Finite-Time Analysis of Adaptive Temporal Difference Learning with Deep Neural Networks

Abstract

Temporal difference (TD) learning with function approximations (linear functions or neural networks) has achieved remarkable empirical success, giving impetus to the development of finite-time analysis. As an accelerated version of TD, the adaptive TD has been proposed and proved to enjoy finite-time convergence under the linear function approximation. Existing numerical results have demonstrated the superiority of adaptive algorithms to vanilla ones. Nevertheless, the performance guarantee of adaptive TD with neural network approximation remains widely unknown. This paper establishes the finite-time analysis for the adaptive TD with multi-layer ReLU networks approximation whose samples are generated from a Markov decision process. Our established theory shows that if the width of the deep neural network is large enough, the adaptive TD using neural network approximation can find the (optimal) value function with high probabilities under the same iteration complexity as TD in general cases. Furthermore, we show that the adaptive TD using neural network approximation, with the same width and searching area, can achieve theoretical acceleration when the stochastic semi-gradients decay fast.

1 Introduction

Temporal difference (TD) learning is a popular and successful iterative algorithm in the area of reinforcement learning (RL) to evaluate a given policy [42, 46], often employed for critic part evaluation in various RL algorithms [25, 37, 38]. Classical TD algorithm adopts the tabular representation for the value function, which stores value estimates on a per-state basis. In large-scale scenarios, the tabular approach becomes intractable due to a large number of states. Thus, the function approximation techniques have been integrated with TD for better scalability and efficiency [4, 45, 33, 50]. The function approximation techniques have achieved remarkable empirical success and are theoretically justifiable when the linear function approximation is used [20]. As a special function approximation approach, deep neural networks (DNNs) have also been integrated with TD [35, 34, 30], achieving phenomenal performance in several applications. However, from the theoretical perspective, establishing theoretical convergence guarantees for training DNNs is much more complicated than that for the linear approximation algorithms, which is still widely open. Some convergence results of TD with DNN approximations have been proved by [18, 7, 48], under some extra assumptions and restrictions. The adaptive methods for DQN have been proposed in [34, 19], inspired by the adaptive stochastic algorithms to accelerate TD. The numerical results show that the adaptive versions of TD can outperform the vanilla ones in many tasks.

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Mathematically, the (projected) TD (with function approximation) can be described as

$$\boldsymbol{\theta}^{k+1} = \mathbb{P}_{\mathbf{V}}(\boldsymbol{\theta}^k - \eta \mathbf{g}^k), \tag{1}$$

where $\mathbf{V}$ is the constrained set, and $\mathbb{P}_{\mathbf{V}}$ denotes the projection onto $\mathbf{V}$, $\boldsymbol{\theta}^k$ is the iterate, and $\mathbf{g}^k$ is the stochastic semi-gradient [1]. Although the update scheme of TD looks similar to the stochastic gradient descent (SGD), it is much more complicated due to the Markov noise, even in the linear function approximation case. Motivated by the adaptive SGD [17, 24], the adaptive TD with linear function approximation is proposed in [47, 41]. The main difference between the adaptive TD and the standard TD lies in the use of adaptive stepsize and momentum. This paper considers the neural adaptive TD. In the kth iteration of neural adaptive TD, it performs

$$\begin{cases} \mathbf{m}^k = \beta \mathbf{m}^{k-1} + (1-\beta)\mathbf{g}^k, \\ v^k = v^{k-1} + \|\mathbf{g}^k\|^2, \\ \boldsymbol{\theta}^{k+1} = \mathbb{P}_{\mathbf{V}}(\boldsymbol{\theta}^k - \eta \mathbf{m}^k/(v^k)^{\frac{1}{2}}), \end{cases} \tag{2}$$

where $\beta > 0$ is the momentum parameter, and $\eta > 0$. The points $\mathbf{m}^k$ and v^k contain past information. The numerical results have demonstrated the advantage of adaptive TD over vanilla TD. The adaptive TD is proved to be convergent with linear function approximation, but the convergence remains unclear when the neural network approximation is used. *This paper proves that the adaptive TD with neural network approximation converges when the width of a ReLU network is sufficiently large. Moreover, we prove that adaptive TD is faster than TD with the ReLU DNN approximation.*

