

The optimizing mode classification stabilization of sampled stochastic jump systems via an improved hill-climbing algorithm based on Q-learning

Guoliang Wang

Abstract—This paper addresses the stabilization problem of stochastic jump systems (SJSs) closed by a generally sampled controller. Because of the controller's switching and state both sampled, it is challenging to study its stabilization. A new stabilizing method deeply depending on the mode classifications is proposed to deal with the above sampling situation, whose quantity is equal to a Stirling number of the second kind. For the sake of finding the best stabilization effect among all the classifications, a convex optimization problem is developed, whose globally solution is proved to be existent and can be computed by an augmented Lagrangian function. More importantly, in order to further reduce the computation complexity but retaining a better performance as much as possible, a novel improved hill-climbing algorithm is established by applying the Q-learning technique to provide an optimal attenuation coefficient. A numerical example is offered so as to verify the effectiveness and superiority of the methods proposed in this study.

Index Terms—Stochastic jump systems; sampled control; mode classification and optimization; Lagrangian function; hill-climbing algorithm; Q-learning.

I. INTRODUCTION

SIGNIFICANTLY different from deterministic systems, stochastic jump system (SJS) can represent physical systems experiencing random structure changes. Due to this system with multiple structures or modes, it makes its stabilization problems quite distinctive whose controller's quantity is not unique. According to the designed controller depending on mode or not, the existing stabilization results are mainly classified into two categories. The first kind needs the mode of controller to keep pace with others and are usually called to be mode-dependent control method such as [1]–[6]. Due to all the modes synchronized with each other at the same time, the effect of stabilization effect will be the best and have the least conservatism. However, this advantage is also its disadvantage which will be limited in practice, since a lot of effort is needed to keep this synchronization all the time. In theoretical research, an ideal assumption for all the mode information available in real time is commonly acquiescent. The second kind is mentioned as mode-independent control approach [7]–[10], in which the mode information is totally removed even if

it is available sometimes. Because the mode information is totally removed in controller, it can stabilize a system regardless of its mode accessible or not and naturally bring more conservatism. In contrast to mode-dependent control being very ideal, mode-independent control is excessively absolute due to mode information neglected completely. For the purpose of bridging the above methods and balancing their advantages and disadvantages, some improved controllers were developed, such as partially mode-dependent controller [11] and partial information controller [12]. However, the drawbacks of mode-dependent and -independent control methods have not been partly solved, and there are still some problems to be further studied. For example, the quantity of controllers denoted as M in the above references is no less than number of modes referred to be N . Consequently, an interesting problem about SJSs can be proposed such that whether one can stabilize an SJS using fewer controllers but more than one. In this situation, some results were found, see, e.g., disordered or unmatched controller [13], [14], and scheduling controller [15], [16]. Though the aim of fewer controllers stabilizing an SJS with more subsystems was realized, it can be found from these references that for each subsystem or mode, at most total M controllers were added once, while the equipment cost and conservatism of stabilization realizing were both increased. The main reason is that the developed controller in these references was designed on the whole modes without considering them separately. When the controller is designed according to the mode classifications, a quantity limited controller method was in proposed in [17]. However, the mode classification method for designing controllers is not fully studied, and many interesting and significant topics need to be further researched. For example, whether there is an optimization classification or not and how to find the best mode classification will be the first problems, and more extensions based on optimizing mode classification method will also be meaningful.

On the other hand, it is well known that networked control system (NCS) is a control system whose components are connected through a communication network. Though NCS has such advantages, there is a precondition that all the transmitted data should be first sampled. Importantly, the appearance of sampled data complicates system analysis and synthesis and causes many unpredictable problems. It can further bring some negative effects such as communication delays, data packet dropouts and/or packet disordering, network attack, some of which are not easy. Thus, it is very important to research the

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G. Wang is with School of Information and Control Engineering, Liaoning Petrochemical University, Fushun 113001, Liaoning, China (e-mail: glwang@lnpu.edu.cn).

sampling phenomenon scientifically. Particularly, when an SJS is connected by a network, some novel but hard issues will encounter such that not only state but also switching signal are sampled. Due to switching being stochastic, its sampling is specific and significantly different from state sampled. Accordingly, some interesting but challenging problems emerge. Up to now, very few results are found to study the stabilization of stochastic systems via a sampled-switching controller. On the one hand, the author in [18] first considered its stabilization problem by applying an auxiliary system approach. More extensions [19], [20] were further obtained base in this method. Since the stabilization problem of original sampling system was transformed to study its auxiliary system without any sampling, it decided that the sampling bound should be very small. On the other hand, based on an augmented system method, the stabilization of Markovian jump systems (MJSs) closed by a sampled switching and state controller was studied in [21], whose sampling effect was modeled to be a time-varying exponential matrix. Unfortunately, the convergency guaranteed by the reference was only asymptotically mean stable and was worse than the common stability concepts such as almost surely exponentially stable, globally asymptotically stable and so on. To summarize, it is necessary to further investigate the stabilization problem of sampled stochastic jump systems. Moreover, new approaches are expected to provide larger sampling bounds and better convergency properties. Meanwhile, many difficulties will encounter in the research, some of which are challenging. For example, how to establish a quantized correlation between the original switching and its sampled value will be the first difficulty to encounter, since no more suitable models are available. Particularly, it will be challenging to achieve the above expected objectives, because both switching and state are sampled simultaneously. Especially so many possible combinations of switching and its sampled values will emerge and bring large difficulties in system analysis and synthesis.

The main contributions of this paper are as follows: 1) For the aim of overcoming the control difficulties of sampling in both state and switching signals, a new stabilizing controller is established to be very closed to the mode classifications. On the one hand, the controller's quantity is smaller than mode-dependent controllers, and its switching is not necessary synchronous to the original one. On the other hand, the conservatism is smaller than mode-independent ones [7]–[10]; 2) So as to further improve the stabilization performance, a convex optimization problem is presented to determine the best classification of sampled stabilization, which is better than [17] without mode optimization. It can be shown that the optimal solution can exist and be obtained by computing some equations coming from an augmented Lagrangian function; 3) Due to the classification quantity being a Stirling number of the second kind and large, a novel method having less complexity but better control performance is proposed based on the hill-climbing algorithm by using the Q-learning technique to ensure an optimal attenuation coefficient. Particularly, not only the monotonicity but also the convergency of method in this

paper is guaranteed; 4) Compared with the existing auxiliary system approach [18] and augmented system method [21], the developed method is less conservative and has larger sampling bounds. Moreover, the key idea in this paper can popularize in many situations as long as the system switching experiences a sampling or mismatching phenomenon.

Notation \mathbb{R} , $\mathbb{R}_{>0}$, $\mathbb{R}_{\geq 0}$ and $\mathbb{N}_{>0}$ represent the sets of real numbers, positive real numbers, non-negative real numbers and positive integers respectively. \mathbb{R}^n denotes the n -dimensional Euclidean space, and $\mathbb{R}_{\geq 0}^N$ is a set of vector whose each element is non-negative. $\mathbb{P}(\cdot)$ and $\mathbb{E}(\cdot)$ are the probability and expectation operators respectively. $\|\cdot\|$ refers to the Euclidean vector norm or spectral matrix norm. $\lambda_{\min}(M)$ and $\lambda_{\max}(M)$ denote the smallest and largest eigenvalues of a square matrix M , $\chi_M \triangleq \lambda_{\max}\left(\frac{M+M^T}{2}\right)$ and $(M)^* \triangleq M + M^T$.

II. PROBLEM FORMULATION

Consider a stochastic jump system described as

$$\dot{x}(t) = A_{\eta(t)}x(t) + B_{\eta(t)}u(t) \quad (1)$$

where $x(t) \in \mathbb{R}^n$, $A_{\eta(t)} \in \mathbb{R}^{n \times n}$ and $B_{\eta(t)} \in \mathbb{R}^{n \times m}$ and $u(t) \in \mathbb{R}^m$. The stochastic switching process $\{\eta(t), t \geq 0\}$ is a piecewise constant function and right-continuous. In detail, it is actual the semi-Markovian switching and takes values from a set $\mathcal{N} \triangleq \{1, 2, \dots, N\}$ such as $\eta(t) = i \in \mathcal{N}$, $\forall t \in [T_n, T_{n+1})$, $n \in \mathbb{N}$, where switching instant T_n satisfies $0 = T_0 < T_1 < \dots < T_n < \dots$. For simplicity, $\eta(T_n) = i$ at the n th switching instant is simply denoted to be $\eta_n = i$.

In contrast to the usual controllers designed for stochastic systems including (1), a general controller with sampling phenomenon is described to be

$$u(t) = K_{\eta(t_k)}x(t_k), \forall t \in [t_k, t_{k+1}), \forall k \in \mathbb{N} \quad (2)$$

where $K_{\eta(t_k)} \in \mathbb{R}^{m \times n}$ is the control gain, t_k is the sampling instant such as $0 = t_0 < t_1 < \dots < t_k < \dots$, and $s_k \triangleq t_k - t_{k-1}$ is denoted as the sampling interval. Obviously, the simultaneous existence of two instant sequences $\{T_n\}_{n \in \mathbb{N}}$ and $\{t_k\}_{k \in \mathbb{N}}$ makes the system analysis and synthesis not easy. Especially, due to $\eta(t)$ being sampled, it will bring large difficulties and also complicates the closed-loop system. The first reason is that so many combinations about original signal $\eta(t)$ and its sampled signal $\eta(t_k)$ encounter and inevitably result in great complexity and large conservatism. The second reason, but not the last, is that the sampled state $x(t_k)$ existing simultaneously also leads to negative effects such that the correlation between the original and sampled states on each subinterval is hard to be done. In order to solve the above mentioned problems, the controller is developed as

$$u(t) = \sum_{\ell=1}^N \alpha_{\ell}^{[\nu(\eta(t_k))]} \bar{K}_{\ell} x(t_k), \forall t \in [t_k, t_{k+1}), \forall k \in \mathbb{N} \quad (3)$$

where $\alpha_{\ell}^{[\nu(\eta(t_k))]} \in \mathbb{R}$ and $\bar{K}_{\ell} \in \mathbb{R}^{m \times n}$ are to be designed. Particularly, the sampling instant t_k of controller (3) is differ-

ent from (2). In detail, the sampling instant of (3) is event-triggered such as

$$t_{k+1} = \min_{t \geq t_k} \{t : \eta(t_k) \in \mathcal{N}_h, \eta(t) \notin \mathcal{N}_h\} \quad (4)$$

Accordingly, function $\nu(\eta(t))$ on interval $\forall t \in [t_k, t_{k+1})$ with property (4) is defined as $\nu(\eta(t)) \equiv \nu(\eta(t_k)) = h$, when $\eta(t) \in \mathcal{N}_h$. Meanwhile, set \mathcal{N}_h is a newly constructed subset based on \mathcal{N} and defined as

$$\mathcal{N}^D = \{\mathcal{N}_1, \mathcal{N}_2, \dots, \mathcal{N}_M\} \quad (5)$$

where $\mathcal{N}_h \subseteq \mathcal{N}$, $h \in \mathcal{M} \triangleq \{1, 2, \dots, M\}$, and constant M is the quantity of elements of set \mathcal{N}^D and denoted as $\text{num}(\mathcal{N}^D) = M \in \mathbb{N}_{>0}$. Moreover, all the elements of set \mathcal{N}^D are mutually exclusive such as

$$\mathcal{N}_h \cap \mathcal{N}_\ell = \emptyset, \forall h \neq \ell \in \mathcal{M} \quad (6)$$

Then, it can be concluded that

$$\mathcal{N} = \bigcup_{h=1}^M \mathcal{N}_h \text{ and } \sum_{h=1}^M n_h = N \quad (7)$$

where $n_h \triangleq \text{num}(\mathcal{N}_h) \in \mathbb{N}_{>0}$. Moreover, subset \mathcal{N}_h can be further expressed as

$$\mathcal{N}_h = \{i_1^{[\mathcal{N}_h]}, i_2^{[\mathcal{N}_h]}, \dots, i_{n_h}^{[\mathcal{N}_h]}\} \quad (8)$$

where $i_\ell^{[\mathcal{N}_h]} \in \mathcal{N}$, $\ell = 1, 2, \dots, n_h$. When the above mentioned classification and event-trigger scheme achieve, controller (3) will be superior to (2), which can be realized easily and of larger significance in theory and practice. Moreover, a mapping between M controllers and N subsystems or modes should be introduced. In the next, it will be seen that this mapping can not only be used to design a new controller for stochastic systems but also bridge the traditionally mode-dependent and -independent controllers very well.

REMARK 1: It is noted that controller (3) includes some existing controllers without any sampling signals as special situations in which neither $x(t)$ or $\eta(t)$ is sampled. First of all, when $\mathcal{M} = \mathcal{N}$ and $\mathcal{N}_h = \{h\}$, $\forall h \in \mathcal{N}$, in addition to $\sum_{\ell=1}^N \alpha_\ell^{[h]} \equiv 1$ with $\alpha_\ell^{[h]} \in \{0, 1\}$ but $\alpha_h^{[h]} \equiv 1$, controller (3) without sampled state $x(t_k)$ but with $x(t)$ will be simplified to the traditionally mode-dependent controller such as $u(t) = \bar{K}_{\eta(t)} x(t)$. Second, under a deterministic classification same to the first situation, and if $\sum_{\ell=1}^N \alpha_\ell^{[h]} \equiv 1$ with $\alpha_\ell^{[h]} \in (0, 1)$, one could get a partial information controller similar to [12]. Moreover, if $\alpha_\ell^{[h]}$ takes discrete values such as $\alpha_\ell^{[h]} \in \{0, 1\}$, a disordered controller similar to [13] will be obtained. Thirdly, when there is only one element in \mathcal{N}^D such as $\mathcal{N}_1 = \mathcal{N}$, there will be only one element in \mathcal{M} such as $\mathcal{M} = \{1\}$. Then, controller (3) with $x(t_k)$ replaced by $x(t)$ will reduce to be the traditionally mode-independent controller $u(t) = K_1 x(t) = \sum_{\ell=1}^N \alpha_\ell \bar{K}_\ell$. When α_ℓ is equal to the stationary distribution, the mode-independent controller $u(t) = \sum_{\ell=1}^N \alpha_\ell \bar{K}_\ell$ will be an optimal estimation of \bar{K}_ℓ similar to [17], [22].

