## STOCHASTIC HYBRID SYSTEMS WITH RENEWAL TRANSITIONS: MOMENT ANALYSIS WITH APPLICATION TO NETWORKED CONTROL SYSTEMS WITH DELAYS\*

DUARTE ANTUNES<sup>†</sup>, JOÃO P. HESPANHA<sup>‡</sup>, AND CARLOS SILVESTRE<sup>§</sup>

Abstract. We consider stochastic hybrid systems (SHSs) for which the lengths of times that the system stays in each mode are independent random variables with given distributions. We propose an approach based on a set of Volterra equations to compute any statistical moment of the state of the SHS. Moreover, we provide a method to compute the Lyapunov exponents of a given degree, i.e., the exponential rate of decrease or increase at which statistical moments converge to zero or to infinity, respectively. We also discuss how, by computing the statistical moments, one can provide information about the probability distribution of the state of the SHS. The applicability of the results is illustrated in the analysis of a networked control problem with independently distributed intervals between data transmissions and delays.

Key words. networked control systems, Volterra integral equations, semi-Markov processes

AMS subject classifications. 60K30, 45D05, 60K15

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1. Introduction and examples. Stochastic hybrid systems (SHSs) combine continuous dynamics and discrete logic. The execution of an SHS is specified by the dynamic equations of the continuous state, a set of rules governing the transitions between discrete modes, and reset maps determining jumps of the state at transition times. As surveyed in [1], various models of SHSs [2], [3], [4] have been proposed, mostly differing in the way randomness enters the different equations governing the evolution of the system. See also [5], [6], [7], [8].

In the present work, we consider SHSs with linear dynamics and linear reset maps and for which the length of time that the system stays in each mode are independent and identically distributed random variables, whose distributions may depend on the discrete mode. The process that combines the transition times and the discrete mode is typically called a Markov renewal process [9], which motivated us to refer to these systems as *stochastic hybrid systems with renewal transitions*. This class of systems can be viewed as a special case of the SHS model in [5], which in turn is a special case of a piecewise deterministic process [4] and also a special case of a state-dependent

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<sup>&</sup>lt;sup>†</sup>Hybrid and Networked Systems Group, Department of Mechanical Engineering, Eindhoven University of Technology, the Netherlands (D.Antunes@tue.nl). Former address: Department of Electrical Engineering and Computer Science, Instituto Superior Técnico, ISR, Lisboa, Portugal.

<sup>&</sup>lt;sup>‡</sup>Department of Electrical and Computer Engineering, University of California, Santa Barbara, CA 93106-9560 (hespanha@ece.ucsb.edu).

<sup>§</sup>Department of Electrical and Computer Engineering of the Faculty of Science and Technology of the University of Macau, Taipa, Macau; on leave from the Department of Electrical Engineering and Computer Science, Instituto Superior Técnico, ISR, 1046-001 Lisboa, Portugal (cjs@isr.ist.utl.pt).

jump-diffusion [6, sect. 5.3]. Alternatively, SHSs with renewal transitions can be viewed as a generalization of a Markov jump linear system (MJLS) [8], in which the length of time that the system stays in each mode follows an exponential distribution, or as a generalization of an impulsive renewal system [10], in which there is only one discrete mode and one reset map.

Our contribution concerns the analyses of the transitory and asymptotic behavior of the statistical moments of an SHS with renewal transitions and is summarized in two main results, made possible by the special structure of SHSs with renewal transitions. In the first main result, we provide expressions for the moments of the SHS at a given time t in terms of a set of Volterra equations. Since [8, Chap. 2] presents a method to obtain the statistical moments of an MJLS, based on the solution of linear differential equations, that does not appear to generalize to SHS with renewal transitions (cf. [8, sect. 8.2]), this first main result shows that for this latter class of systems, one can still compute statistical moments by solving Volterra equations, instead of differential equations. We also discuss how, by computing the moments using these Volterra equations, one can provide information about the probability distribution of the state of the SHS and highlight the advantages of this approach with respect to more general methods available in the literature [4], [5], [6]. The second main result provides a method to obtain the Lyapunov exponents of degree m, which are defined, in accordance with [8, p. 41], as the exponential rate of decrease or increase at which the expected value of the mth power of the norm of the state of the SHS tends to zero or to infinity, respectively. We show that when m is even, Lyapunov exponents can be efficiently determined by finding the zeros of monotonic functions. As a corollary, we provide necessary and sufficient conditions for mean exponential stability, which is defined in terms of the exponential convergence to zero of the expected value of the squared norm of the state of the SHS.

The applicability of our theoretical results is illustrated in a networked control problem with independently distributed intervals between data transmissions of the control signal in a feedback loop, which is a reasonable assumption in networked control systems utilizing CSMA communication protocols [10]. The impulsive renewal systems considered in [10] did not permit us to consider the effect of network induced delays, which is now possible with SHSs with renewal transitions.

A preliminary version of the results presented here appeared in the conference paper [11]. However, in this paper we have been able to reduce significantly the dimension of the Volterra equations needed to compute the statistical moments of the SHS. In fact, the dimension of the Volterra equations in the present paper grows polynomially with m, whereas in [11] it grows exponentially with m. This enables the use of our results to compute statistical moments of much larger order. An additional novelty of the results presented here is that the asymptotic analysis extends both [10] and [11], where the asymptotic analysis is restricted to a special moment, i.e., the squared norm of the state.

The remainder of the paper is organized as follows. SHSs with renewal transitions are defined in section 2. In section 3, we establish and discuss our main results. Section 4 addresses the applicability of the results to a networked control example. In section 5 we draw final conclusions.

Notation and preliminaries. The Kronecker product is denoted by  $\otimes$ . The notation  $x(t_k^-)$  indicates the limit from the left of a function x(t) at the point  $t_k$ . For vectors  $u_1, \ldots, u_m$  in  $\mathbb{R}^n$ , we define  $(u_1, \ldots, u_m) := [u_1^{\mathsf{T}} \ldots u_m^{\mathsf{T}}]^{\mathsf{T}}$ . The spectral radius of a matrix A is denoted by  $r_{\sigma}(A)$ . We denote by  $e_i$  the canonical vector in  $\mathbb{R}^n$ , i.e., the component j of  $e_i \in \mathbb{R}^n$  equals 1 if j = i and 0 otherwise. The indicator function

of a set A is denoted by  $\chi_{x \in A}$ , which equals 1 if  $x \in A$  and 0 otherwise. The number of elements of a finite set C is denoted by |C|.

2. SHS with renewal transitions. A linear SHS with renewal transitions is defined by (i) a linear differential equation

$$\dot{x}(t) = A_{q(t)}x(t), \quad x(0) = x_0, \ q(0) = q_0, \ t_0 = 0,$$

where  $x(t) \in \mathbb{R}^n$  and  $q(t) \in \mathcal{Q} := \{1, \dots, n_q\}$ ; (ii) a family of  $n_\ell$  discrete transition/reset maps

$$(2.2) (q(t_k), x(t_k)) = (\xi_{\ell}(q(t_k^-)), J_{q(t_k^-), \ell}x(t_k^-)), \ell \in \mathcal{L} := \{1, \dots, n_{\ell}\},$$

where  $\xi_{\ell}$  is a map from  $\mathcal{Q}$  to  $\mathcal{Q}$  and the matrix  $J_{q(t_k^-),\ell}$  belongs to a given set  $\{J_{i,\ell} \in \mathbb{R}^{n \times n}, i \in \mathcal{Q}, \ell \in \mathcal{L}\}$ ; and (iii) a family of reset-time measures

Between transition times, denoted by  $t_k$ , the discrete mode q remains constant whereas the continuous state x flows according to (2.1). At transition times, the continuous state and discrete mode of the SHS are reset according to (2.2). The intervals between transition times are independent random variables determined by the reset-time measures (2.3) as follows. A reset-time measure can be either a probability measure or identically zero. In the former case,  $\mu_{i,\ell}$  is the probability measure of the random time that transition  $\ell \in \mathcal{L}$  takes to trigger in the state  $q(t) = i \in \mathcal{Q}$ . The next transition time is determined by the minimum of the triggering times of the transitions associated with state  $q(t) = i \in \mathcal{Q}$ . When  $\mu_{i,\ell}([0,s)) = 0 \ \forall_{s \geq 0}$ , the transition  $\ell$  does not trigger in state  $i \in \mathcal{Q}$ , which allows for some reset maps not to be active in some states.

