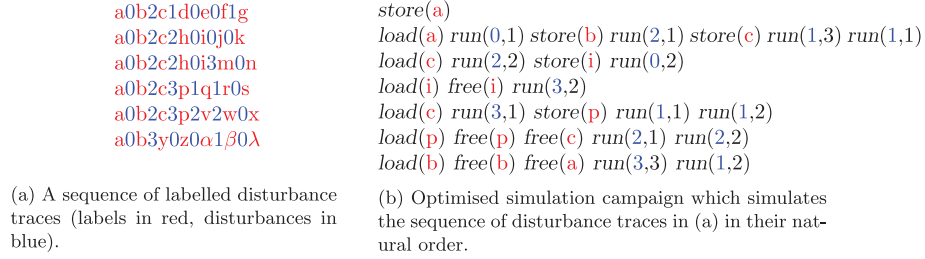


Fig. 5. Labelled disturbance traces and optimised simulation campaign.

2.5.1. Disturbance trace generation and splitting

Our CMurphi-based trace generator (see Section 2.1 and Fig. 4a) works in Depth-First Search (DFS) mode, and, given the set of disturbances, produces a sequence Δ of n disturbance traces. Each generated trace δ in Δ is annotated with labels and is of the form $\delta = (l_0, d_0, l_1, d_1, \dots, l_{h-1}, d_{h-1}, l_h)$, where $\delta = (d_0, \dots, d_{h-1})$ is a sequence of disturbances satisfying the disturbance model and l_0, \dots, l_h belong to a countable set of labels L .

Labels are defined by an injective map λ from finite sequences of disturbances (including the empty sequence) to L . Label l_0 is common to all traces and it is associated to the simulator initial state. Prefixes of disturbance sequences $(\hat{d}_0, \dots, \hat{d}_{p-1})$ common to multiple disturbance traces in Δ are followed by the same label $\hat{l}_p = \lambda(\hat{d}_0, \dots, \hat{d}_{p-1})$. Fig. 5a shows a short sequence of labelled disturbance traces.

Labels identifying common disturbance prefixes are essential in the efficient computation of optimised simulation campaigns. Note that, given that our CMurphi-based generator runs in DFS mode, disturbance traces can be labelled at *no additional computational cost* during generation. Trace labelling during generation greatly increases the efficiency of the simulation campaign optimiser, as shown in [8].

In the following, we will use δ^λ instead of δ (respectively, Δ^λ instead of Δ) when we want to emphasise that a trace δ is annotated (respectively, all traces in set Δ are annotated) with labels, or when we need such labels. In order to enable parallel verification via $k \in \mathbb{N}^+$ processes, we evenly partition the sequence of labelled disturbance traces Δ^λ into $k \in \mathbb{N}^+$ sequences of disturbance traces $\Delta_0^\lambda, \dots, \Delta_{k-1}^\lambda$ (called *slices*).

The splitting process produces slices containing $\lceil n/k \rceil$ traces each, except the last slice, which may contain fewer traces if n is not a multiple of k .

2.5.2. Computation of optimised simulation campaigns

In the next phase of the workflow in Fig. 4, we use the k slices $\Delta_0^\lambda, \dots, \Delta_{k-1}^\lambda$ generated so far to compute, independently and in parallel, k highly optimised simulation campaigns (Fig. 4b). Such simulation campaigns can be simulated, again independently and in parallel, using k simulators, each one running (e.g., on a different core of a bunch of multi-core machines) a model for \mathcal{H} (Fig. 4c–d).

The answer to the SLFV problem is *FAIL* if one of the simulation campaigns raises the simulator output function to 1 (in this case the disturbance trace δ which raised the error is returned as a counterexample, see Fig. 4e). The answer is *PASS* otherwise.

Each simulator accepts four basic commands: *store*, *load*, *free*, *run*. Command *store(l)* stores in memory the current state of the simulator and labels with l such a state. Command *load(l)* loads into the simulator the stored state labelled with l . Command *free(l)* removes from the memory the state labelled with l . Command *run(e, t)* (with $e \in [0, d]$ and $t \in \mathbb{R}^+$) injects disturbance e and then advances the simulation of time t . A *simulation campaign* is thus a sequence of simulator commands.

The simulation campaign χ_i ($0 \leq i < k$) computed from input slice Δ_i^λ steers the simulator as to execute all disturbance traces in Δ_i^λ according to their order in the slice (this is the reason why we refer to this approach as Deterministic SLFV).

Each disturbance trace is executed as if starting from the simulator initial state. However, by using commands *store* and *load*, the optimiser avoids revisiting simulation states as much as possible (as in explicit model checking). Using command *free* the optimiser removes from the simulator memory states that will never be needed in the remaining part of the simulation campaign. This is needed since each state may require many KB of memory (150–300 KB in the case studies presented in this paper).

Fig. 5b shows the optimised simulation campaign computed from the sequence of labelled disturbance traces in Fig. 5a, in the simple case where we ignore any limit on number of states that can be kept simultaneously stored in the simulator memory.

3. Omission Probability

This section formally defines the notion of Omission Probability (OP) (Definitions 7 and 8) and provides an upper bound for it, which can be computed anytime during the parallel verification process from the number of the yet-to-be-simulated traces in each slice (Theorem 1).

Notation 1 (Set of permutations of a set). Let $\Delta = \{\delta_0, \dots, \delta_{n-1}\}$ be a finite non-empty set. We denote with $\text{Perm}(\Delta)$ the set of permutations of elements of Δ :

$$\text{Perm}(\Delta) = \{(\theta_0, \dots, \theta_{n-1}) \mid (\forall i \in [0, n-1] \theta_i \in \Delta) \wedge (\forall i, j \in [0, n-1] i \neq j \rightarrow \theta_i \neq \theta_j)\}$$

If $\hat{\Delta} = (\delta_0, \dots, \delta_{n-1}) \in \text{Perm}(\Delta)$ we write also $\hat{\Delta}(i)$ for δ_i .

A Random Sequence Generator (RSG) models the extraction of a random permutation from a given finite non-empty set (which, in our case, will be the set of admissible disturbance traces Δ). This is formalised in Definition 7.

Definition 7 (Random Sequence Generator). Let Δ be a finite non-empty set. A Random Sequence Generator (RSG) for Δ is a probability space $(\Omega, \mathcal{F}, \text{Pr})$, where:

- Ω (the space of *outcomes*) is the set of permutations of Δ , that is $\Omega = \text{Perm}(\Delta)$.
- \mathcal{F} (the space of *events*) is the set of subsets of Ω , that is: $\mathcal{F} = 2^\Omega = \{E \mid E \subseteq \Omega\}$.
- $\text{Pr} : \mathcal{F} \rightarrow [0, 1]$ is a probability measure such that, for all $\omega \in \Omega$, $\text{Pr}(\omega) = \frac{1}{|\Delta|!}$. That is, permutations of Δ are extracted with uniform probability. Since Ω is countable (actually finite), the probability of any event $E \in \mathcal{F}$ is defined as $\text{Pr}(E) = \sum_{\omega \in E} \text{Pr}(\omega)$.

Let $(\Delta_0, \dots, \Delta_{k-1})$ be a partition of Δ into $k \in \mathbb{N}^+$ disjoint non-empty sets. For any $0 \leq i < k$, let $(\Omega_i, \mathcal{F}_i, \text{Pr}_i)$ be a RSG for Δ_i . A Random Sequence Generator for $(\Delta_0, \dots, \Delta_{k-1})$ is a

probability space $(\Omega, \mathcal{F}, \text{Pr})$, where: $\Omega = \times_{i=0}^{k-1} \Omega_i$, $\mathcal{F} = \times_{i=0}^{k-1} \mathcal{F}_i$ and, for each event $E_0 \times \dots \times E_{k-1} \in \mathcal{F}$ ($E_i \in \mathcal{F}_i$ for each $0 \leq i < k$), $P(E_0 \times \dots \times E_{k-1}) = \prod_{i=0}^{k-1} \text{Pr}_i(E_i)$.

Note that, by [Definition 7](#), a RSG for a partition $(\Delta_0, \dots, \Delta_{k-1})$ of Δ models the extraction of k permutations of, respectively, $\Delta_0, \dots, \Delta_{k-1}$. For all $0 \leq i < k$, the extracted permutation of Δ_i is chosen *uniformly* among all possible permutations of Δ_i . Also, the k permutations are extracted *independently* from each other.

[Definition 8](#) defines the probability of omitting the simulation of a trace $\delta \in \Delta$ containing an error (i.e., $\mathcal{H}(\delta) = 1$) when the verification process has *already* examined q_i disturbance traces, where $q_i \in \{0, \dots, |\Delta_i|\}$, from a random permutation of slice Δ_i , for all $0 \leq i < k$. Thus q_i represents the state of advancement of the computation on the i th slice Δ_i . Hence (q_0, \dots, q_{k-1}) defines the state of advancement (stage) of the computation on all slices. We denote this probability as Omission Probability (OP).