Though some existing controllers are included as special situations of sampled controller (3), their methods cannot be

used to deal with (3). Obviously, controller (3) deeply depends on the mode classification. Meanwhile, so many possible combinations of the division about set \mathcal{N} are involved whose division value is a positive natural number. Particularly, it can be known from [23] that its value is equal to $S_2(N, M)$ which is a Stirling number of the second kind and described to be

$$S_2(N, M) = S_2(N-1, M-1) + M S_2(N-1, M)$$

Its detailed expansion [24] is given to be

$$S_2(N, M) = \frac{1}{M!} \sum_{v=0}^M (-1)^v \binom{M}{v} (M-v)^N \quad (9)$$

When the classification of (3) is not given in advance, controller (3) should deeply depend on the Stirling number of the second kind and be rewritten as

$$u(t) = \sum_{\ell=1}^N \alpha_\ell^{[\nu(\eta(t_k))]}(c) \bar{K}_\ell x(t_k), \forall t \in [t_k, t_{k+1}), \forall k \in \mathbb{N} \quad (10)$$

where c is selected to be only one value from $c \in \mathcal{C} \triangleq \{1, 2, \dots, S_2(N, M)\}$. Accordingly, the related symbols in definitions (5)-(8) will change to be

$$\begin{aligned} \mathcal{N}^D(c) &= \{\mathcal{N}_1(c), \mathcal{N}_2(c), \dots, \mathcal{N}_M(c)\} \\ \mathcal{N}_h(c) &= \{i_1^{[\mathcal{N}_h(c)]}, i_2^{[\mathcal{N}_h(c)]}, \dots, i_{n_h(c)}^{[\mathcal{N}_h(c)]}\} \end{aligned}$$

where $n_h(c) \triangleq \text{num}(\mathcal{N}_h(c)) \in \mathbb{N}_{>0}$, whose properties (6) and (7) satisfy too.

Then, two necessary but not easy problems are proposed as follows:

Problem 1: Whether does the best classification of set \mathcal{N} exist or not?

Problem 2: How to find the best classification to design a sampled stabilizing controller such as (10)?

First of all, a cost function is introduced to be

$$J(\alpha(c)) = \sum_{h=1}^M \left\| \sum_{\ell=1}^N \alpha_\ell^{[h]}(c) \bar{K}_\ell \right\|_2 \quad (11)$$

where matrix $\alpha(c) \triangleq (\alpha_\ell^{[h]}(c)) \in \mathbb{R}^{M \times N}$ is composed of $\alpha_\ell^{[h]}(c)$ but depends on classification parameter c . Then, the related optimization problem is to find the optimal classification and its related optimal classification parameter $\alpha_\ell^{[h]}(c)$ of sampled controller (10) satisfying

$$\begin{aligned} J^* &= \min_{c \in \mathcal{C}} \min_{\alpha_\ell^{[h]}(c) \in \mathbb{R}} \sum_{h=1}^M \left\| \sum_{\ell=1}^N \alpha_\ell^{[h]}(c) \bar{K}_\ell \right\|_2 \\ &= \min_{c \in \mathcal{C}} \sum_{h=1}^M J^{[h]} \left((\alpha^{[h]}(c))^* \right) = \min_{c \in \mathcal{C}} J(\alpha^*(c)) \end{aligned} \quad (12)$$

where the optimal solution $\alpha^*(c)$ satisfies $J(\alpha^*(c)) \triangleq \min_{\alpha_\ell^{[h]}(c) \in \mathbb{R}} \sum_{h=1}^M \left\| \sum_{\ell=1}^N \alpha_\ell^{[h]}(c) \bar{K}_\ell \right\|_2$. Similarly, one can define $(\alpha^{[h]}(c))^*$ such as $J^{[h]}((\alpha^{[h]}(c))^*) \triangleq \min_{\alpha_\ell^{[h]}(c) \in \mathbb{R}} \left\| \sum_{\ell=1}^N \alpha_\ell^{[h]}(c) \bar{K}_\ell \right\|_2$,

where $J^{[h]}((\alpha^{[h]}(c))) \triangleq \left\| \sum_{\ell=1}^N \alpha_{\ell}^{[h]}(c) \bar{K}_{\ell} \right\|_2 > 0$ means there is always a controller added to each subsystem. It can be seen from (12) that a normal method getting the solution to **Problem 2** is to compute cost function (11) one by one whose process will repeat $S_2(N, M)$ times. After doing that, one can select the solution minimizing the value of (11) as the globally optimal solution. Obviously, when N and M become large, the complexity of computation will be very large. Thus, it is necessary to develop a new way which provides a suboptimal solution nearing the globally optimal solution as much as possible but has smaller complexity without computing (11) one by one. In this paper, an improved hill-climbing algorithm (HCA) based on the Q-learning technique will be developed, which can not only avoid traversing all the classifications but also guarantee the proposed algorithm some good properties such as convergence and monotonicity. As a result, the computation complexity will be largely reduced, while the cost function will also have an optimal value. Some fundamentally dealing processes should be established firstly.

First of all, some constants about classification set $\mathcal{N}^D(c)$ are needed and important in designing the optimization method to be presented. In detail, each element $\mathcal{N}_h(c)$ of $\mathcal{N}^D(c)$ is arranged by h such as $\mathcal{N}^D(c) = \{\mathcal{N}_1(c), \mathcal{N}_2(c), \dots, \mathcal{N}_M(c)\}$. Meanwhile, all the elements of $\mathcal{N}_h(c)$ are rewritten as an ascending sequence such as $i_1^{[\mathcal{N}_h(c)]}(c) < i_2^{[\mathcal{N}_h(c)]}(c) < \dots < i_{n_h(c)}^{[\mathcal{N}_h(c)]}(c)$, since they are natural numbers and distinct from each other. Then, two kind parameters referred to be ξ and ζ about subset $\mathcal{N}_h(c)$ can be generated. On the one hand, parameter ξ is generated by arranging every $n_h(c)$ with h selecting $1, 2, \dots, M$ simultaneously and sequently such as $\xi = n_1(c)n_2(c) \cdots n_M(c)$. On the other hand, similar to ξ , parameter ζ is constructed by arranging all the elements of every subset following an ascending sequence such as

$$\zeta = \underbrace{i_1^{[\mathcal{N}_1(c)]} i_2^{[\mathcal{N}_1(c)]} \dots i_{n_1(c)}^{[\mathcal{N}_1(c)]}}_{\text{subset 1}} \dots \underbrace{i_1^{[\mathcal{N}_M(c)]} i_2^{[\mathcal{N}_M(c)]} \dots i_{n_M(c)}^{[\mathcal{N}_M(c)]}}_{\text{subset M}}$$

Based on these parameters, an one-to-one correlation or mapping between sets $\mathcal{D} \triangleq \{\mathcal{N}^D(1), \mathcal{N}^D(2), \dots, \mathcal{N}^D(S_2(N, M))\}$ and \mathcal{C} can be obtained by arranging the value of c orderly along with the values of ξ and ζ increasing successively. To illustrate the above process clear and vividly, an example about $\mathcal{N} = \{1, 2, 3, 4, 5\}$ and $\mathcal{M} = \{1, 2, 3\}$ is given in Table I. There, all the subset classifications about $N = 5$ and $M = 3$ with $S_2(5, 3) = 25$ are listed followed by the above proposed process, while parameters ξ and ζ are computed too. Without loss of generality, only two situations are discussed in detail. The first situation is $\mathcal{N}_1(c) = \{1\}$, $\mathcal{N}_2(c) = \{2\}$ and $\mathcal{N}_3(c) = \{3, 4, 5\}$. Obviously, one knows $n_1(c) = 1$, $n_2(c) = 1$ and $n_3(c) = 3$, while $\xi = 113$ and $\zeta = 12345$. Because of $\zeta = 12345$ being the smallest among $\xi = 113$, and based on the proposed mapping principle, one should select $c = 1$. Similarly, $c = 2$ means the second-smallest value of ζ among $\xi = 113$ such as $\zeta = 13245$. The second situation is $\mathcal{N}_1(c) = \{1\}$, $\mathcal{N}_2(c) = \{2, 3\}$ and $\mathcal{N}_3(c) = \{4, 5\}$. Then, one knows $n_1(c) = 1$, $n_2(c) = 2$ and $n_3(c) = 2$, such that

$\xi = 122$ and $\zeta = 12345$. Though the values of ζ about the two situations are the same, the values of ξ are different. Due to $113 < 122$, the values of parameter c are different and selected to be $c = 1$ and $c = 11$ respectively followed by the above mapping. The other situations are all given in Table I, whose detailed computation processes are omitted.

TABLE I
THE CLASSIFICATION OF COMBINATIONS AND THEIR NUMBERS

$\mathcal{N}_1(c)$	$\mathcal{N}_2(c)$	$\mathcal{N}_3(c)$	ξ	ζ	c
{1}	{2}	{3, 4, 5}	113	12345	1
{1}	{3}	{2, 4, 5}	113	13245	2
{1}	{4}	{2, 3, 5}	113	14235	3
{1}	{5}	{2, 3, 4}	113	15234	4
{2}	{3}	{1, 4, 5}	113	23145	5
{2}	{4}	{1, 3, 5}	113	24135	6
{2}	{5}	{1, 3, 4}	113	25134	7
{3}	{4}	{1, 2, 5}	113	34125	8
{3}	{5}	{1, 2, 4}	113	35124	9
{4}	{5}	{1, 2, 3}	113	45123	10
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{1}	{2, 3}	{4, 5}	122	12345	11
{1}	{2, 4}	{3, 5}	122	12435	12
{1}	{2, 5}	{3, 4}	122	12534	13
{2}	{1, 3}	{4, 5}	122	21345	14
{2}	{1, 4}	{3, 5}	122	21435	15
{2}	{1, 5}	{3, 4}	122	21534	16
{3}	{1, 2}	{4, 5}	122	31245	17
{3}	{1, 4}	{2, 5}	122	31425	18
{3}	{1, 5}	{2, 4}	122	31524	19
{4}	{1, 2}	{3, 5}	122	41235	20
{4}	{1, 3}	{2, 5}	122	41325	21
{4}	{1, 5}	{2, 3}	122	41523	22
{5}	{1, 2}	{3, 4}	122	51234	23
{5}	{1, 3}	{2, 4}	122	51324	24
{5}	{1, 4}	{2, 3}	122	51423	25

Secondly, after introducing the mapping between sets \mathcal{D} and \mathcal{C} , an optimization algorithm based on a local search can be proposed and remain some advantages. In detail, on the one hand, a trajectory only using the selected partial nodes instead of exploiting all the division nodes is generated. In other words, it can lead to a locally even globally optimal solution by an iterating method. On the other hand, a favorable convergence and monotonicity of the proposed algorithm should be satisfied. Particularly, based on the proposed mapping, the normal HCA can be used to realize the former requirement of locally optimal solution. However, the other properties on the convergence and monotonicity described in the latter are not guaranteed. In order to make a brief introduction to the HCA, some definitions are needed to be clarified, while the other details can be found in the existing similar references such as [25]. For convenience, in the following, we will mention the detailed classification set $\mathcal{N}^D(c)$ by using scalar c equivalently. Then, the scalar parameter c in (11) will have a specific meaning, which actually corresponds a detailed classification $\mathcal{N}^D(c)$ and makes it possible to get the optimal solution by applying the HCA. Unfortunately, at least two shortcomings are found in directly applying set \mathcal{C} , when the HCA is to be used. For one thing, the element of \mathcal{C} has a clustering phenomenon. In other words, the clustering phenomenon will likely lead to a worse locally optimal solution, when the search interval radius of the traditionally HCA is not large enough.

