

Perturbation Analysis: A Framework for Data-Driven Control and Optimization of Discrete Event and Hybrid Systems

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Abstract: The history of Perturbation Analysis (PA) is intimately related to that of Discrete Event Dynamic Systems (DEDS), starting with a solution of a long-standing problem in the late 1970s and continuing today with the control and optimization of Hybrid Systems and the emergence of event-driven control methods. We review the origins of the PA theory and how it became part of a broader framework for models, control and optimization of DEDS. We then discuss the theoretical underpinnings of Infinitesimal Perturbation Analysis (IPA) as a data-driven stochastic gradient estimation method and how it has been applied over the past few decades. We explain how IPA offers a basis for general-purpose stochastic optimization of Markov systems through the notion of the performance potential and how it has evolved beyond DEDS and now provides a framework for control and optimization of Hybrid Systems and, more generally, event-driven methodologies.

1. THE ORIGIN OF PERTURBATION ANALYSIS

In pioneering the field of Discrete Event Systems (DES) in the early 1980s, Y.C. Ho and his research group at Harvard University discovered that event-driven dynamics give rise to state trajectories (sample paths) from which one can very efficiently and nonintrusively extract sensitivities of state variables (therefore, various performance metrics as well) with respect to at least certain types of design or control parameters. This eventually led to the development of a theory for *Perturbation Analysis* (PA) in DES (Cassandras and Lafortune (2008), Ho and Cao (1991), Glasserman (1991)), the most successful branch of which is *Infinitesimal Perturbation Analysis* (IPA) due to its simplicity and ease of implementation. In fact, by the early 2000s, IPA was shown to apply to all virtually arbitrary Hybrid Systems (HS) and continues to be today one of the most attractive tools for data-driven control and optimization, especially in stochastic environments where modeling random aspects of a process is prohibitively hard.

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The origin of the key concepts that form the cornerstones of the PA theory are found in a long-standing problem in operations research and industrial engineering known as the *buffer allocation problem*. In its industrial engineering version, it was presented to Ho's research group by the FIAT automobile company in the late 1970s as follows. A typical serial transfer line consists of N workstations in tandem, each with different characteristics in terms of its production rate, failure rate and repair time when failing. In order to accommodate this inhomogeneous behavior, a buffer is placed before the i th workstation, $i = 1, \dots, N$, with B_i discrete slots where production parts can be queued. Since the space within which this transfer line operates is limited, there is an upper bound B to the total number of buffer slots that can be allocated over the N workstations so that $\sum_{i=1}^N B_i = B$. The problem is to allocate these B buffer slots, i.e., determine a vector $[B_1 \dots B_N]$, so as to maximize the throughput of the transfer line while also maintaining a low overall average delay of the parts moving from an entry point before the first workstation to an exit point following the N th workstation. Tackling this problem in a "brute force" manner requires considering all possible buffer allocations, a number given by $\binom{B+N-1}{B}$. For a reasonably small

problem such as $B = 24$ and $N = 6$, this gives 118,755 possible solutions. A direct trial-and-error approach where one is allowed to test each allocation for about a week would require about 2300 years. If one were to reduce the initial solution space to only 1000 “good guesses” and use early 1980s simulation technology requiring about 3 min per trial to estimate the resulting performance, the overall task would take about 250 days of CPU time.

The approach taken by Ho’s group and first reported in Ho et al. (1979) was to study the serial transfer line as a dynamic system whose state includes the integer-valued buffer contents along with real-valued “clocks” associated with each workstation as it processes a part. The question then posed was: “what would happen if in a given allocation a specific value B_i were changed to B_i+1 ?” The “brute force” way to answer this question is to first simulate the system under the nominal allocation with the value B_i and estimate the system’s performance over a (sufficiently long) time period T which may be denoted by $L_T(B_i)$. Then, repeat the simulation under $B_i + 1$ to obtain $L_T(B_i+1)$. The difference $\Delta L_T(B_i) = L_T(B_i+1) - L_T(B_i)$ provides an estimate of the system’s performance sensitivity with respect to B_i . What the research team realized, however, is that this is unnecessary: indeed, the initial simulation alone yielding $L_T(B_i)$ and a simple thought experiment can deliver the value of $\Delta L_T(B_i)$. Moreover, the same thought experiment can deliver the entire vector $[\Delta L_T(B_1), \dots, \Delta L_T(B_N)]$ with minimal extra effort.

The key observation that led to a formal procedure describing this thought experiment is the following. When B_i is replaced by $B_i + 1$, no change in the state of the system can take place unless one of two “events” is observed at time t : (i) The i th buffer content, say $x_i(t)$, reaches its upper limit, i.e., $x_i(t) = B_i$ and a part is ready to leave the $(i - 1)$ th workstation. In this case, this upstream workstation is “blocked” since there is no place for the departing part to go. However, in a perturbed system with B_i replaced by $B_i + 1$ that would not happen and one can simply predict a buffer content perturbation $\Delta x_i(t) = 1$. Moreover, one can record when this blocking occurs at time $t \equiv t_{i,B}$ and the next time that a part departs from the i th workstation, $t_{i,D}$. Then, $t_{i,D} - t_{i,B}$ is the amount of time that would be gained (i.e., no blocking would have occurred) in a perturbed system realization. The important observation here is that $t_{i,D}$, $t_{i,B}$ are directly observed along the nominal system realization. (ii) The $(i + 1)$ th buffer content reaches its lower limit, i.e., $x_{i+1}(t) = 0$ and a part is ready to leave the i th workstation. In this case, if $\Delta x_i(t) = 1$, i.e., the i th workstation has already gained a part from an earlier blocking event, then this gain can now propagate downstream and we can set $\Delta x_{i+1}(t) = \Delta x_i(t) = 1$.

This simple observation leads to the conclusion that estimating the effect of replacing B_i by $B_i + 1$ boils down to observing just a few events along the nominal system realization: *blocking events* (when $x_i(t) = B_i$ and a part departure from $i - 1$ takes place) and *idling events* (when $x_i(t) = 0$ at any $i = 1, \dots, N$). This can be formalized into an “estimator” for buffer perturbations $\Delta x_i(t)$ and event timing perturbations for all part departures at workstations. More generally, this estimator transforms a given

hypothetical perturbation $\Delta B_i(t) = 1$ (or -1) into state perturbations, which can ultimately be used to estimate a performance perturbation $\Delta L_T(B_i)$. Most importantly, this is accomplished without ever having to implement the perturbation $\Delta B_i(t)$, since the estimator depends only on directly observable data from the nominal system realization; in particular, it suffices to observe selected events and associated event times and to perform extremely simple calculations.