The construction of a sample path of the SHS with renewal transitions can then be described as follows:

- 1. Set k = 0,  $t_0 = 0$ ,  $(q(t_k), x(t_k)) = (q_0, x_0)$ .
- 2. For every  $j \in \mathcal{L}$ , obtain  $\bar{h}_k^j$  as a realization of a random variable distributed according to  $\mu_{q(t_k),j}$  if  $\mu_{q(t_k),j}$  is not identically zero, and set  $\bar{h}_k^j = \infty$  otherwise.
- 3. Take  $h_k = \min\{\bar{h}_k^j, j \in \mathcal{L}\}$  and set the next transition time to  $t_{k+1} = t_k + h_k$ . The state of the SHS in the interval  $t \in [t_k, t_{k+1})$  is given by  $(q(t), x(t)) = (q(t_k), e^{A_{q(t_k)}(t-t_k)}x(t_k))$ .
- 4. If  $t_{k+1} < \infty$  (otherwise stop), let  $l_k$  denote the index of the transition that achieves the minimum in step 3, i.e.,  $l_k = j$ :  $h_k = \bar{h}_k^j$ , and update the state according to  $(q(t_{k+1}), x(t_{k+1})) = (\xi_{l_k}(q(t_{k+1}^-)), J_{q(t_{k+1}^-), l_k}x(t_{k+1}^-))$ . Set k = k+1 and repeat the construction from step 2.

We assume that each transition probability measure  $\mu_{i,\ell}$  can be decomposed as  $\mu_{i,\ell} = \mu_{i,\ell}^c + \mu_{i,\ell}^d$  with  $\mu_{i,\ell}^c([0,t)) = \int_0^t f_{i,\ell}(s) ds$  for some density function  $f_{i,\ell}(s) \geq 0$ , and  $\mu_{i,\ell}^d$  is a discrete measure that captures possible point masses  $\{b_{i,\ell}^r, r \geq 1\}$  such that  $\mu_{i,\ell}^d(\{b_{i,\ell}^r\}) = w_{i,\ell}^r$ . The integral with respect to the measure  $\mu_{i,\ell}$  is then defined as

(2.4) 
$$\int_0^t W(s)\mu_{i,\ell}(ds) = \int_0^t W(s)f_{i,\ell}(s)ds + \sum_{r:b_{i,\ell}^r \in [0,t]} w_{i,\ell}^r W(b_{i,\ell}^r).$$

When the discrete measures  $\mu_{i,\ell}^d$  are nonzero measures, different transitions may be triggered at the same time with nonzero probability, leading to an ambiguity in choosing the next state. To avoid ambiguity, we assume the following.

Assumptions 2.1.

- (i) For every  $i \in \mathcal{Q}$ , the reset-time measures  $\mu_{i,\ell}$ ,  $\ell \in \mathcal{L}$ , for mode i have a finite number of point masses.
- (ii) For every  $i \in \mathcal{Q}$ , no two measures  $\mu_{i,\ell_1}$ ,  $\mu_{i,\ell_2}$  for mode i have common point masses, i.e.,  $b_{i,l_1}^j \neq b_{i,l_2}^{\kappa}$  for  $l_1 \neq l_2, l_1 \in \mathcal{L}$ ,  $l_2 \in \mathcal{L} \ \forall_{j \geq 1, \kappa \geq 1}$ . Due to Assumption 2.1(ii), there is zero probability that in step 3 of the construc-

Due to Assumption 2.1(ii), there is zero probability that in step 3 of the construction of a sample path of the SHS the minimum is achieved by two or more indices j.

**3.** Main results. Consider a general mth degree uncentered moment of the state of the SHS with renewal transitions, i.e.,

(3.1) 
$$\mathbb{E}[x_1(t)^{i_1}x_2(t)^{i_2}\dots x_n(t)^{i_n}], \quad \sum_{j=1}^n i_j = m, \ i_j \ge 0.$$

We provide a method to compute (3.1) in subsection 3.1, we obtain the Lyapunov exponents of special moments of the SHS in subsection 3.2, and we discuss how by computing the moments one can reconstruct probability distributions in subsection 3.3. The proofs of the main results are deferred to subsection 3.4.

**3.1.** Moment computation. It is easy to see that there are

(3.2) 
$$p := \frac{(m+n-1)!}{m!(n-1)!}$$

different monomials of degree m and hence p different moments (3.1) of degree m. Let

(3.3) 
$$\{\rho(\kappa) = [i_1(\kappa), \dots i_n(\kappa)], 1 \le \kappa \le p\}$$

be an enumeration of the indices  $i_1 \dots, i_n$  uniquely characterizing such monomials, e.g., for m = n = 2, one such enumeration is  $\rho(1) = [1 \ 1], \ \rho(2) = [2 \ 0], \ \rho(3) = [0 \ 2].$  Then, we use the notation

$$x^{[\kappa]} := x_1^{i_1} \dots x_n^{i_n} \text{ for } \kappa : \rho(\kappa) = [i_1, \dots, i_n], \ 1 \le \kappa \le p,$$

and define the map  $\Gamma^m : \mathbb{R}^{n \times n} \mapsto \mathbb{R}^{p \times p}$  as

$$(3.4) \Gamma^m(A) = B,$$

where  $B = [B_{\iota,\kappa}]$  is uniquely determined by

(3.5) 
$$(Ax)^{[\iota]} = \sum_{\kappa=1}^{p} B_{\iota\kappa} x^{[\kappa]}, 1 \le \iota, \, \kappa \le p.$$

Note that indeed the left-hand side of (3.5) is a linear combination of monomials of degree m with unique coeficients  $B_{\iota,\kappa}$ . The following theorem provides a method to compute any moment of the state of the SHS. Let

(3.6) 
$$\gamma_{\kappa} := \frac{m!}{(i_1! \dots i_n!)} \text{ for } \kappa : \rho(\kappa) = [i_1, \dots, i_n], \sum_{i=1}^n i_i = m,$$

and define the operator

$$\Theta(u^{\kappa}(t)) := (\Theta_1(u^{\kappa}(t)), \dots, \Theta_{n_q}(u^{\kappa}(t))),$$

where  $u^{\kappa}(t) := (u^{\kappa,1}(t), \dots, u^{\kappa,n_q}(t))$  and each  $\Theta_i$  is a convolution operator defined by

(3.7) 
$$\Theta_{i}(u^{\kappa}(t)) := \sum_{\ell=1}^{n_{\ell}} \int_{0}^{t} \Gamma^{m}(e^{A_{i}^{\mathsf{T}}s} J_{i,\ell}^{\mathsf{T}}) u^{\kappa,\xi_{\ell}(i)}(t-s) \frac{r_{i}(s)}{r_{i,\ell}(s)} \mu_{i,\ell}(ds),$$

where  $i \in \mathcal{Q}$ ,  $r_i(s) := \prod_{\ell=1}^{n_\ell} r_{i,\ell}(s)$ ,  $r_{i,\ell}(s) := \mu_{i,\ell}((s,\infty])$ .

THEOREM 3.1. A moment of degree m (3.1) of the SHS (2.1)-(2.3), indexed by  $\kappa : \rho(\kappa) = [i_1 \dots i_n]$ , can be computed as

(3.8) 
$$\mathbb{E}[x(t)^{[\kappa]}] = \sum_{\iota=1}^{p} x_0^{[\iota]} u_{\iota}^{\kappa, q_0}(t),$$

where the  $u^{\kappa} = (u^{\kappa,1}, \dots, u^{\kappa,n_q}), u^{\kappa,i} \in \mathbb{R}^p$ , are the unique solution to the Volterra equation

(3.9) 
$$u^{\kappa}(t) = \Theta(u^{\kappa}(t)) + h^{\kappa}(t), \quad t \ge 0,$$

where  $h^{\kappa}(t) := (h^{\kappa,1}(t), \dots, h^{\kappa,n_q}(t)),$ 

(3.10) 
$$h^{\kappa,i}(t) = \Gamma^m(e^{A_i^{\mathsf{T}}t}) e_{\kappa} \frac{r_i(t)}{\gamma_{\kappa}}, \ i \in \mathcal{Q},$$

where  $e_{\kappa}$  is the canonical vector in  $\mathbb{R}^p$  with nonzero component  $\kappa$ .

An explicit solution to (3.9) takes the form

(3.11) 
$$u^{\kappa}(t) = \sum_{k=0}^{\infty} \Theta^{k}(h^{\kappa}(t)),$$

where  $\Theta^k(h^{\kappa}(t))$  denotes composition, e.g.,  $\Theta^2(h^{\kappa}(t)) = \Theta(\Theta(h^{\kappa}(t)))$  and  $\Theta^0(h^{\kappa}(t)) = h^{\kappa}(t)$ . However, in practice, a numerical method is generally preferable to solve the Volterra equation (3.9). We discuss next the computation complexity of one such method, comparing it also with the complexity of previous methods.