For another, due to the element of \mathcal{C} being a natural number, a big range between two elements exists and complicates the computation of (11). In order to solve the problems coming from \mathcal{C} , another set Ω referred to be “the nominal solution set” is defined as $\Omega \triangleq \{\omega(1), \omega(2), \dots, \omega(S_2(N, M))\}$, where $\omega(c), \forall c \in \mathcal{C}$, is an any real number in the interval $(0, \epsilon]$ with a given real $\epsilon \in \mathbb{R}_{>0}$. Meanwhile, it is further required that $\omega(c), \forall c \in \mathcal{C}$, is distinct such as $\omega(c) \neq \omega(j), \forall c, j \in \mathcal{C}$ and $c \neq j$. More importantly, a corresponding correlation between sets \mathcal{C} and Ω is denoted such that the value of $\omega(c)$ corresponds to the classification set $\mathcal{N}^D(c)$, when the scalar c belongs to set \mathcal{C} . After doing that, another discrete mapping between sets $\mathcal{J} \triangleq \{J(\alpha^*(1)), J(\alpha^*(2)), \dots, J(\alpha^*(S_2(N, M)))\}$ and Ω is introduced to be

$$f: \Omega \rightarrow \mathcal{J} \text{ such as } f(\omega(c)) = J(\alpha^*(c)), \forall c \in \mathcal{C} \quad (13)$$

When some elements' values of set \mathcal{J} are equal, one just needs to remain one and delete the other same values. Based on the characteristic of mapping (13), it can be seen that the deletion of the same values of set \mathcal{J} has nothing to do with the considered problem in this paper. Without loss of generality, all the elements' values of \mathcal{J} are assumed to be distinct in the next. Particularly, due to $f(\omega(c)) = J(\alpha^*(c))$, $\omega(c)$ is said to be “the nominally optimal solution” to (11) with $c \in \mathcal{D}$ given in advance, whose optimization effect is the same as $\alpha^*(c)$. After doing this, the problems mentioned can be done by applying set \mathcal{C} directly. Finally, in order to exploit the HCA successfully, another variable $\hat{\omega}(k) \in \Omega, k = 1, 2, \dots, k_{\max}$, is introduced, where k is the a number of k -th iteration and k_{\max} is the maximum number of iteration and obviously satisfies $k_{\max} \leq S_2(N, M)$. Because of $\hat{\omega}(k) \in \Omega$, it is true that for any iteration number k , there is always only one $\omega(c)$ satisfying $\hat{\omega}(k) = \omega(c), \exists c \in \mathcal{C}$. Meanwhile, when the HCA is mentioned, variable $\hat{\omega}(k)$ is normally simply said to be “a solution” and selected as the k -th iteration value. For the sake of distinction and simplicity, it will be named as “nominal solution” in the next. Then, a neighbour set $\Theta(\hat{\omega}(k))$ about nominal solution $\hat{\omega}(k) \in \Omega$ is defined as

$$\Theta(\hat{\omega}(k)) \triangleq \left\{ \hat{\omega}'(k) \in \Omega \mid |\hat{\omega}'(k) - \hat{\omega}(k)| < L, \hat{\omega}'(k) \neq \hat{\omega}(k) \right\} \quad (14)$$

where $L > 0$ is the radius of search interval. The neighborhood probability mass function $g_{\hat{\omega}(k)}(\hat{\omega}'(k))$ is described as

$$g_{\hat{\omega}(k)}(\hat{\omega}'(k)) \triangleq \mathbb{P}(\hat{\omega}'(k) \in \Theta(\hat{\omega}(k))) = \frac{1}{\text{num}(\Theta(\hat{\omega}(k)))} \quad (15)$$

Meanwhile, the update law of $\hat{\omega}(k)$ is given by

$$\hat{\omega}(k+1) = \hat{\omega}'(k), \text{ if } f(\hat{\omega}'(k)) - f(\hat{\omega}(k)) < 0 \quad (16)$$

It is noted that the above algorithm only ensures the monotonicity of function such as $f(\hat{\omega}(k+1)) < f(\hat{\omega}(k))$ but not the convergence of nominal solution $\hat{\omega}(k)$ such as $|e_{\hat{\omega}}(k+1)| < |e_{\hat{\omega}}(k)|$ where $e_{\hat{\omega}}(k) \triangleq \hat{\omega}(k+1) - \hat{\omega}(k)$. Thus, the convergence speed of nominal solution cannot be guaranteed.

To illustrate the utility of the above proposed algorithm, an example simulation about 100 distinct random points

will be given. In detail, their horizontal and vertical coordinates referred to be $\omega(c)$ and $f(\omega(c))$ respectively are randomly generated on interval $(0, 20]$, which are further denoted as $\Omega = \{\omega(1), \omega(2), \dots, \omega(100)\}$ and $\mathcal{J} = \{f(\omega(1)), f(\omega(2)), \dots, f(\omega(100))\}$. First of all, some simulations are shown in Fig. 1, in which different search radius L such as $L = 20, 10.71, 5.68$ and 2.84 are given. On

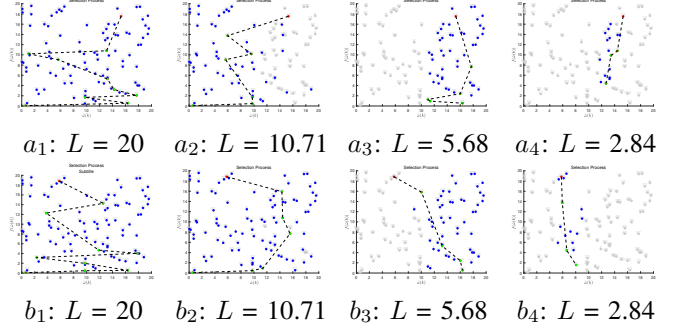


Fig. 1. The iteration effects of (14) under different constants L with different initial points.

the one hand, the red point is the initial iteration point, and the gray points are “useless” ones which are not used at all. Without loss of generality, two situations about different initial iteration points such as Point 16 and Point 83 are considered. Their horizontal and vertical coordinates are selected to be $(\hat{\omega}(1), f(\hat{\omega}(1))) = (\omega(16), f(\omega(16)))$ and $(\hat{\omega}(1), f(\hat{\omega}(1))) = (\omega(83), f(\omega(83)))$ respectively. On the other hand, the blue points are the failed ones during the iteration, while the green ones are the successful ones satisfying the update law (16). By investigating these simulations, it can be found that the bigger the value L takes, the smaller the value $f(\hat{\omega}(k))$ arrives. However, the computation complexity will be larger as L takes bigger values. To the contrary, when the radius is selected to be small enough, it can be concluded from Fig. 1 that the computation complexity will be great reduced. However, the probability of the obtained optimal solution being a locally optimal solution will be higher and have a worse quality. Thus, how to select a suitable radius is also one topic studied in this paper.

The search radius L of (14) is selected to be

$$L_{k+1} \triangleq \delta \cdot |\hat{\omega}(k+1) - \hat{\omega}(k)|, k = 1, 2, \dots \quad (17)$$

where $\delta \in (0, 1]$ is a attenuation coefficient. Particularly, the initial search radius L_1 is also defined as $L_1 \triangleq \delta \cdot |\hat{\omega}(1) - \hat{\omega}(0)|$, where $\hat{\omega}(0)$ is given to be

$$\hat{\omega}(0) = \begin{cases} 0, & \text{if } \hat{\omega}(1) > \frac{\epsilon}{2} \\ \epsilon, & \text{otherwise} \end{cases}$$

Then, one can get a simulation Fig. 2 similar to Fig. 1, where the values of δ about the initial radius are selected to be $\delta = 1, 0.7, 0.4$ and 0.2 respectively. Accordingly, the values of initial radius L_1 about initial value $\hat{\omega}(1) = \omega(16)$ based on (17) are computed to be $L_1 = 15.3, 10.71, 6.12$ and 3.06 respectively. Meanwhile, the values of initial radius L_1 about

initial value $\hat{\omega}(1) = \omega(82)$ are computed to be $L_1 = 14.2, 9.94, 5.68$ and 2.84 respectively. On the one hand, by the simulations of Fig. 2 presented in rows respectively, it can be concluded that the computation complexity is usually reduced along with δ decreasing under the same initial condition. On the other hand, in order to make some comparisons between Fig. 1 based on (14) and Fig. 2 based on (17), it had better to compare them under the same conditions such as the same $\hat{\omega}(1)$ and L_1 . Thus, only a part but not all the subgraphs are necessary to make comparisons such as the comparisons on subgraphs $a_i, i = 1, 2$, and $b_i, i = 1, 3, 4$ between Fig. 1 and Fig. 2 respectively. In detail, not only the fewer iteration times but also the fewer failed points are presented in Fig. 2. More importantly, the convergence of nominal solution $\hat{\omega}(k)$ is guaranteed in Fig. 2, which is not satisfied in Fig. 1. However, there is also a negative effect that the gaps among the locally optimal solutions obtained in Fig. 2 are bigger, some of which are farther away from the globally optimal solution to subgraphs a_i and b_i of Fig. 1. Thus, it is necessary to further study the effect of variable radius in order to give an optimal variation of radius L_k .

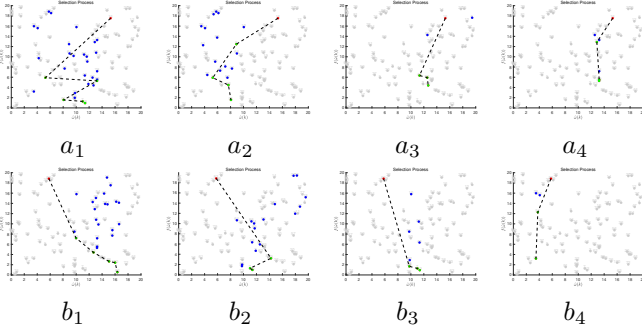


Fig. 2. The iteration effects of (17) under different constants L with different initial points.

For another, based on (17), it can be known that δ plays an important role in generating variable L_k . In order to realize an optimization to radius L_k , expression (17) is revised to be

$$L_{k+1} \triangleq \delta_{k+1} \cdot |\hat{\omega}(k+1) - \hat{\omega}(k)|, k = 1, 2, \dots \quad (18)$$

where $\delta_k \in (0, 1]$. Similar to the definition L_1 in (17), it is defined as $L_1 \triangleq \delta_1 \cdot |\hat{\omega}(1) - \hat{\omega}(0)|$. Then, similar to Fig. 2, similar simulations under (18) are given in Fig. 3. Particularly, the initial values of parameters δ_k and L_k of subgraphs in Fig. 3 are given such as $a_1 : \delta_1 = 1, L_1 = 15.3$; $a_2 : \delta_1 = 0.7, L_1 = 10.71$; $a_3 : \delta_1 = 0.4, L_1 = 6.12$; $a_4 : \delta_1 = 0.2, L_1 = 3.06$; $b_1 : \delta_1 = 1, L_1 = 14.2$; $b_2 : \delta_1 = 0.7, L_1 = 9.94$; $b_3 : \delta_1 = 0.4, L_1 = 5.68$ and $b_4 : \delta_1 = 0.2, L_1 = 2.84$. By making the comparisons of all the subgraphs between Figs. 2 and 3, it can be found that less computation complexity of (18) is presented in Fig. 3, whose iteration time is fewer too. Moreover, the minimum value of $f(\hat{\omega}(k))$ obtained in Fig. 2 based on (17) is larger than one in Fig. 3 by applying (18). In other words, it is necessary and important to study how to select the detailed

value during every iteration to gain a better performance such as smaller value of (12), less computation complexity, but better monotonicity and convergence. All these problems will be considered and finally solved in this paper.

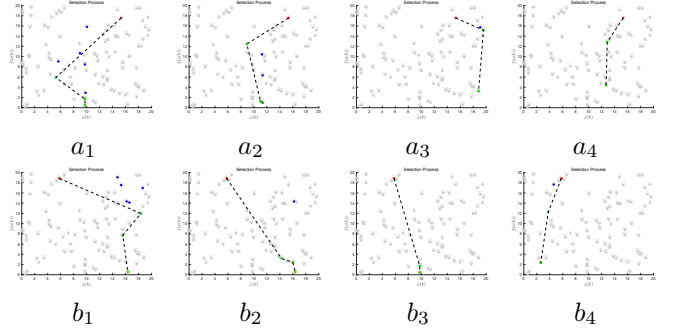


Fig. 3. The iteration effects of (18) under different constants L with different initial points..

DEFINITION 1: A control update sequence $\{t_k, k \in \mathbb{N}\}$ is said to have the finite sampling rate property if there exists a scalar \underline{s} such that $s_k \geq \underline{s} > 0, k \in \mathbb{N}_{>0}$.

Assumption 1: It is assumed that the event-triggered sampling interval $s_k, k \in \mathbb{N}_{>0}$, has an upper bound such as $s_k \leq \min \{\bar{\tau}_1, \bar{\tau}_2, \dots, \bar{\tau}_M\}$, where $\bar{\tau}_h = \sum_{\ell \in \mathcal{N}_h} \hat{\tau}_\ell, \forall h \in \mathcal{M}$, and $\hat{\tau}_\ell \triangleq \mathbb{E}[\tau_\ell(n)], \forall \ell \in \mathcal{N}_h, \forall n \in \mathbb{N}_{>0}$.

DEFINITION 2: System (1) closed by controller (3) is globally asymptotically stable almost surely (GAS a.s.), if the following conditions hold simultaneously.

- (C₁) For any $\epsilon \in (0, 1)$, there is a $\delta = \delta(\epsilon) > 0$ such that $\mathbb{P}(\sup_{t \geq 0} \|x(t)\| < \epsilon) > 1 - \epsilon$, when $\|x_0\| < \delta$.
- (C₂) For any $\gamma > 0$ and $\xi > 0$, there is a positive random variable $\varphi(\gamma, \xi)$ such that $\mathbb{P}(\sup_{t \geq \varphi(\gamma, \xi)} \|x(t)\| < \xi) = 1$, when $\|x_0\| < \gamma$.

