Data-Based Optimal Control for Weakly Coupled Nonlinear Systems Using Policy Iteration

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Abstract—In this paper, a data-based online learning algorithm is established to solve the optimal control problem for weakly coupled continuous-time nonlinear systems with completely unknown dynamics. Using the weak coupling theory, we reformulate the original problem into three reduced-order optimal control problems. We establish an online model-free integral policy iteration algorithm to solve the decoupled optimal control problems without system dynamics. To implement the data-based online learning algorithm, the actor-critic technique based on neural networks and the least squares method are used. Two simulation examples are given to verify the effectiveness of the developed algorithm.

Index Terms—Adaptive dynamic programming (ADP), neural networks (NNs), optimal control, policy iteration (PI), unknown dynamics, weakly coupled systems.

I. INTRODUCTION

N THE real world, many large-scale systems are naturally weakly coupled, such as electrical networks, transportation systems, chemical reactors, and power systems. For these real physical systems, a traditional challenge is the optimal control problem. A common approach is to split this large-scale optimal control problem into some decoupled subproblems using the decentralized control method [1], [2]. While the coupling effects are usually neglected and the obtained control laws may do not have ideal performance. In 1969, Kokotović *et al.* [3] introduced the weakly coupled linear systems to the control systems community. Since then, many theoretical aspects of the optimal control problem for weakly coupled systems have been studied. Gajić and Shen [4], [5] obtained the optimal

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control law through a decoupling transformation which leads to solving two independent reduced-order optimal control problems. For weakly coupled bilinear systems, the optimal control problem has also been solved in a similar way [6], [7]. Jiang and Jiang [8] presented a new approach to decouple the weakly coupled large-scale linear systems and accomplished the stability analysis using the small-gain theory. The optimal control law of the nonlinear systems can be obtained by solving the Hamilton-Jacobi-Bellman (HJB) equations. However, due to the intractable form of the HJB equations, obtaining closed-form optimal controllers by directly solving the HJB equations is difficult. By using the reduced-order scheme and the successive Galerkin approximation (SGA), the optimal control law for the weakly coupled nonlinear system has been constructed based on the solutions of two independent reduced-order HJB equations [9]. Carrillo et al. [10] proposed a learning algorithm to derive the optimal control law using a three-critics/four-actors approximator structure with system dynamics. For large-scale real physical systems, it is difficult to obtain the exact knowledge of system dynamics. Therefore, a kind of data-based algorithms is needed to solve the optimal control problem with unknown system dynamics.

Dynamic programming provides a principled method for determining optimal control laws for dynamical systems in the case of completely known dynamics. While due to the "curse of dimensionality" [11], it is often computationally untenable to obtain the optimal control laws. Among the methods of solving optimal control problems, adaptive dynamic programming (ADP), and reinforcement learning (RL) relax the need for a complete and exact model of the dynamical systems by using compact parameterized approximators. ADP has received increasing attention due to its learning capabilities [12]-[30]. RL is an effective computational method and it can find the optimal policy interactively [31]-[34]. In the existing literature of ADP-based and RL-based optimal control, either policy iteration (PI) or value iteration is utilized to solve the HJB equation. Vrabie and Lewis [35] established an integral RL algorithm to obtain direct adaptive optimal control for nonlinear continuous-time systems with partial system dynamics. Liu et al. [36] developed an online synchronous approximate optimal learning algorithm based on PI to solve a multiplayer nonzero-sum game without the exact knowledge of dynamical systems. Jiang and Jiang [37] presented a novel PI method to solve optimal control problems for linear systems with completely unknown dynamics. Jiang and Jiang [38] presented a novel method of global ADP for the adaptive

optimal control of nonlinear polynomial systems to achieve global asymptotic stability. Without the exact knowledge of system dynamics, Lee *et al.* [39] derived a model-free integral *Q*-learning approach for nonlinear system.

Although ADP-based and RL-based algorithms are widely used, there are few related results which can be used to tackle the optimal control problem for the weakly coupled nonlinear systems. The novelty of this paper is that we establish a data-based learning algorithm to solve this problem with completely unknown dynamics. By partitioning the HJB equation, the original optimal control problem of the weakly coupled systems is reformulated into three reduced-order optimal control problems. We establish the model-free integral PI algorithm to solve the decoupled optimal control problems without system dynamics. The actor-critic technique based on neural networks (NNs) and the least squares method are used to implement the derived online learning algorithm.

The rest of this paper is organized as follows. In Section II, the optimal control problem for the weakly coupled nonlinear systems is described. In Section III, the original problem is reformulated into three reduced-order optimal control problems and a model-free integral PI algorithm using online learning manner with unknown system dynamics is established. Two simulation examples are provided to demonstrate the applicability of the established optimal control policy in Section IV. In Section V, we conclude this paper with a few remarks.

II. PROBLEM FORMULATION

In this paper, we consider the continuous-time nonlinear system with weakly coupled structure

$$\begin{bmatrix} \dot{x}_{1}(t) \\ \dot{x}_{2}(t) \end{bmatrix} = \begin{bmatrix} f_{11}(x_{1}) + \varepsilon f_{12}(x) \\ \varepsilon f_{21}(x) + f_{22}(x_{2}) \end{bmatrix} + \begin{bmatrix} g_{11}(x_{1}) & \varepsilon g_{12}(x) \\ \varepsilon g_{21}(x) & g_{22}(x_{2}) \end{bmatrix} \begin{bmatrix} u_{11}(t) + \varepsilon u_{12}(t) \\ \varepsilon u_{21}(t) + u_{22}(t) \end{bmatrix}$$
(1)

where $x_1(t) \in \mathbb{R}^{n_1}$ and $x_2(t) \in \mathbb{R}^{n_2}$ are the system state vectors, $u_{11}(t), u_{12}(t) \in \mathbb{R}^{m_1}$ and $u_{21}(t), u_{22}(t) \in \mathbb{R}^{m_2}$ are the control input vectors, n_1, n_2, m_1 , and m_2 are positive integers, and ε is a small positive weak coupling parameter. Using the following expressions:

$$x(t) = \begin{bmatrix} x_1(t) \\ x_2(t) \end{bmatrix}, f(x) = \begin{bmatrix} f_{11}(x_1) + \varepsilon f_{12}(x) \\ \varepsilon f_{21}(x) + f_{22}(x_2) \end{bmatrix}$$
$$g(x) = \begin{bmatrix} g_{11}(x_1) & \varepsilon g_{12}(x) \\ \varepsilon g_{21}(x) & g_{22}(x_2) \end{bmatrix}, u(t) = \begin{bmatrix} u_{11}(t) + \varepsilon u_{12}(t) \\ \varepsilon u_{21}(t) + u_{22}(t) \end{bmatrix}$$

the system dynamics (1) can be rewritten as

$$\dot{x}(t) = f(x) + g(x)u(t). \tag{2}$$

We assume that the system (2) is controllable, $f: \mathbb{R}^n \to \mathbb{R}^n$ and $g: \mathbb{R}^n \to \mathbb{R}^{n \times m}$ are Lipschitz continuous on the set $\Omega \subseteq \mathbb{R}^n$, where $n = n_1 + n_2$, $m = m_1 + m_2$, and there must exist a continuous control policy which asymptotically stabilizes the system. Additionally, we let the following assumptions hold through out this paper.

Assumption 1: The state vector x = 0 is the equilibrium of the system.

Assumption 2: The functions $f(\cdot)$ and $g(\cdot)$ are differentiable in their arguments, and f(0) = 0.

Assumption 3: The feedback control vector u(x) = 0 when x = 0.

According to the optimal control theory, we known that solving the optimal control problem is equal to find the optimal control policy $u^*(x(t))$ which minimizes the expenditure of control effort. For this, we define the value function as

$$V(x(t)) = \int_{t}^{\infty} \left[x^{\mathsf{T}}(\tau) Q x(\tau) + u^{\mathsf{T}}(\tau) R u(\tau) \right] d\tau \tag{3}$$

where $Q \in \mathbb{R}^{n \times n}$ and $R \in \mathbb{R}^{m \times m}$ are positive definite symmetric matrices, and $r(x, u) = x^{\mathsf{T}}(t)Qx(t) + u^{\mathsf{T}}(t)Ru(t)$ is the utility function. The matrices Q and R have the following weakly coupled structures:

$$Q = \begin{bmatrix} Q_1 & \varepsilon Q_{\varepsilon} \\ \varepsilon Q_{\varepsilon}^{\mathsf{T}} & Q_2 \end{bmatrix}, \qquad R = \begin{bmatrix} R_1 & 0 \\ 0 & R_2 \end{bmatrix}$$

where $Q_1 \in \mathbb{R}^{n_1 \times n_1}$, $Q_2 \in \mathbb{R}^{n_2 \times n_2}$, $R_1 \in \mathbb{R}^{m_1 \times m_1}$, and $R_2 \in \mathbb{R}^{m_2 \times m_2}$ are positive definite symmetric matrices, and $Q_{\varepsilon} \in \mathbb{R}^{n_1 \times n_2}$. We know that the designed feedback control policy u(x(t)) must not only stabilize the system on Ω , but also guarantee that the value function (3) is finite. That is to say, the control policy must be admissible.