1.1 Related Works

Analysis and the recent development of TD. Leveraging the stochastic approximation techniques, the authors in [22] establish the first convergence results for TD. The limiting convergence of TD with linear function approximation is proved in [45] with the perspective of the dynamics. Since the seminal work of [45], many works have been using the ODE-based method to study the asymptotic convergence of TD since TD update does not follow the (stochastic) gradient direction of any objective function [6, 44]. Some variants of TD have been proposed in [13] with asymptotic convergence guarantees. The first non-asymptotic analysis for the gradient TD, a variant of the TD, has been studied in [31]. Finite-time analysis of TD with independent and identically distributed (i.i.d.) observation assumption has been presented in [12]. The Markov sampling convergence analysis is proved in a subsequent paper [5]. In a concurrent line of research, TD has been studied from the perspective of stochastic linear systems [26]. The finite-time analysis for Markov sampling stochastic linear system has been developed by [39, 21]. The finite-time analysis of multi-agent TD is proved by [14]. A unified analysis for a class of TD learning algorithms with Markov jump is established in [21]. Based on Nesterov's acceleration method, a class of accelerated TD is developed and analyzed in [15]. From the algorithmic viewpoint, the adaptive TD has been recently proposed in [41, 47] to accelerate TD, inspired by the adaptive SGD. In [40], the authors present the finite-time convergence results of decentralized TD with linear approximations.

TD with deep learning. In contrast to the tremendous empirical success of the deep Q-network (DQN), the theory is still relatively weak; until recently, only a handful of papers have studied the theoretical results of TD using neural network approximation. In [18], the authors prove the convergence rates of fitting Q-iteration with a sparse multi-layer ReLU network under i.i.d. observations. The convergence of TD with two-layer neural network approximation is provided by [7] with i.i.d assumption on the samples. The TD-based algorithm with multi-layer ReLU networks under Markov samples is studied in [48]. The theory of the TD with a multi-layer ReLU network relies heavily on the existing results about overparameterized deep networks [23, 11, 16, 2, 1, 3, 49].

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[1] We call it as semi-gradient because its stationary expectation is not the gradient of any fixed objective function.

 1.2 Our Contributions

 In this paper, we consider the adaptive TD with multi-layer ReLU network approximation under
 Markov observations. In contrast to the existing works [18, 7, 48], we study the adaptive variant
 of TD. The scheme of the algorithm is much more complicated, raising tremendous challenges in
 theoretical analysis. Our main theoretical contributions are summarized below:

 • We extend the analysis of adaptive TD with linear function approximation [41, 47] to
 multiple layers neural network approximation under Markovian samplings. The theoretical
 results show that adaptive TD still works for the neural network approximation, even with
 deep networks.

 • We establish the finite-time analyses of adaptive TD with multi-layer ReLU approximation
 under Markov observations. In particular, we show that the adaptive algorithms guarantee
 convergence when the neural network is sufficiently wide, and adaptive TD with neural
 network approximation converges to a projected optimal action-value function. The tech-
 nique required to connect Adam-type algorithms and neural TD is non-trivial since they
 belong to two very different research areas. To this end, we develop a new technique that
 uses expectation with a fixed delay, different from the coupling technique used by [5, 48].

 • We prove that the speed of the adaptive TD can be faster than the vanilla ones with multi-
 layer ReLU approximation. Specifically, we show that adaptive ones use fewer iterations
 to reach the same desired error with the same network widths and searching areas. Our
 theoretical results first explained why Adam-type algorithms perform better than TD in
 DQN, which has been observed in practice.

 ## 2 Preliminaries

 We introduce the notation, some basic concepts, and properties of TD in this section.