III. MAIN RESULTS

THEOREM 1: Consider the system composed of (1) and (3), where the detailed classification of controller (3) is assumed to be given in advance such that \mathcal{N}^D is given beforehand, so do the $\alpha_\ell^{[h]}$ and $\bar{K}_\ell, \forall h \in \mathcal{M}, \forall \ell \in \mathcal{N}$. Then, the related closed-loop system is GAS a.s., if given parameters $\lambda_i \in \mathbb{R}, \mu_i > 1$ and $\sigma_h \in \mathbb{R}_{>0}$, there exists a matrix $P_i > 0$ satisfying

$$\left(P_i \bar{A}_i^{[h]}\right)^* + 2\sigma_h \|P_i B_i\| \sum_{\ell=1}^N \alpha_\ell^{[h]} \bar{K}_\ell \|I\| \leq \lambda_i P_i, \forall h \in \mathcal{M}, \forall i \in \mathcal{N}_h \quad (19)$$

$$P_i \leq \mu_i P_j, \forall i, j \in \mathcal{N} \quad (20)$$

$$\sum_{i=1}^N \pi_i \left(\lambda_i + \frac{\ln \mu_i}{\hat{\tau}_i} \right) < 0 \quad (21)$$

where $\bar{A}_i^{[h]} = A_i + B_i \sum_{\ell=1}^N \alpha_\ell^{[h]} \bar{K}_\ell$, $\pi_i = \frac{\bar{\pi}_i \hat{\tau}_i}{\sum_{j=1}^N \bar{\pi}_j \hat{\tau}_j}$, and $\bar{\pi}_i, \forall i \in \mathcal{N}$, is the stationary distribution of the embedded

chain of semi-Markov process. Meanwhile, the upper sampling bound $\bar{\tau}_h$ is computed by

$$\bar{\tau}_h = \begin{cases} \frac{1}{\|\hat{A}_h\|} \frac{\sigma_h}{1 + \sigma_h}, & \text{if } \hat{\chi}_h \leq 0 \\ \frac{1}{\hat{\chi}_h} \log \left(\frac{\sigma_h \hat{\chi}_h}{\|\hat{A}_h\| (1 + \sigma_h)} + 1 \right), & \text{if } \hat{\chi}_h > 0 \end{cases}, \forall h \in \mathcal{M} \quad (22)$$

where $\|\hat{A}_h\| = \max_{\ell \in \mathcal{N}_h} \{\|\bar{A}_\ell^{[h]}\|\}$ and $\hat{\chi}_h = \max_{\ell \in \mathcal{N}_h} \{\chi_{A_\ell}\}$.

Proof The proof is given in Appendix A. \square

Next, **Problems 1** and **2** will be considered. Before discussing them, a convex optimization problem with some constraints should be proposed first. On the one hand, from Example 3.11 of [27] on Page 82, it is clear that $J^{[h]}((\alpha^{[h]}(c)))$, $\forall h \in \mathcal{M}$ but given $c \in \mathcal{C}$, is convex. Moreover, function $J(\alpha(c)) = \sum_{h=1}^M J^{[h]}((\alpha^{[h]}(c)))$ is also convex. On the other hand, even if the set classification of controllers is fixed, the given conditions of Theorem 1 to be as constraints will be hard to be solved in the optimization problems, since parameters \bar{K}_ℓ and P_i are vector and matrix variables respectively. As a result, different from Theorem 1, not only \bar{K}_ℓ satisfying $K_h \triangleq \sum_{\ell=1}^N \alpha_\ell^{[h]} \bar{K}_\ell$ but also P_i here is given beforehand. Without loss of generality, they will be calculated by following LMIs

$$\begin{bmatrix} (A_i + B_i Y_i)^* - \lambda_i X_i & B_i Y_i & X_i \\ * & -2X_i + I & 0 \\ * & * & -\sigma^{-2} I \end{bmatrix} \leq 0, \forall i \in \mathcal{N} \quad (23)$$

$$\begin{bmatrix} -\mu_i X_j & X_j \\ * & -X_i \end{bmatrix} < 0, \forall i, j \in \mathcal{N} \quad (24)$$

and (21), where $\lambda_i \in \mathbb{R}$, $\sigma \in \mathbb{R}_{>0}$ and $\mu_i > 1$ are given scalars, and $X_i \triangleq P_i^{-1}$ and $Y_i \triangleq \bar{K}_i X_i$ are variables to be computed by solving the above LMIs. Due to page limitation, the detailed calculation process is omitted here.

REMARK 2: It is worth mentioning that the computation way for \bar{K}_ℓ and P_i described in (21), (23) and (24) is not unique. In other words, they can be computed in advance by other ways, as long as the corresponding conditions described by (19)-(21) with given \bar{K}_ℓ and P_i satisfy. Naturally, there exist inevitable differences among the different techniques providing different parameters' values. However, such differences will not lead to essential distinctions to the optimization problems in this paper and will not discuss in detail.

Since the parameters of Theorem 1 such as $\lambda_i \in \mathbb{R}$, $\mu_i > 1$, $\sigma_h \in \mathbb{R}_{>0}$, \bar{K}_ℓ and P_i are given beforehand, conditions (20) and (21) are known and always satisfied. In detail, parameters $\lambda_i \in \mathbb{R}$, $\mu_i > 1$, \bar{K}_ℓ and P_i are defined by (21), (23) and (24), while $\sigma_h \in \mathbb{R}_{>0}$ is a new given parameter and only defined in (19). Then, the constraints described by conditions (19)-(21) are equivalently simplified to be

$$\text{s.t. } \mathcal{H}_i(\alpha(c)) \leq 0, \forall i \in \mathcal{N}_h, c \in \mathcal{C} \quad (25)$$

where

$$\mathcal{H}_i(\alpha(c)) \triangleq \sum_{h=1}^M \left\{ \left[P_i \left(A_i + B_i \sum_{\ell=1}^N \alpha_\ell^{[h]}(c) \bar{K}_\ell \right) \right]^* - \lambda_i P_i + 2\sigma_h \left\| P_i B_i \sum_{\ell=1}^N \alpha_\ell^{[h]}(c) \bar{K}_\ell \right\| I \right\}$$

According to Example 2.10 on Page 38 and Example 3.11 on Page 82 of [27], it is clear that constraint (25) is a convex set. Then, a convex optimization problem described by cost function (11) with constraint (25) is successfully proposed. According to Example 3.10 of [27] on Page 82, it can be concluded that function $\lambda_{\mathcal{H}_i(\alpha(c))}^{\max}$ is also convex. Thus, the feasible domain of (25) is equivalently rewritten to be

$$\text{s.t. } \lambda_{\mathcal{H}_i(\alpha(c))}^{\max} \leq 0, \forall i \in \mathcal{N}_h, c \in \mathcal{C} \quad (26)$$

Since (26) being inequalities, a slack variable such as $\varsigma_i(c)$ is introduced to change (26) be an equation constraint such as

$$\text{s.t. } \lambda_{\mathcal{H}_i(\alpha(c))}^{\max} + (\varsigma_i(c))^2 = 0, \forall i \in \mathcal{N}_h, c \in \mathcal{C} \quad (27)$$

Based on the given results of [28] on Pages 98 and 99, it can be known that such slack variables do not bring any negative effects to (26) but make the equivalent constraint (27) solved easily. Moreover, it can be concluded that constraint (27) is convex. Thus, the original convex optimization problem is transformed into convex cost function (11) with convex constraint (27). In order to solve this convex optimization problem, an augmented Lagrangian function is constructed as

$$\mathcal{L}(\alpha(c), \gamma(c), \varsigma(c), \phi) \triangleq J((\alpha(c))) + L_1(\alpha(c), \gamma(c), \varsigma(c)) + L_2(\alpha(c), \varsigma(c), \phi) \quad (28)$$

where

$$L_1(\alpha(c), \gamma(c), \varsigma(c)) \triangleq \sum_{i=1}^N \gamma_i(c) E_i(\alpha(c), \varsigma_i(c))$$

$$L_2(\alpha(c), \varsigma(c), \phi) \triangleq \frac{\phi}{2} \sum_{i=1}^N E_i^2(\alpha(c), \varsigma_i(c))$$

$$E_i(\alpha(c), \varsigma_i(c)) \triangleq \lambda_{\mathcal{H}_i(\alpha(c))}^{\max} + (\varsigma_i(c))^2$$

$$\gamma(c) \triangleq [\gamma_1(c) \quad \gamma_2(c) \quad \cdots \quad \gamma_N(c)]^T \in \mathbb{R}_{\geq 0}^N$$

$$\varsigma(c) \triangleq [\varsigma_1(c) \quad \varsigma_2(c) \quad \cdots \quad \varsigma_N(c)]^T \in \mathbb{R}_{\geq 0}^N$$

and penalty factor $\phi \in \mathbb{R}_{>0}$. Here, $\varsigma(c)$ is the slack variable and $\gamma(c)$ is the multiplier. Because (11) and (27) are both convex, it is obvious that function (28) is convex too.

Meanwhile, it is obvious that the above computation process of the globally optimal value J^* should be repeated by selecting every $c \in \mathcal{C}$. Based on the above illustrations and simulations about mode classification, the repetition number of computation will equal to the Stirling number of the second kind. Obviously, the computation complexity will be very large especially N and M increase big. In order to deal with this problem and motivated by the phenomena discovered in Figs. 1-3, an optimization algorithm based on an improved HCA

will be developed, which can generate an iteration sequence of solution $\{\hat{\omega}(k)\}_{k=1}^{k_{\max}}$ and so does the corresponding sequence of objective function $\{f(\hat{\omega}(k))\}_{k=1}^{k_{\max}}$. Moreover, the locally optimal solution $\hat{\omega}^*$ and its value $f^* \triangleq f(\hat{\omega}^*)$ can be reached such as $\hat{\omega}^* = \hat{\omega}(k_{\max})$. Moreover, one can conclude from Figs. 1-3 that strategy (18) will be the best if its dynamically attenuation coefficient δ_k is suitably adjusted. In detail, it not only ensures all the advantages of strategy (17) but also improves the iteration speed and the quality of optimal solution $\hat{\omega}^*$. How to dynamically adjust the attenuation coefficient δ_k to make (18) work optimally is an open problem. Here, we will use the Q-learning algorithm to dynamically adjust δ_k , so as to dynamically change the radius of search interval L_k and lead to an improvement in performance. Meanwhile, it is very necessary and important to analyze its mathematical model, which includes the state, action, and reward function.

State: The state S_k is a feedback of the agent at the k th step and denotes the impact of its taking action on the environment. Here, S_k described as the produced solution $\hat{\omega}(k) \in \Omega$ at step k such as $S_k \triangleq \hat{\omega}(k)$. Meanwhile, a termination state S_{end} should be given beforehand but does not belong to the solution set Ω . Its aim is to end the current training round when the state S_k arrives at the termination state S_{end} . Finally, the state space \mathcal{S} is composed of the solution set Ω of HCA and the termination state S_{end} such as $\mathcal{S} = \Omega \cup \{S_{\text{end}}\}$.

Action: The action \mathcal{A}_k is a description of the agent's behavior at the k th step. Here, it is the specific value of δ_k of HCA at the k th iteration step such as $\mathcal{A}_k \triangleq \delta_k$, which can change the neighborhood set $\Theta(\hat{\omega}(k))$ and affect the generation of the next solution. As a result, all values of δ_k or \mathcal{A}_k constitute the action space \mathcal{A} .

Reward: The reward \mathcal{R}_{k+1} is a feedback of the environment in which the agent has taken the action \mathcal{A}_k in state S_k . Here, it is described by

$$\mathcal{R}_{k+1} = \begin{cases} \frac{f(\hat{\omega}(k)) - f(\hat{\omega}'(k))}{|\hat{\omega}'(k) - \hat{\omega}(k)|}, & \text{if } \text{num}(\Theta(\hat{\omega}(k))) \neq 0 \\ -f(\hat{\omega}(k)), & \text{otherwise} \end{cases} \quad (29)$$

where function $f(\cdot)$ has been defined in (13). In detail, two situations will be found if condition $\text{num}(\Theta(\hat{\omega}(k))) \neq 0$ is satisfied. On the one hand, the agent will be penalized and receive a negative reward, if the agent takes an action \mathcal{A}_k in state S_k such that the solution $\hat{\omega}'(k)$ satisfies $f(\hat{\omega}'(k)) \geq f(\hat{\omega}(k))$. On the other hand, the agent will be encouraged and receive a positive reward, if the solution $\hat{\omega}'(k)$ satisfies $f(\hat{\omega}'(k)) < f(\hat{\omega}(k))$. In other words, if the agent wants to obtain a larger positive reward, the solution $\hat{\omega}'(k)$ should satisfy not only $f(\hat{\omega}(k))$ is larger than $f(\hat{\omega}'(k))$ as much as possible, but also the distance between $\hat{\omega}'(k)$ and $\hat{\omega}(k)$ is as small as possible. To the contrary, $\text{num}(\Theta(\hat{\omega}(k))) = 0$ means the neighbor set $\Theta(\hat{\omega}(k))$ of $\hat{\omega}(k)$ is the empty set. It also means that no solution within $\Theta(\hat{\omega}(k))$ whose objective function value is smaller than $f(\hat{\omega}(k))$. In this situation, the agent is still penalized but receives a negative reward $-f(\hat{\omega}(k))$. Then, according to [29] on Page 70, it is well

known that the Q-learning effect of taking action \mathcal{A}_k in state S_k can be depicted by the Q-value, i.e. it can make an evaluation about the variation of δ_k . Accordingly, a specific definition of Q-value similar to (3.11) in [29] is given as

$$Q^\varpi(s, a) \triangleq \mathbb{E}_\varpi \left(\sum_{q=0}^{\infty} \gamma^q \mathcal{R}_{k+q+1} \middle| S_k = s, \mathcal{A}_k = a \right) \quad (30)$$

where $Q^\varpi(s, a)$ is the Q-value and quantizes the effect of taking action $\mathcal{A}_k = a$ followed by the policy ϖ in state $S_k = s$, and $\gamma \in [0, 1]$ is a given discount factor. Particularly, policy ϖ is a mapping from each state $s \in \mathcal{S}$ and action $a \in \mathcal{A}$ to the probability $\mathbb{P}(a|s)$ of taking action a when in state s . As a result, a different taking action $\mathcal{A}_k = a$ can be generated by a different strategy ϖ in state $S_k = s$ and will lead to a different $Q^\varpi(s, a)$. In detail, when the Q-value is larger, the taking action will be better. To the contrary, the action selection will be worse.