This initial procedure pertaining to a very specific type of dynamic system and problem was given the name *Perturbation Analysis* (PA). It soon became clear that it could be extended to any system with a structure similar to that of the serial transfer line and to a perturbation in any system parameter. Thus, one could consider, for instance, speeding up the operation of a workstation and studying the effect of a perturbation Δr_i in the operation rate r_i of the i th workstation. The general procedure is one where some parameter perturbation $\Delta \theta$ generates a state perturbation $\Delta x_i(t)$ when a specific event occurs at time t . Subsequently, the system dynamics dictate how $\Delta x_i(t)$ propagates through the system by affecting $\Delta x_i(t)$ or $\Delta x_j(t)$ for $j \neq i$. Depending on a performance metric of interest, this ultimately yields $\Delta L_T(\Delta \theta)$, the change in performance due to $\Delta \theta$. As for the system structure amenable to this kind of efficient PA, it became obvious that it fits the general class of queueing networks.

An obvious next question was: “Does PA hold for any value of $\Delta \theta$ or do we have to restrict it to “small” $\Delta \theta$ when θ is real-valued?” There was ample empirical evidence collected over the early 1980s that $\Delta \theta$ had to be small but not necessarily “very small”. In other words, the values of $\Delta L_T^{PA}(\Delta \theta)$ obtained through PA were identical to those obtained through the “brute force” finite difference $L_T(\theta + \Delta \theta) - L_T(\theta)$ for “sufficiently small” $\Delta \theta$. This led to the term *Infinitesimal Perturbation Analysis* (IPA) to capture the fact that the methodology was applicable to perturbations which were “infinitesimally” small, although a formal quantification characterizing limits for $\Delta \theta$ was lacking. Moreover, when $\Delta \theta$ became larger, it was still possible to satisfy $\Delta L_T^{PA}(\Delta \theta) = L_T(\theta + \Delta \theta) - L_T(\theta)$ at the expense of observing more “interesting events” and performing a few more calculations. For instance, in the case of the integer-valued buffer size parameter B_i , the minimal feasible perturbation is obviously either $+1$ or -1 . To differentiate these cases, the term *Finite Perturbation Analysis* (FPA) was introduced. FPA reverts to IPA when parameters are real-valued and may be allowed to take “sufficiently small” values $\Delta \theta$.

To illustrate the distinction between IPA and FPA, we consider the case of a simple First-In-First-Out (FIFO) queueing system with a single server preceded by a queue. Let $\{A_k\}$ be the sequence of (generally random) arrival times, $k = 1, 2, \dots$, and $\{D_k\}$ be the corresponding sequence of departure times from the system. If Z_k denotes the service time of the k th entity (customer) processed, then the Lindley equation

$$D_k = \max(A_k, D_{k-1}) + Z_k \quad (1)$$

describes the departure time dynamics with $k = 1, 2, \dots$. Suppose that all (or just some selected subset) of the service times are perturbed by ΔZ_k , $k = 1, 2, \dots$. Let

$I_k = A_k - D_{k-1}$ and observe that when $I_k > 0$ it captures an idle period (since the server must wait until $A_k > D_{k-1}$ to become busy again) and when $I_k < 0$ it captures the waiting time $D_{k-1} - A_k$ of the k th arriving entity in the system. It is easy to obtain from (1) the following departure time perturbation equation:

$$\Delta D_k = \Delta Z_k + \begin{cases} \Delta D_{k-1} & \text{if } I_k \leq 0, \Delta D_{k-1} \geq I_k \\ 0 & \text{if } I_k > 0, \Delta D_{k-1} \leq I_k \\ I_k & \text{if } I_k \leq 0, \Delta D_{k-1} \leq I_k \\ \Delta D_{k-1} - I_k & \text{if } I_k > 0, \Delta D_{k-1} \geq I_k \end{cases} \quad (2)$$

where ΔD_k can be obtained from the generated perturbations ΔZ_k and directly observed data in the form of I_k . This is the FPA procedure for evaluating $\Delta D_k, k = 1, 2, \dots$. Observe, however, that if we select $\Delta Z_k > 0$ to be sufficiently small so that $\Delta D_{k-1} > 0$ can never exceed the finite value of $I_k > 0$, then this reduces to

$$\Delta D_k = \Delta Z_k + \begin{cases} \Delta D_{k-1} & \text{if } I_k \leq 0 \\ 0 & \text{otherwise} \end{cases} \quad (3)$$

which is the much simpler IPA version, requiring only the detection of an idling interval when $I_k > 0$, at which time the departure time perturbation is reset to $\Delta D_k = \Delta Z_k$. Naturally, the question is: ‘‘How small do perturbations need to be before the IPA equation can be used?’’ In a stochastic system, the answer to this question is generally dependent on the specific realization based on which ΔD_k is evaluated. Thus, it is logical to extend the question of estimating $\Delta D_k(\Delta\theta), k = 1, 2, \dots$, the departure time perturbations, to estimating the derivative $\frac{d\Delta D_k}{d\theta}$ by allowing $\Delta\theta \rightarrow 0$. This became one of the two important questions facing PA researchers in the early 1980s:

Q1. When can IPA be used instead of FPA and is it possible to transition the remarkably efficient PA methodology from one limited to sensitivity analysis for finite perturbations to one for gradient estimation of a large class of dynamic systems?

Q2. What is the class of dynamic systems for which the PA methodology applies?

To address **Q2**, it soon became clear that using traditional models for system dynamics of the general form $\dot{x} = f(x, u, t)$ was extremely inefficient and ultimately pointless for systems such as queueing networks. In such systems, at least some state variables (i.e., queueing contents) are discrete and almost always fixed, changing only when specific events occur (i.e., an entity enters or leaves the queue). This led to the realization that traditional dynamic systems described through $\dot{x} = f(x, u, t)$ are *time-driven*, whereas this different class of systems is *event-driven*. The term *Discrete Event Dynamic System* (DEDS) was coined in 1980 and first appeared in the literature in Ho and Cassandras (1980) and Ho and Cassandras (1983), while a first general IPA and FPA framework for queueing networks appeared in Ho et al. (1983). However, the event-driven nature of this class of systems was not formalized until an ‘‘event domain formalism’’ was first proposed in Cassandras and Ho (1985). Over the next several years, it became clear that the class of DEDS is much broader than queueing networks and that PA techniques could be extended to all such systems (Ho and Cao (1991), Glasserman (1991)). In parallel, a novel

control theory for such systems, with the broader term *Discrete Event System* (DES) used, was being developed by Ramadge and Wonham culminating with what has become known as the *supervisory control theory* for DES (Ramadge and Wonham (1980)). It took about a decade before the supervisory control theory and PA were merged into complementary approaches for studying DES and are now viewed as a staple of any study of dynamic systems (Cassandras (1993), Cassandras and Lafortune (2008)).