Definition 1: A control policy u(x) is said to be admissible with respect to (3) on Ω , denoted by $u(x) \in \Psi(\Omega)$ [$\Psi(\Omega)$ is the set of all admissible control laws], if u(x) is continuous on Ω , u(0) = 0, u(x) stabilizes the system (2) on Ω , and V(x(t)) is finite $\forall x_0 \in \Omega$, where x_0 is the initial system state [40].

According to the optimal control theory, the optimal value function is defined as

$$V^*(x(t)) = \min_{u \in \Psi(\Omega)} \int_t^{\infty} \left[x^{\mathsf{T}}(\tau) Q x(\tau) + u^{\mathsf{T}}(\tau) R u(\tau) \right] d\tau.$$

We define the Hamiltonian function of system (2) as

$$H(x, u, V_x) = V_x^{\mathsf{T}} [f(x) + g(x)u] + r(x, u)$$
 (4)

with V(0) = 0, and the term $V_x = \partial V(x)/\partial x$ denotes the partial derivative of the value function with respect to the state. We minimize the Hamiltonian function (4) to obtain the optimal control policy

$$u^*(x) = \arg\min_{u \in \Psi(\Omega)} H(x, u, V_x) = -\frac{1}{2} R^{-1} g^{\mathsf{T}}(x) V_x^*.$$
 (5)

Using the optimal control policy $u^*(x)$, the optimal value function $V^*(x)$ can be described as the unique positive-definite solution of the following HJB equation:

$$0 = V_x^{*T} [f(x) + g(x)u^*(x)] + r(x, u^*(x)).$$
 (6)

Remark 1: In the class of nonlinear systems, the optimal control scheme is based on the solution of the HJB equation (6). Because the solution of HJB equation for nonlinear systems can hardly be found, the SGA method [41], [42] is developed. However, the SGA method has the weakness that the complexity of computation increases rapidly with the order of the system, where the order indicates the dimension of a system, i.e., n. Kim and Lim [9] established the optimal control from two independent reduced-order HJB equations using the

SGA method. Due to the difficulties when obtaining the exact knowledge of the system dynamics, model-free methods are more practical. Motivated by the weak coupling theory and [9], we establish a novel model-free algorithm to derive the optimal control based on three reduced-order HJB equations.

III. COMPUTATIONAL CONTROLLER DESIGN USING DATA-BASED ONLINE LEARNING ALGORITHM

By partitioning the HJB equation, the original problem of the weakly coupled systems is reformulated into three reducedorder optimal control problems. We established the model-free integral PI algorithm to solve the decoupled optimal control problems without system dynamics. The actor-critic technique based on NNs and the least squares method are used to implement the derived online learning algorithm.

A. Problem Transformation

The value function (3) can be partitioned as

$$V(x(t)) = V_1(x_1(t)) + V_2(x_2(t)) + \varepsilon V_{\varepsilon}(x(t))$$

where

$$\begin{split} V_{1}(x_{1}(t)) &= \int_{t}^{\infty} \left[x_{1}^{\mathsf{T}} Q_{1} x_{1} + u_{11}^{\mathsf{T}} R_{1} u_{11} \right] \mathrm{d}\tau \\ V_{2}(x_{2}(t)) &= \int_{t}^{\infty} \left[x_{2}^{\mathsf{T}} Q_{2} x_{2} + u_{22}^{\mathsf{T}} R_{2} u_{22} \right] \mathrm{d}\tau \\ V_{\varepsilon}(x(t)) &= 2 \int_{t}^{\infty} \left[x_{1}^{\mathsf{T}} Q_{\varepsilon} x_{2} + u_{11}^{\mathsf{T}} R_{1} u_{12} + u_{22}^{\mathsf{T}} R_{2} u_{21} \right] \mathrm{d}\tau \\ &+ \varepsilon \int_{t}^{\infty} \left[u_{12}^{\mathsf{T}} R_{1} u_{12} + u_{21}^{\mathsf{T}} R_{2} u_{21} \right] \mathrm{d}\tau. \end{split}$$

According to the reduced-order scheme [9], setting $\varepsilon^2 = 0$, $\varepsilon V_{\varepsilon}(x(t))$ can be represented as

$$\varepsilon V_{\varepsilon}(x(t)) = 2\varepsilon \int_{t}^{\infty} \left[x_{1}^{\mathsf{T}} Q_{\varepsilon} x_{2} + u_{11}^{\mathsf{T}} R_{1} u_{12} + u_{22}^{\mathsf{T}} R_{2} u_{21} \right] d\tau.$$

We give the following definitions to denote the partial derivatives of the value functions $V_1(x_1(t))$, $V_2(x_2(t))$, and $V_{\varepsilon}(x(t))$ with respect to the states x_1 and x_2 , respectively:

$$V_{1x_1} = \frac{\partial V_1}{\partial x_1}, \quad V_{2x_2} = \frac{\partial V_2}{\partial x_2}$$

 $V_{\varepsilon x_1} = \frac{\partial V_{\varepsilon}}{\partial x_1}, \quad V_{\varepsilon x_2} = \frac{\partial V_{\varepsilon}}{\partial x_2}.$

Theorem 1: Partitioning the HJB equation (6), we get an $\mathcal{O}(\varepsilon^2)$ approximation in terms of three reduced-order decoupled HJB equations

$$\begin{split} 0 &= V_{1x_{1}}^{*\mathsf{T}} \big[f_{11}(x_{1}) + g_{11}(x_{1})u_{11}^{*}(x_{1}) \big] \\ &+ x_{1}^{\mathsf{T}} Q_{1}x_{1} + u_{11}^{*\mathsf{T}}(x_{1})R_{1}u_{11}^{*}(x_{1}) \\ 0 &= V_{2x_{2}}^{*\mathsf{T}} \big[f_{22}(x_{2}) + g_{22}(x_{2})u_{22}^{*}(x_{2}) \big] \\ &+ x_{2}^{\mathsf{T}} Q_{2}x_{2} + u_{22}^{*\mathsf{T}}(x_{2})R_{2}u_{22}^{*}(x_{2}) \\ 0 &= V_{1x_{1}}^{*\mathsf{T}} f_{12}(x) + V_{2x_{2}}^{*\mathsf{T}} f_{21}(x) + V_{\varepsilon x_{1}}^{*\mathsf{T}} f_{11}(x_{1}) \\ &+ V_{\varepsilon x_{2}}^{*\mathsf{T}} f_{22}(x_{2}) + 2x_{1}^{\mathsf{T}} Q_{\varepsilon}x_{2} \\ &- 2u_{11}^{*\mathsf{T}}(x_{1})R_{1}u_{12}^{*}(x) - 2u_{22}^{*\mathsf{T}}(x_{2})R_{2}u_{21}^{*}(x). \end{split}$$

The optimal control law (5) can be partitioned as

$$u_{11}^{*}(x_{1}) = -\frac{1}{2}R_{1}^{-1}g_{11}^{\mathsf{T}}(x_{1})V_{1x_{1}}^{*}$$

$$u_{12}^{*}(x) = -\frac{1}{2}R_{1}^{-1}\left[g_{11}^{\mathsf{T}}(x_{1})V_{\varepsilon x_{1}}^{*} + g_{21}^{\mathsf{T}}(x)V_{2x_{2}}^{*}\right]$$

$$u_{21}^{*}(x) = -\frac{1}{2}R_{2}^{-1}\left[g_{22}^{\mathsf{T}}(x_{2})V_{\varepsilon x_{2}}^{*} + g_{12}^{\mathsf{T}}(x)V_{1x_{1}}^{*}\right]$$

$$u_{22}^{*}(x_{2}) = -\frac{1}{2}R_{2}^{-1}g_{22}^{\mathsf{T}}(x_{2})V_{2x_{2}}^{*}.$$
(7)