Definition 8 (Omission Probability). Let $(h, d, \Delta, \mathcal{H})$ be a System Level Formal Verification (SLFV) problem and $(\Delta_0, \dots, \Delta_{k-1})$ be a partition of Δ into $k \in \mathbb{N}^+$ disjoint non-empty sets. Let $(\Omega, \mathcal{F}, \text{Pr})$ be a RSG for $(\Delta_0, \dots, \Delta_{k-1})$, and (q_0, \dots, q_{k-1}) a tuple such that $q_i \in \{0, \dots, |\Delta_i|\}$ for each $0 \leq i < k$.

The *Omission Probability* (OP) for $(\Delta_0, \dots, \Delta_{k-1})$ at stage (q_0, \dots, q_{k-1}) , is defined as:

$$OP_{\mathcal{H}}(|\Delta_0|, \dots, |\Delta_{k-1}|, q_0, \dots, q_{k-1}) \\ = \text{Pr} \left(\left\{ (\omega_0, \dots, \omega_{k-1}) \mid \begin{array}{l} \forall i \in [0, k-1] \ \omega_i \in \Omega_i \wedge \\ AB((\omega_0, \dots, \omega_{k-1}), (q_0, \dots, q_{k-1})) \end{array} \right\} \right)$$

where:

- AB is defined as $A((\omega_0, \dots, \omega_{k-1}), (q_0, \dots, q_{k-1})) \wedge B((\omega_0, \dots, \omega_{k-1}), (q_0, \dots, q_{k-1}))$
- A (After) is: $A((\omega_0, \dots, \omega_{k-1}), (q_0, \dots, q_{k-1})) = \exists i \in [0, k-1] \exists j \in [q_i, |\Delta_i|] \ \mathcal{H}(\omega_i(j)) = 1$;
- B (Before) is: $B((\omega_0, \dots, \omega_{k-1}), (q_0, \dots, q_{k-1})) = \forall i \in [0, k-1] \forall j \in [0, q_i-1] \ \mathcal{H}(\omega_i(j)) = 0$.

In [Definition 8](#), formula A (After) states that there exists a yet-to-be-simulated trace δ (some trace $j \geq q_i$ of some slice i) containing an error, i.e., such that $\mathcal{H}(\delta)$ evaluates to 1. Formula B (Before) states that none of the already simulated traces contains an error, i.e., function \mathcal{H} evaluates to 0 for all of them.

The following [Theorem 1](#) gives an upper bound to the OP, after having simulated q_i randomly extracted traces from slice Δ_i (for each $0 \leq i < k$). In particular, [Theorem 1](#) provides a bound to the probability of omitting the simulation of a trace $\delta \in \Delta$ containing an error when the verification process has already examined q_i disturbance traces from the (the permutation of) slice Δ_i (for all $0 \leq i < k$). Importantly, the bound provided does *not* depend on \mathcal{H} , i.e., it is *independent* of the system model. The proof of the theorem is in [Section 5](#).

Theorem 1. Let $(h, d, \Delta, \mathcal{H})$ be a SLFV problem and $(\Delta_0, \dots, \Delta_{k-1})$ be a partition of Δ into $k \in \mathbb{N}^+$ disjoint non-empty sets. Let $(\Omega, \mathcal{F}, \text{Pr})$ be a Random Sequence Generator for $(\Delta_0, \dots, \Delta_{k-1})$ and (q_0, \dots, q_{k-1}) a tuple such that $q_i \in \{0, \dots, |\Delta_i|\}$ for each $0 \leq i < k$. We have:

$$OP_{\mathcal{H}}(|\Delta_0|, \dots, |\Delta_{k-1}|, q_0, \dots, q_{k-1}) \leq 1 - \min \left\{ \frac{q_i}{|\Delta_i|} \mid 0 \leq i < k \right\} \quad (1)$$

Note that the construction of the slices $\Delta_0, \dots, \Delta_{k-1}$ from Δ is *non-deterministic* (i.e., any partitioning of Δ would work), whereas, for each slice, the selection of a permutation is a *probabilistic* process, modelled as a RSG. Accordingly, [Theorem 1](#) bounds the OP using the *worst case* distribution, i.e., the distribution yielding the

greatest OP. From this stems the min function in the right member of the inequality in [Theorem 1](#).

Finally, we observe that, from [Theorem 1](#), it follows that $OP_{\mathcal{H}}(|\Delta_0|, \dots, |\Delta_{k-1}|, |\Delta_0|, \dots, |\Delta_{k-1}|) = 0$, that is, our verification task terminates after $\max\{|\Delta_i| \mid 0 \leq i < k\}$ parallel steps, having simulated all traces in Δ .

4. Enabling Omission Probability computation: random exhaustive SLFV

Our disturbance trace generator (see [Section 2.1](#)) produces a sequence of disturbance traces Δ^λ , whose order is deterministically chosen by the employed algorithm (in our case, Depth-First Search (DFS)). As a consequence, no information about the Omission Probability (OP) can be inferred *during* the verification process if the simulation campaign computed from each slice verifies the sequence of disturbance traces therein according to their order (as done by the Deterministic SLFV (dSLFV) approach of [Section 2.5](#)), or according to *any deterministic* order, as argued in [Section 1.1](#).

From [Section 3](#) it follows that it is possible to enable OP computation (and hence, *graceful degradation* during the computationally very expensive simulation phase) by simulating the disturbance traces within each slice in an order uniformly chosen at random.

Here we show how we can add support to OP computation in the workflow of [Fig. 4](#) in a *not-too-complicated* and *non-invasive* way, by introducing an additional step, in the parallel portion of the workflow, which implements a Random Sequence Generator (RSG).

The new workflow, which we refer to as random exhaustive SLFV (rSLFV), is shown in [Fig. 6](#). All slices are given as input, in parallel, to instances the new Random Sequence Generator module, shown as [Algorithm 1](#). The Random Sequence Generator module reads a sequence of disturbance traces and computes a *random permutation* of it, uniformly chosen among all possible permutations, thus implementing a RSG ([Definition 7](#)). We argue that the introduction of the new Random Sequence Generator module is non-too-complicated for what concerns both its implementation ([Algorithm 1](#)) and its computational viability (see [Section 6.3](#)). It is also non-invasive, as it seamlessly interoperates with the dSLFV workflow of [Section 2.5](#).

As the input sequence can be too large to be kept in main memory, the Random Sequence Generator module implements a disk-based multi-round algorithm (shown as [Algorithm 1](#)) which takes efficiency into account by using, in each round, as much main memory as possible and by reading/writing the input/output trace files sequentially.

Let $n = |\Delta^\lambda|$ be the number of disturbance traces in the input sequence. Given parameter $z \in \mathbb{N}^+$ for the maximum number of disturbance traces which can be simultaneously stored in main memory, the algorithm, at each round $r \geq 1$, selects the z traces which will have output positions in the interval $[(r-1)z, \min(n, rz-1)]$.

The above selection is performed by computing the first z elements of a random permutation of the traces not yet in the output file $\Delta_{\text{rnd}}^\lambda$, chosen uniformly among all possible permutations. Such a permutation prefix is computed by function `rndPermPrefix()` ([Algorithm 1](#), from line 20). Function `rndPermPrefix()` performs a swap-based computation of a permutation of integers $[0, |\Delta_{\text{in}}^\lambda| - 1]$ (uniformly chosen at random) and interrupts the computation as soon as the first z elements (see variable `result`) have been determined. During its operation, function `rndPermPrefix()` represents the partially computed permutation as an associative array `perm`, which, at any step of the function execution, represents

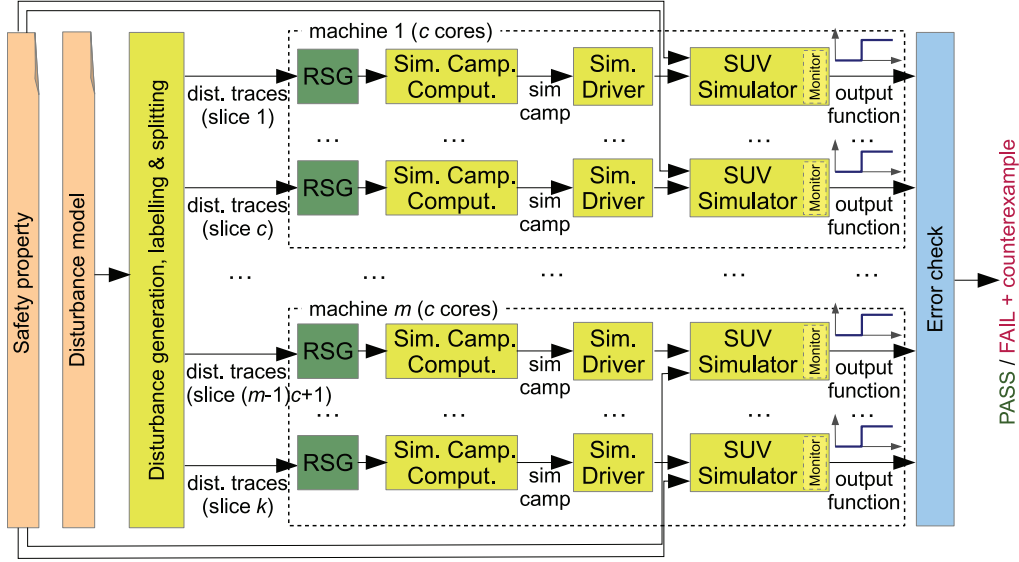


Fig. 6. Parallel HILS based rSLFV.

permutation π of $[0, n - 1]$ such that, for each $j \in [0, n - 1]$:

$$\pi(j) = \begin{cases} v & \text{if } (j, v) \in \text{perm} \\ j & \text{otherwise, i.e., } \exists v (j, v) \in \text{perm} \end{cases}$$

The main algorithm then appends the z randomly selected traces (as chosen by function $\text{rndPermPrefix}()$) to $\Delta_{\text{rnd}}^\lambda$ (according to their output positions), and dumps all the others to a temporary file, which becomes the input of the next round. Algorithm 1 terminates in $\lceil n/z \rceil$ rounds.