**3.1.1. Numerical computation.** One numerical method to solve the Volterra equation (3.9) consists in choosing a set of integration nodes  $\{a_l \in [0, t]\}$  and obtaining  $u^{\kappa}(a_l)$  by iteratively replacing the integrals in (3.9) by quadrature formulas at the nodes  $\{a_j : a_j \in [0, a_l]\}$  (cf. [12]). Suppose that  $a_l = l\bar{h}, l \in \{0, \dots, \bar{L}\}$  are nodes in an interval of interest  $t \in [0, \bar{L}\bar{h}]$ , where  $\bar{L}$  is the number of nodes and  $\bar{h}$  is the spacing between two consecutive nodes. Assuming, for simplicity, that the distributions  $\mu_{i,\ell}$  have no atom points, we have (3.12)

$$\bar{u}^{\kappa,i}(l\bar{h}) = \sum_{\ell=1}^{n_\ell} \sum_{k=0}^l \bar{c}_{l,k} \Gamma^m(e^{A_i^\intercal(k\bar{h})} J_{i,\ell}^\intercal) \bar{u}^{\kappa,\xi_\ell(i)} ((l-k)\bar{h}) \frac{r_i(k\bar{h})}{r_{i,\ell}(k\bar{h})} f_{i,\ell}(k\bar{h}) + \Gamma^m(e^{A_i^\intercal l\bar{h}}) \mathbf{e}_\kappa \frac{r_i(l\bar{h})}{\gamma_\kappa} \mathbf{e}_\kappa \frac{r_i(l\bar{h})}{r_i(l\bar{h})} \mathbf{e}_\kappa \frac{r_i(l\bar{h})}$$

for  $i \in \mathcal{Q}$ , where  $\bar{c}_{l,k}$  denotes the weights of the quadrature formula and  $\bar{u}^{\kappa,i}(l\bar{h})$  approximates  $u^{\kappa,i}(l\bar{h})$ . Considering possible atom points of  $\mu_{i,\ell}$  would add additional additive terms on the right-hand side of (3.12) taking into account the atom points in

the interval  $b_{i,\ell}^r \in [0, l\bar{h}]$  which would not significantly impact in our discussion. The number of matrix multiplications  $\Gamma^m(e^{A_i^\intercal(k\bar{h})}J_{i,\ell}^\intercal)u^{\kappa,\xi_\ell(i)}((l-k)\bar{h})$  in (3.12) is given by

(3.13) 
$$\sum_{\ell=1}^{n_{\ell}} |\{k \in \{0, \dots, l\} : f_{i,\ell}(k\bar{h}) > 0 \text{ and } r_i(k\bar{h}) > 0\}|$$

which is bounded by  $ln_{\ell}$ . Each of these matrix multiplications requires in general  $p^2$  scalar multiplications. Moreover, if  $f_{i,\ell}(0) \neq$  for some  $i \in \mathcal{Q}$  and some  $\ell \in \mathcal{L}$ , a linear system of at most  $n_q p$  equations must in general be solved to obtain  $u^{\kappa}(l\bar{h})$  from the set of equations (3.12),  $i \in \mathcal{Q}$ . Thus, the number of total multiplications needed to iterate (3.12) from l = 0 to  $l = \bar{L} - 1$  is bounded by

(3.14) 
$$p^2 n_q n_\ell \frac{\bar{L}(\bar{L}+1)}{2} + \bar{L}G(n_q p),$$

where  $G(a) = \frac{a^3}{3} + a^2 - \frac{a}{3}$  is the number of multiplications required by the standard Gaussian method to solve a system with  $a \in \mathbb{N}$  variables. Note that the matrices

(3.15) 
$$\Gamma^m(e^{A_i^{\mathsf{T}}(l\bar{h})}J_{i,\ell}^{\mathsf{T}}), \quad \Gamma^m(e^{A_i^{\mathsf{T}}lh}), \quad l \in \{0,\dots,\bar{L}\},$$

can be computed and stored in memory before iterating (3.12). The computation complexity of computing these at most  $2(\bar{L}+1)$  matrices by symbolically manipulating monomials is no larger than the computational complexity of iterating (3.12), which is bounded by (3.14). In fact, suppose that we wish to compute  $C = \Gamma^m(D)$  for  $D \in \mathbb{R}^{n \times n}$  with row vectors  $d^i \in \mathbb{R}^{1 \times n}$ ,  $i \in \{1, \ldots, n\}$ . To compute a row  $\iota$  of C such that  $\rho(\iota) = [i_1 \ldots i_n]$  we recursively compute  $\zeta(y) = (d^1 y)^{i_1} \times (d^2 y)^{i_2} \ldots (d^n y)^{i_n}$ ,  $y \in \mathbb{R}^n$ , according to  $\zeta(y) = \xi_m(y)$ ,

(3.16) 
$$\xi_{k+1}(y) = (d^{r(k)}y)\xi_k(y), \quad \xi_1(y) = 1, \quad k = 1, \dots, m-1,$$

where r(k) = j if  $\sum_{s=0}^{j-1} i_s \le k \le \sum_{s=0}^{j} i_s$ ,  $j \in \{1, \dots, n\}$ ,  $k \in \{1, \dots, m\}$ . Since  $\xi_k(x)$  is a sum of monomials with dimension k and the number of such monomials can be obtained from (3.2), we have that (3.16) requires  $n \times \sum_{k=1}^{m-1} \frac{(k+n-1)!}{k!(n-1)!}$  multiplications to compute each row of C, which is bounded by nmp. Thus, the total number of multiplications to compute C is bounded by  $nmp^2$  and the number of multiplications needed to compute (3.15) is bounded by

$$2(\bar{L}+1)nmp^2.$$

A result similar to Theorem 3.1 was given in a preliminary version of the present paper [11]. However, there exists redundancy both in the state of the Volterra equation proposed in [11] and in the state of the differential equation proposed in [8], which is eliminated by Theorem 3.1. In fact, given an SHS with state-space dimension n and  $n_q$  discrete modes, in [11] one needs to solve a Volterra equation with  $n_q \times n^m$  unknown functions to obtain a moment of degree m. This is also the dimension of the linear differential equation presented in [8] to compute the moments in the special case of exponential reset-time distributions. The Volterra equation (3.9) has dimension  $n_q \times p = n_q \times \frac{(m+n-1)\times \cdots \times (m+1)}{(n-1)!}$ , i.e., the number of unknown functions in (3.9) grows polynomially with m, instead of exponentially.

From a practical point of view our approach (3.12) is feasible in a current personal computer if p is in the order of tens of thousands. For example, in section 4.1 we shall

consider n = 4, for which we can compute up to about m = 10 moments of the state in which case p = 10296. Note that with previous approaches [11], [8] this would require us to manipulate Volterra equations with dimension two orders of magnitude larger  $4^{10} = 1048576$ , which clearly would not be feasible in a current personal computer.

**3.2. Lyapunov exponents.** The following definition of Lyapunov exponent of degree m is adapted from [8, p. 41]. Recall the definition of  $L_p$  norm, for  $p \in \mathbb{N}$ , of a vector x(t)

(3.17) 
$$||x(t)||_p := \left(\sum_{i=1}^n |x_i(t)|^p\right)^{1/p}, \quad p \ge 1.$$

DEFINITION 3.2. Suppose that  $\mathbb{E}[\|x(t)\|_m^m] \neq 0 \ \forall t_{\geq 0}$  and that for every  $x_0 \neq 0$  the following limit exists:

(3.18) 
$$\lambda_L^m(x_0) := \lim_{t \to \infty} \frac{1}{t} \log \mathbb{E}[\|x(t)\|_m^m].$$

Then the Lyapunov exponent  $\lambda_L^m$  of degree m for the SHS (2.1)-(2.3) is defined as

$$\lambda_L^m := \sup_{x_0 \in \mathbb{R}^n} \lambda_L^m(x_0).$$

Moreover, if  $\exists_{b>0}$ :  $\mathbb{E}[\|x(t)\|_m^m] = 0 \ \forall t > b$ , then  $\lambda_L^m := -\infty$ . Let

$$\hat{\Theta}^m(z) := \begin{bmatrix} \hat{\Theta}^m_{1,1}(z) & \dots & \hat{\Theta}^m_{1,n_q}(z) \\ \vdots & \ddots & \vdots \\ \hat{\Theta}^m_{n_q,1}(z) & \dots & \hat{\Theta}^m_{n_q,n_q}(z) \end{bmatrix},$$

where

$$\hat{\Theta}^m_{i,j}(z) := \sum_{\ell: \mathcal{E}_{\ell}(i) = j} \int_0^\infty \Gamma^m(e^{A^\intercal s} J_{i,\ell}^\intercal) e^{-zs} \frac{r_i(s)}{r_{i,\ell}(s)} \mu_{i,\ell}(ds).$$

For  $a \in \mathbb{R}$ , we say that  $\hat{\Theta}^m(a)$  converges absolutely if for every  $i, j \in \mathcal{Q}$ 

$$\sum_{\ell: \xi_\ell(i)=j} \int_0^\infty \|\Gamma^m(e^{A^\intercal s}J_{i,\ell}^\intercal)\|e^{-as}\frac{r_i(s)}{r_{i,\ell}(s)}|\mu_{i,\ell}|(ds) < \infty,$$

where for a matrix X,  $||X||^2 := \operatorname{tr}(X^{\intercal}X)$  and  $|\mu_{i,\ell}|(ds)$  is the total variation norm of the measure  $\mu_{i,\ell}$  (cf. [13]).

