

Online Reinforcement Learning for Dynamic Multimedia Systems

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Abstract—In our previous work, we proposed a systematic cross-layer framework for dynamic multimedia systems, which allows each layer to make *autonomous* and *foresighted* decisions that maximize the system's *long-term performance*, while meeting the application's real-time delay constraints. The proposed solution solved the cross-layer optimization *offline*, under the assumption that the multimedia system's probabilistic dynamics were *known a priori*, by modeling the system as a layered Markov decision process. In practice, however, these dynamics are *unknown a priori* and, therefore, must be learned *online*. In this paper, we address this problem by allowing the multimedia system layers to learn, through repeated interactions with each other, to autonomously optimize the system's long-term performance at run-time. The two key challenges in this layered learning setting are: (i) each layer's learning performance is directly impacted by not only its own dynamics, but also by the learning processes of the other layers with which it interacts; and (ii) selecting a learning model that appropriately balances time-complexity (i.e., learning speed) with the multimedia system's limited memory and the multimedia application's real-time delay constraints. We propose two reinforcement learning algorithms for optimizing the system under different design constraints: the first algorithm solves the cross-layer optimization in a centralized manner and the second solves it in a decentralized manner. We analyze both algorithms in terms of their required computation, memory, and interlayer communication overheads. After noting that the proposed reinforcement learning algorithms learn too slowly, we introduce a complementary accelerated learning algorithm that exploits partial knowledge about the system's dynamics in order to dramatically improve the system's performance. In our experiments, we demonstrate that decentralized learning can perform equally as well as centralized learning, while enabling the layers to act autonomously. Additionally, we show that existing application-independent reinforcement learning algorithms, and existing myopic learning algorithms deployed in multimedia systems, perform significantly worse than our proposed application-aware and foresighted learning methods.

Index Terms—Cross-layer multimedia system optimization, layered Markov decision process, reinforcement learning, resource-constrained multimedia system.

I. INTRODUCTION

STATE-of-the-art multimedia technology is poised to enable widespread proliferation of a variety of life-enhancing applications, such as video conferencing, emergency services,

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surveillance, telemedicine, remote teaching and training, augmented reality, and distributed gaming. However, efficiently designing and implementing such delay-sensitive multimedia applications on resource-constrained, heterogeneous devices and systems is challenging due to the real-time constraints, high workload complexity, and the time-varying environmental dynamics experienced by the system (e.g., video source characteristics, user requirements, workload characteristics, number of running applications, memory/cache behavior etc.).

Cross-layer adaptation is an increasingly popular solution for addressing these challenges in dynamic multimedia systems (DMSs) [1]–[8]. This is because the performance of DMSs (e.g., video rate-distortion costs, delay, and power consumption) can be significantly improved by jointly optimizing parameters, configurations, and algorithms across two or more system layers (i.e., the application, operating system, and hardware layers), rather than optimizing and adapting them in isolation. However, existing cross-layer solutions have several important limitations.

Centralized Cross-Layer Solutions: The majority of these solutions require a central optimizer to coordinate the application, operating system, and hardware adaptations to optimize the performance of one or multiple multimedia applications sharing the DMS's limited resources. To this end, a new interface between the central optimizer and all of the layers is created, requiring extensive modifications to the operating system [4], [5] or the introduction of an entirely new middleware layer [1]–[3], [24], [25]. Unfortunately, these approaches ignore the fact that the majority of modern system's are comprised of layers (components or modules) that are designed by different manufacturers, which makes such tightly integrated designs impractical, if not impossible [8], [18], [19].

Myopic Cross-Layer Solutions: Another limitation of many existing cross-layer solutions for DMSs is that they are *myopic*. In other words, cross-layer decisions are made reactively in order to optimize the *immediate* utility, without considering the impact of these decisions on the future utility. However, in DMSs, it is essential to predict the impact of the current decisions on the long-term utility because the multimedia source characteristics, workload characteristics, OS and hardware dynamics are often correlated across time. Moreover, in DMSs, utility fluctuations across time lead to poor user experience and bad resource planning leads to inefficient resource usage and wasted power [26], [27]. In contrast, *foresighted* (i.e., long-term) optimization techniques take into account the impact of immediate cross-layer decisions on the DMS's expected future utility. Importantly, foresighted policy optimizations have been successfully deployed at the hardware layer to solve the dynamic power management problem [11]–[13]; however,

with the exception of our prior work [9], they have never been used for cross-layer DMS optimization, where *all the layers make foresighted decisions*.

Single-layer Learning Solutions: The various multimedia system layers must be able to adapt at run-time to their experienced dynamics. Most existing learning solutions for multimedia systems are concerned with modeling the unknown and potentially time-varying environment experienced by a single layer. A broad range of single-layer learning techniques have been deployed in multimedia (and general purpose) systems in recent years, which include estimation techniques such as maximum likelihood estimation [4], [5], [10], [12], [13], statistical fitting [7], regression methods [28], adaptive linear prediction [29], and ad-hoc estimation heuristics [3]. However, these solutions ignore the fact that DMSs do not only experience dynamics at a single layer (e.g., time-varying workload), but at multiple (possibly all) layers, which interact with each other. Hence, it is necessary to enable the layers to learn at run-time how to autonomously and asynchronously adapt their processing strategies based on forecasts about (i) their *future dynamics* and, importantly, (ii) *how they impact and are impacted by the other layers*.

In summary, while significant contributions have been made to enhance the performance of DMSs using cross-layer design techniques, no systematic cross-layer framework exists that explicitly considers: (i) the design and implementation constraints imposed by a layered DMS architecture; (ii) the ability of the various layers to autonomously make foresighted decisions in order to jointly maximize the DMS's utility; and, most importantly, (iii) how layers can learn their unknown environmental dynamics and determine how they impact and are impacted by the other layers.

In this paper, similar to [6]–[8], we consider a multimedia system in which (i) a video encoder at the application layer makes rate-distortion-complexity tradeoffs by adapting its configuration and (ii) the operating system layer makes energy-delay tradeoffs by adapting the hardware layer's operating frequency. Our contributions are as follows.

- We propose a centralized reinforcement learning algorithm for online cross-layer DMS optimization based on the well-known Q-learning algorithm. This centralized solution can be easily implemented if a single manufacturer designs all of the layers in the multimedia system.
- We propose a novel layered Q-learning algorithm, which allows the loosely-coupled multimedia system layers to autonomously learn their optimal foresighted policies online, through repeated interactions with each other. This decentralized solution can be implemented if the layers are designed by different manufacturers. We show experimentally that the proposed layered learning algorithm, which adheres to the layered system architecture, performs equally as well as a traditional centralized learning algorithm, which does not.
- We propose a reinforcement learning technique that exploits partial *a priori* knowledge about the system's dynamics in order to accelerate the rate of learning and improve overall learning performance. Unlike existing reinforcement learning techniques [14], [15], which can only

learn about previously visited state-action pairs, the proposed algorithm exploits our partial knowledge about the system's dynamics in order to learn about multiple state-action pairs even before they have been visited. Importantly, the accelerated learning algorithm can be implemented in both the centralized and layered scenarios.

The remainder of this paper is organized as follows. In Section II, we present the system model. In Section III, we formulate the cross-layer problem as a layered Markov decision process (MDP). In Section IV, we propose a centralized and layered (decentralized) Q-learning algorithm for run-time cross-layer optimization. In Section V, we propose a technique to accelerate the rate of learning by exploiting partial *a priori* knowledge that we have about the system's dynamics. In Section VI, we present our experimental results and we conclude the paper in Section VII.

II. SYSTEM SPECIFICATION

In this section, we model a system in which a video encoding application makes rate-distortion-complexity tradeoffs by adapting its configuration and the operating system makes energy-delay tradeoffs by dynamically adapting the DVS-enabled (dynamic voltage scaling) hardware's operating frequency. In Section III, we map this illustrative system into the proposed layered MDP model.

As in [21], we model the video source as a sequence of *video data units* (for example, video macroblocks, groups of macroblocks, or pictures), which arrive at a constant rate into the application's pre-encoding buffer. Similar to [11], we assume that the system operates over discrete time slots. Additionally, we assume that the time slots have variable length, such that one data unit is encoded in each time slot. We interchangeably refer to the time slot during which the n th data unit is encoded as the n th "time slot" or "stage," where $n \in \mathbb{N}$.

A. Dynamic Voltage Scaling Model

The n th data unit is processed at the current operating frequency, which is a member of the state set $\mathcal{S}_f = \{f_i : i = 1, \dots, N_f\}$. The operating frequency can be adapted from one time slot to the next by asserting a command in the action set $\mathcal{A}_u = \{u : u \in \mathcal{S}_f\}$. Similar to [11], we assume that there is a nondeterministic delay associated with the operating frequency transition such that

$$p_f(f^{n+1}|f^n, u^n) = \begin{cases} \beta, & \text{if } f^{n+1} = u^n \neq f^n \\ 1 - \beta, & \text{if } f^{n+1} = f^n \neq u^n \\ 1, & \text{if } f^{n+1} = f^n = u^n \\ 0, & \text{otherwise} \end{cases} \quad (1)$$

where $\{f^{n+1} : n \in \mathbb{N}\}$ is the sequence of operating frequencies and $\beta \in (0, 1]$ ($1 - \beta$) is the probability of a successful (an unsuccessful) operating frequency transition. In Fig. 1, this transition is illustrated by a temporal dependency between the frequency command (action) and operating frequency (state) in the joint OS/HW layer. Note that, regardless of the frequency command u^n , the n th data unit is processed at the operating frequency $f^n \in \mathcal{S}_f$. We define the cost $J_{OS}(f^n) = P(f^n)$ (watts) as the power dissipated at operating frequency $f \in \mathcal{S}_f$,

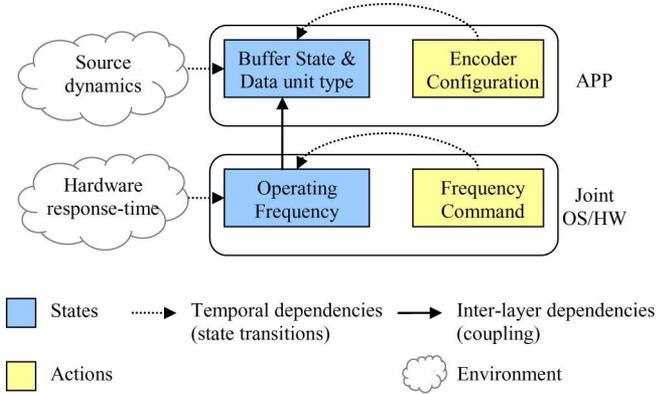


Fig. 1. System diagram showing the states, actions, and environment at each layer.

where $P(\cdot)$ is the system's power-frequency function (see, for example, [5] and [7]).