Based on the optimal control theory, $u_{11}^*(x_1)$ can be seen as the optimal control law for the subsystem 1

$$\dot{x}_1(t) = f_{11}(x_1) + g_{11}(x_1)u_{11}(t)$$

with respect to the value function $V_1(x_1)$. $u_{22}^*(x_2)$ can be seen as the optimal control law for the subsystem 2

$$\dot{x}_2(t) = f_{22}(x_2) + g_{22}(x_2)u_{22}(t)$$

with respect to the value function $V_2(x_2)$. $u_{12}^*(x)$ and $u_{21}^*(x)$ can be solved from the optimal control problem of the virtual subsystem 3 with respect to the value function $V_2^*(x)$

$$V_3^*(x) = 2 \int_t^\infty \left[u_{11}^{*\mathsf{T}}(x_1) R_1 u_{12}^*(x) + u_{22}^{*\mathsf{T}}(x_2) R_2 u_{21}^*(x) - x_1^{\mathsf{T}} Q_{\varepsilon} x_2 \right] d\tau$$

where
$$V_3^*(x) = V_{\varepsilon}^*(x) - 4 \int_t^{\infty} x_1^{\mathsf{T}} Q_{\varepsilon} x_2 d\tau$$
 with $V_3^*(0) = 0$. *Proof:* Refer to the Appendix.

Remark 2: The original optimal control problem with the HJB equation (6) is transformed into three reduced-order HJB equations which should be solved without system dynamics. In the following section, we will derive the data-based online learning algorithm.

B. Model-Free Integral PI Algorithm

The optimal control formulation developed in (7) displays an array of closed-form expressions, which obviates the need to search for the optimal control law via optimization process. To obtain the optimal control law, the existence of $V^*(x)$ satisfying the HJB equation (6) is the necessary and sufficient condition. Instead of directly solving (6), we can successively solve the nonlinear Lyapunov equation (4) and update the control policy based on (7) to obtain the solution $V^*(x)$. This successive approximation is known as the model-based PI algorithm [42]–[45], and it is fundamental for the model-free integral PI algorithm and we describe it as follows.

1) Model-Based PI Algorithm: Step 1: Give a small positive real number ϵ . Let i=0 and start with an initial admissible control policy $u^0(x)$.

Step 2 (Policy Evaluation): Based on the control policy $u^i(x)$, solve $V^i(x)$ from the following nonlinear Lyapunov equation:

$$r(x, u^{i}(x)) + V_{x}^{i\mathsf{T}} [f(x) + g(x)u^{i}(x)] = 0.$$

Step 3 (Policy Improvement): Update the control policy by

$$u^{i+1}(x) = -\frac{1}{2}R^{-1}g^{\mathsf{T}}(x)V_x^i. \tag{8}$$

Step 4: If $||u^{i+1}(x) - u^i(x)|| \le \epsilon$, stop and obtain the approximate optimal control policy $u^{i+1}(x)$; else, set i = i+1 and go to step 2.

In [41], it was shown that on the domain Ω , the cost function $V^i(x)$ uniformly converges to $V^*(x)$ with monotonicity $V^{i+1}(x) \leq V^i(x)$, and the control policy $u^i(x)$ is admissible and converges to $u^*(x)$.

To deal with the optimal control problem without system dynamics, we develop a data-based online learning algorithm called model-free integral PI algorithm. Consider a nonlinear system which is explored by a known bounded piecewise continuous probing signal e(t)

$$\dot{x}(t) = f(x) + g(x)[u(t) + e(t)]$$

where

$$u(t) + e(t) = \begin{bmatrix} [u_{11}(t) + e_1(t)] + \varepsilon u_{12}(t) \\ \varepsilon u_{21}(t) + [u_{22}(t) + e_2(t)] \end{bmatrix}.$$

Now, we consider the subsystem 1 with exploration signal

$$\dot{x}_1(t) = f_{11}(x_1) + g_{11}(x_1)[u_{11}(t) + e_1(t)]. \tag{9}$$

The derivative of the value function $V_1(x_1(t))$ with respect to time along the trajectory of the explored system (9) can be calculated as

$$\dot{V}_1(x_1(t)) = V_{1x_1}^{\mathsf{T}} \left[f_{11}(x_1) + g_{11}(x_1) [u_{11}(t) + e_1(t)] \right]
= -r_1(x_1, u_{11}(x_1)) + V_{1x_1}^{\mathsf{T}} g_{11}(x_1) e_1(t)$$
(10)

where $r_1(x_1, u_{11}(x_1)) = x_1^\mathsf{T} Q_1 x_1 + u_{11}^\mathsf{T}(x_1) R_1 u_{11}(x_1)$ is the utility function for the subsystem 1 given in (9).

We present a lemma which is essential to prove the convergence of the model-free integral PI algorithm.

Lemma 1: Solving for $V_1(x_1)$ in the following equation:

$$V_{1}(x_{1}(t+T)) - V_{1}(x_{1}(t))$$

$$= \int_{t}^{t+T} V_{1x_{1}}^{\mathsf{T}} g_{11}(x_{1}) e_{1}(\tau) d\tau - \int_{t}^{t+T} r_{1}(x_{1}, u_{11}(x_{1})) d\tau$$
(11)

is equivalent to finding the solution of (10).

Proof: Since $u_{11}(x_1) \in \Psi_1(\Omega_1)$ [$\Psi_1(\Omega_1)$ is the set of all admissible control laws for the subsystem 1], the value function $V_1(x_1)$ is a Lyapunov function for the subsystem 1, and it satisfies (10) with $r_1(x_1, u_{11}(x_1)) > 0$, $x_1 \neq 0$. We integrate (10) over the interval [t, t+T] to obtain (11). This means that the unique solution of (10), $V_1(x_1)$, also satisfies (11). To complete the proof, we show that (11) has a unique solution by contradiction.

We assume that there exists another value function $\bar{V}_1(x_1)$ which satisfies (11) with bounding condition $\bar{V}_1(0)=0$. This value function also satisfies $\dot{\bar{V}}_1(x_1)=-r_1(x_1,u_{11}(x_1))+\bar{V}_{1x_1}^\mathsf{T}g_{11}(x_1)e_1(t)$. Subtracting this from (10), we obtain

$$0 = \left(\frac{d[\bar{V}_1(x_1) - V_1(x_1)]^T}{dx_1}\right) \times [\dot{x}_1(t) - g_{11}(x_1)e_1(t)]$$
$$= \left(\frac{d[\bar{V}_1(x_1) - V_1(x_1)]^T}{dx_1}\right) \times [f_{11}(x_1) + g_{11}(x_1)u_{11}(x_1)]$$

Algorithm 1 Model-Free Integral PI Algorithm

- 1: Give a small positive real number ϵ . Let i = 0 and start with an initial admissible control policy $u_{11}^0(x_1)$.
- 2: **Policy Evaluation and Improvement**: Based on the control policy $u_{11}^{i}(x_1)$, solve $V_1^{i}(x_1)$ and $u_{11}^{i+1}(x_1)$ from the integral equation (13).
- integral equation (13).
 3: If ||u₁₁ⁱ⁺¹(x₁) u₁₁ⁱ(x₁)|| ≤ ε, stop and obtain the approximate optimal control policy u₁₁ⁱ⁺¹ for the subsystem 1; else, set i = i + 1 and go to Step 2.

which must hold for any x_1 on the system trajectories generated by the stabilizing policy $u_{11}(x_1)$. According to the above equation, we have $\bar{V}_1(x_1) = V_1(x_1) + c$. As this relation must hold for $x_1(t) = 0$, we know $\bar{V}_1(0) = V_1(0) + c$, c = 0. Thus, $\bar{V}_1(x_1) = V_1(x_1)$, i.e., (11) has a unique solution which is equal to the solution of (10). The proof is complete.