The following theorem asserts that Algorithm 1, when applied to a sequence of disturbance traces Δ^λ , produces a uniformly chosen permutation over the set of all permutations. The proof of the theorem is in Section 5.

Theorem 2. Let Δ^λ be a file containing n disturbance traces. For any $z \geq 1$, Algorithm 1 stores in file $\Delta_{\text{rnd}}^\lambda$ a permutation of the traces in Δ^λ extracted with uniform probability among all possible permutations.

The k output randomised slices (computed in parallel by k instances of the Random Sequence Generator algorithm from k input slices) are then given as input to k parallel instances of the simulation campaign computation module described in Section 2.5.2 which compute k highly optimised simulation campaigns (Fig. 6). As each simulation campaign verifies the traces in its input slice according to their order, the introduction of the Random Sequence Generator module effectively enforces a random order in the disturbance trace verification within each slice, satisfying the requirements stated in Theorem 1 in order to compute, during simulation, an upper bound to the OP.

5. Proof of results

Theorem 1. Let $(h, d, \Delta, \mathcal{H})$ be a System Level Formal Verification (SLFV) problem and $(\Delta_0, \dots, \Delta_{k-1})$ be a partition of Δ into $k \in \mathbb{N}^+$ disjoint non-empty sets. Let $(\Omega, \mathcal{F}, \text{Pr})$ be a Random Sequence Generator for $(\Delta_0, \dots, \Delta_{k-1})$ and (q_0, \dots, q_{k-1}) a tuple such that $q_i \in \{0, \dots, |\Delta_i|\}$ for each $0 \leq i < k$. We have:

$$OP_{\mathcal{H}}(|\Delta_0|, \dots, |\Delta_{k-1}|, q_0, \dots, q_{k-1}) \leq 1 - \min \left\{ \frac{q_i}{|\Delta_i|} \mid 0 \leq i < k \right\} \quad (1)$$

Proof. If, for all $\delta \in \Delta$, $\mathcal{H}(\delta) = 0$, then the left member of (1) is $\text{Pr}(\emptyset) = 0$ and the thesis follows.

Otherwise, let E be a nonempty set containing error traces, that is $E = \{\delta \mid \delta \in \Delta \wedge \mathcal{H}(\delta) = 1\} \neq \emptyset$. $OP_{\mathcal{H}}(|\Delta_0|, \dots, |\Delta_{k-1}|, q_0, \dots, q_{k-1})$ can be rewritten as (see Definition 8):

$$\Pr(\{(\omega_0, \dots, \omega_{k-1}) \mid \forall i \in [0, k-1] \omega_i \in \Omega_i \wedge \forall i \in [0, k-1] \forall j \in [0, q_i-1] \omega_i(j) \notin E\}) \quad (2)$$

because $(\Delta_0, \dots, \Delta_{k-1})$ is a partition of Δ .

Consider any $\bar{\delta} \in E$. Probability (2) is less than or equal to

$$\Pr(\{(\omega_0, \dots, \omega_{k-1}) \mid \forall i \in [0, k-1] \omega_i \in \Omega_i \wedge \forall i \in [0, k-1] \forall j \in [0, q_i-1] \omega_i(j) \neq \bar{\delta}\}) \quad (3)$$

Given that $\bar{\delta}$ belongs to exactly one of $\Delta_0, \dots, \Delta_{k-1}$, say $\Delta_{\bar{i}}$, and given the definition of $\text{Pr}(\omega_0, \dots, \omega_{k-1})$ in Definition 7, expression (3) is equal to:

$$\left(\prod_{\substack{i=0 \\ i \neq \bar{i}}}^{k-1} \Pr(\{\omega_i \in \Omega_i\}) \right) \times \Pr(\{\omega_{\bar{i}} \in \Omega_{\bar{i}} \mid \forall j \in [0, q_{\bar{i}}-1] \omega_{\bar{i}}(j) \neq \bar{\delta}\})$$

which is equal to

$$\Pr(\{\omega_{\bar{i}} \in \Omega_{\bar{i}} \mid \forall j \in [0, q_{\bar{i}}-1] \omega_{\bar{i}}(j) \neq \bar{\delta}\}) \quad (4)$$

as, for each $i \neq \bar{i}$ ($0 \leq i < k$), $\Pr(\{\omega_i \in \Omega_i\}) = 1$.

Probability (4) is the probability of picking a permutation $\omega_{\bar{i}}$ of $\Delta_{\bar{i}}$ which does not have $\bar{\delta}$ in the first $q_{\bar{i}}$ positions, and evaluates to $1 - \frac{q_{\bar{i}}(|\Delta_{\bar{i}}|-1)!}{|\Delta_{\bar{i}}|!} = 1 - \frac{q_{\bar{i}}}{|\Delta_{\bar{i}}|}$ which is less than or equal to $1 - \min\{\frac{q_i}{|\Delta_i|} \mid 0 \leq i < k\}$. The thesis follows. \square

Theorem 2. Let Δ^λ be a file containing n disturbance traces. For any $z \geq 1$, Algorithm 1 stores in file $\Delta_{\text{rnd}}^\lambda$ a permutation of the traces in Δ^λ extracted with uniform probability among all possible permutations.

Proof. We prove the following: (i) all traces in the input sequence (file Δ^λ) will occur in the output sequence (file $\Delta_{\text{rnd}}^\lambda$) exactly once (i.e., the algorithm computes a permutation); (ii) for each δ occurring in Δ^λ , the probability that δ occurs in $\Delta_{\text{rnd}}^\lambda$ at any position is $\frac{1}{n}$ (i.e., the computed permutation is uniformly extracted among all possible permutations).

Algorithm 1 Random Sequence Generator.

Input: Δ^λ , a file holding a labelled sequence of disturbance traces
Output: $\Delta_{\text{rnd}}^\lambda$, a file holding a random permutation of disturbance traces, uniformly chosen among all possible permutations

- 1 **param** z , *max # of traces that can be kept in RAM*
- 2 $\Delta_{\text{rnd}}^\lambda \leftarrow$ a new empty file;
- 3 $\Delta_{\text{tmp}}^\lambda \leftarrow \Delta^\lambda$; /* input of 1st round */
- 4 $r \leftarrow 1$; /* round counter */
- 5 **while** $\Delta_{\text{tmp}}^\lambda$ *is not empty* **do**
- 6 $\Delta_{\text{in}}^\lambda \leftarrow \Delta_{\text{tmp}}^\lambda$; /* tmp file of prev. round */
- 7 $\Delta_{\text{tmp}}^\lambda \leftarrow$ a new temp file; /* next round input */
- 8 **if** $z > |\Delta_{\text{in}}^\lambda|$ **then** $z \leftarrow |\Delta_{\text{in}}^\lambda|$;
- 9 $\text{selected_pos} \leftarrow \text{rndPermPrefix}(|\Delta_{\text{in}}^\lambda|, z)$; /* selected_pos is a mapping from $[0, z - 1]$ to $[0, |\Delta_{\text{in}}^\lambda| - 1]$ */
- 10 $\text{selected_traces} \leftarrow$ empty array of z elements;
- 11 $p \leftarrow 0$;
- 12 **foreach** $\delta \in \Delta_{\text{in}}^\lambda$ **do**
- 13 **if** $\exists i \text{ selected_pos}[i] = p$ **then** $\text{selected_traces}[i] \leftarrow \delta$;
- 14 **else** append trace δ to $\Delta_{\text{tmp}}^\lambda$;
- 15 $p \leftarrow p + 1$;
- 16 **for** $i \leftarrow 0 \dots z - 1$ **do**
- 17 append trace $\text{selected_traces}[i]$ to $\Delta_{\text{rnd}}^\lambda$;
- 18 $r \leftarrow r + 1$;
- 19 **return** $\Delta_{\text{rnd}}^\lambda$;