 Notation: We use $\mathbb{E}[\cdot]$ to denote the expectation with respect to the underlying probability space
 without stochasticity of the initial point, and we use $\|\cdot\|$ to denote the Frobenius norm. $\sigma(\cdot)$ denotes
 the ReLU activation function. We use $\phi : \mathcal{S} \times \mathcal{A} \to \mathbb{R}^d$ to denote the feature mapping. Given a
 closed set $\mathbf{V}$, $\mathbb{P}_{\mathbf{V}}(\mathbf{x})$ represents the projection of $\mathbf{x}$ onto $\mathbf{V}$. The initial point is defined as $\boldsymbol{\theta}^{\text{init}}$. We
 use $a = \tilde{\mathcal{O}}(b)$ to hide the logarithmic factor of b still with the same order. If $a = \mathcal{O}(b)$ and $b = \mathcal{O}(a)$,
 we use the notation $a = \Theta(b)$ and $a = \tilde{\Theta}(b)$ to hide the logarithmic factor. We denote χ^k as the
 sub-algebra generated by $\boldsymbol{\theta}^0, \boldsymbol{\theta}^1, \ldots, \boldsymbol{\theta}^k$, where $\boldsymbol{\theta}^k$ is the value in the kth iteration.

 2.1 Markov Decision Process

For the sake of presentation, we consider the finite state space[2]. Consider a Markov decision process
(MDP) described as a tuple $(\mathcal{S}, \mathcal{A}, \mathcal{P}, \mathcal{R}, \gamma)$, where $\mathcal{S}$ denotes the state space, $\mathcal{A}$ denotes the action
space, $\mathcal{P}_a$ represents the transition matrix associated with action a, and $0 < \gamma < 1$ is the discount
factor. In this case, let $\mathcal{P}_a(s'|s)$ denote the transition probability from state s to state s' under
the action a. The corresponding transition reward is $r(s, a)$. We consider the stochastic policy
$\pi : \mathcal{S} \to \Delta(\mathcal{A})$ that specifies a probability density of all actions given the current state s. $\pi(s, a)$
denotes the probability to choose action a when the current state is s, and $\sum_{a \in \mathcal{A}} \pi(s, a) = 1$. We
consider the *on policy* setting, where both target and behavior policies are π. The corresponding
action-value function $Q_\pi(s, a) : \mathcal{S} \times \mathcal{A} \to \mathbb{R}$ is defined as

$$Q_\pi(s, a) := \mathbb{E}\Big[\sum_{t=0}^{\infty} \gamma^t r(s_t, a_t) | s_0 = s, a_0 = a \Big],$$

and the associated value function $\mathbf{V}_\pi : \mathcal{S} \to \mathbb{R}$ is defined as

$$\mathbf{V}_\pi(s) := \mathbb{E}\Big[\sum_{t=0}^{\infty} \gamma^t r(s_t, a_t) | s_0 = s \Big].$$

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[2]Our results can be extended to infinite state cases, and we consider the finite state for simplicity.

105 It is evident that the restriction on discount $0 < \gamma < 1$ can guarantee the boundedness of $Q_\pi(s, a)$.
106 The Markovian property of MDP yields the following celebrated Bellman equation

$$\mathcal{T}_\pi Q_\pi = Q_\pi, \tag{3}$$

107 where the Bellman operator $\mathcal{T}_\pi$ on a value function Q can be represented as

$$(\mathcal{T}_\pi Q)(s, a) = r(s, a) + \gamma \sum_{s' \in \mathcal{S}} \mathcal{P}_a(s'|s) \mathbf{V}_\pi(s'),$$

108 where $s \in \mathcal{S}, a \in \mathcal{A}$. Directly solving (3) needs $\mathcal{O}(|\mathcal{S}|^3 |\mathcal{A}|^3)$ computations, which is very expensive
109 since $|\mathcal{S}||\mathcal{A}|$ is usually very large. Thus, people turn to use a linear or non-linear approximation as
110 $Q_\pi \approx \Phi(\boldsymbol{\theta}) : \mathbb{R}^D \to \mathbb{R}^{|\mathcal{S}||\mathcal{A}|}$, where $\boldsymbol{\theta} \in \mathbb{R}^D$ and $D \ll |\mathcal{S}||\mathcal{A}|$. In this way, the dimension can be
111 significantly reduced, and we get an approximate solution very fast.