B. Application Model

The n th data unit can be classified as being one of N_z types in the state set $\mathcal{S}_z = \{z_i : i = 1, \dots, N_z\}$. The set of states \mathcal{S}_z depends on the specific video coder being used at the APP layer. For illustration, we assume that $N_z = 3$ because video streams are typically compressed into group of pictures structures containing intrapredicted (I), interpredicted (P), and bi-directionally predicted (B) data units (e.g., MPEG-2, MPEG-4, and H.264/AVC); however, the set of data unit types can be further refined based on, for example, each frame's activity level [7], [20]. It has been shown that transitions among data unit types and frame activity levels in an (adaptive) group of pictures structure can be modeled as a stationary Markov process [7], [20], [30]: i.e.,

$$p_z(z^{n+1}|z^n) \quad (2)$$

where $\{z^n : n \in \mathbb{N}\}$ is the sequence of data unit types. We assume that the choice of data unit type is governed by an algorithm similar to the one in [20]; therefore, the probabilities in (2) depend on the desired ratio of I, P, and B data units (e.g., IBBP...), the source characteristics (i.e., the APP layer's environment illustrated in Fig. 1), and the condition used to decide when to code an I frame [20]. Modeling how the multimedia data evolves over time enables us to utilize the system's resources more efficiently, while still ensuring continuity of the multimedia stream's quality over time.

The n th data unit can be coded using any one of N_h encoding parameter configurations in the action set $\mathcal{A}_h = \{h_i : i = 1, \dots, N_h\}$ (for example, to support its real-time requirements, a video encoder can adapt its choice of quantization parameter, macroblock partition size, deblocking filter, entropy coding scheme, sub-pixel motion accuracy, motion-vector search range, and motion estimation search algorithm). Given the data unit type $z^n \in \mathcal{S}_z$, each configuration $h^n \in \mathcal{A}_h$ achieves different operating points in the rate-distortion-complexity space [6]. We let $b^n(z^n, h^n)$, $d^n(z^n, h^n)$, and $c^n(z^n, h^n)$ represent the encoded bit-rate (bits per data unit), encoded distortion (mean square error), and encoding complexity (cycles), respectively.

We assume that $b^n(z^n, h^n)$, $d^n(z^n, h^n)$, and $c^n(z^n, h^n)$ are instances of the i.i.d. random variables $B(z^n, h^n)$, $D(z^n, h^n)$, and $C(z^n, h^n)$, respectively. Note that, in the proposed learning framework, we do not need to know the distributions of these random variables (see Section IV).

We penalize the application configuration by employing the Lagrangian cost measure $J_{\text{APP}}(z^n, h^n) = d^n(z^n, h^n) + \lambda_{rd} b^n(z^n, h^n)$ used in the H.264/AVC reference encoder for making rate-distortion optimal mode decisions, where $\lambda_{rd} \in [0, \infty)$ is a Lagrangian multiplier, which can be set based on the rate-constraints.

At stage n , the application's pre-encoding buffer contains $q^n \in \mathcal{S}_q = \{q_i : i = 0, \dots, N_q\}$ data units, where N_q is the maximum number of data units that can be stored in the buffer. In this paper, we assume that N_q is not limited by the available memory, but instead it is determined by the specific application's delay-constraints [21]. We assume that data units arrive in the pre-encoding buffer at η data units per second based on a deterministic arrival process from the video capture device (for example, if data units are frames, then η will typically be 15, 24, or 30 frames per second). If the buffer is full when a data unit arrives, then that data unit is discarded. In the following paragraphs, we discuss the buffer model in detail because the accelerated learning algorithm proposed in Section V exploits its structure.

The pre-encoding buffer's state at stage $n + 1$ can be expressed recursively based on its state at stage n : i.e.,

$$\begin{aligned} q^{n+1} &= \min \left\{ [q^n + \lfloor t^n(z^n, h^n, f^n) \cdot \eta \rfloor - 1]^+, N_q \right\} \\ q^0 &= q_{\text{init}} \end{aligned} \quad (3)$$

where

$$t^n(z^n, h^n, f^n) = \frac{c^n(z^n, h^n)}{f^n} \text{ (seconds)} \quad (4)$$

is the n th data unit's processing delay (in Fig. 1, the coupling between the operating frequency and processing delay is represented by the interlayer dependency arrow); $\lfloor x \rfloor$ is the integer part of x ; and $[x]^+ = \max\{x, 0\}$. Note that $c^n(z^n, h^n)$ is an instance of the random variable $C(z^n, h^n)$ with distribution $C(z^n, h^n) \sim p_C(c^n|z^n, h^n)$.

In (3), the -1 indicates that the n th data unit departs the pre-encoding buffer after the nondeterministic encoding delay $T(f^n, z^n, h^n) = C(z^n, h^n)/f^n$ (seconds). Meanwhile, the number of data units that arrive in the pre-encoding buffer during the n th time slot, $\tilde{t}^n(z^n, h^n, f^n) = t^n(z^n, h^n, f^n) \cdot \eta$, is an instance of the random variable $\tilde{T}(f^n, z^n, h^n) = \eta C(z^n, h^n)/f^n$ (data units) with distribution

$$\begin{aligned} \tilde{T}(f^n, z^n, h^n) &\sim p_{\tilde{T}}(\tilde{t}^n|f^n, z^n, h^n) \\ &= \frac{f^n}{\eta} \cdot p_C\left(\frac{f^n}{\eta} \cdot c^n|z^n, h^n\right). \end{aligned} \quad (5)$$

Based on (3) and (5), the pre-encoding buffer's state transition can be modeled as a controllable Markov chain with transition

probabilities [see equation (6), shown at the bottom of the page], where $\Omega = q^{n+1} - q^n + 1$ and $\{q^n : n \in \mathbb{N}\}$ is the sequence of buffer states. Note that, from (3), $q^{n+1} - q^n \geq -1$, with equality when there are no data unit arrivals, i.e., $\tilde{t}^n = 0$.

We define a utility gain function to nonlinearly reward the system for maintaining queuing delays less than the maximum tolerable delay, thereby protecting against overflows that may result from a sudden increase in encoding delay. Formally, we define the utility gain at stage n as

$$g_{\text{APP}}([f^n, q^n, z^n], h^n) = 1 - \left(\frac{q^n + \lfloor \tilde{t}^n(f^n, z^n, h^n) \rfloor - 1}{N_q} \right)^2 \quad (7)$$

which is near its maximum when the buffer is empty (corresponding to zero queuing delay) and is minimized when the buffer is full. In the Appendix, we discuss in detail why (7) is a good definition for the utility gain.

III. FORMULATION AS A LAYERED MDP

In Section III-A, we formulate the cross-layer problem as a layered MDP and map the system described in Section II to the proposed layered MDP model. Then, in Section III-B, we present the system optimization objective.

A. Layered MDP Model

A layered MDP is a tuple $\langle \mathcal{L}, \mathcal{A}, \mathcal{S}, p, R \rangle$, as follows.

- $\mathcal{L} = \{1, \dots, L\}$ is a set of L autonomous layers, which participate in the cross-layer optimization.¹ Each layer is indexed $l \in \{1, \dots, L\}$ with layer 1 corresponding to the lowest participating layer (e.g., the HW layer) and layer L corresponding to the highest participating layer (e.g., the APP layer).
- \mathcal{A} is the global action set $\mathcal{A} = \times_{l \in \mathcal{L}} \mathcal{A}_l$,² where \mathcal{A}_l is the l th layer's local action set.
- \mathcal{S} is the global state set $\mathcal{S} = \times_{l \in \mathcal{L}} \mathcal{S}_l$, where \mathcal{S}_l is the l th layer's local state set.
- p is the joint transition probability function mapping the global state, global action, and global next-state to a value in $[0, 1]$, i.e., $p : \mathcal{S} \times \mathcal{A} \times \mathcal{S} \mapsto [0, 1]$.

¹We note that for a layer to “participate” in the cross-layer optimization, it must be able to adapt one or more of its parameters, configurations, or algorithms (e.g., the APP layer can adapt its source coding parameters); alternatively, if a layer does not “participate”, then it is omitted

² $\times_{l \in \mathcal{L}} \mathcal{A}_l = \mathcal{A}_1 \times \dots \times \mathcal{A}_L$ is the L -ary Cartesian product.

- R is the expected reward function, which maps the global state and global action to a real number representing the system's expected reward, i.e., $R : \mathcal{S} \times \mathcal{A} \mapsto \mathbb{R}$.
- $\gamma \in [0, 1)$ is a discount factor, defined below.

We now map the system described in Section II to the layered MDP model. We consider a system in which two layers participate in the cross-layer optimization (i.e., $\mathcal{L} = \{1, 2\}$). In particular, the *application layer* (APP, $l = 2$) makes rate-distortion-complexity tradeoffs by adapting its configuration and the *operating system layer* (OS, $l = 1$) makes energy-delay tradeoffs by adapting the *hardware layer's* (HW) operating frequency (using, for example, the platform independent open standard known as the Advanced Configuration and Power Interface [11]). Due to the fact that the OS layer controls the HW layer, we combine them into a single decision making layer, which we call the joint OS/HW layer.

Table I summarizes the mapping between the system defined in Section II and the layered MDP model. Since f , q , and z are Markovian, the transition of the global state \mathbf{s} is Markovian with transition probability

$$p(\mathbf{s}' | \mathbf{s}, \mathbf{a}) = \underbrace{p_f(f' | f, u)}_{p_1(s'_1 | s_1, a_1)} \underbrace{p_q(q' | q, f, z, h)}_{p_2(s'_2 | \mathbf{s}, a_2)} p_z(z' | z). \quad (8)$$

Last, the reward function is defined as

$$R(\mathbf{s}, \mathbf{a}) = g_{\text{APP}}(\mathbf{s}, \mathbf{a}) - \omega_{\text{APP}} J_{\text{APP}}(z, h) - \omega_{\text{OS}} J_{\text{OS}}(f) \quad (9)$$

where ω_{APP} and ω_{OS} are positive weights that are assumed to be given. The forms of the factored transition probability function and the additively decomposed reward function are exploited in Section IV-C to derive the proposed layered learning algorithm.

B. System Optimization Objective

Unlike existing cross-layer optimization solutions, which focus on optimizing the myopic (i.e., immediate) utility, the goal in the proposed cross-layer framework is to find the optimal actions at each stage $n \in \mathbb{N}$ that maximize the *discounted sum of future rewards* [9], [11], [14], [15], i.e.,

$$\sum_{n=0}^{\infty} (\gamma)^n R(\mathbf{s}^n, \mathbf{a}^n | \mathbf{s}^0) \quad (10)$$

where the parameter γ ($0 \leq \gamma < 1$) is the “discount factor,” which defines the relative importance of present and future rewards, and \mathbf{s}^0 is the initial state. We refer to the Markov decision policy $\pi^* : \mathcal{S} \mapsto \mathcal{A}$, which maximizes the discounted sum of future rewards from each initial state $\mathbf{s}^0 \in \mathcal{S}$, as the optimal

$$p_{\text{APP}}^{(q)}(q^{n+1} | q^n, f^n, z^n, h^n) = \begin{cases} p_{\tilde{T}} \left\{ \tilde{T} < 2 | f^n, z^n, h^n \right\}, & q^n = q^{n+1} = 0 \\ p_{\tilde{T}} \left\{ \tilde{T} \geq N_q - q + 1 | f^n, z^n, h^n \right\}, & q^{n+1} = N_q \\ p_{\tilde{T}} \left\{ \Omega \leq \tilde{T} < \Omega + 1 | f^n, z^n, h^n \right\}, & \text{otherwise} \end{cases} \quad (6)$$

TABLE I
ABBREVIATED LIST OF NOTATION

f	Operating frequency	u	Frequency command
z	Data unit type	h	Encoding parameter configuration
q	Buffer occupancy	$a_1 = a_{OS} = u$	Joint OS/HW layer action
$s_1 = s_{OS} = f$	Joint OS/HW layer state	$a_2 = a_{APP} = h$	APP layer action
$s_2 = s_{APP} = (z, q)$	APP layer state	$\mathbf{a} = (a_1, a_2)$	Global action
$\mathbf{s} = (s_1, s_2)$	Global state	$g_2(\mathbf{s}, a_2)$	Utility gain
$p_1(s'_1 s_1, a_1),$ $p_2(s'_2 \mathbf{s}, a_2)$	Transition probability functions for the joint OS/HW and APP layers	$J_1(s_1),$ $J_2(s_2, a_2)$	Cost functions for the joint OS/HW and APP layers

foresighted policy. We use a *discounted* sum of rewards instead of, for example, the *average* reward, because: (i) typical video sources have temporally correlated statistics over short time intervals such that the future environmental dynamics cannot be easily predicted without error [7], and, therefore, the system may benefit by weighting its immediate reward more heavily than future rewards; and, (ii) The multimedia session's lifetime is not known *a priori* (i.e., the session may end unexpectedly), and, therefore, rewards should be maximized sooner rather than later.