THEOREM 3.3. Suppose that all the distributions  $\mu_{i,\ell}$  have finite support and let  $b := \inf\{a : \hat{\Theta}^m(a) \text{ converges absolutely}\}$ . Then, if m is even, the spectral radius  $r_{\sigma}(\hat{\Theta}^m(a))$  of  $\hat{\Theta}^m(a)$  is a monotone nonincreasing function of a for a > b and the Lyapunov exponent  $\lambda_L^m$  for the SHS (2.1)-(2.3) is the (unique) real root a to the equation

$$r_{\sigma}(\hat{\Theta}^m(a)) = 1$$

<sup>&</sup>lt;sup>1</sup>The definition in [8, p. 41] does not specify which norm to take for ||x(t)|| (we choose to take (3.17)). Note that in fact (3.18) does not change with the choice of the norm, due to the equivalence of norms in finite-dimensional vector spaces.

if this equation has a real root and  $-\infty$  otherwise.

As we shall see in section 3.4, the proof of Theorem 3.3 relies on establishing that  $\hat{\Theta}^m(z)$  for real z is a positive operator (cf. [14]). This holds only if m is even and this is the main reason for restricting Theorem (3.3) to this case. We do not address in the present paper the case in which m is odd.

We say that the SHS with renewal transitions is mean exponentially stable if there exist constants c > 0 and  $\alpha > 0$  such that for every  $(x_0, q_0)$ ,

$$\mathbb{E}[x(t)^{\mathsf{T}}x(t)] \le ce^{-\alpha t} x_0^{\mathsf{T}} x_0 \ \forall_{t>0},$$

Theorem 3.3 allows us to establish necessary and sufficient conditions for the stability of an SHS with renewal transitions defined as follows.

COROLLARY 3.4. Suppose that all the distributions  $\mu_{i,\ell}$  have bounded support. Then the SHS with renewal transitions is mean exponentially stable if and only if

(3.19) 
$$r_{\sigma}(\hat{\Theta}^2(0)) < 1.$$

3.3. Probability distribution. Let  $\mu$  be the probability distribution of the state of the SHS at time t, i.e.,  $\int_E \mu(ds) := \operatorname{Prob}[x(t) \in E \subseteq \mathbb{R}^n]$ . In this subsection we address the problem of obtaining information on  $\mu$  based on its moments (3.1). It is well known that when the moments are finite and known, the measure  $\mu$  can be uniquely determined [15], while in the case where only a finite number of moments are known, this reconstruction is naturally not unique. In this latter case, an elegant procedure to obtain an explicit expression for an approximating distribution to  $\mu$  can be found in the literature. (We follow closely [15]; see also [16], [17].) Assume, for simplicity, that  $\mu$  has bounded support, which without loss of generality (by proper scaling) can be assumed to be contained in the interval  $D := [0,1]^n = [0,1] \times \cdots \times [0,1]$ . (See [15] for the case of unbounded support.) Given a continuous function  $f: D \mapsto \mathbb{R}$  define the higher dimensional Bernstein polynomials as

$$B_r(f)(y) = \sum_{k:0 < k_1 < r_1} f\left(\frac{k_1}{r_1}, \dots, \frac{k_n}{r_n}\right) P_{r,k}(y), \quad y \in \mathbb{R}^n,$$

where  $r = (r_1, \ldots, r_n)$  is a vector of integers, and  $P_{r,k}(y) := \prod_{i=1}^n {r_i \choose k_i} y_i^{k_i} (1 - y_i)^{r_i - k_i}$ . It is shown in [15] that  $B_r(f)(y)$  converges uniformly to f(y) as  $r_i \to \infty \ \forall_{1 \le i \le n}$ . Thus, if we define  $I_{r,k} := \int_D P_{r,k}(y) \mu(ds)$ , we have that

$$(3.20) \sum_{0 \le k_i \le r_i} f\left(\frac{k_1}{r_1}, \dots, \frac{k_n}{r_n}\right) I_{r,k}$$

is an approximation to  $\mathbb{E}[f(x(t))] = \int_E f(s)\mu(ds)$  that converges to  $\mathbb{E}[f(x(t))]$  as  $r_i \to \infty \ \forall_{1 \le i \le n}$ . This means that the measure  $\mu$  can be approximated by the discrete measure  $\mu_r$  defined by

(3.21) 
$$\sum_{0 \le k_i \le r_i} I_{r,k} \delta\left(\left[\frac{k_1}{r_1}, \dots, \frac{k_n}{r_n}\right]\right),$$

where  $\delta(y)$  denotes the Dirac measure at  $y \in \mathbb{R}^n$ . Since Bernstein polynomials are linear combinations of monomials, we can compute the  $I_{r,k}$  through a linear combination of moments, which can be computed as proposed in section 3.1.

It is important to mention that the problem of recovering a probability distribution from its moments is typically a numerically ill-conditioned problem, and small numerical errors in the moments may lead to large errors in the coefficients  $I_{r,k}$  that define the approximate discrete measure in (3.21). Thus, it is especially important to have a computationally efficient method to compute the moments, as provided by Theorem 3.1, so that one can obtain a good numerical precision in this procedure of obtaining an approximating probability distribution.

One way to interpret the procedure just described is that each integral  $I_{r,n}$  approximates the probability that x lies in a small neighborhood of  $(\frac{k_1}{r_1}, \ldots, \frac{k_r}{r_n})$ . Instead of obtaining an approximation, one can use similar ideas to obtain an upper bound on the probability that x belongs to a given region. That is, suppose, for example, that we wish to determine a bound on the probability that a random variable z = g(x(t)) is greater than a given value. Assume that z takes values in a compact set, which without loss of generality is assumed to be [0,1]. Then we can choose a polynomial  $p_n: [0,1] \to \mathbb{R}_{\geq 0}$  such that  $p_n(w) \geq \chi_{[a,1]}(w), w \in [0,1], a < 1$ , and obtain the bound

(3.22) 
$$\operatorname{Prob}[z \ge a] \le \mathbb{E}[p_n(z)],$$

where the right-hand side of (3.22) can be computed using Theorem 3.1, when g(x(t)) is a sum of monomial functions of the state. Note that when  $p_n(w) = \frac{w}{a}$  and  $p_n(w) = \frac{w^2}{a^2}$ , (3.22) corresponds to the Markov and Chebychev inequalities, respectively. This procedure is used in the networked control example of section 4.

Alternative approaches to estimate the probability distribution of the state of the SHS include those based on the Focker-Plank equation and Dynkin's formula. The Focker-Plank equation of the SHS with renewal transitions can be obtained by specializing the expressions provided in [5], [6, sect. 5.3] to this class of systems, and can be shown to be integro-partial differential equations. Besides the numerical difficulty and computational burden associated with solving these equations, the derivation of these Focker-Plank equations requires the map  $y \mapsto Jy$ ,  $y \in \mathbb{R}^n$ , to be invertible (equivalent to matrix J invertible). The approach based on the Dynkin's formula can be found in [4, Chap. 3, sect. 32.2], where a numerical method is provided to compute the expected value of a given function  $\mathbb{E}[f(x(t))]$  and a fortiori estimating the probability distribution from the relation  $\mathbb{E}[\chi_E(x(t))] = \operatorname{Prob}[x(t) \in E]$  for given  $E \subseteq D$ . Note that the approximation (3.20), which can be obtained with the methods derived in the present paper, provides an alternative to the method in [4, Chap. 3, sect. 32.2] to approximate the expected value of continuous functions of the state. Although we omit the derivations here, it is possible to prove that when specialized to SHSs with renewal transitions and to the case where f is a monomial, the recursive method provided in [4, Chap. 3, sect. 32.2] is equivalent to providing an approximation to (3.8) at each iteration  $n_r$  of the recursive algorithm, taking the form

(3.23) 
$$\sum_{i=1}^{p} x_0^{[\iota]} v_{\iota}^{\kappa, q_0}(t)$$

with

$$(3.24) v^{\kappa}(t) = \sum_{k=0}^{n_{\rm r}} \Theta^k(h^{\kappa}(t)).$$

This last equation converges to (3.11), and hence (3.23) converges to (3.8), but this is clearly an inefficient method for obtaining the solution to the Volterra equation (3.9)

when compared to the method described in subsection 3.1 (cf. [12]). Thus, by exploiting linearity and the special structure of the SHS with renewal transitions, our approach provides an insight that allows us to compute moments more efficiently, than when seeing the SHS with renewal transitions as a piecewise deterministic process and specializing the approach proposed in [4, Chap. 3, sect. 32.2].

**3.4. Proofs of Theorems 3.1 and 3.3.** We start by establishing four preliminary facts stated in the form of propositions.

Let  $A^{(m)}$ , m > 0, denote the mth-fold Kronecker product of a matrix or a vector A with itself, i.e.,

$$A^{(m)} := \underbrace{A \otimes A \cdots \otimes A}_{m},$$

and recall that

$$(3.25) (A \otimes B)(C \otimes D) = (AC) \otimes (BD)$$

(cf. [18]).

Proposition 3.5. The following holds:

(3.26) 
$$\mathbb{E}[x_1(t)^{i_1}x_2(t)^{i_2}\dots x_n(t)^{i_n}] = \mathbb{E}[(x(t)^{\mathsf{T}})^{(m)}c_{\kappa}],$$

where

$$(3.27) c_{\kappa} := \mathbf{e}_{1}^{(i_{1})} \otimes \cdots \otimes \mathbf{e}_{n}^{(i_{n})} \text{ for } \rho(\kappa) = [i_{1}, \dots, i_{n}].$$

*Proof.* The proof follows directly by using (3.25).