20 **function** $\text{rndPermPrefix}(n, z)$
Input: n , positive integer
Input: z , positive integer ($z \leq n$)
Output: result , the first z elements of a permutation of $[0, n - 1]$ uniformly chosen at random

- 21 $\text{result} \leftarrow$ array of z elements;
- 22 $\text{perm} \leftarrow$ empty associative array;
- 23 **for** $i \leftarrow 0 \dots z - 1$ **do**
- 24 $j \leftarrow$ random index between i and $n - 1$;
- 25 **if** *exists* v s.t. $\langle i, v \rangle \in \text{perm}$ **then** $\text{value}_i \leftarrow v$;
- 26 **else** $\text{value}_i \leftarrow i$;
- 27 **if** *exists* v s.t. $\langle j, v \rangle \in \text{perm}$ **then** $\text{value}_j \leftarrow v$;
- 28 **else** $\text{value}_j \leftarrow j$;
- 29 put (i, value_j) into perm ;
- 30 put (j, value_i) into perm ;
- 31 $\text{result}[i] \leftarrow \text{value}_j$;
- 32 **return** result ;
- 33 **end**

Point (i) is immediate: at each round, z traces are appended to $\Delta_{\text{rnd}}^\lambda$, all the others are appended to $\Delta_{\text{tmp}}^\lambda$, which becomes the input of the next round. Also, the algorithm terminates only if the $\Delta_{\text{tmp}}^\lambda$ produced at the previous round is empty.

To prove point (ii), for any p, q in $[0, n - 1]$, we compute the probability that trace δ having position p in the input sequence (file Δ^λ) will have position q in the output sequence (file $\Delta_{\text{rnd}}^\lambda$).

We omit to prove that function $\text{rndPermPrefix}(n, z)$ actually computes the first z elements of a permutation of the integer interval $[0, n - 1]$ uniformly selected at random, as the function interrupts a well-known swap-based permutation algorithm as soon as the first z elements have been determined.

Given that, at each round $r \geq 1$, the main algorithm selects the z input traces which will have output positions $(r - 1)z$ to $\min(n, rz - 1)$, δ is selected only at round $r_\delta = \lceil (q + 1)/z \rceil$. Thus, the probability that δ , having input position p , will have output position q is:

$$\Pr \left(\left(\bigcap_{r=1}^{r_\delta-1} \neg E_r \right) \cap O_{r_\delta}^{q'} \right) \quad (5)$$

where, for all r , $\neg E_r$ denotes the event “ δ is not selected at round r ” and $O_{r_\delta}^{q'}$ denotes the event “ δ is the q' th trace selected in round r_δ ,” where $q' = (q - z(r_\delta - 1))$.

By the chain rule, (5) becomes:

$$\Pr(-E_1) \times \cdots \times \Pr(-E_{r_\delta-1} | -E_{r_\delta-2}, \dots, -E_1) \\ \times \Pr(O_{r_\delta}^{q'} | -E_{r_\delta-1}, \dots, -E_1). \quad (6)$$

For all $1 \leq r \leq r_\delta - 1$:

$$\Pr(-E_r | -E_{r-1}, \dots, -E_1) = \frac{n - z(r-1) - 1}{n - z(r-1)} \times \cdots \\ \times \frac{n - z(r-1) - z}{n - z(r-1) - (z-1)}$$

where, for all $0 \leq i < z$, the i th factor is the number of ways we can choose a trace different from δ out of $n - z(r-1) - i$ (where $n - z(r-1)$ is the number of traces still not selected at the beginning of round r). The expression simplifies to:

$$\Pr(-E_r | -E_{r-1}, \dots, -E_1) = \frac{n - zr}{n - z(r-1)}$$

and the product $\Pr(-E_1) \times \cdots \times \Pr(-E_{r_\delta-1} | -E_{r_\delta-2}, \dots, -E_1)$ is:

$$\Pr(-E_1) \times \cdots \times \Pr(-E_{r_\delta-1} | -E_{r_\delta-2}, \dots, -E_1) = \prod_{r=1}^{r_\delta-1} \frac{n - zr}{n - z(r-1)} \\ = \frac{n - z(r_\delta - 1)}{n}. \quad (7)$$

As for $\Pr(O_{r_\delta}^{q'} | -E_{r_\delta-1}, \dots, -E_1)$, i.e., the probability that δ is the q' th selected trace in round r_δ provided that it has not been selected in previous rounds, it can be computed as:

$$\frac{n - z(r_\delta - 1) - 1}{n - z(r_\delta - 1)} \times \cdots \times \frac{n - z(r_\delta - 1) - (q' - 1)}{n - z(r_\delta - 1) - (q' - 2)} \\ \times \frac{1}{n - z(r_\delta - 1) - (q' - 1)} = \frac{1}{n - z(r_\delta - 1)} \quad (8)$$

where, for all $0 \leq i < q' - 1$, the i th factor is the number of ways we can choose a trace different from δ out of $n - z(r_\delta - 1) - i$ (where $n - z(r_\delta - 1)$ is the number of traces still not selected at the beginning of round r_δ). The last factor is the probability of selecting δ (still unselected) out of $n - z(r_\delta - 1) - (q' - 1)$ traces.

By (7) and (8), probability (6) evaluates to $\frac{1}{n}$, which is independent of p and q . \square

6. Experimental results

In this section we evaluate the effectiveness of our random exhaustive SLFV (rSLFV) approach as follows. First, we evaluate the *overhead* due to the randomisation of disturbance traces needed to enable computation of Omission Probability (OP), by comparing our rSLFV approach with the Deterministic SLFV (dSLFV) approach of [8]. Second, we evaluate the behaviour of the coverage and the OP bound with respect to simulation time. Third, we evaluate speed-up and efficiency of our parallel approach.

6.1. Experimental setting

In this section we describe the case studies and the computational infrastructure we used in our experiments.

6.1.1. Case studies

We experiment with two case studies, using two models included in the Simulink distribution, namely the Inverted Pendulum on a Cart (IPC) and the Fuel Control System (FCS).

6.1.1.1. Inverted Pendulum on a Cart (IPC). The IPC is a control loop system where the controlled system is an inverted pendulum installed on a cart (see Fig. 7). The controller (actually a control software) senses the angular position θ of the pendulum, and computes the force F to be applied to the cart to move it left or right

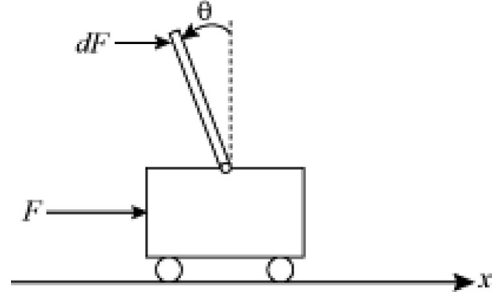


Fig. 7. Inverted Pendulum on a Cart (IPC) case study (from mathworks.com).

along the x axis. The goal is to keep the pendulum in its upright (vertical) unstable position. The physical constraint between the cart and the pendulum gives that both the cart and the pendulum have one degree of freedom each (x and θ , respectively).

The controlled system consists of the cart and the pendulum, whereas the controller consists of the control software computing F from the plant outputs (x and θ). Accordingly, our overall System Under Verification (SUV) model consists of the controlled system and the controller, whose Simulink block diagram is shown in the upper box of Fig. 8.

The system level property that we verify is that after 2 s the pendulum is in upright position, i.e., angle θ is always between $[-0.1, 0.1]$. The monitor checking for this property is shown in the lower box of Fig. 8.

We introduce disturbances by injecting irregularities in the cart rail. We model such irregularities with a modification on the cart weight m with respect to its nominal value of 0.455 kg. For this, we define three disturbances representing normal rail operation ($m = 0.455$ kg), abnormal rail operation ($m = 1.455$ kg), and stressed rail operation ($m = 2.455$ kg).

We consider two disturbance models for the IPC, \mathcal{D}_{IPC}^1 and \mathcal{D}_{IPC}^2 . Model \mathcal{D}_{IPC}^1 has a horizon of $h = 90$ and defines 3 208 276 disturbance traces. Model \mathcal{D}_{IPC}^2 is defined extending \mathcal{D}_{IPC}^1 with more complex operational scenarios and defines 35 641 501 disturbance traces over a horizon of $h = 200$. For both models we set τ (quantum between disturbances) to 0.1 s. A detailed description of \mathcal{D}_{IPC}^1 and \mathcal{D}_{IPC}^2 is not relevant for the evaluation of our experiments below. We only observe that, in our setting, the complexity of answering a System Level Formal Verification (SLFV) problem primarily depends on the number of disturbance traces to be simulated. Thus, the worst case for our approach is when all disturbance traces have to be simulated, i.e., when the answer to the SLFV problem is *PASS*. Indeed, both \mathcal{D}_{IPC}^1 and \mathcal{D}_{IPC}^2 satisfy this requirement.