Throughout this paper, we will find it convenient to work with the optimal *action-value function* $Q^* : \mathcal{S} \times \mathcal{A} \mapsto \mathbb{R}$ [15], which satisfies the following Bellman optimality equation:

$$Q^*(\mathbf{s}, \mathbf{a}) = R(\mathbf{s}, \mathbf{a}) + \gamma \sum_{s' \in \mathcal{S}} p(s' | \mathbf{s}, \mathbf{a}) V^*(s') \quad (11)$$

where $V^*(\mathbf{s}) = \max_{\mathbf{a}} Q^*(\mathbf{s}, \mathbf{a})$, $\forall \mathbf{s} \in \mathcal{S}$, is the optimal *state-value function* [15]. In words, $Q^*(\mathbf{s}, \mathbf{a})$ is the expected sum of discounted rewards achieved by taking action \mathbf{a} in state \mathbf{s} and then following the optimal policy $\pi^* : \mathcal{S} \mapsto \mathcal{A}$ thereafter, where

$$\pi^*(\mathbf{s}) = \arg \max_{\mathbf{a}} Q^*(\mathbf{s}, \mathbf{a}), \quad \forall \mathbf{s} \in \mathcal{S}. \quad (12)$$

IV. LEARNING THE OPTIMAL DECISION POLICY

A. Selecting a Learning Model

As we mentioned before, the multimedia system's dynamics (i.e., $R(\mathbf{s}, \mathbf{a})$ and $p(s' | \mathbf{s}, \mathbf{a})$) are *unknown*, and, therefore, the optimal action-value function Q^* and the optimal policy π^* must be learned online, based on experience. However, it remains to explain how we can learn the optimal policy online despite these unknown quantities.

One option is to directly estimate the reward and transition probability function and then to perform value iteration [15] to determine the optimal policy; however, this solution is too complex because each iteration of the algorithm has complexity $O(|\mathcal{S}|^2 |\mathcal{A}|)$ and the number of iterations required to converge to the optimal policy is polynomial in the discount factor γ .³

³Policy iteration or linear programming could also be used, but these solutions are also too complex.

Additionally, storing the estimated transition probability function incurs a memory overhead of $O(|\mathcal{S}|^2 |\mathcal{A}|)$, which for a large number of states is unacceptably large (e.g., our experimental test-bed in Section VI uses 765 states and 15 actions and, therefore, requires over 33 MBs of memory to directly store the transition probability function using 32 bit single-precision floating point numbers). A second option, proposed in [12], avoids the complex value iteration algorithm, but incurs even greater memory overheads by requiring a large number of policies to be computed and stored offline to facilitate the online decision making process.

Clearly, given the multimedia system's limited memory and the multimedia application's real-time delay constraints, neither of the above solutions is practical. Thus, in our resource-constrained cross-layer setting with many states and actions (and many unknown parameters), we adopt a *model-free* reinforcement learning solution, which can be used to learn the optimal action-value function Q^* and optimal policy π^* online, without estimating the reward and transition probability functions. Specifically, we adopt a low complexity algorithm called Q-learning, which also has limited memory requirements (see Table II in Section IV-D). To the best of our knowledge, Q-learning has never been deployed in a multimedia system, let alone for cross-layer optimization.

B. Proposed Centralized Q-Learning

Central to the conventional centralized Q-learning algorithm is a simple update step performed at the end of each time slot based on the *experience tuple* (ET) $(\mathbf{s}^n, \mathbf{a}^n, r^n, \mathbf{s}^{n+1})$

$$\delta^n = \left[r^n + \gamma \max_{\mathbf{a}' \in \mathcal{A}} Q^n(\mathbf{s}^{n+1}, \mathbf{a}') - Q^n(\mathbf{s}^n, \mathbf{a}^n) \right] \quad (13)$$

$$Q^{n+1}(\mathbf{s}^n, \mathbf{a}^n) \leftarrow Q^n(\mathbf{s}^n, \mathbf{a}^n) + \alpha^n \delta^n \quad (14)$$

where \mathbf{s}^n , \mathbf{a}^n , and $r^n = g_L^n - \sum_{l \in \mathcal{L}} \omega_l J_l^n$ are the state, performed action, and corresponding reward in time slot n , respectively; \mathbf{s}^{n+1} is the resulting state in time slot $n + 1$; \mathbf{a}' is the

TABLE II
COMPARISON OF COMPUTATION, MEMORY, AND COMMUNICATION OVERHEADS (PER TIME SLOT)

	Centralized Q-learning	Layered Q-learning
Computation overheads	Action Selection: $O(\mathcal{A})$ Update: $O(\mathcal{A})$	Action Selection: $O(\mathcal{A} + \mathcal{S}_1 \times \mathcal{A}_2)$ Update: $O(\mathcal{A})$
Memory overheads	$O(\mathcal{S} \times \mathcal{A})$	$O(\mathcal{S} \times \mathcal{A} + \mathcal{S} \times \mathcal{A}_2 \times \mathcal{S}_1)$
Communication overheads	$O(1)$ (8 messages)	$O(1)$ (7 messages)

greedy action⁴ in state \mathbf{s}^{n+1} ; δ^n is the so-called *temporal-difference* (TD) error [14]; and, $\alpha^n \in [0, 1]$ is a time-varying learning rate parameter. We note that the action-value function can be initialized arbitrarily at time $n = 0$.

It is well known that if (i) the rewards and transition probability functions are stationary, (ii) all of the state-action pairs are visited infinitely often, and (iii) α^n satisfies the *stochastic approximation conditions*⁵ $\sum_{n=0}^{\infty} (\alpha^n) = \infty$ and $\sum_{n=0}^{\infty} (\alpha^n)^2 < \infty$, then Q converges with probability 1 to Q^* [16]. Subsequently, the optimal policy can be found using (12).

During the learning process, it is not obvious what the best action is to take in each state. On the one hand, the optimal action-value function can be learned by randomly *exploring* the available actions in each state. Unfortunately, unguided randomized exploration cannot guarantee acceptable performance during the learning process. On the other hand, taking greedy actions, which *exploit* the available information in the action-value function $Q(\mathbf{s}, \mathbf{a})$, can guarantee a certain level of performance. Unfortunately, exploiting what is already known about the system prevents the discovery of new, better, actions. To judiciously trade off exploration and exploitation, we use the so-called ϵ -greedy action selection method [14]:

ϵ -greedy action selection: With probability $1 - \epsilon$, take the greedy action that maximizes the action-value function, i.e., $\mathbf{a}^* = \arg \max_{\mathbf{a}} \{Q(\mathbf{s}, \mathbf{a})\}$; and, with probability ϵ , take an action randomly and uniformly over the action set.

We write $\mathbf{a} = \Phi_{\epsilon}(Q, \mathbf{s})$ to show that \mathbf{a} is an ϵ -greedy joint-action. Note that, if the ϵ -greedy action is taken from the optimal state-value function Q^* , then actions corresponding to the optimal policy will only be taken with probability $1 - \epsilon$; therefore, if ϵ does not decay to 0, then optimal performance will not be achieved.

Fig. 2 illustrates the information exchanges required to deploy the centralized Q-learning algorithm in a two layer multimedia system during one time slot. In Fig. 2, the top and bottom blocks represent the APP layer and joint OS/HW layer, respectively. The center block represents the centralized optimizer, which selects both layers' actions and updates the system's global action-value function Q . As we mentioned be-

⁴A greedy action \mathbf{a}^* is one that maximizes the current estimate of the action-value function, i.e., $\mathbf{a}^* = \arg \max_{\mathbf{a}} \{Q(\mathbf{s}, \mathbf{a})\}$.

⁵For example, $\alpha^n = 1/(n + 1)$ satisfies the stochastic approximation conditions.

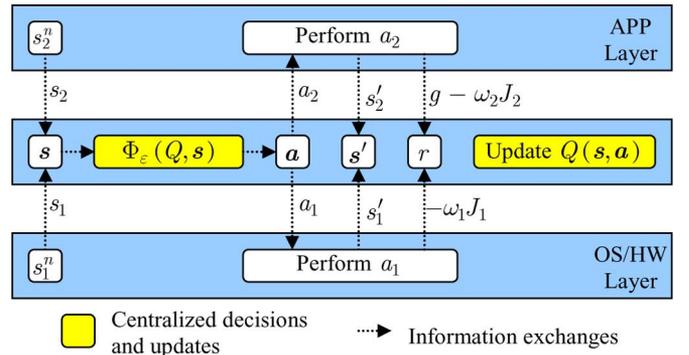


Fig. 2. Information exchanges required to deploy the centralized Q-learning update step in one time slot.

fore, the centralized optimizer may be located at either the APP layer, the joint OS/HW layer, or a separate middleware layer. To best highlight the information exchanges, we illustrate the centralized optimizer as a separate middleware layer. Although we do not explicitly indicate this in Fig. 2, the ET $(\mathbf{s}, \mathbf{a}, r, \mathbf{s}')$ is used by the block labeled “Update $Q(\mathbf{s}, \mathbf{a})$,” which performs the update step defined in (14).

C. Proposed Layered Q-Learning

Thus far, we have discussed how the traditional Q-learning algorithm can be implemented by a centralized optimizer. As we discussed in the introduction, this centralized solution is most appropriate when a single manufacturer designs all of the system's layers. However, we are also interested in developing learning solutions that allow layers designed by different manufacturers to act autonomously, but still cooperatively optimize the system's performance. This is a challenging problem in our *cross-layer* setting because the dynamics experienced at each layer may depend on the learning processes of the other layers. Hence, a key challenge is determining how to coordinate the autonomous learning processes of the various layers to enable them to successfully learn. In this section, we propose a *layered Q-learning* algorithm, which we derive by exploiting the structure of the transition probability and reward functions in order to decompose the global action-value function into local action-value functions at each layer. Then, in Section IV-D, we compare the computation, communication, and memory overheads

incurred by the centralized and the proposed layered learning algorithms.

In this subsection, we first show how the action-value function can be decomposed by exploiting the structure of the considered cross-layer problem. Subsequently, we show how this decomposition leads to a layered Q-learning algorithm, which solves the cross-layer optimization online, in a *decentralized manner*, when the transition probability and reward functions are unknown *a priori*. For notational simplicity, we drop the time slot index n , and denote the next-state \mathbf{s}^{n+1} using \mathbf{s}' .