Returning to the first question **Q1** above, IPA in the form of (3) was successfully used in the early 1980s for many applications that involved stochastic systems with event-driven behaviors, including routing, scheduling and general resource allocation problems in complex manufacturing systems and computer and communication networks (e.g., Cassandras et al. (1990)). The generalization of (3) is to apply it to any performance metric $J(\theta) = E[L(\theta)]$ where $L(\theta)$ is a sample function dependent on θ . IPA is an efficient way to obtain $\nabla L(\theta)$ from observable data on a nominal system realization. However, what is ultimately of interest is $\nabla J(\theta) = \nabla E[L(\theta)]$, and its estimation by $\nabla L(\theta)$ can use in a large class of gradient-based optimization problems. As IPA was applied to harder and harder problems (i.e., systems with event-driven dynamics) much more complex than Lindley equations such as (1), it became clear that IPA estimates $\nabla L(\theta)$ were not accurate compared to $\nabla J(\theta)$ when this could be evaluated through analytical methods in some simple cases or accurately approximated through exhaustive time-consuming simulation methods. Indeed, one could have situations where the signs of $\nabla J(\theta)$ and $\nabla L(\theta)$ were different, resulting in heavily *biased* IPA gradient estimates. It took several years and occasionally controversial debates to realize that the key issue was one of testing the validity of unbiasedness for IPA gradient estimation, i.e., formal conditions under which

$$\nabla E[L(\theta)] = E[\nabla L(\theta)]$$

holds, or in simpler scalar form:

$$\frac{d}{d\theta} E[L(\theta)] = E \left[\frac{dL(\theta)}{d\theta} \right]$$

The way this key issue was addressed is discussed in the next section.

2. INFINITESIMAL PERTURBATION ANALYSIS

By the mid 1980s, it was realized that IPA provided a general framework for computing gradients of sample performance functions defined on the state space of an extensive class of DEDS beyond queueing networks. Furthermore, it was shown to admit especially simple computations by data gathered directly from the sample path of the system. Consequently IPA became the focal point of research in PA, with an eye on potential applications in performance optimization by stochastic gradient-descent algorithms.

The basic thinking about IPA was shaped by the view that it essentially compares two sample paths in slightly changed parameters with a set of common random variables for both paths. A formal mathematical formulation of the IPA framework appeared in Cao (1985). A sample path of a stochastic system is denoted as a pair (θ, ξ) with $\theta \in R^n$ denoting the parameter concerned, and ξ , a set of random variables describing all the randomness involved in

the system. Let $L_T(\theta, \xi)$ be a performance function of the system in a given finite period $[0, T]$. IPA consists of the sample gradient (derivative) $\nabla L_T(\theta, \xi) := \frac{\partial}{\partial \theta}(L_T(\theta, \xi))$, called the *sample derivative*, or *sample gradient* (Cao (1985)). Due to the DEDS structure of the system, it was convenient to think of this derivative as the limit of finite differences with the common sample path, namely

$$\frac{\partial L_T(\theta, \xi)}{\partial \theta} = \lim_{\Delta\theta \rightarrow 0} \frac{L_T(\theta + \Delta\theta, \xi) - L_T(\theta, \xi)}{\Delta\theta}.$$

However, one of the expressed objectives of IPA is to estimate the derivative of the mean performance $J_T(\theta) := E[L_T(\theta, \xi)]$.¹ This naturally raises the question “is the sample derivative given by IPA an *unbiased* estimate of the derivative of the mean performance?” i.e.,

$$E \left[\frac{\partial}{\partial \theta} L_T(\theta) \right] = \frac{\partial}{\partial \theta} E[L_T(\theta)]? \quad (4)$$

The unbiasedness (4) explains why IPA gives accurate derivative estimates for some systems but not others. Roughly speaking, when the sample function $L_T(\theta)$ has a jump (discontinuity) at θ , the interchangeability of expectation and differentiation inherent in Eq. (4) would not hold. Intuitive conditions for this interchangeability were given in Cao (1985); essentially, if a parameter change at θ may cause a change in the order of events in a DEDS, the sample performance function may have a jump at θ , and then the interchangeability in (4) may hold only if the probability of such jumps in $[\theta, \theta + \Delta\theta]$ is of the order $o(\Delta\theta)$. This intuition evolved from the consideration of many engineering problems defined on DEDS. An alternative view of the issue of unbiasedness is through the theory of multivariable calculus. After all, Eq. (4) amounts to an interchangeability of differentiation with respect to θ and integration with respect to ξ (expectation), and this is roughly equivalent to the continuity of the function $L(\theta, \xi)$ in θ with probability one (in ξ). The former view is based on engineering intuition, whereas the latter one, founded upon well-known results in mathematics, often provides a practical way of discerning whether IPA is unbiased.

These realizations stimulated subsequent research in two directions:

1. Identifying classes of systems and problems where IPA is unbiased. Early works include Suri and Zazanis (1988), which proves the unbiasedness for the GI/G/1 queue, and Cao (1988), which proves it for closed Jackson networks. Glasserman (1991) presents IPA in the framework of generalized semi-Markov processes, and extends the results in Cao (1985) and Cao (1988) to a *commuting condition* for the unbiasedness of the derivative estimates. For many other applications in this direction, please see Cassandras and Lafortune (2008).

2. The development of alternative perturbation-analytic techniques that provide unbiased derivative estimates, or reduce the bias, for problems for which IPA estimates are biased. Several such techniques have appeared in the literature; among them is *Smoothed Perturbation Analysis (SPA)* (Gong and Ho (1987), Fu and Hu (1997)). The main idea of this technique is to use the derivative of a

¹ To simplify the presentation we will drop the explicit notational dependence of the sample performance functions on ξ .

conditional mean of the sample function as the estimate of the derivative of the mean performance. While a sample function may have jumps and therefore its derivative is a biased estimate, a conditional mean of the sample function may be smooth enough to provide unbiased estimates. More precisely, suppose that there is a random variable (or vector) denoted as Z such that

$$\begin{aligned} E \left[\frac{\partial}{\partial \theta} E[L_T(\theta)|Z] \right] &= \frac{\partial}{\partial \theta} E[E[L_T(\theta)|Z]] \\ &= \frac{\partial}{\partial \theta} E[L_T(\theta)], \end{aligned} \quad (5)$$

then we can use $\frac{\partial}{\partial \theta} E[L_T(\theta)|Z]$, i.e., the derivative of the conditional-mean sample function, as an unbiased estimate of the performance derivative $\frac{\partial}{\partial \theta} E[L_T(\theta)]$. A potential difficulty with this approach is that the conditional IPA estimator may require a significantly higher computational effort than the basic IPA to the point that it is rendered impractical. In other words, precision and accuracy can be obtained at the expense of high computational complexity.