6.1.1.2. Fuel Control System (FCS). The FCS is a controller for a fault tolerant gasoline engine, which has also been used as a case study in [8,12,15–18].

The FCS has four sensors: throttle angle, speed, EGO (measuring the residual oxygen present in the exhaust gas) and MAP (manifold absolute pressure). The goal of the control system is to maintain the air-fuel ratio (the ratio between the air mass flow rate pumped from the intake manifold and the fuel mass flow rate injected at the valves) close to the stoichiometric ratio of 14.6, which represents a good compromise between power, fuel economy, and emissions.

From the sensor measurements, the FCS estimates the mixture ratio and provides feedback to the closed-loop control, yielding an increase or a decrease of the fuel rate.

The FCS sensors are subject to faults (disturbances), and the whole control system can tolerate single sensor faults. In particular, if a sensor fault is detected, the FCS changes its control law

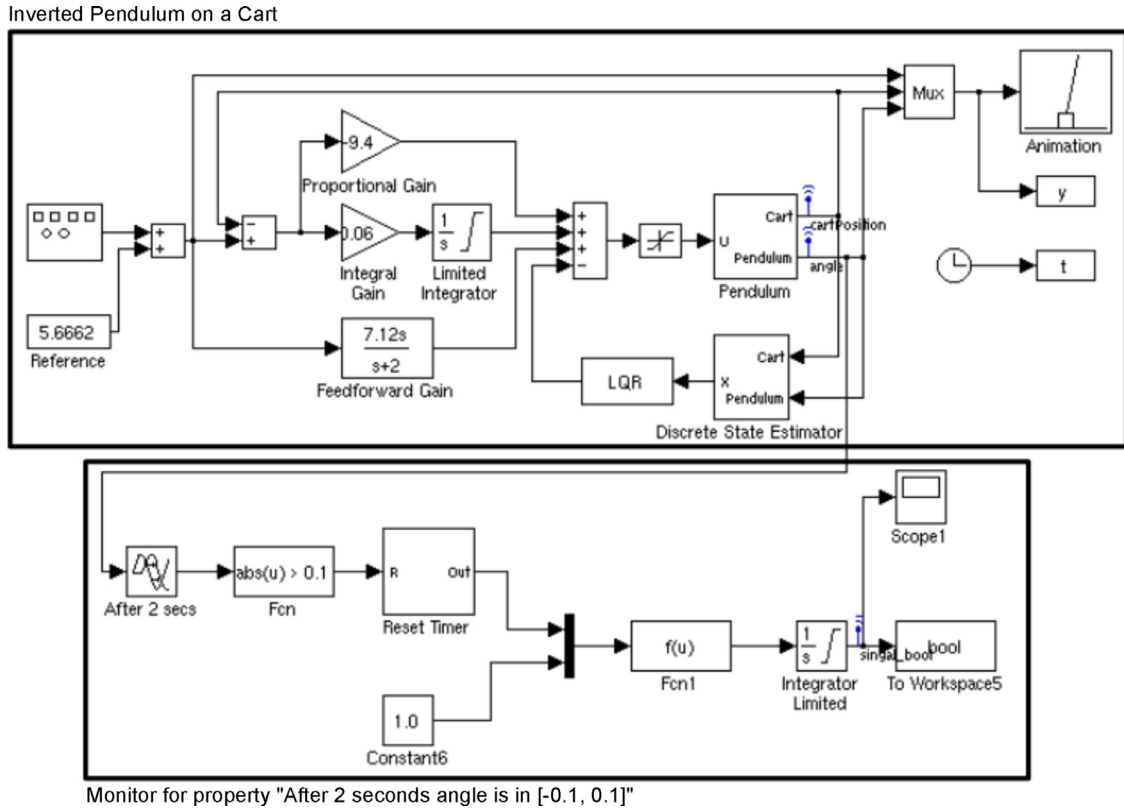


Fig. 8. Simulink block diagram for Inverted Pendulum on a Cart (from mathworks.com) with an embedded property monitor.

by operating the engine with a higher fuel rate to compensate. In case two or more sensors fail, the FCS shuts down the engine, as the air-fuel ratio cannot be controlled.

The control logic of the FCS is implemented by six automata, each one with a number of states ranging from two to five. The signal flow is further subdivided into three subsystems, which exhibit several different Simulink block types, involving arithmetic, lookup tables, integrators, filters and interpolation [19] (see [20] for more details).

We verify one of the system level specifications for such a model, namely: the *fuel_air* model variable is never 0 for more than one second. Accordingly, our SUV consists of the Simulink FCS model along with a monitor for the property under verification (such a model is shown as Fig. 9).

We consider two disturbance models for the FCS, \mathcal{D}_{FCS}^1 and \mathcal{D}_{FCS}^2 . Model \mathcal{D}_{FCS}^1 has a horizon of $h = 100$ and defines 4 023 955 disturbance traces. Model \mathcal{D}_{FCS}^2 is defined extending \mathcal{D}_{FCS}^1 with more complex operational scenarios and defines 12 948 712 disturbance traces over a horizon of $h = 200$. For both models we set τ (quantum between disturbances) to 1 s. A detailed description of \mathcal{D}_{FCS}^1 and \mathcal{D}_{FCS}^2 is not relevant for the evaluation of our experiments below, and can be found in [8,21]. We only observe that, as it happens with our disturbance models for the IPC, for all disturbance traces entailed by \mathcal{D}_{FCS}^1 and \mathcal{D}_{FCS}^2 , the property to be verified is satisfied. This yields the worst scenario to answer our SLFV problem, as all traces need to be simulated.

6.1.2. Computational infrastructure

We ran experiments on multiple Linux PCs, each one equipped with 2 Intel Xeon 3.0 GHz CPUs with 4 cores each and 8 GB RAM. We executed 8 processes (optimisation and simulation) in parallel (one per available core) on each machine.

Table 1

Disturbance trace generation.

SUV	Dist. model	#Traces	Gen. time	File size
IPC	\mathcal{D}_{IPC}^1	3208276	0:9:58	4.6 GB
IPC	\mathcal{D}_{IPC}^2	35641501	7:28:24	107 GB
FCS	\mathcal{D}_{FCS}^1	4023955	0:28:39	3.5 GB
FCS	\mathcal{D}_{FCS}^2	12948712	4:45:47	39 GB

As, in a multi-core setting, the local disk may quickly become a performance bottleneck if heavily used by multiple processes, we have replaced it with 8 RAM disks of 500 MB each per machine, in order to store simulation states. Accordingly, we have used the multi-core version of the dSLFV optimiser of [8] as presented in [12]. Given that, in our case studies, the size of the simulation state files is of about 150–300 KB, this experimental setting allowed our optimiser to count on the possibility, for each simulator, to keep at most 1800 states simultaneously stored.

6.2. Disturbance trace generation and splitting

Along the lines of [8], we use CMurphi to generate the set of labelled admissible disturbance traces entailed by the disturbance model given as input.

Table 1 shows, for each of the four disturbance models we consider (two for IPC and two for FCS) the number of entailed disturbance traces (column #Traces), the time needed by CMurphi to generate them (column Gen. time) and the size of the file computed by our generator to store them (column File size).

We then split the generated sequences of disturbance traces into k slices, with $k = 128, 256, 512$ to enable parallel computation on, respectively, 16, 32, 64 (8-core) machines. Splitting always takes negligible time.