1) *Action-Value Function Decomposition*: The proposed decomposition assumes that (i) each layer has access to the global state \mathbf{s} in each time slot; (ii) the APP layer knows the number of states at the joint OS/HW layer, i.e., $|\mathcal{S}_{\text{OS}}|$; and, (iii) the joint OS/HW layer knows the number of states and actions at the APP layer, i.e., $|\mathcal{S}_{\text{APP}}|$ and $|\mathcal{A}_{\text{APP}}|$.

Given the additive reward function defined in (9) and the factored transition probability function defined in (8), we can rewrite the optimal action-value function defined in (11), as follows:

$$Q^*(\mathbf{s}, \mathbf{a}) = g_2(\mathbf{s}, a_2) - \omega_1 J_1(s_1, a_1) - \omega_2 J_2(s_2, a_2) + \gamma \sum_{s'_1 \in \mathcal{S}_1, s'_2 \in \mathcal{S}_2} p_1(s'_1 | s_1, a_1) p_2(s'_2 | \mathbf{s}, a_2) V^*(\mathbf{s}') \quad (15)$$

where $V^*(\mathbf{s}') = \max_{\mathbf{a}' \in \mathcal{A}} Q^*(\mathbf{s}', \mathbf{a}')$ is the optimal state-value for state $\mathbf{s}' = (s'_1, s'_2)$ (refer to Table I for a review of the notation).

Observing that $g_2(\mathbf{s}, a_2) - \omega_2 J_2(s_2, a_2)$ is independent of s'_1 and that $\sum_{s'_1 \in \mathcal{S}_1} p_1(s'_1 | s_1, a_1) = 1$, we may rewrite the Bellman optimality equation in (15) as follows:

$$Q^*(\mathbf{s}, a_1, a_2) = -\omega_1 J_1(s_1, a_1) + \sum_{s'_1 \in \mathcal{S}_1} \left[p_1(s'_1 | s_1, a_1) \times \left(g_2(\mathbf{s}, a_2) - \omega_2 J_2(s_2, a_2) + \gamma \sum_{s'_2 \in \mathcal{S}_2} p_2(s'_2 | \mathbf{s}, a_2) V^*(s'_1, s'_2) \right) \right]. \quad (16)$$

Given the global state $\mathbf{s} = (s_1, s_2)$ and the optimal state-value function V^* , the APP layer's action-value function can be expressed using the inner Bellman optimality equation:

$$Q_1^*(\mathbf{s}, a_2, s'_1) = \left(g_2(\mathbf{s}, a_2) - \omega_2 J_2(s_2, a_2) + \gamma \sum_{s'_2 \in \mathcal{S}_2} p_2(s'_2 | \mathbf{s}, a_2) V^*(s'_1, s'_2) \right) \quad (17)$$

$\forall a_2 \in \mathcal{A}_2$ and $\forall s'_1 \in \mathcal{S}_1$

which is independent of the immediate action at the joint OS/HW layer (i.e., a_1), but depends on the joint OS/HW layer's potential next-state s'_1 . For each global state $\mathbf{s} = (s_1, s_2)$, the APP layer must compute the set $\{Q_1^*(\mathbf{s}, a_2, s'_1) : a_2 \in \mathcal{A}_2, s'_1 \in \mathcal{S}_1\}$ using (17). Then, given this set from the APP layer, the joint OS/HW layer's action value function can be expressed using the outer Bellman

optimality equation in (16): i.e.,

$$Q^*(\mathbf{s}, a_1, a_2) = \left(-\omega_1 J_1(s_1, a_1) + \sum_{s'_1 \in \mathcal{S}_1} p_1(s'_1 | s_1, a_1) Q_1^*(\mathbf{s}, a_2, s'_1) \right) \quad (18)$$

$\forall a_1 \in \mathcal{A}_1$ and $\forall a_2 \in \mathcal{A}_2$

Intuitively, $Q_1^*(\mathbf{s}, a_2, s'_1)$ can be interpreted as the APP layer's estimate of the expected discounted future rewards; however, there are two important differences between $Q_1^*(\mathbf{s}, a_2, s'_1)$ and the centralized action-value function $Q^*(\mathbf{s}, \mathbf{a})$. First, $Q_1^*(\mathbf{s}, a_2, s'_1)$ does not include the immediate reward at the joint OS/HW layer (i.e., $-\omega_1 J_1(s_1, a_1)$) because it is unknown to the APP layer; and, second, $Q_1^*(\mathbf{s}, a_2, s'_1)$ depends on the joint OS/HW layer's next-state s'_1 (instead of the expectation over of the next-states as it does in the centralized case) because the APP layer does not know the joint OS/HW layer's transition probability function. After receiving $Q_1^*(\mathbf{s}, a_2, s'_1)$ from the APP layer, the joint OS/HW layer is able to "fill in" this missing information. Clearly, from the derivation, $Q^*(\mathbf{s}, a_1, a_2)$ at the joint OS/HW layer is equivalent to the centralized action-value function $Q^*(\mathbf{s}, \mathbf{a})$.

For more information about this decomposition, we refer the interested reader to our prior work [9], in which we use a similar decomposition to solve the cross-layer optimization *offline*, in a decentralized manner, under the assumption that the transition probability and reward functions are *known a priori*.

2) *Layered Q-Learning*: Recall that the centralized Q-learning algorithm described in Section IV-B requires a centralized manager to select actions for both of the layers. In contrast, the layered Q-learning algorithm that we propose in this subsection—made possible by the action-value function decomposition described above—enables each layer to autonomously select its own actions and to update its own action-value function. In the following, we discuss the local information requirements for each layer, how each layer selects its local actions, and how each layer updates its local action-value function.

Local information: The proposed layered Q-learning algorithm requires that the APP layer maintains an estimate of the action-value function on the left-hand side of (17): i.e.,

$$\{Q_1(\mathbf{s}, a_2, s'_1) : \mathbf{s} \in \mathcal{S}, a_2 \in \mathcal{A}_2, s'_1 \in \mathcal{S}_1\}$$

and that the joint OS/HW layer maintains an estimate of the action-value function on the left-hand side of (18): i.e.,

$$\{Q(\mathbf{s}, a_1, a_2) : \mathbf{s} \in \mathcal{S}, (a_1, a_2) \in \mathcal{A}\}.$$

At time slot $n = 0$, these tables can be initialized to 0, i.e., $Q_1^0(\mathbf{s}, a_2, s'_1) = 0$ and $Q^0(\mathbf{s}, a_1, a_2) = 0$. Thus, at initialization, the only information required by the APP layer about the joint OS/HW layer is the number of states at the joint OS/HW layer, i.e., $|\mathcal{S}_{\text{OS}}|$; and, the only information required by the joint OS/HW layer about the APP layer is the number of states and actions at the APP layer, i.e., $|\mathcal{S}_{\text{APP}}|$ and $|\mathcal{A}_{\text{APP}}|$.

Local action selection: At run-time, given the current global state $\mathbf{s} = (s_1, s_2)$, the APP layer selects an ε -greedy action

$a_2 \in \mathcal{A}_2$ as described in Section IV-B, but with the greedy action selected as follows:

$$(a_2^*, s_1') = \arg \max_{a_2 \in \mathcal{A}_2, s_1' \in \mathcal{S}_1} \{Q_1(s, a_2, s_1')\} \quad (19)$$

and the joint OS/HW layer selects its ε -greedy action $a_1 \in \mathcal{A}_1$, but with the greedy action selected as follows:

$$(a_1^*, \hat{a}_2) = \arg \max_{a_1 \in \mathcal{A}_1, a_2 \in \mathcal{A}_2} \{Q(s, a_1, a_2)\}. \quad (20)$$

Note that the APP layer selects action a_2^* under the assumption that the joint OS/HW layer will select its action a_1^* to transition to the next-state s_1' , which maximizes the APP layer's *estimated* discounted future rewards $Q_1(s, a_2, s_1')$. Similarly, the joint OS/HW layer selects action a_1^* under the assumption that the APP layer will select action \hat{a}_2 , which maximizes the joint OS/HW layer's *estimated* discounted future rewards $Q(s, a_1, a_2)$. The message exchanges during the local learning update procedures (described immediately below) serve to “teach” each layer how they impact and are impacted by the other layer. Thus, through the learning process, the layers improve their assumptions about each other, which enables them to improve their greedy actions.

Local learning updates: After executing the ε -greedy action $\mathbf{a}^n = (a_1^n, a_2^n)$ in state $\mathbf{s}^n = (s_1^n, s_2^n)$, the system obtains the reward $r^n = g_2^n - \omega_1 J_1^n - \omega_2 J_2^n$ and transitions to state $\mathbf{s}^{n+1} = (s_1^{n+1}, s_2^{n+1})$. Based on the experience tuple $(\mathbf{s}^n, \mathbf{a}^n, r^n, \mathbf{s}^{n+1})$, each layer updates its action-value function as follows. First, the joint OS/HW layer must forward the scalar $V^n(s_1^{n+1}, s_2^{n+1}) = \max_{a_1' \in \mathcal{A}_1, a_2' \in \mathcal{A}_2} Q^n(s_1^{n+1}, s_2^{n+1}, a_1', a_2')$ to the APP layer. Using this forwarded information, the APP layer can perform its action-value function update based on the form of (17): i.e.,

$$\delta_2^n = [g_2^n - \omega_2 J_2^n + \gamma V^n(s_1^{n+1}, s_2^{n+1}) - Q_1^n(\mathbf{s}^n, a_2^n, s_1^{n+1})] \quad (21)$$

$$Q_1^{n+1}(\mathbf{s}^n, a_2^n, s_1^{n+1}) \leftarrow Q_1^n(\mathbf{s}^n, a_2^n, s_1^{n+1}) + \alpha_2^n \delta_2^n. \quad (22)$$

Then, given the scalar $Q_1^{n+1}(\mathbf{s}^n, a_2^n, s_1^{n+1})$ from the APP layer and the APP layer's selected action a_2^n , the joint OS/HW layer can perform its action-value function update based on the form of (18) as follows:

$$\delta_1^n = [-\omega_1 J_1^n + Q_1^{n+1}(\mathbf{s}^n, a_2^n, s_1^{n+1}) - Q^n(\mathbf{s}^n, a_1^n, a_2^n)] \quad (23)$$

$$Q^n(\mathbf{s}^n, a_1^n, a_2^n) \leftarrow Q^n(\mathbf{s}^n, a_1^n, a_2^n) + \alpha_1^n \delta_1^n. \quad (24)$$

Note that, by using the layered Q-learning algorithm, the layers do not need to directly share their local rewards (i.e., the local learning updates only require the APP layer to know its local reward $g_2^n - \omega_2 J_2^n$ and for the joint OS/HW layer to know its local cost $\omega_1 J_1^n$). This is because of the coordinating message exchanges during the local learning update process (i.e., $V^n(s_1^{n+1}, s_2^{n+1})$ is forwarded from the joint OS/HW layer to the APP layer and $Q_1^{n+1}(\mathbf{s}^n, a_2^n, s_1^{n+1})$ is forwarded from the APP layer to the joint OS/HW layer).