Around the same time, various other techniques were also developed based largely on so-called “cut-and-paste” operations on the sample path in order to smooth out discontinuities resulting from parameter perturbations. Surveys thereof can be found in Ho and Cao (1991) and Cassandras and Lafortune (2008).

The preceding discussion pertains to finite-horizon sample performance functions. Another important class of functions concern long-run (infinite-horizon) averages. Denoted by $J(\theta)$, they have the form

$$J(\theta) = \lim_{T \rightarrow \infty} \frac{1}{T} L_T(\theta),$$

where the system is assumed to be ergodic for the above limit to exist and be independent of the sample path ξ w.p.1. The time T can be either continuous or discrete. IPA gives the sample derivative $\frac{1}{T} \frac{\partial}{\partial \theta} L_T(\theta)$. The issue here is the strong consistency of the IPA derivative, i.e., whether the following limit is in force Cao (1985),

$$\lim_{T \rightarrow \infty} \frac{1}{T} \frac{\partial}{\partial \theta} L_T(\theta) = \frac{\partial J(\theta)}{\partial \theta}; \quad (6)$$

in other words, “are the operators of limit “ $\lim_{T \rightarrow \infty}$ ” and derivative “ $\frac{\partial}{\partial \theta}$ ” interchangeable?”

The study of this issue led to an important concept, the *perturbation realization*, later extended to the *performance potential*, which has been applied to several research areas like Markov decision processes and stochastic control. The main idea is as follows. In queueing networks, the effect of every single perturbation on the performance is finite and can be precisely measured; the total effect of a parameter change on the performance can be decomposed into the sum of the effects of every single perturbation generated (realized) by this parameter change. The performance derivative with respect to this parameter can then be calculated. To illustrate this concept, consider a closed Jackson network consisting of M servers with service rates μ_i , $i = 1, 2, \dots, M$, and let

$$L_T(\theta) := \int_0^T f[\mathbf{n}(t)] dt,$$

where $\mathbf{n}(t) = (n_1, n_2, \dots, n_M)$ is the system state at time t with n_i is the number of customers at server i ,

and $f[\mathbf{n}(t)]$ is a performance function of the state. Now, suppose at $t = 0$ with initial state \mathbf{n} , server i is subjected to a perturbation Δ , meaning, e.g., its completion time is delayed by the amount of Δ . Define the *perturbation realization factor*, (Ho and Cao (1993), Cao (1994)),

$$c(\mathbf{n}, i) = \lim_{T \rightarrow \infty} \lim_{\Delta \rightarrow 0} E \left\{ \frac{1}{\Delta} \left[\int_0^T f(\mathbf{n}'(t)) dt - \int_0^T f(\mathbf{n}(t)) dt \right] \right\}. \quad (7)$$

The perturbed sample path $\mathbf{n}'(t)$ can be simply obtained by the propagation rule on a single sample path. With the help of the realization factors, we can prove that the strong consistency (6) indeed holds, and we can further derive (take $\theta = \mu_i$ as the perturbed parameter)

$$\begin{aligned} \lim_{T \rightarrow \infty} \left[\mu_i \frac{1}{T} \frac{\partial L_T(\mu_i)}{\partial \mu_i} \right] &= \mu_i \frac{dJ(\mu_i)}{d\mu_i} \\ &= \sum_{\text{all } \mathbf{n}} \pi(\mathbf{n}) c(\mathbf{n}, i), \quad \text{w.p.1,} \end{aligned} \quad (8)$$

where $\pi(\mathbf{n})$ is the steady-state probability of state \mathbf{n} . $c(\mathbf{n}, i)$ can be computed by a set of linear equations.

In this approach, the interchangeability of “ $\lim_{T \rightarrow \infty}$ ” and “ $\frac{\partial}{\partial \theta}$ ” is buried in $\lim_{T \rightarrow \infty} \lim_{\Delta \rightarrow 0}$ in (7) because in a strongly connected network, a perturbation can only affect a system in a finite period, and the difference of the two terms in (7) will be almost zero when T is large enough. The interchangeability can be proved along these lines. Many other examples with perturbation realization can be found in Cao (1994).

The notions of perturbation realization and performance potential continue to underscore subsequent applications to large-scale systems. For non-Markovian queueing networks and other DEES, the unbiasedness of IPA and the computational complexity inherent in alternative PA methods designed to circumvent it, led to a partial shift of IPA research from DEES to stochastic hybrid systems. These developments, which have been taking place over the past fifteen years, are the subject of the next section.

3. STOCHASTIC HYBRID SYSTEMS

In 2002 a new approach to IPA emerged, based on Stochastic Flow Models (SFM) (Cassandras et al. (2002)).² Unlike SPA it does not consist of alternative sample-path representations of $J(\theta)$, but rather on an alternative modelling framework that yields approximate estimates for $J(\theta)$, and, more importantly, whose gradients are *unbiased* and provide approximations to $\nabla J(\theta)$. The SFM concept formulated in Cassandras et al. (2002) grew out of the concept of the fluid queue, and subsequently was extended to flow networks and beyond to a general setting of Stochastic Hybrid Systems (SHS); see Cassandras et al. (2010).

Consider Figure 1 for an illustration of the SHS concept. The system in question is the DEES shown in the figure, and the control parameter assigned to it is $\theta \in R^n$. The system with the particular control variable θ generates a sample path which is fed to two places: (i) an algorithm

² To simplify the exposition in the forthcoming discussion, we omit the notational dependence of the sample function $L_T(\theta)$ and its mean $J_T(\theta)$ on T .

which computes $L(\theta)$ and its IPA gradient $\nabla L(\theta)$, and (ii) a continuous-flow modeling artifact, indicated by SHS in the figure. An algorithm which is based on SHS computes the value of another sample performance function, $L_s(\theta)$, and its IPA gradient $\nabla L_s(\theta)$. It is pointed out that the same sample path, associated with the underlying DEES is used to compute the sample performance functions associated with both the discrete system and the hybrid system, as well as their respective IPA gradients.