THEOREM 2: Consider a convex optimization problem described by (11) and (25), where the parameters \bar{K}_ℓ and P_i are computed by solving conditions (21), (23) and (24), and parameters $\lambda_i \in \mathbb{R}$, $\sigma \in \mathbb{R}_{>0}$, $\sigma_h \in \mathbb{R}_{>0}$ and $\mu_i > 1$ are also given in advance. The optimization problem (11) constrained on (25) will have a globally optimal solution, if the following conditions are satisfied

$$\nabla_{\alpha(c)} \mathcal{L}(\alpha(c), \gamma(c), \varsigma(c), \phi) = 0, \forall c \in \mathcal{C} \quad (31)$$

$$\nabla_{\varsigma(c)} \mathcal{L}(\alpha(c), \gamma(c), \varsigma(c), \phi) = 0, \forall c \in \mathcal{C} \quad (32)$$

$$\nabla_{\gamma(c)} \mathcal{L}(\alpha(c), \gamma(c), \varsigma(c), \phi) = 0, \forall c \in \mathcal{C} \quad (33)$$

Then, the globally optimal value of (11) is obtained by

$$J^* \triangleq J(\alpha^*) = \min_{c \in \mathcal{C}} J(\alpha^*(c)) = \min_{c \in \mathcal{C}} \mathcal{L}(\alpha^*(c), \gamma^*(c), \varsigma^*(c), \phi) \quad (34)$$

Meanwhile, the globally optimal solution is computed by

$$\alpha^* \triangleq \arg \min_{c \in \mathcal{C}} J(\alpha^*(c)) = \arg \min_{c \in \mathcal{C}} \mathcal{L}(\alpha^*(c), \gamma^*(c), \varsigma^*(c), \phi) \quad (35)$$

Particularly, in order to get an optimal solution to convex problem (11) with (25) without computing $c \in \mathcal{C}$ one by one, an improved HCA based on Q-learning (Q-HCA) is proposed, whose searching radius iteration rule is given in (18). Meanwhile, δ_k is obtained by the Q-learning with state $S_k = s$, action $\mathcal{A}_k = a$ and reward \mathcal{R}_{k+1} presented in (29). Then, an iteration rule for estimating Q-value (30) is given as

$$Q_{k+1}(S_k, \mathcal{A}_k) = \begin{cases} Q_k(S_k, \mathcal{A}_k) - \beta [Q_k(S_k, \mathcal{A}_k) - (\mathcal{R}_{k+1} \\ + \gamma \max_{\hat{a} \in \mathcal{A}(S_{k+1})} Q_k(S_{k+1}, \hat{a}))], \\ \text{if } (S_k, \mathcal{A}_k) = (s, a); \\ Q_k(S_k, \mathcal{A}_k), \text{ otherwise} \end{cases} \quad (36)$$

where $\beta \in [0, 1]$ is the given learning rate, and $\mathcal{A}(S_{k+1})$ is the set of taking actions in state S_{k+1} . As a result, an optimal attenuation coefficient δ_k^* is obtained. Moreover, under any initial condition $\forall \hat{\omega}(1) \in \Omega$ with $\hat{\omega}(0)$ similarly defined in

(17), a preferably nominal solution sequence $\{\hat{\omega}^{\text{pre}}(k)\}_{k=1}^{k_{\max}}$ with $\hat{\omega}^{\text{pre}}(1) = \hat{\omega}(1)$ can be developed by following rules (36) and (18), where k_{\max} is a maximum value of iteration. Thus, a locally optimal solution to convex problem (11) with (25) satisfying a part of \mathcal{C} is computed by

$$\hat{\omega}^{\text{pre}}(k) \rightarrow \hat{\omega}^* \triangleq \hat{\omega}^{\text{pre}}(k_{\max}) \quad (37)$$

Meanwhile, the corresponding value of (11) is also locally optimal such as $J^* \leq f^*$ and described to be

$$f^* = f(\hat{\omega}^*) = f(\hat{\omega}^{\text{pre}}(k_{\max})) < \dots < f(\hat{\omega}^{\text{pre}}(1)) \quad (38)$$

More importantly, both convergence and monotonicity properties of Q-HCA are ensured. Finally, the upper sampling bound $\bar{\tau}_h$ of the related optimal controllers is also computed by (22).

Proof The proof is given in Appendix B. \square

Specially, when controller (3) has no sampling phenomenon, it will reduce to be

$$u(t) = \sum_{\ell=1}^N \alpha_{\ell}^{[\nu(\eta(t))]} \bar{K}_{\ell} x(t), \forall t \in [T_n, T_{n+1}), \forall n \in \mathbb{N} \quad (39)$$

where $\alpha_{\ell}^{[\nu(\eta(t_k))]} \in \mathbb{R}$ and $\bar{K}_{\ell} \in \mathbb{R}^{m \times n}$ are similar to (3), and function $\nu(\eta(t))$ also has the similar definition but on interval $\forall t \in [T_n, T_{n+1})$. Then, one has the following corollaries.

COROLLARY 1: Consider the system composed of (1) and (39) with mode classification \mathcal{N}^D and $\alpha_{\ell}^{[h]}$ given beforehand. Then, the related closed-loop system is GAS a.s., if given parameters $\lambda_i \in \mathbb{R}$ and $\mu_i > 1$, there exist $X_i > 0$, \bar{Y}_{ℓ} and W satisfying LMIs (21), (24) and

$$\begin{bmatrix} \Gamma_{i1}^{[h]} & \Gamma_{i2}^{[h]} \\ * & (-W)^* \end{bmatrix} < 0, \forall h \in \mathcal{M}, \forall i \in \mathcal{N}_h \quad (40)$$

where $\Gamma_{i1}^{[h]} = (A_i W + B_i \sum_{\ell=1}^N \alpha_{\ell}^{[h]} \bar{Y}_{\ell})^* - \lambda_i X_i$ and $\Gamma_{i2}^{[h]} = A_i W + B_i \sum_{\ell=1}^N \alpha_{\ell}^{[h]} \bar{Y}_{\ell} + X_i - W^{\top}$. Then, the control gains can be computed by

$$\bar{K}_{\ell} = \bar{Y}_{\ell} W^{-1} \quad (41)$$

Proof The proof is established by Theorem 1 and omitted. \square

REMARK 3: When parameter $\alpha_{\ell}^{[h]}$ of controller (39) is selected to be special values such as $\alpha_{\ell}^{[h]} = \frac{\pi_{\ell}}{\sum_{j \in \mathcal{N}_h} \pi_j}$, if $\ell \in \mathcal{N}_h$; $\alpha_{\ell}^{[h]} = 0$, otherwise. Accordingly, controller (39) can be simply rewritten as $u(t) = \sum_{\ell \in \mathcal{N}_h} \frac{\pi_{\ell}}{\sum_{j \in \mathcal{N}_h} \pi_j} \bar{K}_{\ell} x(t)$ and similar to [17] which is related to the stationary distribution probability. In this situation, similar results can be obtained and omitted for page limitation. Particularly, if all the values of $\sum_{\ell=1}^N \alpha_{\ell}^{[h]} \bar{Y}_{\ell}$ are equal such as $\bar{Y} = \sum_{\ell=1}^N \alpha_{\ell}^{[h]} \bar{Y}_{\ell}$, $\forall h \in \mathcal{N}$ and $\mathcal{M} = \mathcal{N}$, a mode-independent controller $u(t) = Kx(t)$ with $\bar{K} = \bar{Y} W^{-1}$ will be obtained and similar to [17], [22]. In other words, Corollary 1 contains some existing results as special situations, which is more general and less conservative.

Accordingly, when the mode classification is not given in advance and similar to (10), controller (39) is described as

$$u(t) = \sum_{\ell=1}^N \alpha_{\ell}^{[\nu(\eta(t))]}(c) \bar{K}_{\ell}(c) x(t), \forall t \in [T_n, T_{n+1}) \quad (42)$$

After giving parameters $\lambda_i \in \mathbb{R}$ and $\mu_i > 1$, one can compute \bar{K}_{ℓ} and P_i by solving conditions (21), (24) and

$$(A_i + B_i Y_i)^* - \lambda_i X_i \leq 0, \forall i \in \mathcal{N} \quad (43)$$

Then, similar to constraint (25), the constraint of the above conditions is equivalently simplified to be

$$\text{s.t. } \bar{\mathcal{H}}_i(\alpha(c)) \leq 0, \forall i \in \mathcal{N}_h(c), c \in \mathcal{C} \quad (44)$$

where

$$\bar{\mathcal{H}}_i(\alpha(c)) \triangleq \sum_{h=1}^M \left\{ \left[P_i \left(A_i + B_i \sum_{\ell=1}^N \alpha_{\ell}^{[h]}(c) \bar{K}_{\ell} \right) \right]^* - \lambda_i P_i \right\}$$

A similar augmented Lagrangian function is constructed as

$$\begin{aligned} \bar{\mathcal{L}}(\alpha(c), \gamma(c), \varsigma(c), \phi) &\triangleq J((\alpha(c))) + T_1(\alpha(c), \gamma(c), \varsigma(c)) \\ &\quad + T_2(\alpha(c), \varsigma(c), \phi) \end{aligned} \quad (45)$$

where

$$T_1(\alpha(c), \gamma(c), \varsigma(c)) \triangleq \sum_{i=1}^N \gamma_i(c) \bar{E}_i(\alpha(c), \varsigma_i(c))$$

$$T_2(\alpha(c), \varsigma(c), \phi) \triangleq \frac{\phi}{2} \sum_{i=1}^N \bar{E}_i^2(\alpha(c), \varsigma_i(c))$$

$$\bar{E}_i(\alpha(c), \varsigma_i(c)) \triangleq \lambda_{\mathcal{H}_i(\alpha(c))}^{\max} + (\varsigma_i(c))^2$$

and parameters $\gamma(c)$ and $\varsigma(c)$ are defined in (28). Then, one can obtain the following corollary.

COROLLARY 2: Consider a convex optimization problem described by (11) and (44), where the parameters \bar{K}_{ℓ} and P_i are computed by solving conditions (21), (24) and (43) under giving parameters $\lambda_i \in \mathbb{R}$ and $\mu_i > 1$. The optimization problem (11) constrained on (44) will have a globally optimal solution, if the following conditions are satisfied

$$\nabla_{\alpha(c)} \bar{\mathcal{L}}(\alpha(c), \gamma(c), \varsigma(c), \phi) = 0, \forall c \in \mathcal{C} \quad (46)$$

$$\nabla_{\varsigma(c)} \bar{\mathcal{L}}(\alpha(c), \gamma(c), \varsigma(c), \phi) = 0, \forall c \in \mathcal{C} \quad (47)$$

$$\nabla_{\gamma(c)} \bar{\mathcal{L}}(\alpha(c), \gamma(c), \varsigma(c), \phi) = 0, \forall c \in \mathcal{C} \quad (48)$$

Then, the globally optimal value of (11) is obtained by

$$J^* = \min_{c \in \mathcal{C}} \bar{\mathcal{L}}(\alpha^*(c), \gamma^*(c), \varsigma^*(c), \phi) \quad (49)$$

Meanwhile, the globally optimal solution is computed by

$$\alpha^* = \arg \min_{c \in \mathcal{C}} \bar{\mathcal{L}}(\alpha^*(c), \gamma^*(c), \varsigma^*(c), \phi) \quad (50)$$

Moreover, based on the developed Q-HCA algorithm, one also gets a locally optimal solution and its locally optimal value such as (37) and (38) without computing $c \in \mathcal{C}$ one by one.

Proof Its proof is similar to Theorem 2 and omitted. \square

REMARK 4: It can be seen that some parameters such as \bar{K}_{ℓ} and P_i in main results need to be given in advance. Particularly, in order to give them easily and directly, they are solved by some LMIs such as (23) and (24). It is worth noted that these convenient techniques will lead to some conservatism in terms of some of constraints such as (25)

unsatisfied. In this situation, some possible ways may be used to deal with them, and only two possible ways are mentioned here. On the one hand, one can select different values of the scalars given in the presented LMIs time and time again, until the dissatisfied constraints hold. On the other hand, other different conditions providing \bar{K}_ℓ and P_i in advance can be developed so as to make the dissatisfied constraints satisfying.