Based on the model-based PI algorithm and using the representations $V_1^i(x_1(t))$ and $u_{11}^i(x_1)$, the policy improvement (8) for the subsystem 1 can be written as

$$u_{11}^{i+1}(x_1) = -\frac{1}{2}R_1^{-1}g_{11}^{\mathsf{T}}(x_1)V_{1x_1}^{i}$$
 (12)

where i is the iteration index. Integrating (10) from t to t+T with a time period T>0, and using the policy improvement (12), we have

$$V_1^i(x_1(t)) - V_1^i(x_1(t+T)) = \int_t^{t+T} r_1(x_1, u_{11}^i(x_1)) d\tau + 2 \int_t^{t+T} \left(u_{11}^{i+1}(x_1)\right)^\mathsf{T} R_1 e_1(\tau) d\tau.$$
(13)

Since the dynamics $f_{11}(x_1)$ and $g_{11}(x_1)$ are not in the integral equation (13), the integral PI algorithm can be implemented using the data generated from the system instead of the system dynamics. Thus, we obtain the model-free integral PI algorithm (Algorithm 1).

Theorem 2: Give an initial admissible control policy $u_{11}^0(x_1)$ for the subsystem 1. Using the model-free integral PI algorithm established in Algorithm 1, the value function and the control law converge to the optimal value function and the optimal control law as $i \to \infty$, that is

$$V_1^i(x_1) \to V_1^*(x_1), \quad u_{11}^i(x_1) \to u_{11}^*(x_1).$$

Proof: Based on the results in [35], we known that all the subsequent control policies will be admissible during the algorithm implementation if $u_{11}^0(x_1)$ is admissible. Considering the model-based PI algorithm and the formation process of (13), the value function sequence generated in Algorithm 1 will converge to the solution of the HJB equation. So we can conclude that the value function $V_1^i(x_1)$ and the control policy $u_{11}^i(x_1)$ obtained from the proposed model-free integral PI algorithm will converge to the solution of the optimal control problem for the subsystem 1. The proof is complete.

To solve the optimal control policy $u_{22}^*(x_2)$ for the subsystem 2, we can apply Algorithm 1 with some simply replacements. Using the expressions of the optimal control

laws $u_{11}^*(x_1)$ and $u_{22}^*(x_2)$, we derive the following equation which will be used to solve $u_{12}^*(x)$ and $u_{21}^*(x)$:

$$\begin{split} V_3^i(x(t+T)) - V_3^i(x(t)) &= -2 \int_t^{t+T} x_1^\mathsf{T} Q_\varepsilon x_2 \mathrm{d}\tau \\ &+ 2 \int_t^{t+T} \left[u_{11}^{*\mathsf{T}}(x_1) R_1 u_{12}^i(x) + u_{22}^{*\mathsf{T}}(x_2) R_2 u_{21}^i(x) \right] \mathrm{d}\tau \\ &+ 2 \int_t^{t+T} \left[\left(u_{12}^{i+1}(x) \right)^\mathsf{T} R_1 e_1(\tau) + \left(u_{21}^{i+1}(x) \right)^\mathsf{T} R_2 e_2(\tau) \right] \mathrm{d}\tau. \end{split}$$

Using this formulation to replace the integral equation (13) in Algorithm 1, we can calculate $u_{12}^*(x)$ and $u_{21}^*(x)$ iteratively.

C. Algorithm Implementation

For the subsystem 1, we represent $V_1^i(x_1)$ and $u_{11}^{i+1}(x_1)$ by single-layer NNs on a compact set Ω_1 as

$$V_1^i(x_1) = \sum_{j=1}^{N_{1c}} \omega_{1j}^i \phi_{1j}(x_1) + \delta_{1c}^i(x_1)$$

$$u_{11,p}^{i+1}(x_1) = \sum_{j=1}^{N_{1a}} v_{1j,p}^i \psi_{1j}(x_1) + \delta_{1a,p}^i(x_1)$$

where $p=1,2,\ldots,m_1,\,\omega_{1j}^i\in\mathbb{R}$, and $\nu_{1j,p}^i\in\mathbb{R}$ are bounded ideal weight parameters which will be determined by the developed data-based integral PI algorithm, $\phi_{1j}(x_1)\in\mathbb{R}$ and $\psi_{1j}(x_1)\in\mathbb{R}$, $\{\phi_{1j}\}_{j=1}^{N_{1c}}$ and $\{\psi_{1j}\}_{j=1}^{N_{1a}}$ are the sequences of real-valued activation functions which are linearly complete and independent, and $\delta_{1c}^i(x_1)\in\mathbb{R}$ and $\delta_{1a,p}^i(x_1)\in\mathbb{R}$ are the bounded NN approximation errors. Since the ideal weights are unknown, the outputs of the critic network and the action network are

$$\hat{V}_{1}^{i}(x_{1}) = \sum_{j=1}^{N_{1c}} \hat{\omega}_{1j}^{i} \phi_{1j}(x_{1}) = \hat{\omega}_{1}^{i\mathsf{T}} \phi_{1}(x_{1})$$
 (14)

$$\hat{u}_{11,p}^{i+1}(x_1) = \sum_{i=1}^{N_{1a}} \hat{v}_{1j,p}^i \psi_{1j}(x_1) = \hat{v}_{1,p}^{i\mathsf{T}} \psi_1(x_1)$$
 (15)

where $\hat{\omega}_1^i$ and $\hat{v}_{1,p}^i$ are the current estimated weights, and

$$\begin{aligned} \phi_{1}(x_{1}) &= \left[\phi_{11}(x_{1}), \phi_{12}(x_{1}), \dots, \phi_{1N_{1c}}(x_{1})\right]^{\mathsf{T}} \in \mathbb{R}^{N_{1c}} \\ \psi_{1}(x_{1}) &= \left[\psi_{11}(x_{1}), \psi_{12}(x_{1}), \dots, \psi_{1N_{1a}}(x_{1})\right]^{\mathsf{T}} \in \mathbb{R}^{N_{1a}} \\ \hat{\omega}_{1}^{i} &= \left[\hat{\omega}_{11}^{i}, \hat{\omega}_{12}^{i}, \dots, \hat{\omega}_{1N_{1c}}^{i}\right]^{\mathsf{T}} \in \mathbb{R}^{N_{1c}} \\ \hat{v}_{1,p}^{i} &= \left[\hat{v}_{11,p}^{i}, \hat{v}_{12,p}^{i}, \dots, \hat{v}_{1N_{1a},p}^{i}\right]^{\mathsf{T}} \in \mathbb{R}^{N_{1a}} \\ \hat{v}_{1}^{i\mathsf{T}} &= \left[\hat{v}_{1,1}^{i}, \hat{v}_{1,2}^{i}, \dots, \hat{v}_{1,m_{1}}^{i}\right]^{\mathsf{T}} \in \mathbb{R}^{m_{1} \times N_{1a}}. \end{aligned}$$

Define $\text{col}\{\hat{v}_1^{i\mathsf{T}}\} = [\hat{v}_{1.1}^{i\mathsf{T}}, \hat{v}_{1.2}^{i\mathsf{T}}, \dots, \hat{v}_{1.m_1}^{i\mathsf{T}}]^\mathsf{T} \in \mathbb{R}^{m_1N_{1a}}$. Then

$$\begin{aligned} \left(\hat{u}_{11}^{i+1}(x_1)\right)^{\mathsf{T}} R_1 e_1(t) &= \left(\hat{v}_1^{i\mathsf{T}} \psi_1(x_1)\right)^{\mathsf{T}} R_1 e_1(t) \\ &= \left[\psi_1(x_1) \otimes (R_1 e_1(t))\right]^{\mathsf{T}} \mathrm{col} \left\{\hat{v}_1^{i\mathsf{T}}\right\} \end{aligned}$$

where \otimes represents the Kronecker product. Using the real outputs of the networks (14) and (15), the integral equation (13)

has the following general form:

$$\lambda_{1k}^{\mathsf{T}} \begin{bmatrix} \hat{\omega}_1^i \\ \operatorname{col} \left\{ \hat{\nu}_1^{i\mathsf{T}} \right\} \end{bmatrix} = \theta_{1k} \tag{16}$$

with

$$\theta_{1k} = \int_{t+(k-1)T}^{t+kT} \left[x_1^\mathsf{T} Q_1 x_1 + \hat{u}_{11}^{i\mathsf{T}}(x_1) R_1 \hat{u}_{11}^i(x_1) \right] d\tau$$

$$\lambda_{1k} = \left[(\phi_1(x_1(t+(k-1)T)) - \phi_1(x_1(t+kT)))^\mathsf{T} - 2 \int_{t+(k-1)T}^{t+kT} (\psi_1(x_1) \otimes (R_1 e_1(\tau)))^\mathsf{T} d\tau \right]^\mathsf{T}$$

where T is the period of time to measure the data. Since the general form (16) is a 1-D equation, we cannot find the unique weight vector. The least squares method [39] can be used to guarantee the uniqueness of the weights over the compact set Ω_1 . For any positive integer K_1 , we denote $\Lambda_1 = [\lambda_{11}, \lambda_{12}, \dots, \lambda_{1K_1}]$ and $\Theta_1 = [\theta_{11}, \theta_{12}, \dots, \theta_{1K_1}]^T$. Then, we have the following K_1 -dimensional equation

$$\Lambda_1^{\mathsf{T}} \begin{bmatrix} \hat{\omega}_1^i \\ \operatorname{col} \left\{ \hat{v}_1^{i\mathsf{T}} \right\} \end{bmatrix} = \Theta_1.$$