112 We collect necessary assumptions for MDP below.

Assumption 1 *The transition rewards are uniformly bounded by 1, that is,*

$$|r(s, a)| \leq 1, \ \forall \ s, s' \in \mathcal{S}, \ a \in \mathcal{A}.$$

For any $s, \in \mathcal{A}$, it holds

$$\mu(s) = \lim_{t \to \infty} \mathcal{P}(s_t = s | s_0 = s', a_0 = a') > 0.$$

There exists a constant $0 \leq \rho < 1$ and $\bar{\kappa} > 0$ such that

$$\sum_{s \in \mathcal{S}} |\mathcal{P}(s_t = s | s_0 = s', a_0 = a') - \mu(s)| \leq \bar{\kappa} \rho^t.$$

113 The boundedness of $(r(s, s'))_{s,s' \in \mathcal{S}}$ comes from the finiteness of $\mathcal{S}$. In Assumption 1, the uniform
114 boundedness assumption can be replaced by non-uniform boundedness in the finite state case. In this
115 paper, it is used for simplicity. The rest part of Assumption 1 is standard for the Markovian property.
116 It is well-known that irreducible and aperiodic Markov chains can always follow Assumption 1
117 [27]. For Assumption 1, the time that $(s_t, a_t)_{t \geq 0}$ needed for its current state distribution to match
118 the stationary one with ϵ error in total variation distance is $\mathcal{O}(\log \frac{1}{\epsilon} / \log \frac{1}{\rho})$. Thus, the constant ρ
119 represents the speed of the process accessing to the stationary distribution. In finite-time cases, it is
120 easy to prove that ρ is the second largest eigenvalue of $\mathcal{P}$. We can see that the smaller ρ is, the faster
121 the process will converge to stationary states.

122 ## 2.2 Neural Temporal Difference Learning

Although we can get Q by solving the Bellman equation induced by the given policy π, in practice S
may contain a very large number of different states and actions, and it is hard to solve the Bellman
equation directly. Thus, alternative methods have been proposed, including leveraging the linear [43]
or non-linear approximations (e.g., kernels and neural networks [35]). This paper is devoted to the
approximation using a L-hidden-layer ReLU neural network defined as

$$f(\boldsymbol{\theta}; \mathbf{x}) = \sqrt{m} \mathbf{W}_L \sigma(\mathbf{W}_{L-1} \cdots \sigma(\mathbf{W}_1 \mathbf{x}) \cdots),$$

123 where $\mathbf{x} \in \mathbb{R}^d$ is the input data, $\mathbf{W}_1 \in \mathbb{R}^{m \times d}$, $\mathbf{W}_L \in \mathbb{R}^{1 \times m}$ and $\mathbf{W}_l \in \mathbb{R}^{m \times m}$ for $l = 2, \dots, L-1$,
124 $\boldsymbol{\theta} := [\mathbf{W}_1, \dots, \mathbf{W}_L]$ denotes all the weights. In neural TD, we use the following approximation

$$Q_\pi(s, a) \approx f(\boldsymbol{\theta}; \phi(s, a)),$$

125 where $\boldsymbol{\theta} \in \mathbb{R}^{(L-2)m^2 + md + m}$ is a parameter vector, m is usually set such that $(L-2)m^2 + md + m$
126 is smaller than $|\mathcal{S}||\mathcal{A}|$ to reduce the difficulties caused by the high dimensionality.