IV. NUMERICAL EXAMPLE

Example 1: Considering a system of form (1) with $\mathcal{N} = \{1, 2, 3, 4, 5\}$ whose matrices are described as

$$\begin{aligned} A_1 &= \begin{bmatrix} -2 & 2 \\ 0.1 & 0.14 \end{bmatrix}, B_1 = \begin{bmatrix} 0.32 \\ -1.7 \end{bmatrix}; A_2 = \begin{bmatrix} -3.6 & 0.2 \\ 1 & 0.2 \end{bmatrix} \\ B_2 &= \begin{bmatrix} 0.32 \\ -2 \end{bmatrix}; A_3 = \begin{bmatrix} -0.7 & 2 \\ 0 & 0.32 \end{bmatrix}, B_3 = \begin{bmatrix} 0.25 \\ -0.45 \end{bmatrix}; \\ A_4 &= \begin{bmatrix} -3.3 & 0.1 \\ 0.8 & 0.03 \end{bmatrix}, B_4 = \begin{bmatrix} 0.22 \\ -0.61 \end{bmatrix}; \\ A_5 &= \begin{bmatrix} -2.5 & 0 \\ 0.1 & 0.95 \end{bmatrix}, B_5 = \begin{bmatrix} 0.12 \\ -0.53 \end{bmatrix} \end{aligned}$$

The transition probabilities of the embodied chain is given as

$$P = \begin{bmatrix} 0 & 0.1 & 0.25 & 0.25 & 0.4 \\ 0.1 & 0 & 0.2 & 0.4 & 0.3 \\ 0.2 & 0.35 & 0 & 0.3 & 0.15 \\ 0.15 & 0.3 & 0.2 & 0 & 0.35 \\ 0.3 & 0.15 & 0.25 & 0.3 & 0 \end{bmatrix}$$

It has a unique stationary distribution computed as $\bar{\pi} = [0.21 \ 0.19 \ 0.2 \ 0.15 \ 0.25]$. In addition, the expectations of sojourn time for modes are given to be $\hat{\tau}_1 = 0.03$, $\hat{\tau}_2 = 0.05$, $\hat{\tau}_3 = 0.04$, $\hat{\tau}_4 = 0.09$ and $\hat{\tau}_5 = 0.07$. Firstly, based on conditions (23) and (24), after letting $\lambda_1 = -0.8$, $\lambda_2 = -0.9$, $\lambda_3 = -0.5$, $\lambda_4 = -0.6$, $\lambda_5 = 0.4$, $\mu_1 = 1.005$, $\mu_2 = 1.001$, $\mu_3 = 1.009$, $\mu_4 = 1.002$, $\mu_5 = 1.001$ and $\sigma = 0.069$, one can obtain that $\sum_{i=1}^N \pi_i \left(\lambda_i + \frac{\ln \mu_i}{\hat{\tau}_i} \right) = -0.2549 < 0$, and the control gains are computed such as $\bar{K}_1 = [0.1921 \ 0.1857]$, $\bar{K}_2 = [0.2087 \ 0.9061]$, $\bar{K}_3 = [1.0049 \ 4.9363]$, $\bar{K}_4 = [0.6863 \ 3.1342]$ and $\bar{K}_5 = [0.2835 \ 3.7523]$, while matrix P_i is computed as

$$\begin{aligned} P_1 &= \begin{bmatrix} 0.0047 & 0.0034 \\ 0.0034 & 0.0236 \end{bmatrix}, P_2 = \begin{bmatrix} 0.0047 & 0.0033 \\ 0.0033 & 0.0235 \end{bmatrix}, \\ P_3 &= \begin{bmatrix} 0.0047 & 0.0034 \\ 0.0034 & 0.0236 \end{bmatrix}, P_4 = \begin{bmatrix} 0.0047 & 0.0033 \\ 0.0033 & 0.0235 \end{bmatrix}, \\ P_5 &= \begin{bmatrix} 0.0047 & 0.0033 \\ 0.0033 & 0.0235 \end{bmatrix} \end{aligned}$$

As we know, based on the commonly mode-dependent methods, see, e.g., [1]–[4], [6], there will be 5 controllers since the mode quantity of this example is 5. More importantly, an ideal assumption that the switching signals between controller and other system matrices should be synchronous is also needed. To the contrary, the method in this paper can provide fewer effective controllers whose modes are not necessary synchronous to system matrices', which has larger application

scope. Without loss of generality, set \mathcal{N} is separated into 3 classes such as $\mathcal{M} = \{1, 2, 3\}$. Consequently, it can be computed that $S_2(5, 3) = 25$ with $N = 5$ and $M = 3$, and the classification parameter c is the same to Table I. Then, after applying the above designed controllers in addition to some related parameters, the optimal value $J(\alpha^*(c))$ with its optimal solution $\alpha^*(c)$ of each classification is given in Table II, where $\bar{\tau} = \min_{h \in \mathcal{M}} \{\bar{\tau}_h\}$ is denoted as the ultimate upper sampling bound. It can be seen from this table that there is no optimal solution to some classifications. The reason is that the given parameters \bar{K}_ℓ and P_i computed by (23) and (24) dissatisfy constraint (25), whose possibly solvable algorithms are presented in Remark 4. Moreover, some comparisons between this paper and some similar existing references can be done in the next. In references [18]–[20], a method based on an auxiliary system approach was found and can be used to discuss stochastic systems having a sampling phenomenon in either state or switching or both of them. Since their methods about sampling are the same, without loss of generality, only comparisons between this paper and [18] are enough. Under the above given parameters and by using the method in [18], the maximum sampling upper bound can be computed as $\bar{\tau} = 0.000047$. It is far less than $\bar{\tau} = 0.0174$ which results from the globally optimal solution such as $\alpha_1^* = 0.4581$, $\alpha_2^* = 0.4668$, $\alpha_3^* = -0.0713$, $\alpha_4^* = 1.6570$, $\alpha_5^* = -0.6935$ and $J^* = 3.3820$ and also corresponds to the situation of Table II with $c = 1$. It can be further seen from Table II, all the upper sampling bounds are larger than 0.000047, so long as there exists a suitable solution. Based on these comparisons, it can be concluded that the developed mode classification and optimization method is superior to the above auxiliary system approach for providing a larger sampling interval and improving the efficiency of communication network. Moreover, under the initial condition $x_0 = [2 \ -2]^\top$ and selecting any effective classification such as $\mathcal{N}_1 = \{1\}$, $\mathcal{N}_2 = \{2, 3\}$, $\mathcal{N}_3 = \{4, 5\}$, corresponding to $c = 11$ in Table II, one can compute its controller such as $K_{\mathcal{N}_1} = [0.0880 \ 0.5432]$, $K_{\mathcal{N}_2} = [0.8654 \ 2.2401]$ and $K_{\mathcal{N}_3} = [-0.2397 \ 1.5940]$. Then, the state curves of the closed-loop system are shown in Fig. 4 (a), and the schematic diagram of mode signals is displayed in Fig. 4 (b). Although both state and mode are sampled, the presented controller is still useful in stabilizing a stochastic system.

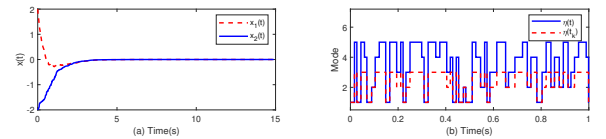


Fig. 4. (a) The state responds of the closed-loop system by controller (3). (b) The simulations of signals $\eta(t)$ and $\eta(t_k)$.

Furthermore, based on the developed Q-HCA in Theorem 2, a suboptimal solution or even an optimal solution can be obtained. In detailed, the iteration processes under different initial points are shown in Fig. 5 with $\epsilon = 10$, where

TABLE II
THE OPTIMIZATION RESULTS OF THEOREM 2 FOR DIFFERENT MODE CLASSIFICATIONS.

c	$\alpha_1^*(c)$	$\alpha_2^*(c)$	$\alpha_3^*(c)$	$\alpha_4^*(c)$	$\alpha_5^*(c)$	$J(\alpha^*(c))$	$\bar{\tau}$
1	0.46	0.47	-0.07	1.66	-0.69	3.3820	0.0174
2	—	—	—	—	—	—	—
3	0.46	2.70	0.05	0.24	-0.12	3.5891	0.0146
4	0.46	1.50	-0.18	0.57	0.45	4.5422	0.0142
5	—	—	—	—	—	—	—
6	0	11.83	-1.70	0.24	0.45	4.8947	0.0136
7	0.68	0.47	-0.09	1.23	0.45	6.4733	0.0123
8	—	—	—	—	—	—	—
9	—	—	—	—	—	—	—
10	0	11.83	-1.70	0.24	0.45	4.8947	0.0137
11	0.46	16.92	-2.65	-0.80	1.09	4.5637	0.0146
12	0.46	-4.97	1.08	1.68	-0.82	3.7408	0.0172
13	0.46	-0.51	0.10	0.75	0.57	5.1091	0.0153
14	-1.88	0.47	0.99	-0.80	1.09	9.2897	0.0123
15	0	0.47	0.67	0.24	-0.18	3.9675	0.0174
16	0	0.47	-0.08	1.01	0.45	4.9755	0.0161
17	—	—	—	—	—	—	—
18	—	—	—	—	—	—	—
19	—	—	—	—	—	—	—
20	0.40	0.08	0.67	0.24	-0.18	4.0838	0.0173
21	-1.88	-4.01	0.99	0.24	1.39	9.4961	0.0123
22	0	16.92	-2.65	0.24	0.45	4.8516	0.0146
23	0	0.59	-0.10	1.06	0.45	5.0880	0.0174
24	-1.88	-1.43	0.99	0.65	0.45	9.6922	0.0091
25	0	16.92	-2.65	0.26	0.45	4.8516	0.0146

the points in the horizontal axis denote as infeasible points. Particularly, the infeasible points in blue means that they have been selected to participate in calculation, while the black ones are only infeasible points but without being selected during the iteration. Moreover, from this simulation, it can be seen that when the initial points are Points 24 and 7 respectively, the final point is Point 12 corresponding to $c = 12$ in Table II, whose cost function is computed as $f^* = 3.7408$. Similarly, one can further know that when the initial point is Point 14, the final point is Point 3 with $f^* = 3.5891$. And when the initial point is Point 7, the final point is Point 1. Especially, in this situation, its cost function is computed as $f^* = 3.3820$ and equals to the globally optimal value $J^* = 3.3820$. Obviously, by comparing Table II and Fig. 5, one can obtain that the computation complexity and time of Fig. 5 by the Q-HCA method is both largely reduced and without computing mode classification one by one. Moreover, it also can be seen from Fig. 5 that the convergence of the Q-HCA method is also ensured, which is asymptotically convergent. In a word, all the statements about the effectiveness and superiority of the developed method in this paper have been shown by the simulations and comparisons.

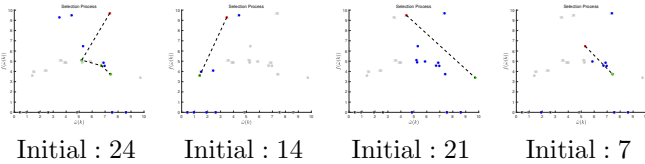


Fig. 5. The effects of Q-HCA in Theorem 2 under different initial points.

Next, some simulations and comparison about Corollary 2

will be done, where set \mathcal{N} is also separated into 3 classes such as $\mathcal{M} = \{1, 2, 3\}$. First, letting $\lambda_1 = -0.8$, $\lambda_2 = -0.9$, $\lambda_3 = -0.5$, $\lambda_4 = -0.6$, $\lambda_5 = 0.4$, $\mu_1 = 1.005$, $\mu_2 = 1.001$, $\mu_3 = 1.009$, $\mu_4 = 1.002$ and $\mu_5 = 1.001$, it can be known from Corollary 2 that $\sum_{i=1}^N \pi_i \left(\lambda_i + \frac{\ln \mu_i}{\bar{\tau}_i} \right) = -0.2549 < 0$. Meanwhile, the control gains computed by conditions (24) and (43) are listed as follows: $\bar{K}_1 = \begin{bmatrix} 1.9651 & 1.6991 \end{bmatrix}$, $\bar{K}_2 = \begin{bmatrix} 1.8758 & 1.1729 \end{bmatrix}$, $\bar{K}_3 = \begin{bmatrix} 4.6357 & 4.8861 \end{bmatrix}$, $\bar{K}_4 = \begin{bmatrix} 6.0120 & 4.4209 \end{bmatrix}$ and $\bar{K}_5 = \begin{bmatrix} 4.7489 & 5.3510 \end{bmatrix}$. Now, more comparisons will be done. In reference [17], a quantity-limited controller was proposed based on mode separation. However, the optimization problem about such mode separation was not considered in the reference. Moreover, it has been shown there that the method in [17] is less conservative than mode-independent ones [7]–[10]. In this situation, only comparisons between Corollary 2 and [17] will be done. Based on [17], it is assumed that modes $\{1\}$ and $\{2\}$ are known, and the rest modes are unknown. In this situation, the above assumption is the same as the following mode classification such as $\mathcal{N}_1 = \{1\}$, $\mathcal{N}_2 = \{2\}$ and $\mathcal{N}_3 = \{3, 4, 5\}$. Then, by using the method in [17], one can obtain the control gains as $K_1 = \begin{bmatrix} 1.7897 & 2.0866 \end{bmatrix}$, $K_2 = \begin{bmatrix} 2.1638 & 4.7343 \end{bmatrix}$ and $K = \begin{bmatrix} -0.1417 & -9.2833 \end{bmatrix}$, and the same cost function (11) about the above designed controllers of [17] is equal to $J = \|K_1\| + \|K_2\| + \|K\| = 17.2387$. Investigating all the optimal function value J^* computed similar to Table II and omitted due to page limitation, it can be seen that the global cost function value after optimization is much smaller and less conservative. In order to further illustrate the superiority, more comparisons are done such that a positive perturbation Δ is only added in block (1,1) of matrix A_1 such as $-2 + \Delta$, and the others remain unchanged. Then, the respective maximum allowable bound about Corollary 2 and [17] under the same λ_1 is given in Table III, which is denoted as Δ_{\max} . It is obvious that Corollary 2 by optimizing the mode classifications is less conservative than [17]. Similar to Fig. 5, some simulations about Q-HCA of Corollary 2 under different initial points are given in Fig. 6 with $\epsilon = 10$. It can be found that when the initial points are Points 21, 22 and 23 respectively, the final point is Point 2 corresponding to $c = 2$ similar to Table II, whose cost function is computed as $f^* = 6.7184$ and equals to the globally optimal value $J^* = 6.7184$. Also, when the initial point is Point 10, the final point is Point 1 with $f^* = 7.1112$ which is a locally optimal value. As a result, similar conclusions about the given Q-HCA in this paper can be obtained and omitted here.