Fig. 3 illustrates the information exchanges required to deploy the layered Q-learning algorithm in a two layer mul-

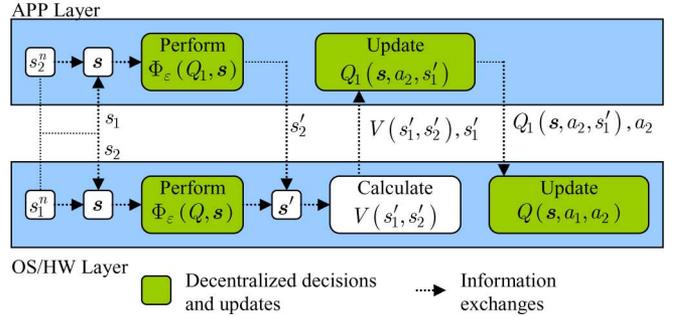


Fig. 3. Information exchanges required to deploy the layered Q-learning update step in one time slot.

timedia system during one time slot. Although we do not explicitly indicate this in Fig. 3, the ETs $(\mathbf{s}, a_2, g_2 - \omega_2 J_2, s_1')$ and $(\mathbf{s}, a_1, -\omega_1 J_1, s_1')$ are used by the blocks labeled “Update $Q_1(s, a_2, s_1')$ ” at the APP layer and “Update $Q(s, a_1, a_2)$ ” at the joint OS/HW layer, respectively.

D. Computation, Communication, and Memory Overheads

In this subsection, we compare the computation, communication, and memory overheads associated with the centralized and layered Q-learning algorithms. Figs. 2 and 3 illustrate the exact information exchanges required for each algorithm and Table II lists the per time slot computation, communication, and memory overheads associated with each algorithm.

Interestingly, the layered Q-learning algorithm is more complex and requires more memory than the centralized algorithm. Note that, in our setting, $|\mathcal{S}_1| = |\mathcal{A}_1|$; hence, the layered Q-learning algorithm incurs approximately twice the computational and memory overheads as the centralized algorithm. The increased overheads can be interpreted as the cost of optimal decentralized learning (optimal in the sense that it performs equally as well as the centralized learning algorithm). Nevertheless, the proposed centralized and layered Q-learning algorithms are far less complex and use less memory than the alternative learning solutions discussed in Section IV-A. We also observe that, in both cases, the communication overheads per time slot are $O(1)$ because they are independent of the size of the state and action sets at each layer; however, the precise number of messages exchanged depends on the deployed learning algorithm as illustrated in Figs. 2 and 3, and noted in Table II.

V. ACCELERATED LEARNING USING VIRTUAL EXPERIENCE

The large number of buffer states at the APP layer significantly limits the system's learning speed because the Q-learning update step defined in (14) updates the action-value function for only one state-action pair in each time slot. Several existing variants of Q-learning adapt the action-value function for multiple state-action pairs in each time slot. These include model-free temporal-difference- λ updates, and model-based algorithms such as Dyna and prioritized sweeping [14], [15]; however, these existing solutions are not system specific and assume no *a priori* knowledge of the problem's structure. Consequently, they can only update previously visited state-action pairs, leaving the learning algorithm to act blindly (or with

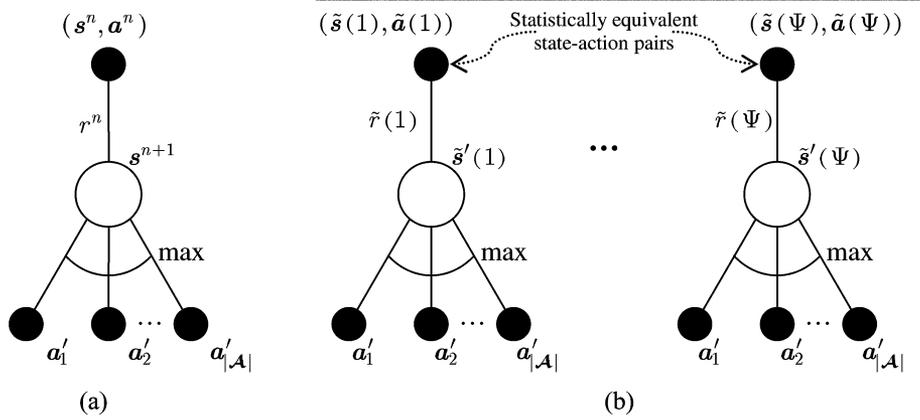


Fig. 4. Backup diagrams. (a) Conventional Q-learning; (b) Q-learning with virtual updates. The virtual update algorithm applies a backup on the actual ET and an additional Ψ backups on virtual ETs in $\Sigma(\sigma^n)$ indexed by $1 \leq \psi \leq \Psi$. (a) Q-learning backup for actual ET $\sigma^n = (s^n, a^n, r^n, s^{n+1})$. (b) Q-learning backup for Ψ virtual ETs $\sigma^n = (\tilde{s}^n, \tilde{a}^n, \tilde{r}^n, \tilde{s}' \in \Sigma(\sigma^n))$.

very little information) the first few times that it visits each state, thereby resulting in suboptimal learning performance (our experimental results in Section VI-D strongly support this conclusion).

In this subsection, we propose a new Q-learning variant, which can be used in addition to the abovementioned variants. Unlike conventional reinforcement learning algorithms, which assume that no *a priori* information is available about the environmental dynamics,⁶ the proposed algorithm exploits the *form* of the transition probability and reward functions (defined in Section II) in order to update the action-value function for multiple *statistically equivalent*⁷ state-action pairs in each time slot, including those that have never been visited. In our specific setting, we exploit the fact that the data unit arrival distribution $p_{\tilde{T}}(\tilde{t}|f, z, h)$ (defined in (5)) is conditionally independent of the current buffer state q^n (given z^n, h^n , and f^n). This allows us to extrapolate the experience obtained in each time slot to other buffer states and action pairs. Thus, in stationary (non-stationary) environments, the proposed algorithm improves convergence time (adaptation speed) at the expense of increased computational complexity. For ease of exposition, we discuss the algorithm in terms of the centralized system; however, it can be easily extended to work with the layered Q-learning algorithm proposed in Section IV-C.

Let $\sigma^n = (s^n, a^n, r^n, s^{n+1})$ represent the ET at stage n , where $s^n = (z^n, q^n, f^n)$, $a^n = (u^n, h^n)$, $r^n = g_2^n - \omega_1 J_1^n - \omega_2 J_2^n$, and $s^{n+1} = (z^{n+1}, q^{n+1}, f^{n+1})$. Given q^n and the number of data unit arrivals $[\tilde{t}^n] = [\tilde{t}^n(z^n, h^n, f^n)]$, the next buffer state q^{n+1} can be easily determined from (3). By exploiting our partial knowledge about the system's dynamics,

⁶In conventional reinforcement learning [14], [15], the actors (i.e., the system layers in our setting) are assumed to have no *a priori* information about the form of the transition probability and reward functions beyond possible high-level structural knowledge about the factored transition and reward dynamics [17], [23]. In other words, in a conventional reinforcement learning framework, the additive decomposition structure of the reward function defined in (9) and the factored transition probability structure defined in (8) may be known *a priori*, but the actual form of the utility gain defined in (7) and the actual form of the transition probability function defined in (6) cannot be known *a priori*.

⁷We say that a state-action pair (\tilde{s}, \tilde{a}) is *statistically equivalent* to the pair (s, a) if $p(s'|s, a) = p(s'|\tilde{s}, \tilde{a})$, $\forall s' \in \mathcal{S}$ and $R(\tilde{s}, \tilde{a})$ can be determined from $R(s, a)$.

we can use the statistical information provided by $[\tilde{t}^n]$ to generate *virtual experience tuples* (virtual ETs) that are statistically equivalent to the actual ET. This statistical equivalence allows us to perform the Q-learning update step for the virtual ETs using information provided by the actual ET.

We let $\tilde{\sigma}^n = (\tilde{s}, \tilde{a}, \tilde{r}, \tilde{s}')$ $\in \Sigma(\sigma^n)$ represent one virtual ET in the set of virtual ETs $\Sigma(\sigma^n)$. In order to be statistically equivalent to the actual ET, the virtual ETs in $\Sigma(\sigma^n)$ must satisfy the following two conditions.

1. The data unit arrival distribution $p_{\tilde{T}}(\tilde{t}|f, z, h)$ defined in (5) must be the same for the virtual ETs as it is for the actual ET. In other words, the virtual operating frequency \tilde{f} , the virtual type \tilde{z} , and the virtual configuration \tilde{h} must be the same as the actual operating frequency f^n , the actual type z^n , and the actual configuration h^n , respectively. This also implies that the virtual costs at the APP and joint OS/HW layers are the same as the actual costs at these layers, i.e., $\tilde{J}_1 = J_1^n$ and $\tilde{J}_2 = J_2^n$.
2. The next virtual buffer state $\tilde{q}' \in \mathcal{S}_q$ must be related to the current virtual buffer state $\tilde{q} \in \mathcal{S}_q$ through the buffer evolution equation defined in (3): i.e.,

$$\tilde{q}' = \min \left\{ [\tilde{q} + [\tilde{t}^n] - 1]^+, N_q \right\} \quad (25)$$

where $[\tilde{t}^n]$ is the number of data unit arrivals under the actual ET.

Any virtual ET that satisfies the two above conditions can have its reward determined using information embedded in the actual ET. Specifically, from the first condition, we know that the virtual ET's local costs are $\tilde{J}_1 = J_1^n$ and $\tilde{J}_2 = J_2^n$. Then, based on the second condition, we can compute the virtual ET's utility gain as $\tilde{g}_2 = 1 - (\tilde{q} + [\tilde{t}^n] - 1/N_q)^2$. Finally, the Q-learning update step defined in (14) can be performed on every virtual ET $\tilde{\sigma}^n = (\tilde{s}, \tilde{a}, \tilde{r}, \tilde{s}')$ $\in \Sigma(\sigma^n)$ as if it is the actual ET.