Let  $\mathcal{T}(m,n)$  be the set of symmetric tensors, i.e., multilinear functions R on the m-fold  $\mathbb{R}^n \times \cdots \times \mathbb{R}^n$  (cf. [19, Chap. 4]) such that

$$R(w_1, w_2, \dots, w_m) = R(w_{\sigma(1)}, w_{\sigma(2)}, \dots, w_{\sigma(m)})$$

for every  $w_i \in \mathbb{R}^n, 1 \leq i \leq m$ , and every one-to-one permutation of indices  $\sigma$ :  $\{1, 2, \ldots, m\} \mapsto \{1, 2, \ldots, m\}$ . We note that there is a natural identification between monomials of degree m in  $\mathbb{R}^n$  and  $\mathcal{T}(m, n)$ . In fact, to every m-degree monomial  $x^{[\kappa]} = x_1^{i_1} x_2^{i_2} \dots x_n^{i_n}$ , indexed by  $\kappa$ , as in (3.3), we can associate an element  $R^{\kappa}$  of the following orthogonal basis of symmetric tensors defined by

$$R^{\kappa}(w_1, w_2, \dots, w_m) := b^{\kappa} w_1 \otimes w_2 \otimes \dots \otimes w_m,$$

where

$$(3.28) b^{\kappa} := \sum_{j \in \mathcal{I}_{n}} (\mathbf{e}_{j_{1}}^{\mathsf{T}}) \otimes (\mathbf{e}_{j_{2}}^{\mathsf{T}}) \otimes \ldots \otimes (\mathbf{e}_{j_{m}}^{\mathsf{T}}), \ 1 \leq \kappa \leq p,$$

and the vector j belongs a set of  $\gamma_{\kappa}$  permutations of indices defined as

(3.29) 
$$\mathcal{J}_{\kappa} := \{ j = (j_1, \dots, j_m) : x_{j_1} x_{j_2} \dots x_{j_m} = x^{[\kappa]} \ \forall_{x \in \mathbb{R}^n} \},$$

where  $\kappa : \rho(\kappa) = [i_1, \dots, i_n].$ 

Proposition 3.6. The following holds:

$$(3.30) (x(t)^{(m)})^{\mathsf{T}} c_{\kappa} = (x(t)^{(m)})^{\mathsf{T}} d_{\kappa},$$

where  $d_{\kappa} := \frac{(b^{\kappa})^{\intercal}}{\gamma_{\kappa}}$  and

(3.31) 
$$x(t)^{(m)} = \sum_{\kappa=1}^{p} x^{[\kappa]} (b^{\kappa})^{\mathsf{T}},$$

where p is given by (3.2).

*Proof.* The proof follows directly from definitions (3.28) and (3.29).

Let  $T_i(t)$  denote the transition matrix of the SHS starting at the discrete mode  $q_0 = i$ , i.e.,  $x(t) = T_{q_0}(t)x_0$ , where

$$T_i(t) = e^{A_{q(t)}(t-t_r)} \dots J_{\xi_{l_0}(i),l_1} e^{A_{\xi_{l_0}(i)}(t_1-t_0)} J_{i,l_0} e^{A_i t_1},$$

 $r = \max\{k \in \mathbb{Z}_{\geq 0} : t_k \leq t\}.$ 

PROPOSITION 3.7. The following holds:

(3.32) 
$$\mathbb{E}[(x(t)^{\mathsf{T}})^{(m)}c_{\kappa}] = (x_0^{\mathsf{T}})^{(m)}w^{\kappa,q_0}(t),$$

where the  $w^{\kappa,i}(t)$ ,  $i \in \mathcal{Q}$ , are defined as

(3.33) 
$$w^{\kappa,i}(t) := \mathbb{E}[(T_i(t)^{\mathsf{T}})^{(m)}d_{\kappa}], \quad i \in \mathcal{Q}.$$

*Proof.* Using (3.25), (3.26), and (3.30) we obtain

$$\mathbb{E}[(x(t)^{\mathsf{T}})^{(m)}c_{\kappa}] = \mathbb{E}[(x(t)^{\mathsf{T}})^{(m)}d_{\kappa}] = (x_0^{\mathsf{T}})^{(m)}\mathbb{E}[(T_{q_0}(t)^{\mathsf{T}})^{(m)}d_{\kappa}] = (x_0^{\mathsf{T}})^{(m)}w^{\kappa,q_0}(t).$$

For a matrix  $A \in \mathbb{R}^{n \times n}$ , we define a map  $T_A : \mathbb{R}^p \mapsto \mathbb{R}^p, y \mapsto v = T_A(y)$  by

(3.34) 
$$\sum_{\kappa_2=1}^p v_{\kappa_2}(b^{\kappa_2})^{\intercal} = A^{(m)} \sum_{\kappa_2=1}^p y_{\kappa_2}(b^{\kappa_2})^{\intercal}.$$

PROPOSITION 3.8. For a matrix  $A \in \mathbb{R}^{n \times n}$ , the map  $T_A : \mathbb{R}^p \mapsto \mathbb{R}^p, y \mapsto v = T_A(y)$  defined by (3.34) can be described by

$$(3.35) v = \Gamma^m(A)y,$$

where  $\Gamma^m$  is given by (3.4)

*Proof.* If we denote the elements and row of A by  $A_{ij}$  and  $a^i$ , respectively, i.e.,  $A = [A_{ij}] = [(a^1)^{\intercal} (a^2)^{\intercal} \dots (a^n)^{\intercal}]^{\intercal}$ , then using (3.25), (3.28), (3.29), and (3.31), we have that for  $\kappa_1 : \rho(\kappa_1) = [i_1 i_2, \dots i_n]$ ,

$$(Ax)^{[\kappa_1]} = (a^1)^{(i_1)} \otimes (a^2)^{(i_2)} \otimes \ldots \otimes (a^n)^{(i_n)} x^{(m)}$$

$$= \frac{1}{\gamma_{\kappa_1}} \sum_{j \in \mathcal{J}_{\kappa_1}} a^{j_1} \otimes a^{j_2} \otimes \ldots \otimes a^{j_m} x^{(m)}$$

$$= \frac{1}{\gamma_{\kappa_1}} \sum_{j \in \mathcal{J}_{\kappa_1}} a^{j_1} \otimes a^{j_2} \otimes \ldots \otimes a^{j_m} \sum_{\kappa_2 = 1}^p (b^{\kappa_2})^{\mathsf{T}} x^{[\kappa_2]}$$

$$= \sum_{\kappa_2 = 1}^p B_{\kappa_1 \kappa_2} x^{[\kappa_2]},$$

where

(3.36) 
$$B_{\kappa_1 \kappa_2} = \frac{1}{\gamma_{\kappa_1}} \sum_{j \in \mathcal{J}_{\kappa_1}} \sum_{i \in \mathcal{J}_{\kappa_2}} A_{j_1 i_1} A_{j_2 i_2} \dots A_{j_n i_n}.$$

Applying the tensors  $b^{\kappa_1}, 1 \leq \kappa_1 \leq p$ , on both sides of (3.34) one also obtains that the linear map (3.34) is described by (3.36) in the basis  $(b^j)^{\mathsf{T}}, 1 \leq j \leq p$ .  $\square$ Proof of Theorem 3.1. We start by showing that

(3.37) 
$$w^{\kappa}(t) := (w^{\kappa,1}(t), \dots, w^{\kappa,n_q}(t)),$$

where the  $w^{\kappa,i}$  are defined in (3.33), satisfies a Volterra equation.

Consider an initial condition  $q_0 = i, i \in \mathcal{Q}$ , and a given time t and partition the probability space into the events  $[t_1 \leq t] \cup [t_1 > t]$ . We can further partition  $[t_1 \leq t]$  into  $[t_1 \leq t] = \bigcup_{\ell=1}^{n_\ell} B_\ell(t) \cup B_0(t)$ , where  $B_0(t)$  is the event of two transitions triggering at the same time in the interval [0,t], which has probability zero due to Assumption 2.1, and  $B_\ell(t)$  is the event of the transition  $\ell$  being the first to trigger in the interval [0,t], i.e.,  $B_\ell(t) = [\min\{\bar{h}_0^j, j \in \mathcal{L}\} = \bar{h}_0^\ell = t_1 \leq t] \wedge [\bar{h}_0^j > \bar{h}_0^\ell, \ell \neq j]$ , where  $\ell$  denotes intersection. Notice that since the initial state is  $\ell$ 0 =  $\ell$ 1,  $\ell$ 2 is distributed according to  $\ell$ 2, for a given  $\ell$ 3 for which  $\ell$ 4, is the nonzero measure. When transition  $\ell$ 3 does not trigger in state  $\ell$ 6 for which  $\ell$ 7, is empty, in which case  $\ell$ 6, is the zero measure. Using this partition we can write

$$(3.38) \qquad \mathbb{E}[(T_i(t)^{\intercal})^{(m)}d_{\kappa}] = \mathbb{E}[(T_i(t)^{\intercal})^{(m)}d_{\kappa}\chi_{[t_1>t]}] + \sum_{\ell=1}^{n_{\ell}} \mathbb{E}[(T_i(t)^{\intercal})^{(m)}d_{\kappa}\chi_{B_{\ell}(t)}],$$

where we denote by  $\chi_{x \in A}$  the characteristic function of a set A, i.e.,  $\chi_{x \in A}$  equals 1 if  $x \in A$  and 0 otherwise. The first term on the right-hand side of (3.38) is given by