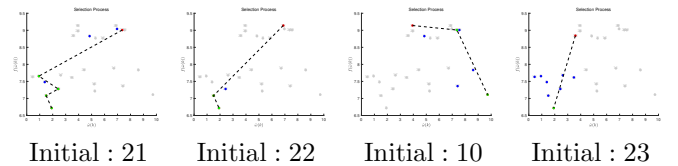


Fig. 6. The effects of Q-HCA in Corollary 2 under different initial points.

Finally, more comparisons will be done to further demon-

TABLE III
THE MAXIMUM ALLOWABLE BOUND Δ_{\max} UNDER λ_1 TAKING DIFFERENT VALUES.

λ_1	-0.1	-0.2	-0.3	-0.4	-0.5
Corollary 2	5.595	5.562	5.529	5.496	5.463
[17]	4.341	4.302	4.263	4.224	4.185
	-0.6	-0.7	-0.8	-0.9	-1
	5.430	5.397	5.364	5.331	5.298
	4.146	4.108	4.069	4.030	3.991

strate the advantages of the method proposed here. As we know, a sampled system can be discussed by transforming the sampled term into a time-varying delay term. Similarly, when both state and switching are sampled, a input delay approach was used in [33]. Unfortunately, its method has some limits such that it cannot be used to stabilize an MJS having non-negative eigenvalues. Meanwhile, an augmented system approach was presented in [21] and can be used to deal with MJSs with sampled state and switching simultaneously. Due to it considering about MJSs instead of semi-Markovian jump systems (SMJSs) and in order to make some comparisons, it is assumed that the related system in this example is reduced to be an MJS. In detail, all the system parameters are the same as the above SMJS, except the transition rate matrix of MJSs is given as

$$\Lambda = \begin{bmatrix} -3 & 1 & 0.5 & 0.2 & 1.3 \\ 1 & -3 & 0.5 & 1.3 & 0.2 \\ 1 & 0.5 & -2 & 0.25 & 0.25 \\ 1.5 & 0.2 & 0.9 & -4 & 1.4 \\ 0.4 & 0.3 & 0.2 & 0.1 & -1 \end{bmatrix}$$

Then, based on [21] with given control gains such as $K_1 = [0.1710 \ 1.1218]$, $K_2 = [0.1853 \ 0.8848]$, $K_3 = [0.9524 \ 4.8487]$, $K_4 = [0.6209 \ 2.8994]$ and $K_5 = [0.2302 \ 3.7272]$, one can get the upper sampling bound is $\bar{\tau} = 0.0028$. On the other hand, it can be known that similar results about MJSs can be easily obtained from the results about SMJSs. Thus, under the same control gains given above, and selecting $\alpha_1 = 0.335$, $\alpha_2 = 29.5068$, $\alpha_3 = -4.8934$, $\alpha_4 = -0.6157$ and $\alpha_5 = 0.9425$ with a specific mode classification such as $\mathcal{N}_1 = \{1\}$, $\mathcal{N}_2 = \{2, 3\}$ and $\mathcal{N}_3 = \{4, 5\}$, the upper sampling bound based on Theorem 1 without any optimization is equal to $\bar{\tau} = 0.01324$. It shows that the method in this paper can provide a larger sampling bound and is less conservative than [21].

V. CONCLUSIONS

In this study, the sampled stabilization for stochastic jump systems has been investigated by a mode classification optimization method. A new mode-classification-based controller has been proposed to handle the general sampling situation that both controller's state and switching are sampled. So as to find the best classification for stabilization, an optimization problem has been proposed to provide a globally optimal solution among such classifications. An Q-HCA algorithm has been established to provide an optimal attenuation coefficient,

whose monotonicity and convergency have been both satisfied. Finally, the utility and advantage of this paper have been obviously illustrated and shown by a numerical example.

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VI. APPENDIX

A. Proof of Theorem 1

First of all, $N_\eta(t_0, t)$, $\forall t \geq 0$, is defined to be the switching quantity of $\eta(t)$ on interval $[t_0, t]$. For any $N_\eta(t_0, t) = n$, $\forall t \geq 0$ and $n \in \mathbb{N}$, one concludes that $[t_0, t] = \bigcup_{i=0}^{n-1} [T_i, T_{i+1}) \cup [T_n, t]$ and $t \in [T_n, T_{n+1})$. Consequently, one gets $[t_0, t] = \bigcup_{j=0}^{k-1} [t_j, t_{j+1}) \cup [t_k, t]$ and $t \in [t_k, t_{k+1})$, $k \in \mathbb{N}$, from event-triggered mechanism (4) under Assumption 1. Moreover, one can further obtain $k \leq n$ such that every subinterval of $\bigcup_{i=0}^{n-1} [T_i, T_{i+1}) \cup [T_n, t]$ such as $[T_i, T_{i+1})$ or $[T_n, t]$ certainly belongs to one subinterval of $\bigcup_{j=0}^{k-1} [t_j, t_{j+1}) \cup [t_k, t]$ such as $[t_j, t_{j+1})$ or $[t_k, t]$ uniquely. Alternatively, $\{t_k\}_{k \in \mathbb{N}}$ established based on principle (4) is a subsequence of $\{T_n\}_{n \in \mathbb{N}}$. As a result, and without loss of generality, it is assumed that

$$[t_j, t_{j+1}) = \bigcup_{q=0}^{N_\eta(t_j, t_{j+1})-1} [T_{N_\eta(t_0, t_j)+q}, T_{N_\eta(t_0, t_j)+q+1}) \quad (51)$$

where $t_j = T_{N_\eta(t_0, t_j)}$ and $t_{j+1} = T_{N_\eta(t_0, t_{j+1})}$. Particularly, $[t_k, t]$ is denoted as

$$[t_k, t] = \bigcup_{q=0}^{N_\eta(t_k, t)-1} [T_{N_\eta(t_0, t_k)+q}, T_{N_\eta(t_0, t_k)+q+1}) \cup [T_{N_\eta(t_0, t)}, t] \quad (52)$$

In addition, for $\forall t \in [t_k, t_{k+1})$, $\forall k \in \mathbb{N}$, one can conclude that $\eta(t) \in \mathcal{N}_{\eta(t_k)}$, if $\eta(t_k) \in \mathcal{M}$. Then, after applying (3), the system (1) closed by an event-triggered controller described by (3)-(8) is equal to

$$\begin{cases} \dot{x}(t) = A_{\eta(t)}x(t) + B_{\eta(t)} \sum_{\ell=1}^N \alpha_\ell^{[\nu(\eta(t_k))]} \bar{K}_\ell x(t_k) \\ x(t_k) = x(t_k^-) \end{cases} \quad (53)$$

Accordingly, choose the Lyapunov function as $V(x(t), \eta(t)) = x^T(t)P_{\eta(t)}x(t)$ and based on (51)-(53) and condition (20), for any $t \in [T_{N_\eta(t_0, t)}, T_{N_\eta(t_0, t+1)})$, one gets that

$$\begin{aligned} & V(x(T_{N_\eta(t_0, t_k)+1}), \eta_{N_\eta(t_0, t_k)}) \\ & \times e^{-\lambda_{N_\eta(t_0, t_k)}(T_{N_\eta(t_0, t_k)+1} - T_{N_\eta(t_0, t_k)})} \\ & = V(x(T_{N_\eta(t_0, t_k)}), \eta_{N_\eta(t_0, t_k)}) \\ & + \int_{T_{N_\eta(t_0, t_k)}}^{T_{N_\eta(t_0, t_k)+1}} d(V(x(s), \eta_{N_\eta(t_0, t_k)})) \\ & \times e^{-\lambda_{N_\eta(t_0, t_k)}(s - T_{N_\eta(t_0, t_k)})} \\ & \leq V(x(T_{N_\eta(t_0, t_k)}), \eta_{N_\eta(t_0, t_k)}) \\ & \leq \mu_{N_\eta(t_0, t_k)-1} V(x(T_{N_\eta(t_0, t_k)}), \eta_{N_\eta(t_0, t_k)-1}) \end{aligned} \quad (54)$$

\vdots

$$\begin{aligned} & V(x(T_{N_\eta(t_0, t)}), \eta_{N_\eta(t_0, t)-1}) \\ & \times e^{-\lambda_{N_\eta(t_0, t)-1}(T_{N_\eta(t_0, t)} - T_{N_\eta(t_0, t)-1})} \\ & = V(x(T_{N_\eta(t_0, t)-1}), \eta_{N_\eta(t_0, t)-1}) \\ & + \int_{T_{N_\eta(t_0, t)-1}}^{T_{N_\eta(t_0, t)}} d(V(x(s), \eta_{N_\eta(t_0, t)-1})) \\ & \times e^{-\lambda_{N_\eta(t_0, t)-1}(s - T_{N_\eta(t_0, t)-1})} \\ & \leq V(x(T_{N_\eta(t_0, t)-1}), \eta_{N_\eta(t_0, t)-1}) \\ & \leq \mu_{N_\eta(t_0, t)-2} V(x(T_{N_\eta(t_0, t)-1}), \eta_{N_\eta(t_0, t)-2}) \\ & V(x(t), \eta_{N_\eta(t_0, t)}) e^{-\lambda_{N_\eta(t_0, t)}(t - T_{N_\eta(t_0, t)})} \\ & = V(x(T_{N_\eta(t_0, t)}), \eta_{N_\eta(t_0, t)}) \\ & + \int_{T_{N_\eta(t_0, t)}}^t d(V(x(s), \eta_{N_\eta(t_0, t)}) e^{-\lambda_{N_\eta(t_0, t)}(s - T_{N_\eta(t_0, t)})}) \\ & \leq V(x(T_{N_\eta(t_0, t)}), \eta_{N_\eta(t_0, t)}) \\ & \leq \mu_{N_\eta(t_0, t)-1} V(x(T_{N_\eta(t_0, t)}), \eta_{N_\eta(t_0, t)-1}) \end{aligned} \quad (55)$$

Particularly, the above inequalities holding also need the following condition

$$\begin{aligned} & \left[x^\top(t)P_{\eta(t)} \left(A_{\eta(t)}x(t) + B_{\eta(t)} \sum_{\ell=1}^N \alpha_\ell^{[h]} \bar{K}_\ell x(t_k) \right) \right]^* \\ & \leq \lambda_{\eta(t)} x^\top(t)P_{\eta(t)}x(t) \end{aligned} \quad (56)$$

where $\forall t \in [t_k, t_{k+1})$, $\eta(t_k) = h \in \mathcal{M}$ and $\eta(t) \in \mathcal{N}_h$. On the other hand, let $\delta(t) = x(t) - x(t_k)$, $\forall t \in [t_k, t_{k+1})$, it can be concluded from (53) that

$$\begin{cases} \dot{\delta}(t) = A_{\eta(t)}\delta(t) + \bar{A}_{\eta(t)}^{[h]}x(t_k) \\ \delta(t_k) = 0 \end{cases}, \forall t \in [t_k, t_{k+1}), \forall k \in \mathbb{N} \quad (57)$$

Its solution is computed as

$$\delta(t) = e^{\int_{t_k}^t A_{\eta(s)}ds} \delta(t_k) + \int_{t_k}^t e^{\int_{\theta}^t A_{\eta(s)}ds} \bar{A}_{\eta(\theta)}^{[h]}x(t_k)d\theta \quad (58)$$

which implies

$$\begin{aligned} \|\delta(t)\| & \leq \int_{t_k}^t \|e^{\int_{\theta}^t A_{\eta(s)}ds}\| \|\bar{A}_{\eta(\theta)}^{[h]}\| \|x(t_k)\| d\theta \\ & \leq \max_{\ell \in \mathcal{N}_h} \left\{ \|\bar{A}_\ell^{[h]}\| \right\} \|x(t_k)\| \int_{t_k}^t e^{\max_{\ell \in \mathcal{N}_h} \{\chi_{A_\ell}\}(t-\theta)} d\theta \\ & = \|\hat{A}_h\| \int_{t_k}^t e^{\hat{\chi}_h(t-\theta)} d\theta \|x(t_k)\| \end{aligned} \quad (59)$$

It can be known that function $\zeta_h(t - t_k) \triangleq \int_{t_k}^t e^{\hat{\chi}_h(t-\theta)} d\theta$, $\forall t \in [t_k, t_{k+1})$, $\forall k \in \mathbb{N}$, is monotonically increasing with t and $\zeta(0) = 0$. Based on Assumption 1, one has

$$\zeta_h(t - t_k) \leq \zeta_h(s_{k+1}) \leq \zeta_h(\bar{\tau}_h) \quad (60)$$

Then, one obtains that

$$\|\delta(t)\| \leq \|\hat{A}_h\| \zeta_h(\bar{\tau}_h) (\|\delta(t)\| + \|x(t)\|) \quad (61)$$