Performing the Q-learning update step on every virtual ET in $\Sigma(\sigma^n)$ incurs a computational overhead of approximately $O(|\Sigma(\sigma^n)| \times |\mathcal{A}|)$ in time slot n (note that $|\Sigma(\sigma^n)|$ is typically as large as the buffer state set $\mathcal{S}_q = \{q : i = 0, \dots, N_q\}$ if you

TABLE III
 ACCELERATED LEARNING USING VIRTUAL EXPERIENCE TUPLES

1.	Initialize $Q(s, a)$ arbitrarily for all $(s, a) \in \mathcal{S} \times \mathcal{A}$;	
2.	Initialize state s^0 ;	
3.	For $n = 0, 1, \dots$	
4.	Take action a^n using ε -greedy action selection on $Q(s^n, \cdot)$;	
5.	Obtain experience tuple $\sigma^n = (s^n, a^n, r^n, s^{n+1})$ and record $[\tilde{t}^n]$;	
6.	For all $\tilde{q} \in \mathcal{S}_q$	% Generate the virtual ET set $\Sigma(\sigma^n)$
7.	$\tilde{q}' = \min \{ [\tilde{q} + [\tilde{t}^n] - 1]^+, N_q \}$;	% virtual ET next buffer state
8.	$\tilde{s} = (\tilde{f}, \tilde{z}, \tilde{q}) = (f^n, z^n, \tilde{q})$;	% virtual ET state
9.	$\tilde{a} = (\tilde{u}, \tilde{h}) = (u^n, h^n)$;	% virtual ET action
10.	$\tilde{r} = \left[1 - \left(\frac{\tilde{q} + [\tilde{t}^n] - 1}{N_q} \right)^2 \right] - \omega_1 J_1^n - \omega_2 J_2^n$;	% virtual ET reward
11.	$\tilde{s}' = (\tilde{f}', \tilde{z}', \tilde{q}') = (f^{n+1}, z^{n+1}, \tilde{q}')$;	% virtual ET next state
12.	$\tilde{\sigma} = (\tilde{s}, \tilde{a}, \tilde{r}, \tilde{s}')$ $\in \Sigma(\sigma^n)$;	% store virtual ET
13.	For Ψ virtual ETs $\tilde{\sigma} \in \Sigma(\sigma^n)$	% Update virtual ETs
14.	$\tilde{\delta} = \left[\tilde{r} + \gamma \max_{a' \in \mathcal{A}} Q(\tilde{s}', a') \right] - Q(\tilde{s}, \tilde{a})$;	% TD error for virtual ET $\tilde{\sigma}^n$
15.	$Q(\tilde{s}, \tilde{a}) \leftarrow Q(\tilde{s}, \tilde{a}) + \alpha \tilde{\delta}$;	% Q-learning update for virtual ET $\tilde{\sigma}^n$

include the actual ET). Unfortunately, it may be impractical to incur such large overheads in every time slot, especially if the data unit granularity is small (e.g., one macroblock). Hence, in our experimental results in Section VI, we show how the learning performance is impacted by updating $\Psi \leq |\Sigma(\sigma^n)|$ virtual ETs in each time slot by selecting them randomly and uniformly from the virtual ET set $\Sigma(\sigma^n)$.

Table III describes the virtual ET based learning procedure in pseudo-code and Fig. 4 compares the backup diagram [14] of the conventional Q-learning algorithm to the proposed algorithm with virtual ET updates. Importantly, because the virtual ETs are statistically equivalent to the actual ET, the virtual ET based learning algorithm is *not* an approximation of the Q-learning algorithm; in fact, the proposed algorithm accelerates Q-learning by exploiting our partial knowledge about the structure of the considered problem.

VI. EXPERIMENTS

In this section, we compare the performance of the proposed learning algorithms against the performance of several existing algorithms in the literature using the cross-layer DMS described in Section II. Table IV details the parameters used in our DMS simulator, which we implemented in Matlab. In our simulations, we use actual video encoder trace data, which we obtained by profiling the H.264 JM Reference Encoder (version 13.2) on a Dell Pentium IV computer. Our traces comprise measurements of the encoded bit-rate (bits/MB), reconstructed distortion (MSE), and encoding complexity (cycles) for each video MB of the *Foreman* sequence (30 Hz, CIF resolution, quanti-

zation parameter 24) under three different encoding configurations. The chosen parameters are listed in Table IV. We use a data unit granularity of one macroblock. As in [7], we assume that the power frequency function is of the form $P(f) = \kappa f^\theta$, where $\kappa \in \mathbb{R}_+$ and $\theta \in [1, 3]$. Since real-time encoding is not possible with the available encoder, we set the data unit arrival rate to $\eta = 44$ DUs/s, which corresponds to 1/9 frames per second.

For the simulations in Sections VI-B and C, we use stationary data traces, which we generate from the measured (nonstationary) data traces by assuming that the rate, distortion, and complexity samples are drawn from i.i.d. random variables with distributions equivalent to the distributions of the quantities over the entire nonstationary data traces. In Section VI-D, we perform simulations using both the measured (nonstationary) and generated (stationary) data traces.

A. Evaluation Metrics

We deploy two different metrics to compare the performance of the various learning algorithms discussed in this paper. The first metric is the *weighted estimation error*, which we define as

$$\text{Weighted estimation error} \triangleq \sum_{\mathbf{s} \in \mathcal{S}} \mu^*(\mathbf{s}) \left| \frac{V^*(\mathbf{s}) - V^n(\mathbf{s})}{V^*(\mathbf{s})} \right| \quad (26)$$

where μ^* is the stationary distribution under the optimal policy π^* , $V^*(\mathbf{s}) = \max_{\mathbf{a}} Q^*(\mathbf{s}, \mathbf{a})$ is the optimal state-value function, and $V^n(\mathbf{s}) = \max_{\mathbf{a}} Q^n(\mathbf{s}, \mathbf{a})$ is the state-value function

TABLE IV
SIMULATION PARAMETERS (FOREMAN SEQUENCE, 30 HZ, CIF RESOLUTION, QUANTIZATION PARAMETER 24)

Layer	Parameter	Value
Application Layer (APP)	Data Unit Granularity	1 Macroblock
	Buffer State Set	$\mathcal{S}_q = \{0, \dots, N_q\}$, $N_q = 50$ DUs
	Data Unit Type Set	$\mathcal{S}_q = \{z_1, z_2, z_3\}$ $z_1 = P$, $z_2 = B$, $z_3 = I$
	Parameter Configuration	$\mathcal{A}_{APP} = \{h_1, h_2, h_3\}$ h_1 : Quarter-pel MV, 8x8 block ME h_2 : Full-pel MV, 8x8 block ME h_3 : Full-pel MV, 16x16 block ME
	APP Cost Weight	$\omega_{APP} = 22/1875$
	Rate-Distortion Lagrangian	$\lambda_{rd} = 1/16$
	Data Unit Arrival Rate	$\eta = 44$ (Data Units/Sec)
Operating System / Hardware Layer (Joint OS/HW)	Power-Frequency Function	$P(f) = \kappa f^\theta$ $\kappa \in 1.5 \times 10^{-27}$ and $\theta = 3$.
	Operating Frequency Set	$\mathcal{A}_{OS} = \{200, 400, 600, 800, 1000\}$ Mhz
	Joint OS/HW Cost Weight	$\omega_{OS} = 22/125$

estimate at stage n . In (26), we weight the estimation error $|V^*(\mathbf{s}) - V^n(\mathbf{s})|$ by $\mu^*(\mathbf{s})$ because the optimal policy will often only visit a small subset of the states, and the remaining states will never be visited; thus, if we were to estimate $V^*(\mathbf{s})$ while following the optimal policy, a *nonweighted* estimation error would not converge to 0, but the proposed weighted estimation error would.

Although the weighted estimation error tells us *how quickly an algorithm can learn the optimal policy*, it does not tell us *how well the algorithm performs while learning*. To see why this is, consider an (exaggerated) scenario in which $|V^*(\mathbf{s}) - V^n(\mathbf{s})| = 0$, but the exploration rate in the greedy action selection procedure is $\varepsilon = 1$ (i.e., always explore). In this scenario, the greedy actions are optimal, but the random policy implemented by always exploring will be far from optimal. For this reason, we define a second evaluation metric using the *average reward*, which more accurately measures the system's performance. For a simulation of duration N , the average reward metric is defined as

$$\text{Average reward} \triangleq \frac{1}{N} \sum_{n=0}^{N-1} r^n \quad (27)$$

where r^n is the reward sample obtained at stage n . If the exploration rate ε decays appropriately to 0 as $n \rightarrow \infty$, then the average reward will increase as the weighted estimation error decreases. Determining an optimal decay strategy for ε is an important problem, but is beyond the scope of this paper. We refer the interested reader to [14], [15] [31] for more information on the exploitation-exploration trade off.

We note that the weighted estimation error cannot be evaluated under nonstationary dynamics because the optimal state value function cannot be determined. Hence, for nonstationary dynamics, we rely solely on the average reward to evaluate

the learning performance. Nevertheless, the rate at which the weighted estimation error converges under stationary dynamics for a particular learning algorithm is indicative of how well that algorithm will perform when the dynamics are nonstationary.

B. Single-Layer Learning Results

Before presenting our cross-layer learning results, we believe it is informative to see how well a single layer can learn (say layer l) when the other layer (say layer $-l$) deploys a static policy. The learning layer deploys a simple adaptation of the centralized Q-learning algorithm in which it knows the global state and global reward, but does not know the actions implemented by the other layer. The details of the single-layer learning algorithm are omitted due to space constraints; however, we note that it is essential that the learning layer knows the global state and global reward so that it can learn how it impacts the other system layers. Otherwise, if the joint OS/HW layer is unaware of its impact on the application's delay and quality, for example, it will always selfishly minimize its own costs by operating at its lowest frequency and power, which can adversely impact the real-time application's performance.

For illustration, we assume that the static layer deploys its optimal local policy π_{-l}^* corresponding to the global optimal policy $\pi^* = (\pi_l^*, \pi_{-l}^*)$; hence, layer l attempts to learn π_l^* . Fig. 5 illustrates the optimal policies at the APP layer (π_2^*) and the joint OS/HW layer (π_1^*) for each buffer state and data unit type when the current operating frequency is $f = 600$ MHz. In Fig. 5, the application actions (i.e., parameter configurations) are as defined in Table IV.

Fig. 6 compares the cumulative average reward obtained using single-layer learning to the optimal achievable reward, and Table V shows the corresponding power, rate-distortion costs, utility gain, and buffer overflows, for a simulation of duration $N = 192,000$ time slots (approximately 485 frames

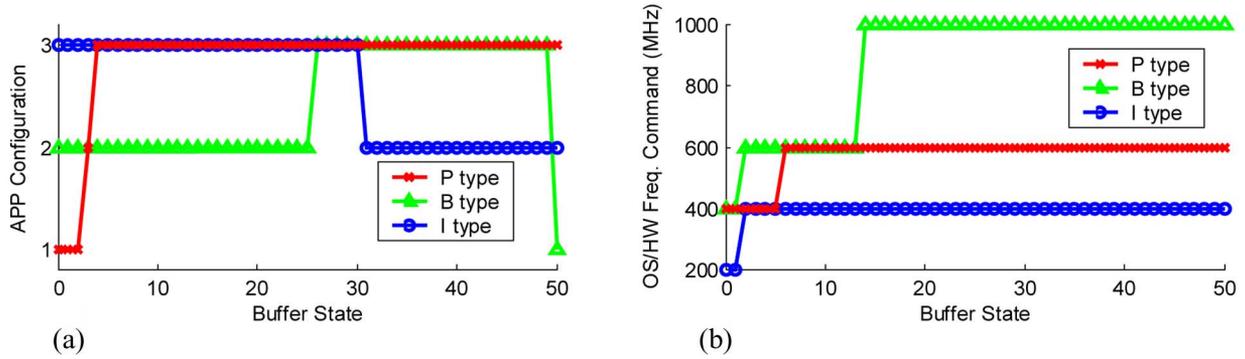


Fig. 5. Optimal policies for each data unit type. (a) APP parameter configurations. (b) Joint OS/HW layer frequency command. The current operating frequency is set to be $f = 600$ MHz.