127 **Assumption 2** *For any state-action pair $(s, a) \in \mathcal{S} \times \mathcal{A}$, we assume the feature vector is uniformly*
128 *bounded such that $\|\phi(s, a)\| = 1$.*

129 A simple normalization can make Assumption 2 hold. This assumption is used to simplify coefficients
130 in the subsequent proofs.

131 Next, we present the scheme of neural temporal difference learning. With $(s_k, a_k)_{k \geq 0}$ be-
132 ing a trajectory sampled from π and the stepsize $\eta > 0$, the neural TD updates with $\mathbf{g}^k \leftarrow$
133 $\overline{\mathbf{g}}(\boldsymbol{\theta}^k; s_k, a_k, s_{k+1}, a_{k+1})$ in (1), where the stochastic semi-gradient is defined as

$$
\begin{aligned}
\overline{\mathbf{g}}(\boldsymbol{\theta}; s_k, a_k, s_{k+1}, a_{k+1}) := &\nabla_{\boldsymbol{\theta}} f(\boldsymbol{\theta}; \phi(s_k, a_k) \\
&\times [f(\boldsymbol{\theta}; \phi(s_k, a_k)) - r(s_k, a_k) - \gamma f(\boldsymbol{\theta}; \phi(s_{k+1}, a_{k+1}))].
\end{aligned}
\tag{4}
$$

134 The projection in scheme (2) is used to ensure the boundedness of $\boldsymbol{\theta}^k$ and simplify the convergence
135 analysis. In the deep neural network approximation, the searching area $\mathbf{V}$ is chosen as a ball around
136 the initialization $\boldsymbol{\theta}^{\text{init}} = [\mathbf{W}_1^{\text{init}}, \ldots, \mathbf{W}_L^{\text{init}}]$ for the ease of analysis. The detailed expression of $\mathbf{V}$ is
137 set as

$$
\mathbf{V} := \{\boldsymbol{\theta} = [\mathbf{W}_1, \ldots, \mathbf{W}_L] \big| \|\mathbf{W}_l - \mathbf{W}_l^{\text{init}}\|_F \leq \omega\},
\tag{5}
$$

138 where $1 \leq l \leq L$. It is assumed that $\boldsymbol{\theta}^{\text{init}}$ is chosen randomly from the Gaussian distribution, and
139 $\mathbf{W}_l^{\text{init}}$ is drawn from $\mathcal{N}(0, 1/m)$ with $l = 1, \ldots, m$.

140 The collection of all local linearization of $f(\boldsymbol{\theta}; \phi(s, a))$ at the initial point $\boldsymbol{\theta}_0$ is defined as

$$
\mathcal{F}_{\mathbf{V}, m} := \{f(\boldsymbol{\theta}^{\text{init}}; \phi(s, a)) + \langle \nabla_{\boldsymbol{\theta}} f(\boldsymbol{\theta}^{\text{init}}; \phi(s, a)), \boldsymbol{\theta} - \boldsymbol{\theta}^{\text{init}} \rangle : \boldsymbol{\theta} \in \mathbf{V}\}.
$$

141 If the function f is linear, i.e.,

$$
f(\boldsymbol{\theta}; \phi(s, a)) = \langle \boldsymbol{\theta}, \phi(s, a) \rangle,
\tag{6}
$$

142 we can see that $\mathcal{F}_{\mathbf{V}, m} = \{\langle \boldsymbol{\theta}, \phi(s, a) \rangle : \boldsymbol{\theta} \in \mathbf{V}\}$. The approximate stationary point of neural TD
143 associated with $\mathcal{F}_{\mathbf{V}, m}$ is defined as follows.