The weight vector can be solved by the following equation when Λ_1^T has full column rank:

$$\begin{bmatrix} \hat{\omega}_1^i \\ \text{col} \left\{ \hat{v}_1^{i\mathsf{T}} \right\} \end{bmatrix} = \left(\Lambda_1 \Lambda_1^{\mathsf{T}} \right)^{-1} \Lambda_1 \Theta_1. \tag{17}$$

Therefore, we need to make sure $(\Lambda_1 \Lambda_1^\mathsf{T})^{-1}$ exists; that is to say, the number of collected points K_1 should satisfy $K_1 \ge \mathrm{rank}(\Lambda_1) = N_{1c} + m_1 N_{1a}$. By collecting enough data points of the explored system (9), the weight parameters in (17) can be obtained in real time. Using the same implementation procedures for the subsystem 1, we can solve the optimal control problems of the subsystems 2 and 3.

IV. NUMERICAL SIMULATION

We provide two simulation examples in this section to demonstrate the applicability of the established data-based integral PI algorithm for weakly coupled nonlinear systems.

Example 1: In this example, we consider the system (1) with the following parameters:

$$f_{11}(x_1) = \begin{bmatrix} -1.93x_{11}^2 \\ -1.394x_{11}x_{12} \end{bmatrix}$$

$$f_{12}(x) = \begin{bmatrix} 0 \\ -4.26x_{21}x_{22} \end{bmatrix}$$

$$f_{21}(x) = \begin{bmatrix} -1.3x_{12}^2 \\ 0.95x_{11}x_{21} - 1.03x_{12}x_{22} \end{bmatrix}$$

$$f_{22}(x_2) = \begin{bmatrix} -0.63x_{21}^2 \\ 0.413x_{21} - 0.426x_{22} \end{bmatrix}$$

$$g_{11}(x_1) = \begin{bmatrix} -1.274x_{11}^2 \\ 0 \end{bmatrix}, \qquad g_{12}(x) = \begin{bmatrix} 0 \\ -6.5x_{22} \end{bmatrix}$$

$$g_{21}(x) = \begin{bmatrix} 0.75x_{11} \\ 0 \end{bmatrix}, \qquad g_{22}(x_2) = \begin{bmatrix} -0.718x_{21} \\ 0 \end{bmatrix}$$

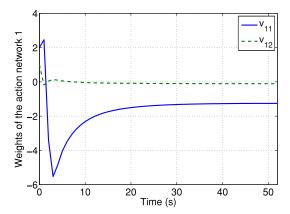


Fig. 1. Evolution of the action network 1's weights.

In the above, $x_1 = [x_{11}, x_{12}]^T \in \mathbb{R}^2$ and $u_{11}(x_1) \in \mathbb{R}$ are the state and control vectors of the subsystem 1, and $x_2 = [x_{21}, x_{22}]^T \in \mathbb{R}^2$ and $u_{22}(x_2) \in \mathbb{R}$ are the state and control vectors of the subsystem 2. The initial system state is $x(0) = [3.4, 2.7, 4.3, 1.2]^T$. The weak coupling parameter is equal to $\varepsilon = 0.05$. The matrices Q and R are chosen as

$$Q_1 = Q_2 = R = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, \qquad Q_{\varepsilon} = \begin{bmatrix} 1 & 0 \\ 0 & 0.05 \end{bmatrix}.$$

We assume that the exact knowledge of the system dynamics is completely unknown during the simulation. We adopt the integral PI algorithm to derive the optimal control law.

For the subsystem 1

$$\dot{x}_1 = \begin{bmatrix} -1.93x_{11}^2 \\ -1.394x_{11}x_{12} \end{bmatrix} + \begin{bmatrix} -1.274x_{11}^2 \\ 0 \end{bmatrix} u_{11}(x_1)$$

the weight parameters of the critic network and the action network are

$$\hat{\omega}_{1} = [\hat{\omega}_{11}, \hat{\omega}_{12}, \hat{\omega}_{13}]^{\mathsf{T}}$$
$$\hat{\nu}_{1} = [\hat{\nu}_{11}, \hat{\nu}_{12}]^{\mathsf{T}}.$$

The activation functions are chosen as

$$\phi_1(x_1) = \begin{bmatrix} x_{11}^2, x_{11}x_{12}, x_{12}^2 \end{bmatrix}^\mathsf{T}$$
$$\psi_1(x_1) = \begin{bmatrix} x_{11}x_{12}, x_{12}^2 \end{bmatrix}^\mathsf{T}.$$

From the activation functions, we have $N_{1c} = 3$ and $N_{1a} = 2$ and we select $K_1 = 10$ to conduct the simulation. The initial weights are chosen as $\hat{\omega}_1 = [0, 0, 0]^{\mathsf{T}}$ and $\hat{v}_1 = [2, 1]^{\mathsf{T}}$. During the online learning process, the time period T = 0.1[s] and the exploration signal $e_1(t) = 3\sin(2\pi t) + 3\cos(2\pi t)$ are used. The least squares problem is solved after K_1 samples are acquired, thus the weights of the NNs are updated every 1[s]. The evolution of the action network 1's weights is illustrated in Fig. 1. After 52 iterations, the precision $\epsilon = 10^{-4}$ is achieved. At time t = 52[s], $\hat{v}_1^* = [-1.2557, -0.1067]^{\mathsf{T}}$.

For the subsystem 2, the activation functions are chosen as

$$\phi_2(x_2) = \begin{bmatrix} x_{21}^2, x_{21}x_{22}, x_{22}^2 \end{bmatrix}^\mathsf{T}$$
$$\psi_2(x_2) = \begin{bmatrix} x_{21}x_{22}, x_{22}^2 \end{bmatrix}^\mathsf{T}.$$

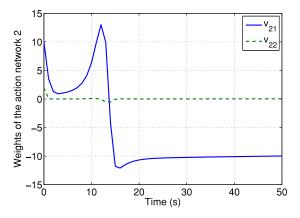


Fig. 2. Evolution of the action network 2's weights.

As $N_{2c} = 3$ and $N_{2a} = 2$, we conduct the simulation with $K_2 = 10$. The initial weights are chosen as $\hat{\omega}_2 = [0, 0, 0]^{\mathsf{T}}$ and $\hat{\nu}_2 = [10, 2]^{\mathsf{T}}$. During the online learning process, the time period T = 0.1[s] and the exploration signal $e_2(t) = 5\sin(2\pi t) + 5\cos(2\pi t)$ are used. The evolution of the action network 2's weights is illustrated in Fig. 2. After 50 iterations, the precision ϵ is achieved. At time t = 50[s], $\hat{\nu}_2^* = [-9.9814, 0.0367]^{\mathsf{T}}$.

For the virtual subsystem 3, the weight parameters of the critic network and the action network are

$$\hat{\omega}_3 = \left[\hat{\omega}_{31}, \hat{\omega}_{32}, \hat{\omega}_{33}, \hat{\omega}_{34}, \hat{\omega}_{35}, \hat{\omega}_{36}\right]^\mathsf{T}$$
$$\hat{\nu}_3 = \left[\hat{\nu}_{31}, \hat{\nu}_{32}, \hat{\nu}_{33}, \hat{\nu}_{34}\right]^\mathsf{T}.$$

The activation functions are chosen as

$$\phi_3(x) = \begin{bmatrix} x_{11}^2, x_{11}x_{12}, x_{12}^2, x_{21}^2, x_{21}x_{22}, x_{22}^2 \end{bmatrix}^\mathsf{T}$$

$$\psi_3(x) = \begin{bmatrix} x_{11}x_{12}, x_{12}^2, x_{21}x_{22}, x_{22}^2 \end{bmatrix}^\mathsf{T}.$$

From the activation functions, we have $N_{3c} = 6$ and $N_{3a} = 4$ and we select $K_3 = 10$ to conduct the simulation. The initial weights are chosen as $\hat{\omega}_3 = [0, 0, 0, 0, 0, 0]^T$ and $\hat{\nu}_3 = [0, -2, 2, 3]^T$. During the online learning process, the time period T = 0.1[s] and the exploration signals $e_1(t)$ and $e_2(t)$ are used. The least squares problem is solved after K_3 samples are acquired, and the weights are updated every 1[s]. After 20 iterations, the precision $\epsilon = 10^{-4}$ is achieved. At time t = 20[s], $\hat{\nu}_3^* = [0.3830, 0.0533, -0.0899, -0.9548]^T$.