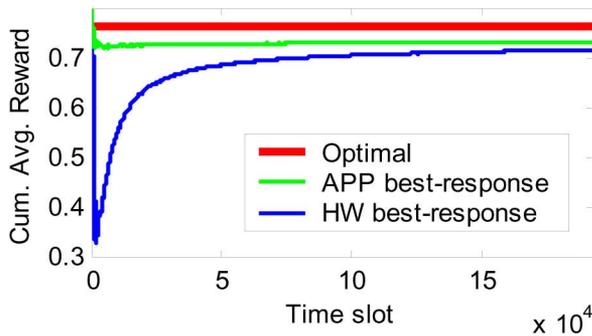


Fig. 6. Cumulative average reward for single-layer learning with stationary trace ($N = 192,000$).

drawn from the Foreman sequence, CIF resolution, by repeating the sequence from the beginning after 300 frames). We observe that the total average reward obtained when the joint OS/HW layer learns in response to the APP layer’s optimal local policy π_2^* is lower than when the APP layer learns in response to the joint OS/HW layer’s optimal local policy π_1^* . This is because there are more actions to explore at the joint OS/HW layer (5 compared to 3), and the joint OS/HW layer’s policy is more important to the system’s overall performance than the APP layer’s policy for the chosen simulation parameters (see Table IV). For instance, given the action sets defined in Table IV, the joint OS/HW layer can significantly impact the application’s experienced delay (i.e., a factor of 5 change from 200 MHz to 1000 MHz), while the APP layer cannot (i.e., its actions impact the delay by less than a factor of 2). Consequently, when the APP layer learns in response to the joint OS/HW layer’s optimal policy [Fig. 5(b)], the system is initially better off than when the joint OS/HW layer learns in response to the APP layer’s optimal policy [Fig. 5(a)].

We note that, in Fig. 6, the saturated performance of the APP and joint OS/HW layers’ learning algorithms is due to the finite exploration probability (i.e., $\epsilon > 0$ in the ϵ -greedy action selection procedure), which forces the layers to occasionally test suboptimal actions.

C. Cross-Layer Learning Results

In this subsection, we evaluate the performance of the centralized and layered Q-learning algorithms proposed in Sections IV-B and C, respectively. Fig. 7 compares the cumulative

TABLE V
SINGLE-LAYER LEARNING PERFORMANCE STATISTICS FOR STATIONARY TRACE
($N = 192,000$)

	APP Layer Learns	Joint OS/HW Layer Learns	Optimal
Avg. Reward	0.7337	0.7172	0.7631
Avg. Power (W)	0.2835	0.4512	0.2435
Avg. R-D	14.93	15.08	14.75
Avg. Gain	0.9588	0.9736	0.9790
No. Overflows	0	61	0

average reward obtained using layered Q-learning to the performance of the centralized learning algorithm, the optimal achievable reward, and the performance of the myopic learning algorithm deployed in the state-of-the-art cross-layer coordination framework called GRACE-1⁸ [5]; Fig. 8 compares the weighted estimation error for the centralized and layered Q-learning algorithms; and, Table VI shows the corresponding power, rate-distortion costs, utility gain, and buffer overflows. As in the previous simulations, the simulation duration is $N = 192,000$ time slots. Recall from Table II that the algorithms compared in this subsection have roughly the same computational complexity (note that the complexity of the GRACE-1 algorithm is roughly $O(|\mathcal{A}|)$ in our setting because it requires finding the action that will most closely meet a data unit’s deadline).

As expected, the layered Q-learning algorithm performs equally as well as the centralized Q-learning algorithm, but also allows the layers to act autonomously through decentralized decisions and updates. Meanwhile, the myopic learning algorithm deployed by GRACE-1 performs very poorly in terms of overall reward. This is because, for each data unit, GRACE-1 myopically selects the lowest processing frequency that can successfully encode it before its deadline; consequently, the buffer quickly fills, which makes the system susceptible to delay deadline violations due to the time-varying workload (this is similar to what happens when using the “conventional gain” function described in the Appendix).

⁸In GRACE-1, the statistical cycle demand (complexity) of a video encoding application is measured using an exponential average of the 95th-percentile complexity of recently completed jobs (processed data units) in a sliding window. Based on the statistical cycle demand, GRACE-1 myopically encodes jobs at the lowest frequency which will meet the job’s deadline. If slack remains after processing a job, then it is reclaimed for future jobs.

TABLE VI
CROSS-LAYER LEARNING PERFORMANCE STATISTICS FOR STATIONARY TRACE ($N = 192,000$)

	Optimal	Centralized Q-learning	Layered Q-Learning	GRACE-1
Avg. Reward	0.7631	0.4727	0.4721	-0.1730
Avg. Power (W)	0.2435	0.3815	0.4061	0.4982
Avg. Rate-Distortion	14.75	15.00	14.99	15.30
Avg. Utility Gain	0.9790	0.7158	0.7195	0.0942
No. Overflows	0	1231	1251	3040

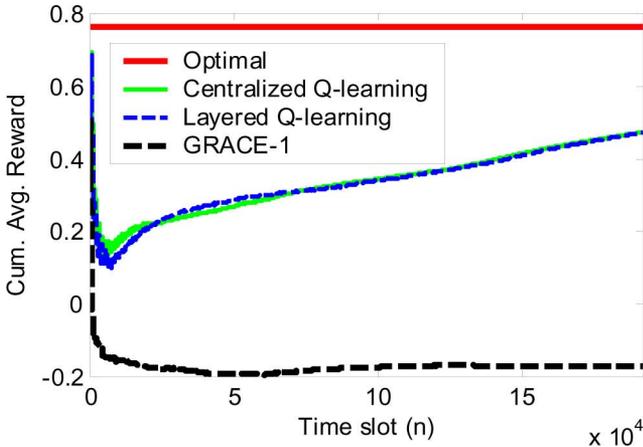


Fig. 7. Cumulative average reward for cross-layer learning algorithms with stationary trace ($N = 192,000$).

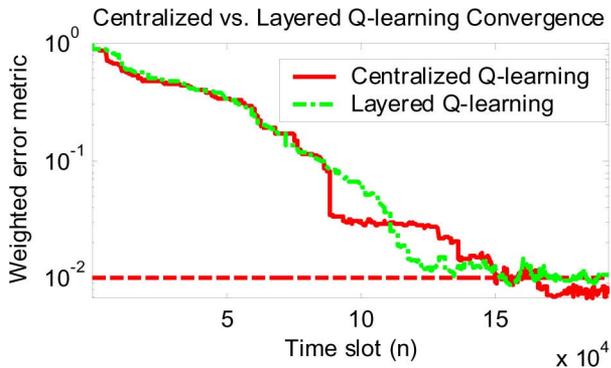


Fig. 8. Weighted estimation error metric for cross-layer learning algorithms with stationary trace ($N = 192,000$).

It is important to note that, even though the proposed learning algorithms outperform the existing myopic learning algorithm, there is still significant improvement to be made. In particular, the centralized and layered Q-learning algorithms learn too slowly because they only update one state-action pair in each time slot. Consequently, the number of buffer overflows and the power consumption are both higher, and the utility gain is lower, than in the optimal case. In Section VI-D, we show that the learning performance (including the weighted estimation error and the average reward) can be dramatically improved by *smartly* updating multiple state-action pairs in each time slot.

D. Accelerated Learning With Virtual ETs

In this subsection, we compare the performance of the proposed virtual ET based learning algorithm (see Section V) to an

existing reinforcement learning algorithm called temporal-difference- λ learning⁹ (i.e., $TD(\lambda)$ [15]), which has comparable complexity (roughly $O((\Psi + 1)|\mathcal{A}|)$ in each time slot for Ψ virtual ET or $TD(\lambda)$ updates in addition to the actual ET update). Recall that, in this subsection, we simulate the system with the measured (nonstationary) trace data and the generated (stationary) trace data.

Fig. 9 illustrates the cumulative average reward achieved when we allow $\Psi \in \{0, 1, 15, 30, 45\}$ virtual ET or $TD(\lambda)$ updates in each time slot (the case when $\Psi = 0$ is equivalent to the conventional centralized Q-learning algorithm in which only the actual ET is updated); Fig. 10 compares the weighted estimation error for the same learning algorithms; and, Table VII illustrates corresponding detailed simulation results.

Interestingly, the $TD(\lambda)$ updates are completely ineffectual at improving the system's performance. In particular, it is obvious that increasing the number of $TD(\lambda)$ updates in each time slot does not guarantee higher average rewards or lower weighted estimation errors. Meanwhile, increasing the number of virtual ET updates in each time slot dramatically improves the weighted estimation error metric and the average reward. The reason for this stark difference in performance is because the virtual ET based learning algorithm can learn about state-action pairs *without visiting them*, while the $TD(\lambda)$ learning algorithm can only learn about *previously visited state-action pairs*. The ability to learn about unvisited states is very important in our problem setting because after the buffer initially fills from early exploration (resulting in the initial drop in rewards illustrated in Figs. 6, 7, and 9), the learning algorithm must learn to efficiently drain the buffer without consuming too much power or unnecessarily reducing the application's quality. This is simple when using virtual ETs because the information obtained from experience in a "near full" buffer state (e.g., $q = N_q - 1$) can be extrapolated to a "near, near full" buffer state (e.g., $q = N_q - 2$) and so on. When using $TD(\lambda)$ updates, however, the algorithm learns very slowly about "near, near full" buffer states because they are visited only infrequently from the "near full" buffer state. For this reason, $TD(\lambda)$ (even for a large number of updates) has trouble emptying the buffer, and performs no better than the conventional centralized Q-learning algorithm, which updates only one state-action pair in each time slot.

⁹In our implementation, $TD(\lambda)$ applies the update step defined in (14) to the Ψ most frequently visited states in the recent past. These recently visited states are updated using the TD error obtained for the current state weighted by the previously visited states' "eligibility" as defined in [15]. Informally, the "eligibility" is the discounted frequency with which states have been visited in the past.

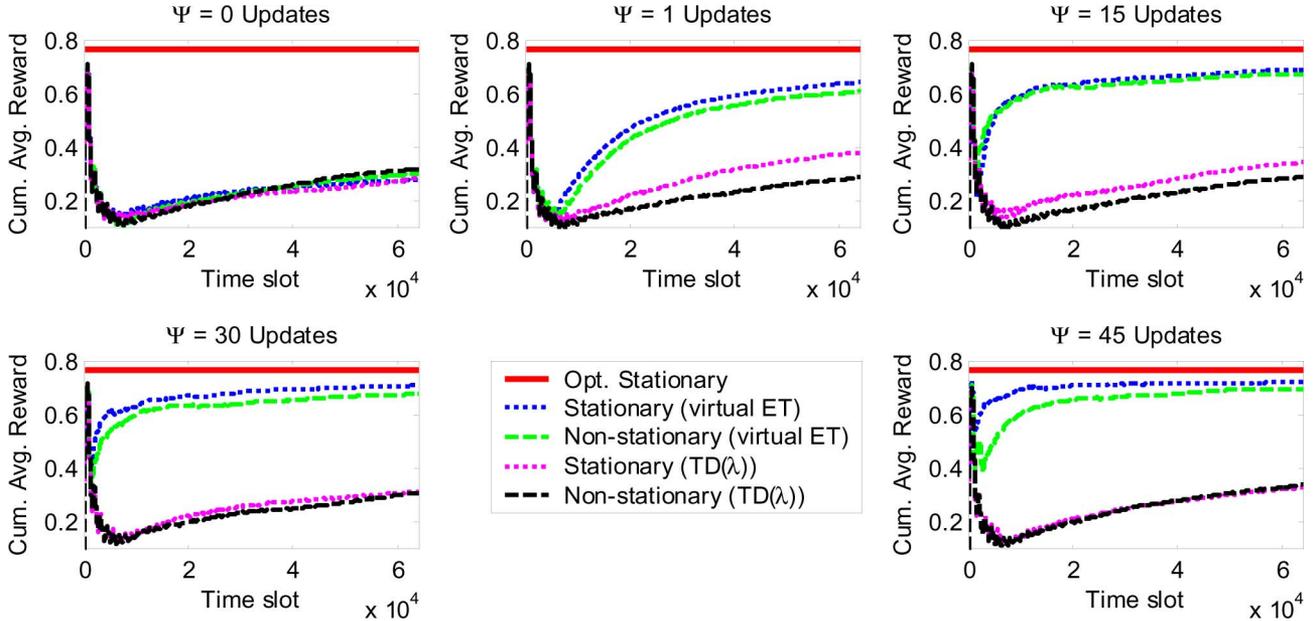


Fig. 9. Cumulative average reward achieved with virtual ET and $TD(\lambda)$ updates for a stationary trace and a nonstationary trace, compared to the optimal achievable reward under the stationary trace ($N = 64,000$).