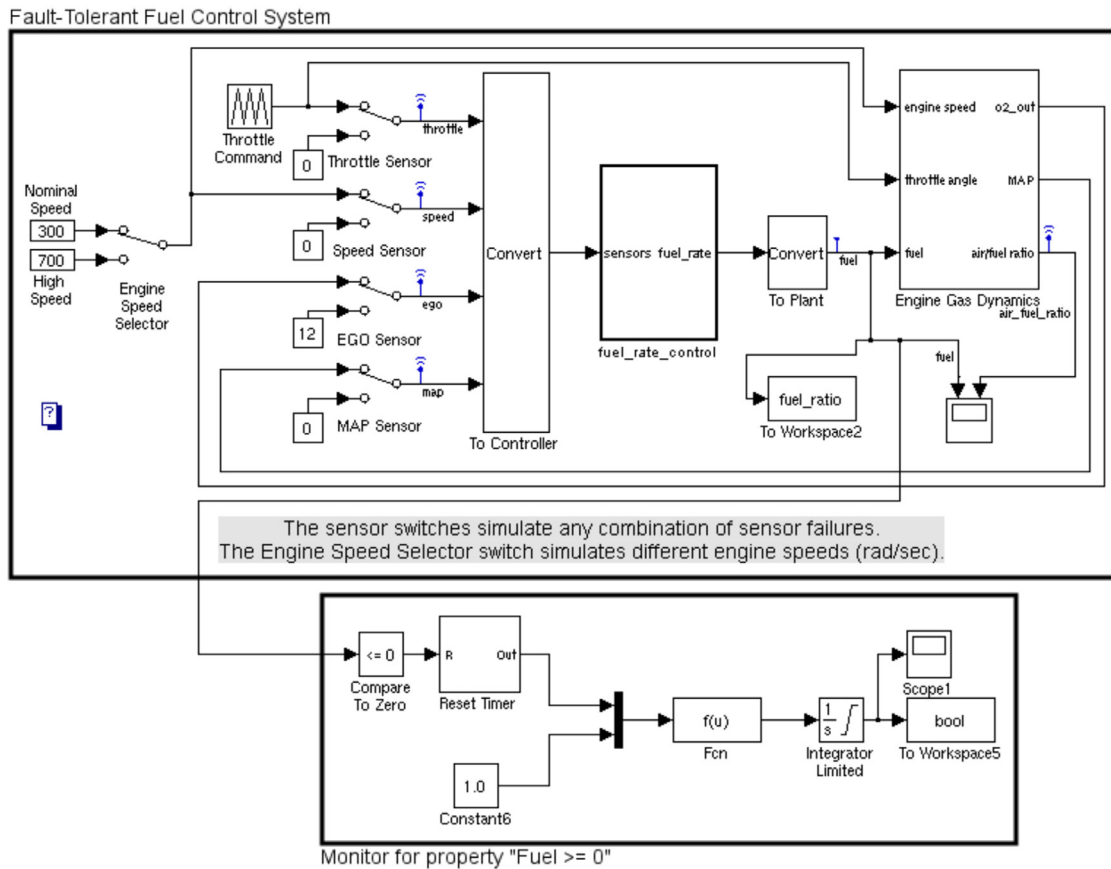


Fig. 9. Simulink/Stateflow representation of the Fuel Control System (from mathworks.com) with an embedded property monitor.

6.3. Computation of optimised simulation campaigns

Table 2 shows, for each of the disturbance models we consider, the time needed by our Random Sequence Generator module to enable anytime OP computation during the verification process (rSLFV) and the time needed to compute optimised simulation campaigns with (rSLFV) and without (dSLFV) trace order randomisation.

Column #Mach gives the number of machines we used in parallel. Column #Slices gives the overall number of slices in which the sequence of admissible disturbance traces has been partitioned (one per available core, i.e., 8 slices per machine). Column #Traces per slice shows the number of traces in any single slice (except the last slice, which has fewer traces when the overall number of traces is not a multiple of #Slices). Column group rSLFV shows the maximum overall time to compute the simulation campaign starting from a slice. This time is split into two parts: the time needed to execute the Random Sequence Generator module (column RSG) and the time needed to compute the simulation campaign starting from the randomised slice (column sim. camp. comp.). Column dSLFV shows the maximum time to compute the simulation campaign starting from a slice when no RSG is performed, thus no anytime OP computation is enabled. Column rSLFV overhead shows the difference between the rSLFV and dSLFV times.

It can be observed that our parallel approach to computation of optimised randomised simulation campaigns is able to effectively exploit parallelism in order to handle disturbance models entailing tens of millions of operational scenarios.

Also, Table 2 shows that the random sequence generation phase makes the rSLFV simulation campaign computation process longer than that for dSLFV (i.e., overhead is positive). The difference

is, however, negligible with respect to the whole verification time, which takes many hours, even for the massively parallel simulation of operational scenarios entailed by our smallest disturbance models, i.e., \mathcal{D}_{IPC}^1 for the IPC and \mathcal{D}_{FCS}^1 for the FCS (see Section 6.4).

6.4. Execution of the simulation campaigns

Table 3 shows the execution time of the simulation campaigns generated by dSLFV and rSLFV.

As the exhaustive simulation of the traces generated by disturbance models \mathcal{D}_{IPC}^2 , \mathcal{D}_{FCS}^2 (entailing, respectively, 35 641 501 and 12 948 712 disturbance traces) would require a prohibitively long time, from now on we restrict ourselves to the simulation of the traces generated by \mathcal{D}_{IPC}^1 and \mathcal{D}_{FCS}^1 (which entail, respectively, 3 208 276 and 4 023 955 disturbance traces).

By enabling the computation of the OP during the simulation activity we have a quite significant increase of simulation time. However, such an overhead can be drastically reduced, or even neutralised, by using more parallel processes (higher values for $k = \#Slices$). This behaviour is due to the fact that the rSLFV optimiser needs to compute a simulation campaign under the restriction that the number of states that the simulator can keep simultaneously stored is limited (to fit within the 500 MB of disk space available to each simulator instance). For high values of $k = \#Slices$ (e.g., $k = 512$), this is not a big obstacle. On the other hand, for low values of k , the number of traces in each slice is higher and they share shorter common prefixes on average. Hence, a fully-optimised random order execution of them would need a too high number of simulation states to be simultaneously kept stored. As a consequence, the optimiser is forced to post free commands for many simulation states which would be needed again in

Table 2

Overhead of enabling OP (rSLFV) with respect to dSLFV in the computation of simulation campaigns (time in h:m:s).

#Mach.	#Slices	#Traces per slice	rSLFV			dSLFV		rSLFV overhead
			RSG	sim.camp.comp.	Total	sim.camp.comp.	overhead	
1	8	401 035	0:0:21	1:8:52	1:9:13	0:8:53	1:0:20	
2	16	200 518	0:0:11	0:11:57	0:12:9	0:4:36	0:7:32	
4	32	100 259	0:0:5	0:6:33	0:6:38	0:2:23	0:4:14	
8	64	50 130	0:0:3	0:2:50	0:2:53	0:1:14	0:1:39	
16	128	25 065	0:0:2	0:1:23	0:1:26	0:0:38	0:0:47	
32	256	12 533	0:0:1	0:1:19	0:1:21	0:0:20	0:1:1	
64	512	6267	0:0:1	0:0:21	0:0:23	0:0:11	0:0:12	
(a) Inverted Pendulum on a Cart (IPC), disturbance model \mathcal{D}_{IPC}^1 : 3 208 276 traces with horizon $h = 90$								
#Mach.	#Slices	#Traces per slice	rSLFV			dSLFV		rSLFV overhead
			RSG	sim.camp.comp.	total	sim.camp.comp.	overhead	
16	128	278 450	0:1:40	2:41:28	2:43:8	0:35:0	2:8:8	
32	256	139 225	0:1:6	1:9:12	1:10:18	0:17:40	0:52:38	
64	512	69 613	0:0:34	0:21:26	0:22:0	0:8:57	0:13:3	
(b) Inverted Pendulum on a Cart (IPC), disturbance model \mathcal{D}_{IPC}^2 : 35 641 501 traces with horizon $h = 200$								
#Mach.	#Slices	#Traces per slice	rSLFV			dSLFV		rSLFV overhead
			RSG	sim.camp.comp.	total	sim.camp.comp.	overhead	
1	8	502 995	0:1:12	0:4:52	0:6:4	0:5:27	0:0:37	
2	16	251 498	0:0:24	0:2:34	0:2:58	0:2:8	0:0:50	
4	32	125 749	0:0:15	0:2:24	0:2:39	0:0:57	0:1:42	
8	64	62 875	0:0:8	0:1:20	0:1:28	0:0:29	0:0:59	
16	128	31 438	0:0:7	0:1:19	0:1:26	0:0:17	0:1:9	
32	256	15 719	0:0:7	0:0:32	0:0:39	0:0:8	0:0:31	
64	512	7860	0:0:6	0:0:5	0:0:11	0:0:4	0:0:7	
(c) Fuel Control System (FCS), disturbance model \mathcal{D}_{FCS}^1 : 4 023 955 traces with horizon $h = 100$								
#Mach.	#Slices	#Traces per slice	rSLFV			dSLFV		rSLFV overhead
			RSG	sim.camp.comp.	total	sim.camp.comp.	overhead	
16	128	101 162	0:0:45	0:20:36	0:21:21	0:8:18	0:13:3	
32	256	50 581	0:0:24	0:15:35	0:15:59	0:4:31	0:11:28	
64	512	25 291	0:0:14	0:6:43	0:6:57	0:2:4	0:4:53	
(d) Fuel Control System (FCS), disturbance model \mathcal{D}_{FCS}^2 : 12 948 712 traces with horizon $h = 200$								

Table 3

Parallel execution of simulation campaigns by dSLFV and rSLFV (time in h:m:s).