When function $\zeta_h(\bar{\tau}_h)$ satisfies

$$\zeta_h(\bar{\tau}_h) \leq \frac{1}{\|\hat{A}_h\|} \frac{\sigma_h}{1 + \sigma_h} \quad (63)$$

where $\sigma_h > 0$ is a suitable design parameter, inequality (62) becomes to be

$$\|\delta(t)\| \leq \sigma_h \|x(t)\|, \forall t \in [t_k, t_{k+1}), \forall k \in \mathbb{N}, \eta(t_k) = h \in \mathcal{M} \quad (64)$$

Under condition $x(t_k) = x(t) - \delta(t)$, inequality (57) is equivalent to

$$\begin{aligned} & x^T(t) \left[\left(P_{\eta(t)} \bar{A}_{\eta(t)}^{[h]} \right)^* - \lambda_{\eta(t)} P_{\eta(t)} \right] x(t) \\ & - 2x^T(t) P_{\eta(t)} B_{\eta(t)} \sum_{\ell=1}^N \alpha_{\ell}^{[\nu(\eta(t_k))]} \bar{K}_{\ell} \delta(t) \leq 0 \end{aligned} \quad (65)$$

Based on condition (19) and (64), inequality (65) is guaranteed. Thus, conditions (54)-(56) are satisfied. They finally yield that for $\forall t \geq 0$,

$$\begin{aligned} V(x(t), \eta(t)) & \leq \mu_{\eta_{N_{\eta}(t_0, t)-1}} V(x(T_{N_{\eta}(t_0, t)}), \eta_{N_{\eta}(t_0, t)-1}) \\ & \quad \times e^{\lambda_{\eta_{N_{\eta}(t_0, t)}}(t - T_{N_{\eta}(t_0, t)})} \\ & \leq \mu_{\eta_{N_{\eta}(t_0, t)-2}} \mu_{\eta_{N_{\eta}(t_0, t)-1}} \\ & \quad \times V(x(T_{N_{\eta}(t_0, t)-1}), \eta_{N_{\eta}(t_0, t)-2}) \\ & \quad \times e^{\lambda_{\eta_{N_{\eta}(t_0, t)}}(t - T_{N_{\eta}(t_0, t)})} \\ & \quad \times e^{\lambda_{\eta_{N_{\eta}(t_0, t)-1}}(T_{N_{\eta}(t_0, t)} - T_{N_{\eta}(t_0, t)-1})} \\ & \quad \vdots \\ & \leq V(x_0, \eta_0) \prod_{k=1}^{N_{\eta}(t_0, t)} \mu_{\eta_k} e^{\int_0^t \lambda_{\eta(s)} ds} \\ & = V(x_0, \eta_0) \prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \end{aligned} \quad (66)$$

Then, one could obtain that

$$\|x(t)\| \leq \frac{\max_{i \in \mathcal{N}} \{\lambda_{\max}(P_i)\}}{\min_{i \in \mathcal{N}} \{\lambda_{\min}(P_i)\}} \|x_0\| \prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \quad (67)$$

It is obvious from (21) that there is always a scalar $\epsilon_1 > 0$ such that

$$\sum_{i=1}^N \pi_i \lambda_i + \sum_{i=1}^N \left(\frac{\pi_i}{\hat{\tau}_i} + \epsilon_1 \right) \ln \mu_i < 0 \quad (68)$$

Then, similar to the proof of Theorem 1 in [26], one knows that if $\forall i \in \mathcal{N}$ and ϵ_1 defined in (68), there exists a positive constant $H(\epsilon_1)$ such that if $t \geq H(\epsilon_1)$, then

$$N_i(t_0, t) \leq \left(\frac{\pi_i}{\hat{\tau}_i} + \epsilon_1 \right) t, \text{ a.s.} \quad (69)$$

Accordingly, for $t \geq H(\epsilon_1)$, one has

$$\begin{aligned} \prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} & = e^{\sum_{i=1}^N N_i(t_0, t) \ln \mu_i + \int_0^t \lambda_{\eta(s)} ds} \\ & \leq e^{\sum_{i=1}^N \left(\frac{\pi_i}{\hat{\tau}_i} + \epsilon_1 \right) (\ln \mu_i) t + \int_0^t \lambda_{\eta(s)} ds}, \text{ a.s.} \end{aligned} \quad (70)$$

Based on the strong law of large numbers, one obtains that

$$\lim_{t \rightarrow \infty} \frac{1}{t} \int_0^t \lambda_{\eta(s)} ds = \sum_{i=1}^N \lambda_i \pi_i, \text{ a.s.} \quad (71)$$

Then, it can be concluded from (21) that

$$\lim_{t \rightarrow \infty} \sum_{i=1}^N \left(\frac{\pi_i}{\hat{\tau}_i} + \epsilon_1 \right) (\ln \mu_i) t + \int_0^t \lambda_{\eta(s)} ds = -\infty, \text{ a.s.} \quad (72)$$

Consequently, one gets that

$$\prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} = 0, \text{ a.s.} \quad (73)$$

Based on the above equation, it can be known that for any given scalars $\gamma > 0$ and $\xi > 0$, there exists a random variable $\varphi(\gamma, \xi)$ such that

$$\begin{aligned} & \mathbb{P} \left(\sup_{t \geq \varphi(\gamma, \xi)} \left(\prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \right) \right. \\ & \quad \left. < \frac{\min_{i \in \mathcal{N}} \{\lambda_{\min}(P_i)\} \xi}{\max_{i \in \mathcal{N}} \{\lambda_{\max}(P_i)\} \gamma} \right) = 1 \end{aligned} \quad (74)$$

Then, for any $\|x_0\| < \gamma$ and combining conditions (67) and (74), one can conclude condition C_2 . Moreover, it can be concluded from (74) that

$$\sup_{t \geq 0} \left(\prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \right) < \infty, \text{ a.s.} \quad (75)$$

Obviously, for any $\epsilon \in (0, 1)$, there is always $\rho > 0$ such that

$$\mathbb{P} \left(\sup_{t \geq 0} \left(\prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \right) < \rho \right) > 1 - \epsilon. \quad (76)$$

By selecting $\delta = \frac{\min_{i \in \mathcal{N}} \{\lambda_{\min}(P_i)\} \epsilon}{\max_{i \in \mathcal{N}} \{\lambda_{\max}(P_i)\} \rho}$ and letting $\|x_0\| < \delta$, and considering inequality (67) under condition (76), one has

$$\begin{aligned} & \mathbb{P} \left(\sup_{t \geq 0} \|x(t)\| \leq \frac{\max_{i \in \mathcal{N}} \{\lambda_{\max}(P_i)\}}{\min_{i \in \mathcal{N}} \{\lambda_{\min}(P_i)\}} \|x_0\| \right. \\ & \quad \times \sup_{t \geq 0} \left(\prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \right) \\ & \quad \left. < \frac{\epsilon}{\rho} \sup_{t \geq 0} \left(\prod_{i=1}^N \mu_i^{N_i(t_0, t)} e^{\int_0^t \lambda_{\eta(s)} ds} \right) < \epsilon \right) \\ & > 1 - \epsilon \end{aligned} \quad (77)$$

which implies condition C_1 . Due to conditions C_1 and C_2 both satisfying, the resulting closed-loop system is finally GAS a.s..

As for condition (63) with $\eta(t_k) = h \in \mathcal{M}, \forall t \in [t_k, t_{k+1})$, it is implied by

$$\int_{t_k}^t e^{\hat{\chi}_h(t-\theta)} d\theta \leq \frac{1}{\|\hat{A}_h\|} \frac{\sigma_h}{1 + \sigma_h} \quad (78)$$

When $\hat{\chi}_h \leq 0$, it could be guaranteed by

$$s_k \leq \bar{\tau}_h \leq \frac{1}{\|\hat{A}_h\|} \frac{\sigma_h}{1 + \sigma_h} \quad (79)$$

which is actually the former condition of (22). To the contrary, condition (63) under $\hat{\chi}_h > 0$ is ensured by

$$\frac{1}{\hat{\chi}_h} e^{\hat{\chi}_h(t-s)} \Big|_{t_k}^{t_k} \leq \frac{1}{\hat{\chi}_h} (e^{\hat{\chi}_h \bar{\tau}_h} - 1) \leq \frac{1}{\|\hat{A}_h\|} \frac{\sigma_h}{1 + \sigma_h} \quad (80)$$

It is equivalent to the latter condition of (22). This completes the proof. \square

B. Proof of Theorem 2

On the one hand, first of all, it is obvious that the augmented Lagrangian function (28) is convex. Then, based on the result given in [30] on Page 108, it can be concluded that there exists a globally optimal solution $(\alpha^*(c), \gamma^*(c), \varsigma^*(c))$ to (28) which can be computed by solving conditions (31)-(33) and defined as $\mathcal{L}(\alpha^*(c), \gamma^*(c), \varsigma^*(c), \phi) \triangleq \min_{\alpha(c), \gamma(c), \varsigma(c)} \mathcal{L}(\alpha(c), \gamma(c), \varsigma(c), \phi)$. Particularly, it can be seen that (28) is strictly convex about $\alpha(c)$ such as $\nabla_{\alpha(c)}^2 \mathcal{L}(\alpha(c), \gamma(c), \varsigma(c), \phi) > 0$. In other words, there is only one globally optimal solution $\alpha^*(c)$ satisfying (31)-(33). Meanwhile, according to reference [31] on Page 97, it can be known that the optimization problem (11) constrained by (27) has the same optimal value as the optimization problem (28). That is

$$J(\alpha^*(c)) = \mathcal{L}(\alpha^*(c), \gamma^*(c), \varsigma^*(c), \phi) \quad (81)$$

Finally, one can easily obtain the globally optimal value of (11) with constraint (27) described by (34), whose corresponding optimal solution is obviously computed by (35).

On the other hand, it has been shown in [32] that $Q_k(s, a)$ generated by (36) can theoretically converge to $Q^*(s, a)$ with probability 1 when $k \rightarrow \infty$. As a result, an optimal taking action \mathcal{A}_k^* , $k = 1, 2, \dots$, maximizing Q-value (30) can be established, which is actually an optimal attenuation coefficient δ_k^* . Then, a preferable searching radius L_k^{pre} , $k = 1, 2, \dots$, can be gotten based on the iterative update rule (18), while the initial condition $\hat{\omega}(1)$ is given to be $\forall \hat{\omega}(1) \in \Omega$ but $\hat{\omega}(0)$ should be given such as one defined in (17). Based on the obtained L_k^{pre} , a preferably nominal solution referred as $\hat{\omega}^{\text{pre}}(k)$, $k = 2, 3, \dots$, can be gotten by using HCA. Finally, by repeating the above process along with iteration number k having a maximum value k_{\max} , two optimal sequences $\{\mathcal{A}_k^*\}_{k=1}^{k_{\max}}$ and $\{\delta_k^*\}_{k=1}^{k_{\max}}$ can be established based on (36), while two preferable sequence $\{L_k^*\}_{k=1}^{k_{\max}}$ and $\{\hat{\omega}^{\text{pre}}(k)\}_{k=1}^{k_{\max}}$

with $\hat{\omega}^{\text{pre}}(1) = \hat{\omega}(1)$ can also be developed by (18). The detailed repeating process is given as follows:

$$\begin{aligned} \hat{\omega}(1) &\rightarrow \delta_1^* \xrightarrow{\hat{\omega}(1)} L_1 \rightarrow \hat{\omega}(2) \rightarrow \dots \rightarrow \\ \hat{\omega}(k_{\max}) &\rightarrow \delta_{k_{\max}}^* \xrightarrow{\hat{\omega}(k_{\max})} L_{k_{\max}} \end{aligned}$$

Consequently, based on the characteristics of Q-HCA, it can be obviously known that sequence $\{\hat{\omega}^{\text{pre}}(k)\}_{k=1}^{k_{\max}}$ will have a good convergence characteristic such as (37). Meanwhile, the monotonicity of (38) can be guaranteed. Moreover, since all the elements of $\{\hat{\omega}^{\text{pre}}(k)\}_{k=1}^{k_{\max}}$ belong to Ω but is only a part of Ω , it is obviously concluded that the obtained optimal value f^* will be no less than J^* obtained by (34). In this situation, it can be said that the above obtained $\hat{\omega}^*$ is only a locally optimal solution. Finally, the iteration process of the developed Q-HCA is given below:

Algorithm 1: Improved hill climbing algorithm based on Q-learning

Initial Q-values with Q-HCA, parameter ϵ , iterative index $k = 1$,
 Select a starting state $S_1 = \hat{\omega}(1)$, calculate L_1 by the parameter ϵ
while ($k \leq$ maximum number of iterations)
 while S_k is not a termination state S_{end}
 Choose \mathcal{A}_k according to the $\max Q(S_k)$
 Take action \mathcal{A}_k
 Return reward \mathcal{R}_{k+1} by the equation (29)
 Update Q-value by the equation (36)
 Update L_k by the equation (19)
 Update $\Theta(\hat{\omega}(k))$ of $\hat{\omega}(k)$ corresponding to S_k
 by the equation (14)
 if $\Theta(\hat{\omega}(k))$ is not empty set
 Generate a new state by $g_{\hat{\omega}(k)}(\hat{\omega}'(k))$ of $\hat{\omega}(k)$
 corresponding to S_k
 Move to new state
 else if $\Theta(\hat{\omega}(k))$ is empty set
 Move to the termination state S_{end}
 end if
end while
end while

This completes the proof. \square