TABLE VII

VIRTUAL EXPERIENCE LEARNING STATISTICS FOR STATIONARY (NON-STATIONARY) DATA TRACE OUTSIDE (INSIDE) PARENTHESES ($N = 64,000$)

	Number of Virtual ET Updates				
	$\Psi = 0$	$\Psi = 1$	$\Psi = 15$	$\Psi = 30$	$\Psi = 45$
Avg. Reward	0.2785 (0.3015)	0.6414 (0.6078)	0.6893 (0.6710)	0.7076 (0.6759)	0.7200 (0.6944)
Avg. Power (W)	0.4212 (0.4632)	0.4164 (0.4314)	0.3817 (0.3864)	0.3432 (0.3657)	0.3108 (0.3381)
Avg. Rate-Distortion	15.00 (14.89)	14.92 (14.77)	15.06 (14.84)	14.93 (14.81)	14.87 (14.71)
Avg. Utility Gain	0.5287 (0.5577)	0.8898 (0.8570)	0.9332 (0.9131)	0.9432 (0.9140)	0.9492 (0.9264)
No. Overflows	1196 (1410)	1035 (1271)	502 (437)	226 (378)	103 (438)
	Number of $TD(\lambda)$ Updates				
	$\Psi = 0$	$\Psi = 1$	$\Psi = 15$	$\Psi = 30$	$\Psi = 45$
Avg. Reward	0.2835 (0.3188)	0.3807 (0.2899)	0.3449 (0.2903)	0.3141 (0.3088)	0.3287 (0.3388)
Avg. Power (W)	0.4494 (0.4707)	0.4649 (0.4732)	0.4551 (0.4672)	0.4530 (0.4644)	0.4544 (0.4875)
Avg. Rate-Distortion	15.07 (14.88)	15.14 (14.92)	15.14 (14.88)	15.03 (14.84)	15.15 (14.91)
Avg. Utility Gain	0.5395 (0.5763)	0.6402 (0.5483)	0.6026 (0.5472)	0.5703 (0.5647)	0.5866 (0.5997)
No. Overflows	1255 (1630)	1230 (1491)	1267 (1511)	1234 (1491)	1250 (1317)

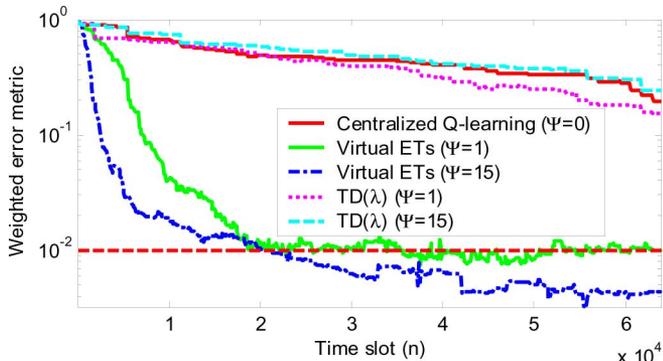


Fig. 10. Weighted estimation error metric for virtual ETs and $TD(\lambda)$ with stationary trace ($N = 64,000$).

One might expect that violating the stationarity assumption, which is widely used in the reinforcement learning literature (see [14]–[16]), would be disastrous to the system’s performance. We observe in Fig. 9 and Table VII, however, that the

proposed virtual ET based learning algorithm is robust to the nonstationary dynamics in the measured data traces. This is primarily due to its ability to quickly propagate new information obtained at one state-action pair throughout the state-action space. From the data in Table VII, we observe that the percent difference between the reward obtained in the stationary case and the nonstationary case is less than 6%.

As in the single-layer learning case, the learning performance with virtual ETs saturates due to the finite exploration probability (i.e., $\epsilon > 0$), which prevents the learning algorithm from converging to the optimal policy.

It is important to note that, in these experiments, we have assumed that the update complexity is negligible. This would be true if the updates were performed infrequently (e.g., per video frame) or if the updates were performed in parallel (e.g., using a vector processor); however, it is not a valid assumption for the very frequent and serial updates deployed here (i.e., per video macroblock, without a vector processor). Hence, performing 15,

30, or 45 updates in each time slot is not reasonable, but performing 1 or 2 virtual ET updates is. Despite this technicality, we have shown the relative performance improvements that can be achieved by updating multiple virtual experience tuples in each time slot. An interesting trade off to investigate in future research is the impact of a coarser data unit granularity with a simultaneous increase in the number of virtual updates (such that the average learning complexity remains constant). Last, we note that the learning performance could be further improved (for the same number of virtual ET updates) by directing the virtual ET updates to states that are most likely to be visited in the near future (e.g., states near the observed next-state), instead of randomly updating the virtual experience tuples.

VII. CONCLUSION

Foresighted decisions are required to optimize the performance of resource-constrained multimedia systems. However, in practical settings, in which the system's dynamics are unknown *a priori*, efficiently learning the optimal foresighted decision policy can be a challenge. Specifically, selecting an improper learning model can lead to an unacceptably slow learning speed, incur excessive memory overheads, and/or adversely impact the application's real-time performance due to large computational overheads. In this paper, we choose reinforcement learning to appropriately balance time-complexity, memory complexity, and computational complexity. We propose a centralized and a layered reinforcement learning algorithm and extend these algorithms to exploit our partial knowledge of the system's dynamics. The proposed accelerated learning algorithm, which is based on updating multiple statistically equivalent state-action pairs in each time slot, allows us to judiciously tradeoff per-stage computational complexity and time-complexity (i.e., learning speed), while incurring low memory overheads.

In our experimental results, we verify that the layered learning solution performs equally as well as the centralized solution, which may be impractical to implement if the layers are designed by different manufacturers or operate at different time-scales. We also illustrate that our proposed foresighted learning algorithms outperform the myopic learning algorithm deployed in an existing state-of-the-art cross-layer optimization framework and that, by exploiting knowledge about the system's dynamics, we can significantly outperform an existing application-independent reinforcement learning algorithm that has comparable computational overheads.

A fruitful direction for future research is the application of the techniques introduced in this paper to cross-layer optimization of the network protocol stack.

APPENDIX

In this appendix, we discuss the form of the utility gain function defined in (7). Conventionally, if a multimedia buffer does not overflow (i.e., the buffer constraints are not violated), then there is no penalty (see, for example, [21] and [22]). Within the proposed MDP-based framework, where it is difficult to explicitly impose constraints because the dynamics are not known *a*

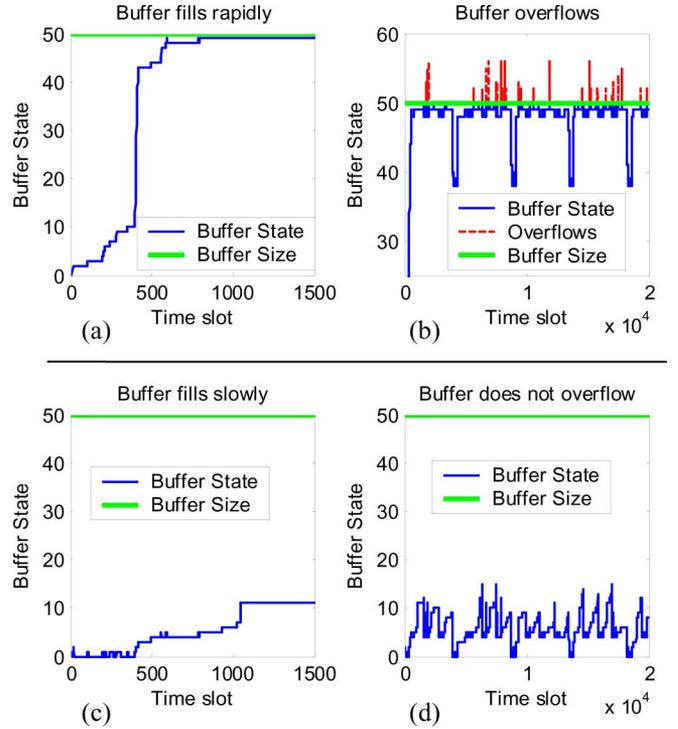


Fig. 11. Buffer evolution in $N = 20,000$ time slot simulation. (a) Conventional utility gain results in the buffer filling rapidly; (b) conventional utility gain leads to overflows; (c) proposed utility gain keeps the buffer occupancy low; (d) proposed utility gain prevents overflows.

priori, the conventional buffer model must be integrated into the system's reward function. Specifically, it can be integrated into the reward through a utility gain function of the form

$$g_{\text{conv}}(s, a_L) = \begin{cases} 1, & q + \lfloor \tilde{t} \rfloor - 1 \leq N_q \\ N_q - (q + \lfloor \tilde{t} \rfloor - 1), & \text{otherwise} \end{cases}$$

which provides a reward of 1 if the buffer does not overflow, and a penalty proportional to the number of overflows otherwise.

In contrast to the proposed continuous utility gain function defined in (7), $g_{\text{conv}}(s, a_L)$ is disjoint. As a result, the optimal foresighted policy obtained using $g_{\text{conv}}(s, a_L)$ will initially fill the buffer rapidly in an attempt to minimize rate-distortion and power costs (because it is not penalized for filling the buffer) as illustrated in Fig. 11(a). Subsequently, the policy will attempt to keep the buffer nearly full in order to balance the cost of overflow with the rate-distortion and power costs required to reduce the buffer's occupancy. Unfortunately, with the buffer nearly full, any sudden burst in the complexity of a data unit will immediately overflow the buffer as illustrated in Fig. 11(b). In contrast, the proposed utility gain function is robust against bursts in complexity because it encourages the buffer occupancy to remain low as illustrated in Fig. 11(c) and (d).

The data in Table VIII shows that the proposed utility gain function not only prevents buffer overflows, but it also achieves comparable power consumption as the conventional utility gain function. Thus, we have verified that our choice of utility gain function is good. An added benefit of the proposed utility gain function is that it aids in the learning process. This is because

TABLE VIII
PERFORMANCE STATISTICS USING THE CONVENTIONAL UTILITY GAIN
FUNCTION AND THE PROPOSED UTILITY GAIN FUNCTION

	Form of the utility gain	
	Conventional	Proposed
Avg. Power (W)	0.2421	0.2427
Avg. Rate-Distortion	14.95	14.84
No. Overflows	394	0

actions are *immediately* rewarded (or penalized) based on how they impact the buffer state.

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