#Mach.	#Slices	Min	Max	Avg	Approach
16	128	2:14:2	4:10:52	3:44:52	dSLFV
		25:14:28	85:10:36	58:9:43	rSLFV
		23:0:26	80:59:44	54:24:51	overhead
32	256	1:6:49	2:5:40	1:50:28	dSLFV
		4:26:16	23:59:0	13:0:24	rSLFV
		3:19:27	21:53:20	11:9:56	overhead
64	512	0:30:30	1:3:10	0:53:45	dSLFV
		0:59:26	1:51:10	1:43:12	rSLFV
		0:28:56	0:48:0	0:49:27	overhead
(a) Inverted Pendulum on a Cart (IPC), disturbance model \mathcal{D}_{IPC}^1					
#Mach.	#Slices	Min	Max	Avg	Approach
16	128	70:6:4	100:17:53	87:49:56	dSLFV
		216:42:13	348:51:47	308:46:18	rSLFV
		146:36:9	248:33:54	220:56:22	overhead
32	256	44:0:27	57:57:27	48:34:6	dSLFV
		63:53:54	136:18:14	108:14:19	rSLFV
		19:53:27	78:20:47	59:40:13	overhead
64	512	18:32:36	26:49:4	23:2:19	dSLFV
		22:9:19	29:23:33	26:43:31	rSLFV
		3:36:43	2:34:29	3:41:12	overhead
(b) Fuel Control System (FCS), disturbance model \mathcal{D}_{FCS}^1					

yet-to-be-simulated traces. Such traces will then be simulated from the simulator initial state, thus yielding performance degradation.

6.5. Overall verification time and scalability

In this section we evaluate the overall impact of enabling any-time OP computation and the scalability of our parallel approach to SLFV.

Table 4 shows the *overall time* needed to carry out our SLFV tasks (IPC with disturbance model \mathcal{D}_{IPC}^1 and FCS with disturbance model \mathcal{D}_{FCS}^1) with k parallel processes, for the values of k already used in the previous tables.

In particular, column *Gen. & comp. sim. camp.* reports the sum of the disturbance trace generation and splitting time, parallel randomisation (only for rSLFV) and parallel simulation campaign computation time (from Table 2). Column *Simulation* reports the parallel simulation time when using k parallel processes (i.e., the maximum simulation time over all the $k = \#Slices$ slices as in column *Max* of Table 3). Column *Overall* reports the overall time to carry out each dSLFV and each rSLFV task, as the sum of the previous two columns.

We observe that our approach takes *negligible* time to generate optimised simulation campaigns with respect to the time needed to actually simulate them (e.g., minutes vs. hours).

6.5.1. Estimation of sequential verification time

In order to evaluate the scalability of our parallel approach to SLFV, Table 4 also reports (in the first two rows of each sub-table),

Table 4

Overall performance overhead (including disturbance trace generation and splitting, trace randomisation, computation of simulation campaigns and Simulink simulations) of rSLFV with respect to dSLFV (time in h:m:s).

#Mach.	#Slices	Gen. & comp. sim. camp.	Simulation	Overall	Speedup	Efficiency	Approach
1	1	0:56:11 26:59:57	458:40:0 ^a	459:36:11 ^a	1.00 ×	100.00%	dSLFV
			880:38:24 ^a	907:38:21 ^a	1.00 ×	100.00%	rSLFV
				+97.48%	+0.00%	+0.00%	overhead
16	128	0:10:36 0:11:24	4:10:52	4:21:28	105.46 ×	82.39%	dSLFV
			85:10:36	85:22:0	10.63 ×	8.30%	rSLFV
				+1858.83%	+89.92%	+74.09%	overhead
32	256	0:10:18 0:11:19	2:5:40	2:15:58	202.80 ×	79.22%	dSLFV
			23:59:0	24:10:19	37.55 ×	14.67%	rSLFV
				+966.57%	+81.48%	+64.55%	overhead
64	512	0:10:9 0:10:21	1:3:10	1:13:19	376.04 ×	73.45%	dSLFV
			1:51:10	2:1:31	448.13 ×	87.53%	rSLFV
				+65.71%	−19.17%	−14.08%	overhead

(a) Inverted Pendulum on a Cart (IPC), disturbance model \mathcal{D}_{IPC}^1

#Mach.	#Slices	Gen. & comp. sim. camp.	Simulation	Overall	Speedup	Efficiency	Approach
1	1	0:35:55 3:4:56	11242:31:28 ^a	11243:7:23 ^a	1.00 ×	100.00%	dSLFV
			13683:20:32 ^a	13686:25:28 ^a	1.00 ×	100.00%	rSLFV
				+21.73%	+0.00%	+0.00%	overhead
16	128	0:28:56 0:30:5	100:17:53	100:46:49	111.56 ×	87.16%	dSLFV
			348:51:47	349:21:52	39.18 ×	30.61%	rSLFV
				+246.66%	+64.88%	+56.55%	overhead
32	256	0:28:47 0:29:18	57:57:27	58:26:14	192.40 ×	75.16%	dSLFV
			136:18:14	136:47:32	100.05 ×	39.08%	rSLFV
				+134.08%	+48.00%	+36.08%	overhead
64	512	0:28:43 0:28:50	26:49:4	27:17:47	411.89 ×	80.45%	dSLFV
			29:23:33	29:52:23	458.15 ×	89.48%	rSLFV
				+9.44%	−11.23%	−9.03%	overhead

(b) Fuel Control System (FCS), disturbance model \mathcal{D}_{FCS}^1

^a Estimated value.

the overall dSLFV and rSLFV times when using only one parallel process (*sequential time*). Unfortunately, as for our case studies a sequential simulation would be prohibitively long, we have estimated the sequential simulation time to carry out both dSLFV and rSLFV as follows.

Let t_k^{avg} be the average time to simulate a slice where $k = \#Slices$ parallel processes are used (row *#slices* = k , column *avg*, for either dSLFV or rSLFV). For any value of k , the sequential simulation time could be estimated as $k \times t_k^{\text{avg}}$. As this value changes a little bit for different values of k , Table 4 estimates sequential simulation time as $\min\{128t_{128}^{\text{avg}}, 256t_{256}^{\text{avg}}, 512t_{512}^{\text{avg}}\}$. Such huge values (weeks of computation) make clear that estimation is the only viable way to compute the simulation sequential times. Note that in our computation we are slightly overestimating the sequential time, since we are assuming that some traces of each slice must be simulated from the initial state. In an actual 1-process execution of a simulation campaign, the optimiser may exploit stored simulator states to avoid simulation of such traces from the initial state. As the time to simulate a single trace is of a few seconds and the simulator can keep only a limited number of stored states, this is negligible with respect to the value of the sequential simulation time.

6.5.2. Speedup and efficiency

Sequential simulation time for both dSLFV and rSLFV is used in Table 4 to compute the *speedup* and the *efficiency* of our parallel approach to SLFV, as typically done in the evaluation of parallel algorithms. In particular, for each $k = \#Slices$, column *Speedup* shows the ratio t_1/t_k , where t_1 is the estimated overall sequential verification time and t_k is the overall verification time when k parallel processes are used. Column *Efficiency* is computed by dividing the speedup by the number of parallel processes $k = \#Slices$.

Table 4 also shows the overhead (see bold values) due to randomisation of the verification task (which is the price to pay in order to enable anytime computation of OP), both in terms of overall verification time increase and in terms of reduction of speedup and efficiency. We observe that such an overhead is significant, but it can be drastically reduced by increasing the number k of parallel processes.

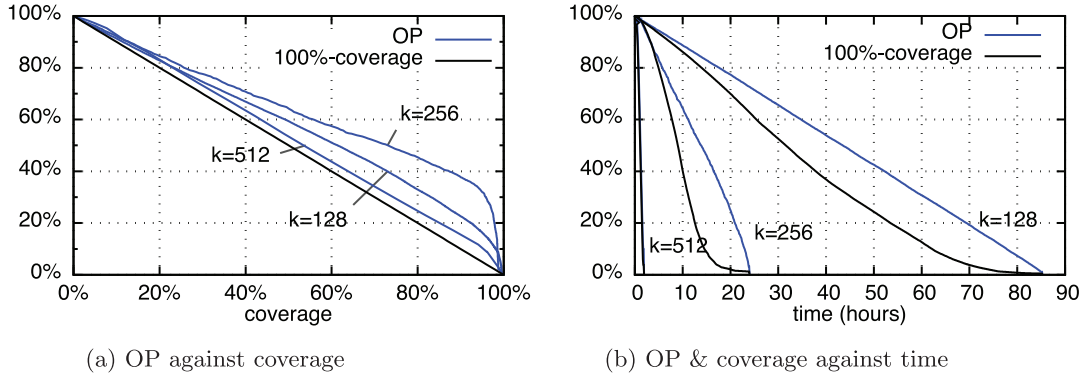
6.6. Omission probability

Figs. 10a and c show how our upper bound to the OP decreases as a function of the coverage (i.e., the ratio of admissible traces simulated) during the parallel execution of the k simulation campaigns (IPC with disturbance model \mathcal{D}_{IPC}^1 and FCS with disturbance model \mathcal{D}_{FCS}^1), for $k = 128, 256, 512$. It can be observed that our OP bound is always *very close* to the ratio of yet-to-be-simulated traces (curves named “100%-coverage”, i.e., 100% minus coverage), which is the best one can do (using only one parallel process) without any assumption on the number of error traces.