(3.39) 
$$\mathbb{E}[(T_i(t)^\intercal)^{(m)}d_\kappa\chi_{[t_1>t]}] = (e^{A_i^\intercal t})^{(m)}d_\kappa \ \mathbb{E}[\chi_{[t_1>t]}]$$
$$= (e^{A_i^\intercal t})^{(m)}d_\kappa \ r_i(t),$$
$$= (e^{A_i^\intercal t})^{(m)}\frac{(b^\kappa)^\intercal}{\gamma_\kappa} \ r_i(t),$$

where we used the fact that  $\mathbb{E}[\chi_{[t_1>t]}] = \operatorname{Prob}([t_1>t]) = \prod_{j=1}^{n_\ell} \operatorname{Prob}[h_0^j>t] = \prod_{j=1}^{n_\ell} r_{i,j}(t) = r_i(t)$ . To obtain an expression for the second term on the right-hand side of (3.38), notice first that for a function of the first jump time  $G(t_1)$ ,

$$\mathbb{E}[G(t_1)\chi_{B_{\ell}(t)}] = \int_0^t \mathbb{E}[G(s)\chi_{[h_0^j > s, j \neq \ell]} | \bar{h}_0^{\ell} = s] \mu_{i,\ell}(ds)$$

$$= \int_0^t G(s) \Pi_{j=1, j \neq \ell}^{n_{\ell}} r_{i,j}(s) \mu_{i,\ell}(ds) = \int_0^t G(s) \frac{r_i(s)}{r_{i,\ell}(s)} \mu_{i,\ell}(ds).$$

Notice also that  $T_i(t) = \hat{T}_{\xi_{\ell}(i)}(t-t_1)(E_{i,\ell}(t_1))$  when the transition  $\ell$  is first triggered, where  $\hat{T}_{\xi_{\ell}(i)}(t-t_1)$  is the transition matrix of the SHS from  $t_1$  to t starting the process at  $q_1 = \xi_{\ell}(i)$  and  $E_{i,\ell}(s) := J_{i,\ell}e^{As}$ ,  $s \ge 0$ . Each of the terms of the summation on the right-hand side of (3.38) can then be expressed as (3.40)

$$\mathbb{E}[(T_i(t)^{\mathsf{T}})^{(m)}d_{\kappa}\chi_{B_{\ell}(t)}] = \int_0^t (E_{i,\ell}(s)^{\mathsf{T}})^{(m)}\mathbb{E}[(\hat{T}_{\xi_{\ell}(i)}(t-s)^{\mathsf{T}})^{(m)}d_{\kappa}]\frac{r_i(s)}{r_{i,\ell}(s)}\mu_{i,\ell}(ds).$$

By construction of the process (or in other words using the Markov property of the process at transition times)  $\mathbb{E}[(\hat{T}_{\xi_{\ell}(i)}(t-s)^{\intercal})^{(m)}d_{\kappa}] = \mathbb{E}[(T_{\xi_{\ell}(i)}(t-s)^{\intercal})^{(m)}d_{\kappa}] = w^{\kappa,\xi_{\ell}(i)}(t-s)$ . Replacing (3.39) and (3.40) in (3.38) and noticing that  $q_0 = i \in \mathcal{Q}$  is arbitrary we obtain that

(3.41) 
$$w^{\kappa,i}(t) = W_i(w^{\kappa}(t)) + (e^{A_i^{\mathsf{T}}t})^{(m)} \frac{(b^{\kappa})^{\mathsf{T}}}{\gamma_{\kappa}} r_i(t), \quad i \in \mathcal{Q},$$

where

$$W_i(w^{\kappa}(t)) := \sum_{\ell=1}^{n_{\ell}} \int_0^t (E_{i,\ell}(\tau)^{\mathsf{T}})^{(m)} w^{\kappa,\xi_{\ell}(i)} (t-\tau) \frac{r_i(\tau)}{r_{i,\ell}(\tau)} \mu_{i,\ell}(d\tau).$$

An explicit solution to the set of equations (3.41) takes the form

$$(3.42) w^{\kappa}(t) = \sum_{k=0}^{\infty} W^k \left( (e^{A_i^{\mathsf{T}} t})^{(m)} \frac{(b^{\kappa})^{\mathsf{T}}}{\gamma_{\kappa}} r_i(t) \right),$$

where  $w^{\kappa} = (w^{\kappa,1}, \dots, w^{\kappa,n_q})$ ,  $W = (W_1, \dots, W_{n_q})$ , and  $W^k$  denotes composition. From (3.42) it is possible to conclude that  $w^{\kappa,i}(t)$  belongs to the dual vector space of the symmetric tensors for every  $i \in \mathcal{Q}$  and therefore can be written as

(3.43) 
$$w^{\kappa,i}(t) = \sum_{j=1}^{p} u_j^{\kappa,i}(t) (b^j)^{\mathsf{T}}.$$

Replacing (3.43) in (3.41) we get (3.44)

$$\sum_{j=1}^p u_j^{\kappa,i}(t)(b^j)^\intercal = (e^{A_i^\intercal t})^{(m)}(b^\kappa)^\intercal \frac{1}{\gamma_\kappa} r_i(t)$$

$$+ \sum_{\ell=1}^{n_{\ell}} \int_{0}^{t} (E_{i,\ell}(\tau)^{\mathsf{T}})^{(m)} \sum_{j=1}^{p} u_{j}^{\kappa,\xi_{\ell}(i)} (t-\tau) (b^{j})^{\mathsf{T}} \frac{r_{i}(\tau)}{r_{i,\ell}(\tau)} \mu_{i,\ell}(d\tau), \quad i \in \mathcal{Q}.$$

Multiplying both sides of (3.44) by  $b^{\iota}, \iota \in \{1, ..., p\}$ , and using the fact that the maps (3.34) and (3.35) are equivalent, we obtain the set of equations (3.9).  $\square$  Proof of Theorem 3.3. Similarly to (3.26) we can write

(3.45) 
$$\mathbb{E}\left[\sum_{j=1}^{n} (x_j(t))^m\right] = \sum_{j=1}^{n} \mathbb{E}[(x(t)^{\mathsf{T}})^{(m)} e_j^{(m)}].$$

Each term of the summation  $\mathbb{E}[(x(t)^{\intercal})^{(m)}e_j^{(m)}]$  can be obtained from Theorem 3.1, where the index  $\kappa$  in (3.8), (3.9), (3.10) should be taken as  $\kappa_j$ , defined by

$$\kappa_i : \rho(\kappa_i) = m \mathbf{e}_i, \ 1 \le j \le n,$$

where  $e_j$  is a canonical vector in  $\mathbb{R}^n$ . Due to the linearity of the Volterra equation we can obtain

(3.46) 
$$\sum_{i=1}^{n} \mathbb{E}[(x_j(t))^m] = \sum_{i=1}^{p} x_0^{[i]} u_i^{m,q_0}(t),$$

where  $u^m = (u^{m,1}, \dots, u^{m,n_q})$  is uniquely determined by the Volterra equation

(3.47) 
$$u^{m}(t) = \Theta(u^{m}(t)) + h^{m}(t), \quad t \ge 0,$$

where  $h^m(t) := (h^{m,1}(t), \dots, h^{m,n_q}(t)),$ 

$$h^{m,i}(t) = \Gamma^m(e^{A_i^{\mathsf{T}}t}) \sum_{j=1}^n e_{\kappa_j} \frac{r_i(t)}{\gamma_{\kappa}}, \ i \in \mathcal{Q},$$

and  $e_{\kappa_j} \in \mathbb{R}^p$ . It is then clear that the limit  $\lambda_L(x_0) = \lim_{t \to \infty} \frac{1}{t} \log(\mathbb{E}[\|x(t)\|_m^m])$  exists for every  $x_0$  if and only if the limit

(3.48) 
$$\lambda_V := \lim_{t \to \infty} \frac{1}{t} \log(\|u^m(t)\|)$$

exists, and  $\lambda_V = \sup_{x_0} \lambda_L(x_0) = \lambda_L$ . Note also that  $\mathbb{E}[\|x(t)\|_m^m] = 0 \ \forall_{x_0}, t > b > 0$ , if and only if  $u^m(t) = 0, t > b > 0$ , in which case both  $\lambda_L$  and  $\lambda_V$  equal  $-\infty$  according to our definitions.