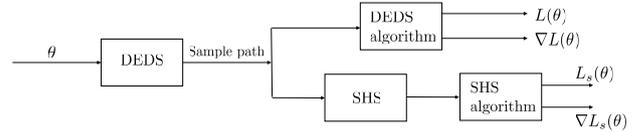


Fig. 1. Relationship between a DEES and its SHS approximation

Suppose that $\nabla L(\theta)$ is a biased estimator of $\nabla J(\theta)$, and hence it does not provide a good approximation to it. The SHS framework is useful as long as $\nabla L_s(\theta)$ provides a close approximation to $\nabla J(\theta)$, which often is the case when the IPA derivative $\nabla L_s(\theta)$ is unbiased. It must be pointed out that approximating $J(\theta)$ usually is a major concern in performance evaluation or estimation, whereas approximations of $\nabla J(\theta)$ is a concern in optimization and control. Since our interest is in the latter but not the former, we have no intrinsic interest in the quality of approximations provided by $L_s(\theta)$ to $J(\theta)$. In fact, extensive testing by simulation has shown that such approximations can be inadequate for the purpose of performance evaluation, whereas the approximation of $\nabla J(\theta)$ provided by $\nabla L_s(\theta)$ suffice to achieve optimization and control objectives. Examples of this point can be found in Cassandras et al. (2002), Cassandras et al. (2010).

The first example analyzed in this context concerns a continuous-flow single-server queue with a finite buffer, where fluid arriving at a full queue is being discarded as a matter of overflow. The basic stochastic-flow modelling construct, depicted in Figure 2, is defined as follows: Given a time-horizon $[0, T]$, let $\{\alpha(t)\}$ and $\{\beta(t)\}$ be the instantaneous fluid-arrival rate to the queue, and the instantaneous fluid service rate at the queue, respectively. These are assumed to be exogenous stochastic processes defined on a common probability space (Ω, \mathcal{F}, P) . Let $b > 0$ denote the size of the buffer. Denote by $X(t)$ and $\gamma(t)$ the instantaneous buffer-contents (amount of fluid in the buffer) and instantaneous spillover rate from the queue due to overflow. Then the processes $\{X(t)\}$ and $\{\gamma(t)\}$ are defined as follows (see Cassandras et al. (2002)):

$$\frac{dX(t)}{dt^+} = \begin{cases} 0, & \text{if } X(t) = 0, \text{ and } \alpha(t) \leq \beta(t) \\ 0, & \text{if } X(t) = b \text{ and } \alpha(t) \geq \beta(t), \\ \alpha(t) - \beta(t), & \text{otherwise,} \end{cases} \quad (9)$$

and

$$\gamma(t) = \begin{cases} \alpha(t) - \beta(t), & \text{if } X(t) = b \\ 0, & \text{if } X(t) < b. \end{cases} \quad (10)$$

A typical control variable consists of a parameter of the arrival-rate process, the service-rate process, or the buffer size. For example, $\beta(\theta; t) = \theta\beta(t)$, where θ represents a controlled flow parameter and $\{\beta(t)\}$ is an exogenous process depending on time t but not on θ . This can arise in communication networks where θ represents the total

transmission rate of a channel and $\beta(t)$ is the fraction of it which is allocated to a particular session. Observe the notation $\beta(\theta; t)$, indicating that the instantaneous service rate depends on θ . As a result, the buffer-occupancy and spillover-rate processes depend on θ as well via (9) and (10), and hence denoted by $\{X(\theta; t)\}$ and $\{\gamma(\theta; t)\}$. Ref. Cassandras et al. (2002) considered various sample performance functions of such control variables. We next present the first SFM analyzed from the standpoint of IPA, which exhibits the salient features of an extensive suite of fluid queueing networks.

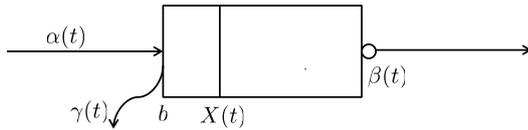


Fig. 2. Elemental fluid-queue model

Consider the case where the control parameter is the buffer capacity, and the performance function is the amount of fluid which is discarded during a given horizon interval $[0, T]$. Thus, $\theta = b$, and the performance function, denoted by $L_\gamma(\theta)$, is

$$L_\gamma(\theta) := \int_0^T \gamma(\theta; t) dt. \quad (11)$$

Note that $L_\gamma(\theta)$ is related to the fraction of discarded fluid from the total arrival volume during the time $t \in [0, T]$, which is $L_\gamma(\theta) / \int_0^T \alpha(t) dt$.

Regarding the IPA derivative $\frac{dL_\gamma(\theta)}{d\theta}$, we note that it is unbiased since the function $L_\gamma(\theta)$ is continuous. The following formula for it was obtained in Cassandras et al. (2002): Define N_T as the number of lossy busy periods in the horizon interval $[0, T]$, namely the number of busy periods during which the buffer becomes full at some time. Then,

$$\frac{dL_\gamma(\theta)}{d\theta} = -N_T. \quad (12)$$

As an example, consider the realization of the state trajectory $\{X(\theta; t)\}$ shown in Figure 3. It is evident that the first and third busy periods are lossy while the second in not, therefore $\frac{dL_\gamma(\theta)}{d\theta} = -2$.

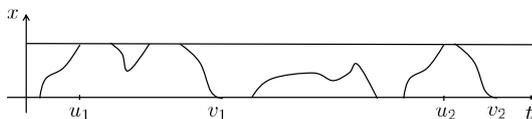


Fig. 3. Typical state trajectory $\{X(\theta; t)\}$

As another example, consider the cumulative workload as a function of the buffer capacity. Thus, $\theta = b$ as in the previous example, and let the performance function, denoted by $L_x(\theta)$, be defined as

$$L_x(\theta) = \int_0^T X(\theta; t) dt. \quad (13)$$

Note that the term $L_x(\theta) / \int_0^T \alpha(t) dt$ serves to approximate the average delay of fluid “molecules” by Little’s law. As for the IPA derivative, fix $\theta > 0$, and let B_m , $m = 1, \dots, M$ denote the lossy busy periods in the interval $[0, T]$ in

increasing order. For every $m = 1, \dots, M$, let $u_m \in B_m$ denote the first time the buffer becomes full in B_m , and let v_m be the end-time of B_m . Then $\frac{dL_x(\theta)}{d\theta}$ has the following form,

$$\frac{dL_x(\theta)}{d\theta} = \sum_{m=1}^M (v_m - u_m). \quad (14)$$

For example, in Figure 3, $\frac{dL_x(\theta)}{d\theta} = \sum_{m=1}^2 (v_m - u_m)$.