According to the results in Section III, the optimal control law of the weakly coupled system can be derived as

$$u^*(x) = \begin{bmatrix} u_{11}^*(x_1) + \varepsilon u_{12}^*(x) \\ \varepsilon u_{21}^*(x) + u_{22}^*(x_2) \end{bmatrix}.$$

Using the optimal control $u^*(x)$ to control the weakly coupled system for 20[s], we obtain the evolution process of the state trajectory and control trajectory shown in Figs. 3 and 4. Obviously, these simulation results have verified the effectiveness of the developed model-free integral PI algorithm.

Example 2: In this example, we use the established modelfree integral PI algorithm to balance a bicycle riding at a constant speed on a horizontal surface. The steering column of the bicycle is vertical, which means that the bicycle is not

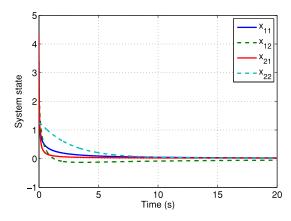


Fig. 3. State trajectory of the weakly coupled system under the derived optimal control.

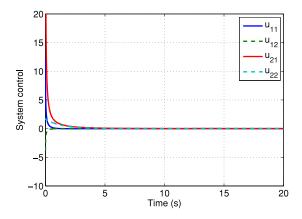


Fig. 4. Control trajectory of the weakly coupled system under the derived optimal control.

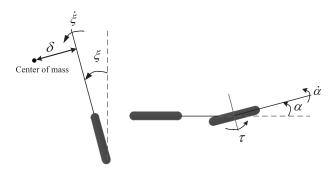


Fig. 5. Schematic representation of the bicycle, as seen from behind (left) and from the top (right).

self-stabilizing, but must be actively stabilized to prevent it from falling [47]. This is a variant of a bicycle balancing and riding problem which is widely used as a benchmark for RL algorithms [46].

The schematic representation of the bicycle is provided in Fig. 5, which includes the system state and control variables. The system state variables are the roll angle ξ [rad] of the bicycle measured from the vertical axis, the angle α [rad] of the handlebar, and the respective angular velocities $\dot{\xi}$, $\dot{\alpha}$ [rad/s]. The control variables are the displacement δ [m] of the bicyclerider common center of mass perpendicular to the plane of the bicycle, and the torque τ [Nm] applied to the handlebar.

TABLE I PARAMETERS OF THE BICYCLE

Symbol	Value	Units	Symbol	Value	Units
M_c M_d M_r g v	15 1.7 60 9.81 10/3.6	kg kg kg m/s ² m/s	$\left egin{array}{c} h \\ l \\ r \\ d_{CM} \\ c \end{array} \right $	0.94 1.11 0.34 0.3 0.66	m m m m

Therefore, the state vector is $x = [\xi, \dot{\xi}, \alpha, \dot{\alpha}]^T$, and the control vector is $u = [\delta, \tau]^T$.

The dynamics of the bicycle can be represented as [46]

$$\ddot{\xi} = \frac{1}{J_{bc}} \left[\sin \beta (M_c + M_r) g h - \cos \beta \left(\frac{J_{dc} v}{r} \dot{\alpha} \right) + \operatorname{sign}(\alpha) \frac{M_d r v^2}{l} (|\sin \alpha| + |\tan \alpha|) \right]$$

$$\ddot{\alpha} = \frac{1}{J_{dl}} \left(\tau - \frac{J_{dv} v}{r} \dot{\xi} \right)$$

where

$$J_{bc} = \frac{13}{3} M_c h^2 + M_r (h + d_{CM})^2, \quad J_{dc} = M_d r^2$$

$$J_{dv} = \frac{3}{2} M_d r^2, \quad J_{dl} = \frac{1}{2} M_d r^2, \quad \beta = \xi + \arctan \frac{\delta}{h}.$$

Table I shows the values of the parameters in the bicycle model. The meanings of these parameters are the same as those in [46]. Using the notations $x_1 = [\xi, \dot{\xi}]^T$, $x_2 = [\alpha, \dot{\alpha}]^T$, $u_1 = \delta$, and $u_2 = \tau$, we rewrite the bicycle dynamics as

$$\begin{bmatrix} \dot{x}_{11} \\ \dot{x}_{12} \\ \dot{x}_{21} \\ \dot{x}_{22} \end{bmatrix} = \begin{bmatrix} x_{12} \\ 4.62x_{11} + 0.054x_{21} + 0.011x_{22} + 4.62u_{1} \\ x_{22} \\ 24.51x_{12} + 10.18u_{2} \end{bmatrix}.$$

Compared with the system (1) with weakly coupled structure, we have the following system dynamics:

$$f_{11}(x_1) = \begin{bmatrix} x_{12} \\ 4.62x_{11} \end{bmatrix}, \quad f_{12}(x) = \begin{bmatrix} 0 \\ 0.54x_{21} + 0.11x_{22} \end{bmatrix}$$

$$f_{21}(x) = \begin{bmatrix} 0 \\ 245.1x_{12} \end{bmatrix}, \quad f_{22}(x_2) = \begin{bmatrix} x_{22} \\ 0 \end{bmatrix}$$

$$g_{11}(x_1) = \begin{bmatrix} 0 \\ 4.62 \end{bmatrix}, \quad g_{12}(x) = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$$

$$g_{21}(x) = \begin{bmatrix} 0 \\ 0 \end{bmatrix}, \quad g_{22}(x_2) = \begin{bmatrix} 0 \\ 10.18 \end{bmatrix}.$$

The initial system state is $x(0) = [0.1, -0.1, 0.1, -0.1]^T$. The weak coupling parameter is $\varepsilon = 0.1$. The matrices Q and R are chosen as

$$Q_1 = Q_2 = R = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, \qquad Q_{\varepsilon} = \begin{bmatrix} 1 & 0 \\ 0 & 0.1 \end{bmatrix}.$$

Assume that the exact knowledge of the bicycle is completely unknown during the simulation. We adopt the model-free integral PI algorithm to solve the bicycle balancing and riding problem.

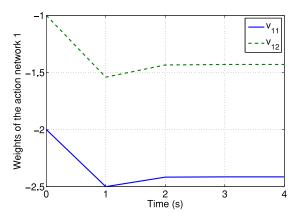


Fig. 6. Evolution of the action network 1's weights.

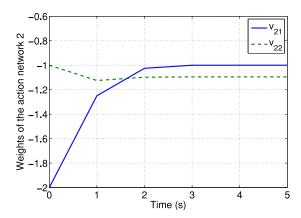


Fig. 7. Evolution of the action network 2's weights.

As in Example 1, for the subsystem 1 the weight parameters of the critic network and the action network are

$$\hat{\omega}_{1} = [\hat{\omega}_{11}, \hat{\omega}_{12}, \hat{\omega}_{13}]^{\mathsf{T}}$$
$$\hat{\nu}_{1} = [\hat{\nu}_{11}, \hat{\nu}_{12}]^{\mathsf{T}}.$$

The activation functions are chosen as

$$\phi_1(x_1) = \left[x_{11}^2, x_{11}x_{12}, x_{12}^2 \right]^\mathsf{T}$$

$$\psi_1(x_1) = \left[x_{11}, x_{12} \right]^\mathsf{T}.$$

From the activation functions, we have $N_{1c} = 3$ and $N_{1a} = 2$ and we select $K_1 = 10$ to conduct the simulation. We set the initial weights as $\hat{\omega}_1 = [0,0,0]^\mathsf{T}$ and $\hat{v}_1 = [-2,-1]^\mathsf{T}$. During the online learning process, the time period T = 0.1[s] and the exploration signal $e_1(t) = 0.05 \sin(2\pi t) + 0.05 \cos(2\pi t)$ are used. The least squares problem is solved after K_1 samples are acquired, and thus the weights of the NNs are updated every 1[s]. The evolution of the action network 1's weights is illustrated in Fig. 6. After 4 iterations, the precision $\epsilon = 10^{-4}$ is achieved. At time t = 4[s], $\hat{v}_1^* = [-2.4142, -1.4301]^\mathsf{T}$.