Definition 1 ([7]) *A point $\boldsymbol{\theta}^* \in \mathbf{V}$ is said to be the approximate stationary point if*

$$
\langle \mathbf{h}(\boldsymbol{\theta}^*), \boldsymbol{\theta} - \boldsymbol{\theta}^* \rangle \geq 0, \ \ \forall \boldsymbol{\theta} \in \mathbf{V},
\tag{7}
$$

where

$$
\mathbf{h}(\boldsymbol{\theta}) := \mathbb{E}\left[\widehat{\Delta}(\boldsymbol{\theta}) \nabla_{\boldsymbol{\theta}} \widehat{f}(\boldsymbol{\theta}; \phi(s, a))\right],
$$

145 *and the temporal difference error $\hat{\Delta}$ is defined as*

$$
\hat{\Delta}(\boldsymbol{\theta}) = \hat{\Delta}(s, a, s', a'; \boldsymbol{\theta}) := \hat{f}(\boldsymbol{\theta}; \phi(s, a)) - r(s, a) - \gamma \hat{f}(\boldsymbol{\theta}; \phi(s', a')),
\tag{8}
$$

146 *and $\hat{f}(\boldsymbol{\theta}; \phi(s, a)) \in \mathcal{F}_{\mathbf{V}, m}$.*

147 The convergence of neural TD is dependent on $\boldsymbol{\theta}^*$, and so is our result. In [7], it has been proved
148 that such a definition is well-defined; the approximate stationary point exists, minimizing the mean
149 squared projected Bellman error (MSPBE). Such a fact is more straightforward if the function f is
150 linear: the approximate stationary point of TD is identical to the unique solution to the projected
151 Bellman equation [45].

2.3 Neural Adaptive Temporal Difference Learning

153 This paper considers the neural adaptive TD. In the kth iteration of neural adaptive TD, we sample
154 $\mathbf{g}^k = \overline{\mathbf{g}}(\boldsymbol{\theta}^k; s_k, a_k, s_{k+1}, a_{k+1})$ in (2) with stochastic semi-gradient (4). The set $\mathbf{V}$ is chosen as (5)
155 with $\boldsymbol{\theta}^{\text{init}}$ also being sampled from the Gaussian distribution.

2.4 Some Useful Properties

157 For neural adaptive TD, several bounds and properties are presented in the following lemma (c.f.
158 Lemma 6.1,[48]; Lemma 4.4, [8])

Algorithm 1 Neural Adaptive Temporal Difference

Require: Parameters: $\eta, \gamma, \beta, L, m, \omega$
    Initialization: $\mathbf{g}^0 = \mathbf{0}, \mathbf{m}^0 = \mathbf{0}, v^0 = 0, \mathbf{W}_l^{\text{init}} \sim \mathcal{N}(0, 1/m)$ with $l = 1, \ldots, m$, $\mathbf{V}$ is set as (5)
    For $k = 1, 2, \ldots$
    1.sample the trajectory $s_0, a_0, s_1, a_1, \ldots$ from π
    2.calculate $\mathbf{g}^k = \overline{\mathbf{g}}(\boldsymbol{\theta}^k; s_k, a_k, s_{k+1}, a_{k+1})$ as (4)
    3.update the parameter $\boldsymbol{\theta}^k$ as (2)
    End for

Lemma 1 *Assume that $(\boldsymbol{\theta}^k)_{1 \leq k \leq K} \in \boldsymbol{B}(\boldsymbol{\theta}^*, \omega)$, given any $0 < \delta < 1$ and $K \in \mathbb{Z}^+$, m, ω and L satisfy*

$$m \geq C \max\{dL^2 \log \frac{m}{\delta}, \omega^{-4/3} L^{-8/3} \log \frac{m}{\omega\delta}\},$$
$$\frac{C_1 d^{3/2}}{L m^{3/4}} \leq \omega \leq \frac{C_2}{L^6 (\log m)^3}, \tag{9}$$

then it holds that

$$\|\nabla_{\boldsymbol{\theta}} f(\boldsymbol{\theta}; \phi(s, a))\| \leq C_3 \sqrt{m}, \forall s \in \mathcal{S}, a \in \mathcal{A}, \tag{10}$$

and

$$|f(\boldsymbol{\theta}^k; \mathbf{x})| \leq C_4 \sqrt{\log(K/\delta)}, 1 \leq k \leq K. \tag{11}$$