Let

$$\mathcal{U}: \{T \in \mathcal{T}(m,n): T(x,x,\ldots,x) \ge 0 \ \forall_{x \in \mathbb{R}^n} \}$$

and for  $y \in \mathbb{R}^p$ 

$$T_y(w_1, w_2, \dots, w_m) = \left(\sum_{\iota=1}^p y_{\iota} b^{\iota}\right) (w_1 \otimes w_2 \otimes \dots \otimes w_m),$$

and consider the set

$$\mathcal{K}: \{y \in \mathbb{R}^p : T_y \in \mathcal{U}\},\$$

where the  $b^{\iota}$  are described by (3.28). The set  $\mathcal{K}$  is a cone, i.e.,  $\mathcal{K}$  is a closed convex set such that if  $y, z \in \mathcal{K}$ , then  $\alpha_1 y + \alpha_2 z \in \mathcal{K}$  for  $\alpha_1 \geq 0$ ,  $\alpha_2 \geq 0$  and is such that the set  $-\mathcal{K} := \{-y : y \in \mathcal{K}\}$  intersects  $\mathcal{K}$  only at the zero vector. Moreover  $\mathcal{K}$  is a solid cone, which in finite-dimensional spaces is equivalent to being reproducing, i.e., any element  $y \in \mathbb{R}^p$  can be written as  $y = y_1 - y_2$ , where  $y_1, y_2 \in \mathcal{K}$  (cf. [14, p. 10]). In fact, let  $z \in \mathcal{K}$  be defined as  $\sum_{\iota=1}^p \beta_{\iota} b^{\iota}$ , where  $\beta_{\iota} = \delta$  if  $\iota : \rho(\iota) = m \mathbf{e}_i$ ,  $\mathbf{e}_i \in \mathbb{R}^n$ , for some  $i \in \{1,\ldots,n\}$ , and  $\beta_i = 0$  otherwise. Then, from (3.28),  $z = \delta \sum_{j=1}^n (e_j)^{(m)}$ . Given any  $y \in \mathbb{R}^p$ , then  $y = y_1 - y_2$  for  $y_1 = z + y$  and  $y_2 = z$ , which belong to  $\mathcal{K}$  for sufficiently large  $\delta$ . Likewise one can also prove that  $\mathcal{K}^{n_q} := \mathcal{K} \times \cdots \times \mathcal{K} \subset \mathbb{R}^p \times \mathbb{R}^p \times \cdots \times \mathbb{R}^p$ ,  $y = (y_1, \ldots, y_{n_q}) \in \mathcal{K}^{n_q}$  if and only if  $y_i \in \mathcal{K} \ \forall_{1 \leq i \leq n_q}$  is a solid cone. We prove that the Volterra equation (3.47) has a positive kernel with respect to the solid cone  $\mathcal{K}^{n_q}$  in the sense of [10], and therefore we can directly apply [10, Thm. 13] to conclude that  $r_{\sigma}(\Theta^{m}(a))$  is nondecreasing and that the root with largest real part of  $\det(I-\Theta^{m}(z))$ is real and coincides with the unique value a such that  $r_{\sigma}(\hat{\Theta}(a)^m) = 1$ . As stated in [10, Thm. 13], the zero a equals  $\lambda_V$ , given by (3.48), which in turn equals the Lyapunov exponent  $\lambda_L$ , provided this zero is not a removable singularity of a given complex function. The proof of this latter statement follows steps similar to the ones provided in [10, Thm. 4] and is therefore omitted.

To prove that the kernel of the Volterra equation (3.47) is a positive operator, we need to prove (cf. [10, sect. IV.B]) that if  $y = (y^1, \ldots, y^{n_q}) \in \mathcal{K}^{n_q}$ ,  $y^i \in \mathbb{R}^n$ ,  $1 \le i \le n_q$ ,

then  $M(s)y \in \mathcal{K}^{n_q}$  for every s > 0, where

$$M(s) := \begin{bmatrix} M_{1,1}(s) & \dots & M_{1,n_q}(s) \\ \vdots & \ddots & \vdots \\ M_{n_q,1}(s) & \dots & M_{n_q,n_q}(s) \end{bmatrix}$$

and

(3.49) 
$$M_{i,j}(s) := \sum_{\ell: \mathcal{E}_{\ell}(i) = j} \Gamma^{m}(e^{A^{\mathsf{T}}s} J_{i,\ell}^{\mathsf{T}}) \frac{r_{i}(s)}{r_{i,\ell}(s)}.$$

Note that if  $(z^1, z^2, \ldots, z^{n_q}) = M(s)y$ , then for a given s, each  $z^i \in \mathbb{R}^p$ ,  $1 \le i \le n_q$ , can be written as a sum of terms taking the form  $\Gamma^m(e^{A^\intercal s}J_{i,\ell}^\intercal)y^j$  multiplied by positive scalars. Thus to prove that  $z^i \in \mathcal{K}$  and hence  $z \in \mathcal{K}^{n_q}$ , it suffices to prove that  $w = \Gamma^m(C)y \in \mathcal{K}$  for an arbitrary  $1 \le j \le n_q$ ,  $y \in \mathcal{K}$ , and  $C \in \mathbb{R}^{n \times n}$ . To this effect, using the fact that the map (3.34) can be written as (3.35), we have that

$$w: \sum_{\iota=1}^p w_\iota(b^\iota)^\intercal = C^{(m)} \sum_{\iota=1}^p y_\iota(b^\iota)^\intercal$$

belongs to  $\mathcal{K}$  because  $\sum_{\iota=1}^{p} w_{\iota}(b^{\iota})$  belongs to  $\mathcal{U}$  since

$$\begin{split} \sum_{\iota=1}^p w_\iota(b^\iota) x^{(m)} &= \sum_{\iota=1}^p y_\iota(b^\iota) (C^\intercal)^{(m)} x^{(m)} \\ &= \sum_{\iota=1}^p y_\iota(b^\iota) (C^\intercal x)^{(m)} \geq 0, \end{split}$$

where in the last equality we used the fact that  $\sum_{i=1}^{p} y_i(b^i) \in \mathcal{U}$ .

**4. Application to networked control.** We consider the following simplified version of the networked control setup in [10]. Suppose that we wish to control a linear plant

$$\dot{x}_{P}(t) = A_{P}x_{P}(t) + B_{P}\hat{u}(t)$$

using a state feedback controller taking the form  $K_C x_P(t)$  that needs to be implemented digitally and suppose that the actuation is held constant  $\hat{u}(t) = \hat{u}(s_{\kappa}), t \in [s_{\kappa}, s_{\kappa+1})$ , between actuation update times denoted by  $\{s_{\kappa}, \kappa \geq 0\}$ .

The controller has direct access to the state measurements but communicates with the plant actuators through a network possibly shared by other users. The controller attempts to do periodic transmissions of data at a desired sampling period  $T_s$ , but these regular transmissions may be perturbed by the medium access protocol. For example, users using CSMA for medium access may be forced to back off for a typically random amount of time until the network becomes available. We assume these random back-off times to be independent and identically distributed and denote by  $\mu_s$  the associated measure.

We consider two different cases.

Case I. After waiting to obtain network access, the controller (re)samples the sensor, computes the control law, and transmits this most recent data. Assuming that the transmission delays are negligible, and defining  $x := (x_P, \hat{u})$ , we have

$$\dot{x} = Ax, \quad A = \begin{bmatrix} A_P & B_P \\ 0 & 0 \end{bmatrix},$$

$$x(s_k) = Jx(s_k^-), \quad J = \begin{bmatrix} I & 0 \\ K_C & 0 \end{bmatrix}.$$

Since the intervals  $\{s_{k+1} - s_k, k \ge 0\}$  result from the controller waiting a fixed time  $T_s$  plus a random amount of time with a measure  $\mu_s(s)$ , these intervals are independent and identically distributed according to

$$\mu([0,\tau)) = \begin{cases} \mu_s(\tau - T_s) & \text{if } \tau \ge T_s, \\ 0, \quad s \in [0, T_s). \end{cases}$$

Note that the system (4.2) is a special case of an SHS with a single state and a single reset map.

Case II. After waiting to obtain access to the network, the controller does not resample the sensor and simply transmits the data that it had collected at the time of the first attempt to transmit the sensor data. We model this by an SHS with the following two discrete modes  $(n_q = 2)$ :

- State q(t) = 1: The controller waits for a fixed time  $T_s$ .
- State q(t) = 2: The controller waits a random time to gain access to the network.

Let  $r_k = s_k + T_s$ ,  $x := (x_P, \hat{u}, v)$ , where  $v(t) := u_\kappa, t \in [r_\kappa, r_{\kappa+1})$ , is a variable that holds the last computed control value. The transitions between the two discrete modes can be modeled by a single transition  $(n_\ell = 1)$  which is a function of the two discrete modes and takes the form (2.2), specified as follows. When in state 1 the SHS transits to state 2  $(\xi_1(1) = 2)$  at times  $r_\kappa$ . The corresponding state jump models the update of the variable  $v(r_\kappa) = u_\kappa$  that holds the last computed control value and is described by

$$x(r_{\kappa}) = J_{1,1}x(r_{\kappa}^{-}), \quad J_{1,1} = \begin{bmatrix} I & 0 & 0 \\ 0 & I & 0 \\ K_{C} & 0 & 0 \end{bmatrix}.$$

When in state 2 the SHS transits to state 1 ( $\xi_1(2) = 1$ ) at actuation update times  $s_{\kappa}$ . The state jump models the actuation update  $\hat{u}(s_{\kappa}) = v(s_{\kappa}^{-})$  and is described by

$$x(s_{\kappa}) = J_{2,1}x(s_{\kappa}^{-}), \quad J_{2,1} = \begin{bmatrix} I & 0 & 0 \\ 0 & 0 & I \\ 0 & 0 & I \end{bmatrix}.$$

The reset-time measures are given by the following:

- $\mu_{1,1}(\tau) = \delta(\tau T_s)$  is a discrete measure that places all mass  $w_i = 1$  at time  $T_s$ .
- $\mu_{2,1}(\tau) = \mu_s(\tau)$ .