We point out that Eqs. (12) and (14) were derived (in Cassandras et al. (2002)) under minimal assumptions on the system. In fact, the only assumption made is that the processes $\{\alpha(t)\}$ and $\{\beta(t)\}$ be piecewise continuous and of bounded variation in $t \in [0, T]$ w.p.1. Therefore, for the purpose of computing Eqs. (12) and (14), these processes can be generated not only from an SFM but also from a DEDS. Such DEDS can be a modeling artifact of the system or the system itself. Furthermore, the formulas (12) and (14) do not depend on observations of the detailed dynamics associated with arrivals or departures of each customer at the queue, nor on the instantaneous values of $\alpha(t)$ or $\beta(t)$. Instead, they require only the observations of macro events like the beginning and end of busy periods and full-buffer periods. Thus, comparing and contrasting these formulas with the analogous equations derived for IPA in the traditional, discrete-queueing setting (e.g., Cassandras and Lafortune (2008)), we see that the SFM-based formulas are unbiased, simpler, and do not require any details of the probability laws underscoring the arrival and service processes. Due to the last property we say that the IPA derivative is *nonparametric*. Furthermore, as mentioned above, the IPA derivatives can be computed from data generated from DEDS as well as SFM. In fact, Cassandras et al. (2002) and subsequent papers on IPA in the SFM setting (e.g., Zhang and Cassandras (2002), Sun et al. (2004), Cassandras et al. (2010) and references therein) report on successful solutions of DEDS optimization problems where the IPA gradients are computed from SFM-derived formulas. All of this suggests that the SFM setting provides an alternative framework to DEDS for the application of IPA, which holds out promise of real-time optimization via closed-loop control.

As we mentioned earlier, following the basic formulation of the SFM and derivation of the aforementioned IPA gradients, there were several efforts to extend the model and results to fluid queueing systems and other multiflow networks. The main results concern the development of IPA gradients for prototypical problems, and their implementation in optimization environments. By and large the simplicity and unbiasedness of the IPA gradients are maintained. The nonparametric property and reliance only on observations of macro events has been almost maintained, but held ground for close approximations where, practically, the errors can be neglected. Moreover, the key structure of perturbation propagation, which rendered IPA attractive from its onset, is maintained in fluid-flow networks. This is all summarized in Cassandras et al. (2010) and formulated in a general framework of *Stochastic Hybrid Systems* (SHS) based on a formal calculus (referred to as the *IPA Calculus*) of event-driven propagations. We close this section by mentioning that the SHS framework can serve not only as an approximation of a DEDS but also as a primary model of systems with discrete and

continuous dynamics that arise in various application areas of current interest.

4. RECENT AND CURRENT TRENDS

This section presents some of the main research directions in IPA which emerged during the past decade. In particular we discuss applications to large-scale Markov processes and stochastic hybrid systems, performance regulation of systems, and the use of event-driven (as opposed to time-driven) methods for control and optimization.

4.1 IPA of Markov Systems and Stochastic Optimization

Until the mid 1990s, IPA was largely limited to “infinitesimal” perturbations, and could not be applied to perturbations with finite size. Around this time, however, it was realized that the perturbation realization principle applies to finite jumps of states as well. This made it possible to develop IPA algorithms for Markov processes.

Consider an irreducible and aperiodic Markov chain $\mathbf{X} = \{X_n : n \geq 0\}$ on a finite state space $\mathcal{S} = \{1, 2, \dots, M\}$ with transition probability matrix $P = [p(j|i)] \in [0, 1]^{M \times M}$. Let $\pi = (\pi_1, \dots, \pi_M)$ be the vector representing its steady-state probabilities, and $f = (f_1, f_2, \dots, f_M)^T$ be the performance vector, where “T” represents its transpose. We have $Pe = e$, where $e = (1, 1, \dots, 1)^T$ is an M -dimensional vector whose all components equal 1, and $\pi = \pi P$. The performance measure is the long term average defined as

$$\begin{aligned} \eta &= \sum_{i=1}^M \pi_i f_i = \pi f = \lim_{L \rightarrow \infty} \frac{1}{L} \sum_{l=0}^{L-1} f(X_l) \\ &= \lim_{L \rightarrow \infty} \frac{F_L}{L}, \quad w.p.1, \end{aligned} \quad (15)$$

where $F_L := \sum_{l=0}^{L-1} f(X_l)$.

Let P' be another irreducible transition probability matrix on the same state space. Suppose P changes to $P(\delta) = P + \delta Q = \delta P' + (1-\delta)P$, with $\delta > 0$, $Q = P' - P = [q(j|i)]$, and the reward function f keeps the same. We have $Qe = 0$. The performance measure will change to $\eta(\delta) = \eta + \Delta\eta(\delta)$. The derivative of η in the direction of Q is defined as

$$\frac{d\eta(\delta)}{d\delta} = \lim_{\delta \rightarrow 0} \frac{\Delta\eta(\delta)}{\delta}. \quad (16)$$

In this system, a perturbation means that the system is perturbed from one state i to another state j . Following the same idea as in (7), we study two independent Markov chains $\mathbf{X} = \{X_n; n \geq 0\}$ and $\mathbf{X}' = \{X'_n; n \geq 0\}$ with $X_0 = i$ and $X'_0 = j$; both of them have the same transition matrix P . The *realization factor* is defined as Cao (2007):

$$\begin{aligned} d(i, j) &= \lim_{L \rightarrow \infty} E \left[\sum_{l=0}^{L-1} (f(X'_l) - f(X_l)) \middle| X_0 = i, X'_0 = j \right], \\ & \quad i, j = 1, \dots, M. \end{aligned} \quad (17)$$

Thus, $d(i, j)$ represents the average effect of a jump from i to j on F_L in (15). From (17), it is easy to see that $d(i, j)$ satisfies the conservation law as in physics:

$$g(i, k) = g(i, j) + g(j, k), \quad i, j, k \in \mathcal{S},$$

Thus, we can define a vector $g = (g(1), g(2), \dots, g(M))^T$, called *performance potential*, such that

$$d(i, j) = g(j) - g(i), \quad i, j \in \mathcal{S},$$

and we can verify that it satisfies the *Poisson equation*

$$(I - P + e\pi)g = f, \quad (18)$$

where I is the $M \times M$ identity matrix.