For the subsystem 2, the activation functions are chosen as

$$\phi_2(x_2) = \begin{bmatrix} x_{21}^2, x_{21}x_{22}, x_{22}^2 \end{bmatrix}^\mathsf{T}$$
$$\psi_2(x_2) = \begin{bmatrix} x_{21}, x_{22} \end{bmatrix}^\mathsf{T}.$$

As $N_{2c} = 3$ and $N_{2a} = 2$, we conduct the simulation with $K_2 = 10$. The initial weights are chosen as $\hat{\omega}_2 = [0, 0, 0]^T$

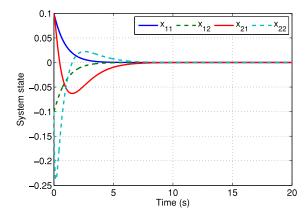


Fig. 8. State trajectory of the bicycle under the derived optimal control.

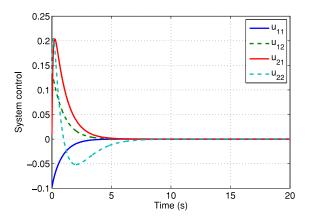


Fig. 9. Control trajectory of the bicycle under the derived optimal control.

and $\hat{v}_2 = [-2, -1]^T$. During the online learning process, the time period T = 0.1[s] and the exploration signal $e_2(t) = 0.05 \sin(2\pi t) + 0.05 \cos(2\pi t)$ are used. The evolution of the action network 2's weights is illustrated in Fig. 7. After 5 iterations, the precision $\epsilon = 10^{-4}$ is achieved. At time t = 5[s], $\hat{v}_2^* = [-1.0000, -1.0955]^T$.

For the virtual subsystem 3, the weight parameters of the critic network and the action network are

$$\hat{\omega}_3 = \begin{bmatrix} \hat{\omega}_{31}, \hat{\omega}_{32}, \hat{\omega}_{33}, \hat{\omega}_{34}, \hat{\omega}_{35}, \hat{\omega}_{36} \end{bmatrix}^\mathsf{T} \\ \hat{\nu}_3 = \begin{bmatrix} \hat{\nu}_{31}, \hat{\nu}_{32}, \hat{\nu}_{33}, \hat{\nu}_{34} \end{bmatrix}^\mathsf{T}.$$

The activation functions are chosen as

$$\phi_3(x) = \begin{bmatrix} x_{11}^2, x_{11}x_{12}, x_{12}^2, x_{21}^2, x_{21}x_{22}, x_{22}^2 \end{bmatrix}^\mathsf{T}$$

$$\psi_3(x) = [x_{11}, x_{12}, x_{21}, x_{22}]^\mathsf{T}.$$

From the activation functions, we have $N_{3c} = 6$ and $N_{3a} = 4$ and we select $K_3 = 10$ to conduct the simulation. The initial weights are chosen as $\hat{\omega}_3 = [0, 0, 0, 0, 0, 0]^T$ and $\hat{\nu}_3 = [0, -2, 2, 3]^T$. During the online learning process, the time period T = 0.1[s] and the exploration signals $e_1(t)$ and $e_2(t)$ are used. The least squares problem is solved after K_3 samples are acquired, and the weights of the NNs are updated every 1[s]. After 23 iterations, the precision $\epsilon = 10^{-4}$ is achieved. At time t = 23[s], $\hat{\nu}_3^* = [0.5420, -0.8267, 0.6952, -0.6516]^T$.

$$0 = \begin{bmatrix} V_{1x_{1}}^{*} + \varepsilon V_{\varepsilon x_{1}}^{*} \\ \varepsilon V_{\varepsilon x_{2}}^{*} + V_{2x_{2}}^{*} \end{bmatrix}^{\mathsf{T}} \left(\begin{bmatrix} f_{11}(x_{1}) + \varepsilon f_{12}(x) \\ \varepsilon f_{21}(x) + f_{22}(x) \end{bmatrix} \right) + \begin{bmatrix} g_{11}(x_{1}) & \varepsilon g_{12}(x) \\ \varepsilon g_{21}(x) & g_{22}(x) \end{bmatrix} \begin{bmatrix} u_{11}^{*}(x_{1}) + \varepsilon u_{12}^{*}(x) \\ \varepsilon u_{21}^{*}(x) + u_{22}^{*}(x) \end{bmatrix} \right)$$

$$+ \begin{bmatrix} x_{1} \\ x_{2} \end{bmatrix}^{\mathsf{T}} \begin{bmatrix} Q_{1} & \varepsilon Q_{\varepsilon} \\ Q_{\varepsilon} & Q_{2} \end{bmatrix} \begin{bmatrix} x_{1} \\ x_{2} \end{bmatrix} + \begin{bmatrix} u_{11}^{*}(x_{1}) + \varepsilon u_{12}^{*}(x) \\ \varepsilon u_{21}^{*}(x) + u_{22}^{*}(x) \end{bmatrix}^{\mathsf{T}} \begin{bmatrix} R_{1} & 0 \\ 0 & R_{2} \end{bmatrix} \begin{bmatrix} u_{11}^{*}(x_{1}) + \varepsilon u_{12}^{*}(x) \\ \varepsilon u_{21}^{*}(x) + u_{22}^{*}(x_{2}) \end{bmatrix}$$

$$= \begin{bmatrix} V_{1x_{1}}^{*} + \varepsilon V_{\varepsilon x_{1}}^{*} \\ V_{1x_{1}}^{*} + \varepsilon V_{\varepsilon x_{1}}^{*} \end{bmatrix}^{\mathsf{T}} \begin{bmatrix} f_{11}(x_{1}) + g_{11}(x_{1}) u_{11}^{*}(x_{1}) + \varepsilon [f_{12}(x) + g_{11}(x_{1}) u_{12}^{*}(x) + g_{12}(x) u_{22}^{*}(x_{2}) \end{bmatrix}$$

$$+ x_{1}^{\mathsf{T}} Q_{1}x_{1} + x_{2}^{\mathsf{T}} Q_{2}x_{2} + 2\varepsilon x_{1}^{\mathsf{T}} Q_{\varepsilon}x_{2} + u_{11}^{*\mathsf{T}}(x_{1}) R_{1}u_{11}^{*}(x_{1}) + u_{22}^{\mathsf{T}}(x_{2}) R_{2}u_{22}^{*}(x_{2})$$

$$+ 2\varepsilon u_{11}^{*\mathsf{T}}(x_{1}) R_{1}u_{12}^{*}(x) + 2\varepsilon u_{22}^{*\mathsf{T}}(x_{2}) R_{2}u_{21}^{*}(x)$$

$$+ 2V_{1x_{1}}^{*\mathsf{T}} [f_{11}(x_{1}) + g_{11}(x_{1}) u_{11}^{*}(x_{1})] + x_{1}^{\mathsf{T}} Q_{1}x_{1} + u_{11}^{*\mathsf{T}}(x_{1}) R_{1}u_{11}^{*}(x_{1})$$

$$+ W_{1x_{1}}^{\mathsf{T}} [f_{11}(x_{1}) + g_{11}(x_{1}) u_{11}^{*}(x_{1})] + x_{1}^{\mathsf{T}} Q_{1}x_{1} + u_{11}^{*\mathsf{T}}(x_{1}) R_{1}u_{11}^{*}(x_{1})$$

$$+ W_{1x_{1}}^{\mathsf{T}} [f_{11}(x_{1}) + g_{11}(x_{1}) u_{11}^{*}(x_{1})] + x_{1}^{\mathsf{T}} Q_{2}x_{2} + u_{22}^{*\mathsf{T}}(x_{2}) R_{2}u_{22}^{*}(x_{2})$$

$$+ 2\varepsilon u_{11}^{*\mathsf{T}} [f_{11}(x_{1}) + g_{11}(x_{1}) u_{11}^{*}(x_{1})] + \varepsilon V_{1x_{1}}^{\mathsf{T}} [f_{12}(x) + g_{11}(x_{1}) u_{12}^{*}(x) + g_{12}(x) u_{22}^{*}(x_{2})]$$