In both discrete modes, the continuous-time dynamics are described by  $\dot{x} = A_i x$ ,  $i \in \{1, 2\}, A_1 = A_2 = A$ , where

$$A = \begin{bmatrix} A_P & B_P & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}.$$

## **4.1.** Numerical example. Suppose that the plant (4.1) is described by

$$A_P = \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}, \ B_P = \begin{bmatrix} 0 \\ 1 \end{bmatrix},$$

which by properly scaling the state and input can be viewed as a linearized model of a damp-free inverted pendulum. Moreover, suppose that the measure  $\mu_s$  is uniform with support on the interval  $[0,\tau]$ , and fix  $T_s=0.1s$  and  $x_P(0)=[1\,0]^\intercal$ . A continuous-time state feedback controller is synthesized using a linear quadratic regulator (LQR) and is given by  $\hat{u}(t)=K_Cx(t)$ ,  $K_C=-\left[1+\sqrt{2}-1+\sqrt{2}\right]$ , which is the solution to the problem  $\min_{\hat{u}(t)}\int_0^\infty [x_P(t)^\intercal x_P(t)+\hat{u}(t)^2]dt$ , yielding the closed-loop eigenvalues  $\lambda_i(A_P+B_PK_C)=\{-1,-\sqrt{2}\}$ . We wish to investigate the stability and performance of the closed loop when instead of the ideal network-free case we consider the scenarios of Cases I and II. To this effect we define the quantity

$$e(t) = x_P(t)^{\mathsf{T}} x_P(t) + \hat{u}(t)^2$$

which can be written as  $e(t) = x^{\mathsf{T}} P x$ , where (i) in the network-free case  $P = I_2 + K_C^{\mathsf{T}} K_C$  and  $x = x_P$ ; (ii) in Case I,  $P = I_3$  and  $x = (x_P, \hat{u})$ ; and (iii) in Case II,  $P = \operatorname{diag}(I_2, 1, 0)$ , and  $x = (x_P, \hat{u}, v)$ . Note that, in the network-free case, e(t) is the quantity whose integral is minimized by LQR control synthesis and e(t) decreases exponentially fast at a rate  $\alpha = 2$ , since the dominant closed-loop eigenvalue equals  $\lambda_i(A_P + B_P K_C) = -1$ . In Cases I and II,  $\mathbb{E}[e(t)]$  converging to zero is equivalent to mean square stability ( $\tau$  and  $\tau_s$  are finite). Corollary 3.4 can be used to determine whether the closed loop is MSS in Cases I and II. Moreover, when the closed loop is MSS, we can determine the exponential decay constant of  $\mathbb{E}[e(t)]$  from Theorem 3.3. The results are summarized in Table 4.1 for different values of the support  $\tau$  of the uniform measure  $\mu_s$  of the back-off time.

The fact that closed-loop stability is preserved for larger values of  $\tau$  in Case I confirms what one would expect intuitively, i.e., Case I is more appropriate when transmitting dynamic data, since the most recent sampling information is sent through the network.

Using the state moment expressions provided by Theorem 3.1, we can perform a more detailed analysis by plotting the moments of e(t), which can be expressed in terms of the moments of the state. For example, the two first moments take the form

$$\begin{split} \mathbb{E}[e(t)] &= \mathbb{E}[x(t)^\intercal P x(t)] = \mathbb{E}[(x(t)^\intercal)^{(2)}] \nu(P), \\ \mathbb{E}[e(t)^2] &= \mathbb{E}[(x(t)^\intercal P x(t))^2] = \mathbb{E}[(x(t)^\intercal)^{(4)}] (\nu(P) \otimes \nu(P)). \end{split}$$

In Figure 4.1, we plot the expected value of the error  $\mathbb{E}[e(t)]$  and its  $2-\sigma$  confidence interval  $\mathbb{E}[e(t)] \pm 2\mathbb{E}[(e(t) - \mathbb{E}[e(t)])^2]^{1/2}$  for a network measure support  $\tau = 0.4$ . Note that from the Chebyshev inequality, we conclude that

 $\begin{array}{c} \text{Table 4.1} \\ \text{Exponential decay rates } \mathbb{E}[e(t)] \leq c e^{-\alpha t} \end{array}$ 

	(a) Case I											
au	0.4	0.6	0.8	1.0	1.2	> 1.211						
$\alpha$	2.000	2.000	1.969	0.477	$7.63 \times 10^{-5}$	Not MSS						
(b) Case II												
$\tau$	0.1	0.2	0.3	0.4	0.5	> 0.521						
$\alpha$	2.000	2.000	2.000	0.849	0.118	Not MSS						

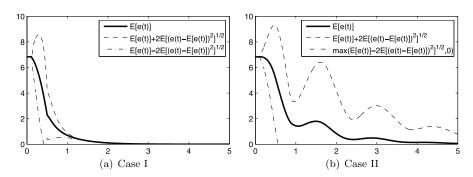


Fig. 4.1. Plot of  $\mathbb{E}[e(t)]$ , where e(t) is quadratically state dependent. For a fixed t,  $\mathbb{E}[e(t)]$  lies between the dashed curves with probability  $> \frac{3}{4}$ .

$$\text{Prob}[|e(t) - \mathbb{E}[e(t)]| > a(t)] \le \frac{\mathbb{E}[(e(t) - \mathbb{E}[e(t)])^2]}{a(t)^2},$$

and therefore one can guarantee that for a fixed t, e(t) lies between the curves  $\mathbb{E}[e(t)] \pm a(t)$ ,  $a(t) = 2\mathbb{E}[(e(t) - \mathbb{E}[e(t)])^2]^{1/2}$  with a probability greater than  $\frac{3}{4}$ . The numerical method used to compute the solution of the Volterra equation is the one described in section 3.1 for which we used a trapezoidal rule for the integration method. In Case I, the expected value of the quadratic state function e(t) tends to zero much faster, and with a much smaller variance than in Case II, confirming once again that Case I is more appropriate when transmitting dynamic data.

The plots of Figure 4.1 can be confirmed through Monte Carlo simulation, although this latter method is in general more computationally demanding. For illustration, in Table 4.2 we compare the numerical error  $\sum_{k=0}^{L} \mathbb{E}[|e(k\underline{h}) - \hat{e}(k\underline{h})|]$ ,  $\underline{L} = 100$ ,  $\underline{h} = 0.01$ , obtained with our proposed method (with trapezoidal rule for integration) and the Monte Carlo method, for Case I, where  $e(k\underline{h})$  is a very accurate approximation of e(t) at time  $k\underline{h}$  and  $\hat{e}(k\underline{h})$  is the approximation obtained with one of the methods presented in Table 4.2. For the Monte Carlo method, we show the numerical errors obtained with a different numbers of Monte Carlo runs, whereas for our proposed method we show the numerical errors obtained with different discretization steps or equivalently with a different number of discretization points L = 1 equally spaced in the interval [0,1]. The computational times were obtained with a computer with the following specifications: Intel Core 2 Duo, 3 GHz processor speed, and 3.5 GB RAM. From the values in Table 4.2 it is clear that our proposed method provides better accuracy with fewer computations.

 $\begin{array}{c} \text{Table 4.2} \\ \textit{Numerical errors } \sum_{k=0}^{\underline{L}} \mathbb{E}[|e(k\underline{h}) - \hat{e}(k\underline{h})|] \textit{ for case I, } \underline{L} = 100, \ \underline{h} = 0.01. \end{array}$ 

(a) Proposed method based on Volterra equations										
Number of discretization points $L$	500	1000	2000	4000	8000	16000				
Computational time (sec)	6.04	7.44	12.17	29.37	103.19	347.52				
Num. error	5.0187	2.4482	1.1796	0.5493	0.2352	0.0783				
(b) Monte Carlo simulations										
Number of Monte Carlo runs	300	600	1200	2400	4800	9600				
Computational time (sec)	10.03	19.47	42.90	90.80	235.64	701.22				
Num. error	3.73	2.26	3.07	2.02	0.79	0.77				

5. Conclusions and future work. We proposed an approach based on Volterra renewal-type equations to analyze SHSs for which the lengths of time that the system stays in each mode are independent random variables with given distributions. We showed that any statistical mth degree moment of the state can be computed using this approach, and we provided a number of results characterizing the asymptotic behavior of a second-degree moment of the system. Due to the large number of problems that fit the stochastic hybrid systems framework, finding more applications where the results can be applied is a topic for future work.

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