Multiplying both sides of the Poisson equation with π on the left, we get

$$\pi g = \pi f = \eta. \quad (19)$$

Multiplying both sides of the Poisson equation with π' on the left yields

$$\pi' Qg = \pi'(P' - P)g = \pi'(I - P)g = \pi' f - \pi g = \pi' f - \eta.$$

That is,

$$\eta' - \eta = \pi' Qg. \quad (20)$$

Setting $P(\delta) = P + \delta Q$ and $\eta' = \eta(\delta)$ and letting $\delta \rightarrow 0$ in (20), we get the desired performance derivative along the direction Q :

$$\frac{d\eta(\delta)}{d\delta} = \pi Qg. \quad (21)$$

The derivatives can be used in performance optimization (Marbach and Tsitsiklis (2001)). Efficient algorithms can be derived for estimating g and estimating the derivative $\frac{d\eta(\delta)}{d\delta}$ directly (Cao (2007)). Now, there is a new subarea in reinforcement learning, called policy gradients, devoted to developing algorithms for the performance derivative (21), see e.g., Baxter and Bartlett (2001), Baxter et al. (2001), etc.

If the reward function also changes from f to f' , let $h := f' - f$. It is easy to check that

$$\eta' - \eta = \pi'(Qg + h) = \pi'\{(P'g + f') - (Pg + f)\}. \quad (22)$$

This is the Performance Difference Formula (PDF); it initiates a new direction in performance optimization, the *direct-comparison based approach*. In fact, it is observed that the PDF contains all the information in comparing the performance of any two policies, and an optimality condition can be simply derived from this equation without dynamic programming or discounting for long-run average performance. For example, because $\pi' > 0$, from (22), we conclude

$$\text{If } P'g + f' \leq Pg + f, \text{ then } \eta' \leq \eta. \quad (23)$$

This leads to the optimality condition: a policy (P^*, f^*) with potential g^* is optimal if and only if $Pg^* + f \leq P^*g^* + f^*$ for all policies P . Policy iteration algorithms can also be developed from (22).

The Direct-Comparison (DC) based approach is an alternative to dynamic programming (DP) to performance optimization of dynamic systems. As illustrated above, this approach is very simple and intuitive for long-run average performance; in fact, a complete theory based on n th bias optimality for long-run average performance can be developed with no discounting (Cao (2007)). Next, the PDF provides global information to performance comparison in the entire period; while dynamic programming works backwards in time at a particular time instant (continuous or discrete), and hence it only provides local information. Therefore, the DC-based approach opens a new horizon for problems requiring global considerations.

For example, the approach naturally solves a long existing issue in time non-homogenous Markov systems, the under selectivity, which means that in performance optimization of time non-homogenous systems, where the transition probabilities and reward functions are different at different time $k = 1, 2, \dots$, the long-run average performance, and for that matter its optimal policy, does not depend the actions (transition probabilities and rewards) in any finite periods (Cao (2015)). The approach has also been applied to stochastic control problems with diffusion processes (continuous time and continuous states); it solves the control problem with non-smooth value functions without resorting to viscosity solutions (Cao (2017)).

Equations (21) and (22) provide a sensitivity-based view to performance optimization. For problems where the performance is not additive, (21) may be used. The DC-based approach links both together naturally, Research in this direction is ongoing and the influence of the sensitivity-based view extends beyond the area of DEDS.

Last but not the least, combined with the aggregation technique, the DC-based approach leads to the theory of *event-based optimization (control)*, in which control actions depend on events rather than the states. This may dramatically reduce the computation since the number of events (event space) is much smaller than that of states (state space). Conditions have been derived under which Hamilton-Jacobi-Bellman (HJB) type of optimality equation holds for event-based control; because the sequence of events is not Markov and aggregation is used, approximation is usually involved, see Cao (2007); Xia et al. (2014); other related topics will be discussed in Section 4.3.

4.2 Performance Regulation

Emerging applications of IPA concern the tracking of a reference input to a dynamical system by its output process. An abstract, discrete-time single-input-single-output system is depicted in Figure 4, where k denotes time, $r \in R$ is the reference input, y_k is the system's output, u_k is the control input to the plant, and $e_k := r - y_k$ is the error signal. The plant generally is a time-varying dynamical system lacking an accurate model and subjected to unpredictable variations.

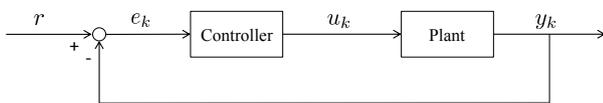


Fig. 4. Control System

As an example of interest, it is desirable to regulate the instructions' throughput of a computer processor by adjusting its clock rate, or frequency. More specifically, the time axis is divided into contiguous periods called *control cycles*, during each of which the frequency is set (fixed) and the average throughput is measured. At the end of the control cycle the frequency is changed by the controller according to the difference between the given target-throughput (setpoint) and the average throughput. In this setting the time counter k indicates the index of the control cycle denoted henceforth by C_k , u_k is the value of the clock frequency during C_k , and y_k is the average instruction-throughput computed during C_k .

The plant in Figure 4 is the processor, and any model thereof would describe the frequency-to-throughput relationship. Since we enact a real-time control, there is no need for a model to close the loop since the output y_k is measured. However, we shall see that a model is needed in order to implement the controller that we have in mind. An established model of an out-of-order architecture is provided by a queueing system (see Hennessey and Patterson (2012), or a simplified exposition in Wardi et al. (2016)) which defies analysis. We use it nonetheless in an effective way, as described below.

Returning to the abstract system in Figure 4, the objective is to design a controller which can deliver the desired tracking without a detailed knowledge of the plant-model while facing wide-ranging variations in the system's input-output relationships. Moreover, the controller has to achieve that in very short time-frames and hence by simple computations. Tracking typically involves an integrator in the loop, and to have the controller be as simple as possible we first considered a standalone integrator. Now it is well known that a standalone integrator may destabilize the closed-loop system and otherwise have poor stability margins. Furthermore, to be effective its gain may have to be determined by data, gathered off line, concerning the system's response. However, we cannot obtain such meaningful data due to the unpredictable variability in a processor's workload during program executions. Therefore we adopted a *variable-gain integrator*, whose gain is recomputed at the beginning of each control cycle as a part of the control loop, hence based only on measurements, in a way that extends the stability margins and provides the desirable tracking. In fact, simulation testing showed that this obviates the addition of a proportional element to the controller.