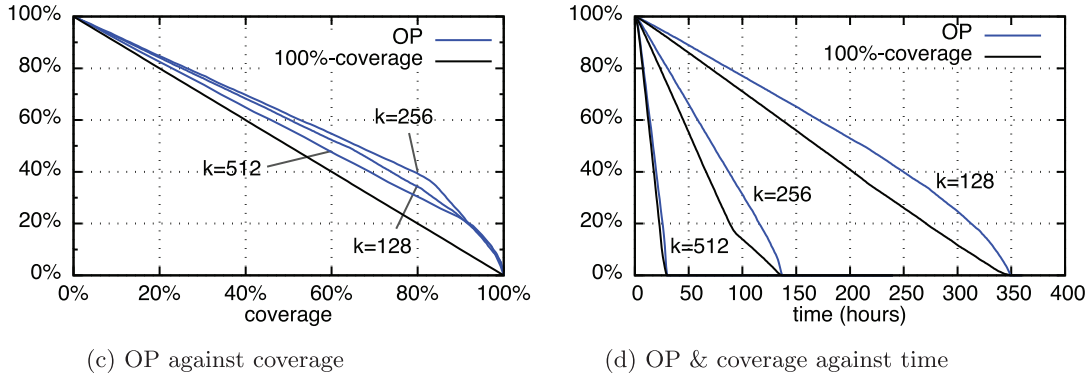
6.7. Completion time estimation

Figs. 10b and d show that OP bound, computed during the parallel execution of the simulation campaigns (IPC with disturbance model \mathcal{D}_{IPC}^1 and FCS with disturbance model \mathcal{D}_{FCS}^1), decreases nearly *linearly* in time. The same happens with the coverage, which can thus be used as a reliable *estimator* for the completion time of the whole verification process.

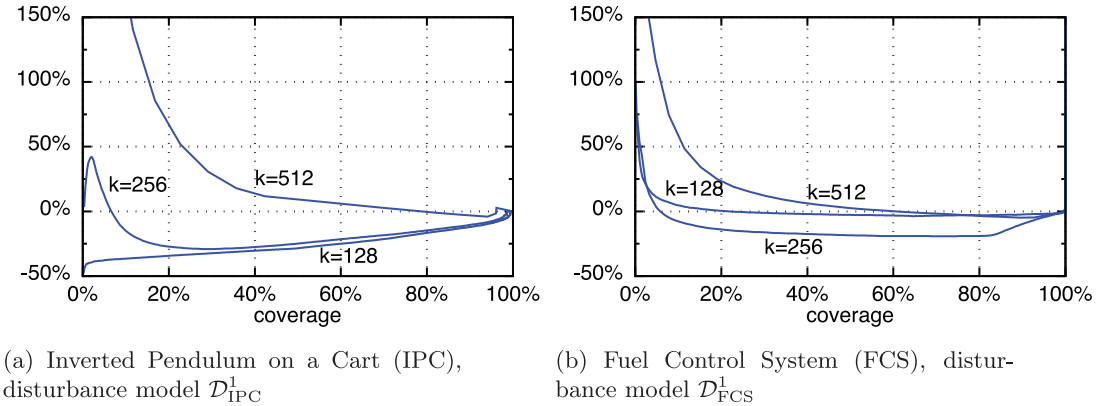
Fig. 11 shows the error percentage (on the true completion time) made by a completion time estimation based on the coverage. For each value x of the coverage, the error is computed as



(a) OP against coverage
(b) OP & coverage against time
Inverted Pendulum on a Cart (IPC), disturbance model \mathcal{D}_{IPC}^1



(c) OP against coverage
(d) OP & coverage against time
Fig. 10. Omission Probability (OP) computation during the parallel execution of the simulation campaigns.



(a) Inverted Pendulum on a Cart (IPC), disturbance model \mathcal{D}_{IPC}^1
(b) Fuel Control System (FCS), disturbance model \mathcal{D}_{FCS}^1
Fig. 11. Completion time estimation error against coverage.

$((t_x/x) - t_c)/t_c$ where t_x is the time elapsed to reach coverage x and t_c is the true completion time. It can be observed that such a completion time estimation becomes accurate quickly (e.g., when the coverage is $\geq 30\%$, the error is within 30%).

7. Related work

A parallel exhaustive Hardware In the Loop Simulation based hybrid system model checking similar to the one described in this work is presented in [8]. The main differences of the present work with respect to [8] are the following. (i) Our simulation campaign optimiser and the one in [8] both take as input the admissible disturbance traces (simulation scenarios). However, the simulation campaigns computed in [8] schedule scenarios according to their order, whereas in this work we introduce an intermediate step which enables simulation of *all* scenarios, *exactly once*, in a uniform

random order. (ii) During the verification process, the approach in [8] only outputs the attained coverage, whereas, in this work also the attained Omission Probability (OP) is computed, by exploiting the randomisation of the order with which scenarios are scheduled.

The work in [9] considers a finite state (digital hardware verification) setting and presents an algorithm to estimate the coverage achieved during SAT based bounded model checking. Since computation paths are not selected uniformly at random, [9] does not provide any information about the OP.

Random model checking is a formal verification approach closely related to our setting. A random model checker provides, at *any time* during the verification process, an upper bound to the OP. Upon detection of an error, a random model checker returns a counterexample. Random model checking algorithms have been investigated, e.g., in [10,22,23]. The main differences with

respect to our approach are the following. (i) All random model checkers generate simulation scenarios using a sort of Monte-Carlo based random walk. As a result, unlike our algorithm, none of them is exhaustive (within a finite time horizon). (ii) Random model checkers (e.g., see [10]) assume availability of a lower bound to the probability of selecting (with a random-walk) an error trace. Of course, being exhaustive, we do not have any such assumption.

The coverage yielded by random sampling a set of test cases has been studied by mapping it to the Coupon Collector's Problem CCP (see, e.g., [24]). In CCP elements are randomly extracted (uniformly and with replacement) from a finite set of n test cases (disturbance traces in our context). Known results (see, e.g., [25]) tell us that the probability distribution of the number of test cases to be extracted in order to collect *all* n elements has *expected* value $\Theta(n \log n)$, and a small variance with known bounds. This allows us to bound the OP during the verification. Differently from such CCP-based approaches, here we not only bound the OP, but also grant the completion of our verification task within *just* n trials. This is made possible by the fact that we first generate all disturbance traces.

Monte-Carlo based robustness analysis of CPSs has been investigated in [26]. We note that, within a finite time bound, we are exhaustive whereas the approach in [26] is not. On the other hand, unlike our approach, [26] also evaluates how *robustly* the given property holds.

Probabilistic (e.g., see [27,28]) and, more specifically, *simulation-based statistical model checking* approaches (e.g., see [15,16,29–34]) are closely related to our work. In particular, [16] addresses statistical model checking of Simulink models and presents experimental results on one of the Simulink case studies we use here. The main differences between such approaches and ours are the following. (i) Probabilistic model checking is a *white-box* approach (a model is available), whereas we are in a *black-box* setting (only a simulator is available). Thus, only simulation-based statistical model checking approaches can be used in our context. (ii) Statistical model checking is not exhaustive (within a finite time horizon), whereas we are. (iii) Both probabilistic and statistical model checking use a stochastic model for the SUV, whereas in our setting the SUV is deterministic and disturbances are nondeterministic. The probability measure in our context, as in random model checking, stems from the randomisation of the verification process itself. (iv) None of the available simulation-based statistical model checking approaches addresses the problem of the optimisation of the simulation campaign, which is an essential step to make our *parallel random exhaustive* (HILS) based model checking viable.

Formal verification of Simulink models has been widely investigated, examples are in [35–37]. Such methods however focus on discrete time models (e.g., Stateflow or Simulink restricted to discrete time operators) with small domain variables. Therefore they are well suited to analyse critical subsystems, but cannot handle complex system level verification tasks (e.g., our case studies). This is indeed the motivation for the development of statistical model checking methods as those in [15,16], for the exhaustive HILS based approach in [8], and for our present parallel random exhaustive HILS based approach.

8. Conclusions

We presented a *parallel random exhaustive* (HILS) based model checker for hybrid systems that, while being *exhaustive* with respect to the disturbance model given as input, provides at *any time* during the verification process an *upper bound* to the probability that the SUV exhibits an error in a yet-to-be-simulated scenario (Omission Probability, OP).

Our experimental results on real world case studies from the Simulink distribution (namely: Inverted Pendulum on a Cart and

Fuel Control System) show that, by exploiting parallelism, our approach to the computation of optimised simulation campaigns is feasible even for disturbance models entailing tens of millions of operational scenarios.

Also, simulation results show that, by exploiting parallelism, our simulation campaign optimiser effectively counteracts the simulation time overhead stemming from randomisation.

Finally, we have shown that our bound to the OP decreases about *linearly* with the coverage, which is as good as it can be even in the worst case scenario (just one error trace). Furthermore, resting on randomisation, we can use the coverage as a reliable estimator for the time needed to complete the verification process.

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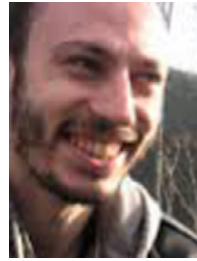
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