$$+ W_{2x_{1}}^{\mathsf{T}} [f_{22}(x_{2}) + g_{22}(x_{2}) u_{22}^{*}(x_{2})] + \varepsilon V_{1x_{1}}^{\mathsf{T}} [f_{12}(x) + g_{11}(x_{1}) u_{12}^{*}(x) + g_{12}(x) u_{22}^{*}(x_{2})]$$

$$+ \varepsilon V_{\varepsilon x_{1}}^{\mathsf{T}} [f_{21}(x) + g_{21}(x) u_{22}^{*}(x_{2})] + \varepsilon V_{2x_{1}}^{\mathsf{T}} [f_{21}(x) + g_{22}(x_{2}) u_{21}^{*}(x) + g_{21}(x) u_{11}^{*}(x_{1})]$$

$$+ \varepsilon V_{\varepsilon x_{1}}^{\mathsf{T}} [f_{22}(x_{2}) + g_{22}(x_{2}) u_{22$$

$$u^{*}(x) = -\frac{1}{2}R^{-1}g^{\mathsf{T}}(x)V_{x}^{*} = -\frac{1}{2}\begin{bmatrix} R_{1} & 0 \\ 0 & R_{2} \end{bmatrix}^{-1}\begin{bmatrix} g_{11}(x_{1}) & \varepsilon g_{12}(x) \\ \varepsilon g_{21}(x) & g_{22}(x_{2}) \end{bmatrix}^{\mathsf{T}}\begin{bmatrix} V_{1x_{1}}^{*} + \varepsilon V_{\varepsilon x_{1}}^{*} \\ \varepsilon V_{\varepsilon x_{2}}^{*} + V_{2x_{2}}^{*} \end{bmatrix}$$

$$= \begin{bmatrix} u_{11}^{*}(x_{1}) + \varepsilon u_{12}^{*}(x) \\ \varepsilon u_{21}^{*}(x) + u_{22}^{*}(x_{2}) \end{bmatrix} = -\frac{1}{2}\begin{bmatrix} R_{1}^{-1}g_{11}^{\mathsf{T}}(x_{1})V_{1x_{1}}^{*} + \varepsilon R_{1}^{-1}\left[g_{11}^{\mathsf{T}}(x_{1})V_{\varepsilon x_{1}}^{*} + g_{21}^{\mathsf{T}}(x)V_{2x_{2}}^{*} \right] \\ \varepsilon R_{2}^{-1}\left[g_{22}^{\mathsf{T}}(x_{2})V_{\varepsilon x_{2}}^{*} + g_{12}^{\mathsf{T}}(x)V_{1x_{1}}^{*} \right] + R_{2}^{-1}g_{22}^{\mathsf{T}}(x_{2})V_{2x_{2}}^{*} \end{bmatrix}$$
(A3)

According to the results in Section III, the optimal control law of the weakly coupled system can be derived as

$$u^*(x) = \begin{bmatrix} u_{11}^*(x_1) + \varepsilon u_{12}^*(x) \\ \varepsilon u_{21}^*(x) + u_{22}^*(x_2) \end{bmatrix}.$$

Using the optimal control $u^*(x)$ to control the weakly coupled system for 20[s], we obtain the evolution process of the state trajectory and control trajectory shown in Figs. 8 and 9. Obviously, these simulation results have verified the effectiveness of the developed model-free integral PI algorithm.

Remark 3: In Figs. 1 and 2, one weight parameter is largely dominated by the other one. While in Figs. 6 and 7, we can find that the weight parameters have the same order of magnitude. Selecting different activation functions may result in different converged weight vector.

V. CONCLUSION

In this paper, a data-based online learning algorithm for weakly coupled nonlinear systems is established. The optimal control law is derived by the optimal controllers of the reduced-order subsystems. We use the model-free integral PI algorithm with an exploration to solve the HJB equations related to the subsystems. We use the actor-critic technique and the least squares method to implement the constructed algorithm. The effectiveness of the developed optimal control law is demonstrated by two simulation examples.

APPENDIX PROOF OF THEOREM 1

Using the notation

$$V_x^* = \begin{bmatrix} V_{1x_1}^* + \varepsilon V_{\varepsilon x_1}^* \\ \varepsilon V_{\varepsilon x_2}^* + V_{2x_2}^* \end{bmatrix}$$

and setting $\varepsilon^2 = 0$, the HJB equation (6) can be rewritten as (A1), shown at the top of the page, which consists of three parts, i.e., HJB1, HJB2, and the last term which will be simplified as HJB3 in (A2), shown at the top of the page. The optimal control law $u^*(x)$ can be calculated as (A3), shown at the top of the page. Then we have the expressions of $u_{11}^*(x_1)$, $u_{12}^*(x)$, $u_{21}^*(x)$, and $u_{22}^*(x_2)$ as in (7). According to the optimal control theory, HJB1 = 0 is the HJB equation for the subsystem 1

$$\dot{x}_1(t) = f_{11}(x_1) + g_{11}(x_1)u_{11}(t)$$

and the optimal control law is $u_{11}^*(x_1) = -(1/2)R_1^{-1}g_{11}^\mathsf{T}(x_1)V_{1x_1}^*$. HJB2 = 0 is the HJB equation for the subsystem 2

$$\dot{x}_2(t) = f_{22}(x_2) + g_{22}(x_2)u_{22}(t)$$

and the optimal control law is $u_{22}^*(x_2) = -(1/2)R_2^{-1}g_{22}^\mathsf{T}(x_2)V_{2x_2}^*$.

To simplify the last term in (A1) besides HJB1 and HJB2, we give the following equations according to (A3):

$$V_{\varepsilon x_{1}}^{*\mathsf{T}} g_{11}(x_{1}) + V_{2x_{2}}^{*\mathsf{T}} g_{21}(x) = -2u_{12}^{*\mathsf{T}}(x) R_{1}$$

$$V_{\varepsilon x_{2}}^{*\mathsf{T}} g_{22}(x_{2}) + V_{1x_{1}}^{*\mathsf{T}} g_{12}(x) = -2u_{21}^{*\mathsf{T}}(x) R_{2}$$

$$V_{1x_{1}}^{*\mathsf{T}} g_{11}(x_{1}) = -2u_{11}^{*\mathsf{T}}(x) R_{1}$$

$$V_{2x_{2}}^{*\mathsf{T}} g_{22}(x_{1}) = -2u_{22}^{*\mathsf{T}}(x) R_{2}. \tag{A4}$$

Based on (A4), we obtain HJB3 = 0 as (A2). To solve $u_{12}^*(x)$ and $u_{21}^*(x)$ from HJB3, we integrate both sides of (A2) from t to ∞ , and obtain

$$\int_{t}^{\infty} \left[V_{1x_{1}}^{*\mathsf{T}} f_{12}(x) + V_{2x_{2}}^{*\mathsf{T}} f_{21}(x) + V_{\varepsilon x_{1}}^{*\mathsf{T}} f_{11}(x_{1}) + V_{\varepsilon x_{2}}^{*\mathsf{T}} f_{22}(x_{2}) \right] d\tau$$

$$= 2 \int_{t}^{\infty} \left[u_{11}^{*\mathsf{T}}(x_{1}) R_{1} u_{12}^{*}(x) + u_{22}^{*\mathsf{T}}(x_{2}) R_{2} u_{21}^{*}(x) - x_{1}^{\mathsf{T}} Q_{\varepsilon} x_{2} \right] d\tau.$$

Using $V_{\varepsilon}^*(x)$, we have

$$\begin{split} V_3^*(x) &= \int_t^\infty & \Big[V_{1x_1}^{*\mathsf{T}} f_{12}(x) + V_{2x_2}^{*\mathsf{T}} f_{21}(x) \\ &\quad + V_{\varepsilon x_1}^{*\mathsf{T}} f_{11}(x_1) + V_{\varepsilon x_2}^{*\mathsf{T}} f_{22}(x_2) \Big] \mathrm{d}\tau \end{split}$$

where $V_3^*(x) = V_{\varepsilon}^*(x) - 4 \int_t^{\infty} x_1^{\mathsf{T}} Q_{\varepsilon} x_2 d\tau$ with $V_3^*(0) = 0$. The proof is complete.

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