The controller has the following form,

$$u_k = u_{k-1} + A_k e_{k-1},$$

where u_k is the control variable set at the start of C_k , A_k is computed from measurements made during C_{k-1} and hence available at the start of C_k , $e_{k-1} = r - y_{k-1}$, and y_{k-1} is computed from measurements during C_{k-1} . When the gain A_k is independent of $k = 1, 2, \dots$ we recognize this as an adder, a discrete-time equivalent of an integrator. The gain A_k is computed by the following formula,

$$A_k = A_{k-1} + \frac{1}{\frac{\partial y_{k-1}}{\partial u_{k-1}} + \eta_{k-1}} e_{k-1}. \quad (24)$$

The output y_{k-1} depends not only on u_{k-1} but possibly also on noise and other exogenous processes as well as past output like y_{k-2} , etc. These variables are not factored in the term $\frac{\partial y_{k-1}}{\partial u_{k-1}}$ which hence literally stands for the partial derivative.

The term $\frac{\partial y_{k-1}}{\partial u_{k-1}}$ has to be estimated in real time during C_{k-1} as a part of the control loop. However, in the computer application described above, and in other DEDS and SHS, we were unable to compute it due to the absence of analytical models for the plant. Therefore, in Equation (24), we allow for an additive error, η_{k-1} , in its computation. Convergence results of the resulting tracking algorithm, derived in Wardi et al. (2016), account for the presence of such error terms. Simulation tests verify these results with substantial relative errors, which can be 30%

or higher. In other words, the performance of the regulation technique is robust with respect to computational errors in the loop. Leveraging this robustness, we have estimated $\frac{\partial y_{k-1}}{\partial u_{k-1}}$ by IPA. But unlike unbiasedness and exact computation, which have been principal concerns in the use of IPA throughout much its development, we were primarily concerned with fast computations while allowing for substantial bias and computational errors.

Results of simulation experiments can be found in Wardi et al. (2016) and references therein. These include queueing networks, Petri nets, transportation models, and other DEDES. Of a particular interest is the case where the plant is a queueing system with biased IPA. For the original problem of interest, namely instruction-throughput regulation in computer processors, we used a detailed system-level simulation platform for computer architectures, called *Manifold* (Yalamancili et al. (2016)). IPA is biased, and we also induced further errors deliberately in order to simplify the computations of its derivative. Lately we implemented the regulation technique on Intel's fourth-generation microarchitecture, Haswell, and tested it on industry-benchmark programs (Hammarlund et al. (2014)). In this implementation we actually adopted a simpler computation of $\frac{\partial y_{k-1}}{\partial u_{k-1}}$ than IPA can provide, one that is based on linear approximation. Various results can be found in (Chen et al. (2016)).

To summarize, the technique described in this subsection extends the research area in IPA in two directions. First, it explores applications to systems' performance regulation rather than optimization. Second, it does not pursue the objectives of unbiased gradient estimates and their precise computations, but rather seeks simple control laws with fast computations in the loop.

4.3 Event-Driven Control and Optimization

The emergence of DEDES in the 1980s brought to the forefront an alternative viewpoint to the traditional *time-driven* paradigm in which time is an independent variable and, as it evolves, so does the state of a dynamic system. The *event-driven* paradigm offers an alternative, complementary look at modeling, control, communication, and optimization (Miskowicz (2015), Cassandras (2014)). The key idea is that a clock should not be assumed to dictate actions simply because a time step is taken; rather, an action should be triggered by an "event" specified as a well-defined condition on the system state or as a consequence of environmental uncertainties that result in random state transitions. Observing that such an event could actually be defined to be the occurrence of a "clock tick", it follows that this framework may in fact incorporate time-driven methods as well. On the other hand, defining the proper "events" requires more sophisticated techniques compared to simply reacting to time steps. In the development of DEDES, such events were seen as the natural means to drive the dynamics of a large class of systems including computer networks, manufacturing systems, and supply chains among many. By the early 1990s, however, it became evident that many interesting dynamic systems are in fact "hybrid" in nature, i.e., at least some of their state transitions are caused by (possibly controllable) events. This has been reinforced by technological advances through which

sensing and actuating devices are embedded into systems allowing physical processes to interface with such devices which are inherently event-driven. More recently, the term *Cyber-Physical System* (CPS) has emerged to describe the hybrid structure of systems where some components operate as physical processes modeled through time-driven dynamics, while other components (mostly digital devices empowered by software) operate in event-driven mode.

Moreover, many systems of interest are now networked and spatially distributed. In such settings, especially when energy-constrained wireless devices are involved, frequent communication among system components can be inefficient, unnecessary, and sometimes infeasible. Thus, rather than imposing a rigid time-driven communication mechanism, it is reasonable to seek instead to define specific events which dictate when a particular node in a network needs to exchange information with one or more other nodes. When, in addition, the environment is stochastic, significant changes in the operation of a system are the result of random event occurrences, so that, once again, understanding the implications of such events and reacting to them is crucial. In distributed systems, event-driven mechanisms have the advantage of significantly reducing communication among networked components without affecting desired performance objectives. In multi-agent systems where the goal is for networked components to cooperatively maximize (or minimize) a given objective, it is shown in Zhong and Cassandras (2010) that an event-driven scheme can still achieve the optimization objective while drastically reducing communication (hence, prolong the lifetime of a wireless network), even when delays are present (as long as they are bounded). Event-driven approaches are also attractive in receding horizon control, where it is computationally inefficient to re-evaluate a control value over small time increments as opposed to event occurrences defining appropriate planning horizons for the controller. Finally, as already pointed out in Section 4.1, the use of event-driven optimization methods has the benefit of scaling with the size of the event-space and not the (generally much larger) state space of a system.

In Section 3, we discussed how IPA is used in the control and optimization of SHS based on the general-purpose IPA Calculus (Cassandras et al. (2010)). However, even when a hybrid system is studied in a deterministic setting, IPA proves extremely useful in evaluating performance gradients *on line* that can be used for the purpose of optimizing the operation of complex multi-agent systems. These are commonly modeled as hybrid systems with time-driven dynamics describing the motion of the agents or the evolution of physical processes in a given environment, while event-driven behavior characterizes events that may occur randomly (e.g., an agent failure) or in accordance with control policies (e.g., an agent stopping to sense the environment or to change directions). As such, a multi-agent system can be studied in the context of the IPA Calculus with parameterized controllers aiming to meet certain specifications or to optimize a given performance metric. In some cases, the solution of a multi-agent dynamic optimization problem is reduced to a policy which is naturally parametric. Therefore, IPA may be used to evaluate on line performance gradients through which one can drive the system towards optimal (at least locally)

points; recent examples of this approach may be found in Cassandras et al. (2013), Zhou et al. (2016).

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