Denote that $\mathbf{h}^k := \widehat{\Delta}(\boldsymbol{\theta}^k) \nabla_{\boldsymbol{\theta}} \widehat{f}(\boldsymbol{\theta}^k)$ ($\widehat{\Delta}(\boldsymbol{\theta}^k)$ is given by (8)), it follows that

$$\|\mathbf{g}^k - \mathbf{h}^k\| \leq C_3(2 + \gamma)\omega^{1/3} L^3 \sqrt{m \log m \log(K/\delta)}$$
$$+ C_4 \omega^{4/3} L^4 \sqrt{m \log m} + C_5 \omega^2 L^4 m, \tag{12}$$

where $\mathbf{g}^k$ is the stochastic semi-gradient in neural adaptive TD, and

$$\|\mathbf{g}^k\| \leq (2 + \gamma) C_7 \sqrt{m \log(K/\delta)} \tag{13}$$

holds with probability at least $1 - 2\delta - 3L^2 \exp(-C_6 m \omega^{2/3} L)$ over the randomness of the initial point, where $\{C_i > 0\}_{i=1,\ldots,7}$ and $C > 0$ are universal constants.

Lemma 1 has been proved by [Lemma 6.1, [48]] and [Lemma 4.4, [8]]. The bounds (10) and (11) are only dependent on the structure of the neural networks. From Lemma 1, with high probabilities, $v^k \leq (2 + \gamma)^2 C_7^2 m \log(K/\delta) k$ when $k \leq K$. But this is the worst-case bound and may not always be achievable. From (4), we can see that $\mathbf{g}^k$ keeps the sparsity of $\nabla_\theta f(\boldsymbol{\theta}^k; \phi(s_k, a_k))$ because

$$[f(\theta^k; \phi(s_k, a_k)) - r(s_k, a_k) - \gamma f(\theta^k; \phi(s_{k+1}, a_{k+1}))] \in \mathbb{R}.$$

Thus, the stochastic semi-gradient $\mathbf{g}^k$ can be very sparse, based on which the following bound is proposed.

Condition 1 *In neural adaptive temporal difference learning, the following bound holds*

$$v^k \leq C_0 [m \log(K/\delta) k]^\alpha, \quad 0 < \alpha \leq 1, \quad 1 \leq k \leq K. \tag{14}$$

We can see that $\alpha = 1$ directly holds for (14) due to the bound (13); while $0 < \alpha < 1$ indicates the stochastic semi-gradients decay fast. In applications, the sparse gradients can always obey such a condition. We stress that such assumption is standard in the analysis of the Adam-type algorithms; see, e.g., references [29, 17, 36, 10, 32].

Now, we turn to the neural approximation case. As mentioned above, the convergence analysis uses the approximate stationary point $\boldsymbol{\theta}^*$. One core part of our proofs is investigating $\mathbf{h}$ versus the iteration. Due to the Markovian observation, $s_k, a_k, s_{k+1}, a_{k+1}$ may miss visiting several states in a single sampling, i.e., choosing some states and actions with probability 0. That indicates $\mathbb{E}[\mathbf{h}^k]$ may be

a biased expectation of $\mathbf{h}$. To get controlled difference bound, we consider $\mathbb{E}(\overline{\mathbf{h}}(\boldsymbol{\theta}^{k-T}) \mid \chi^k)$ and $\mathbf{h}(\boldsymbol{\theta}^{k-T})$, where

$$\overline{\mathbf{h}}(\boldsymbol{\theta}^{k-T}) := \hat{\Delta}(s_k, a_k, s_{k+1}, a_{k+1}; \boldsymbol{\theta}^{k-T}) \nabla_{\boldsymbol{\theta}} \widehat{f}(\boldsymbol{\theta}^{k-T}; \phi(s_k, a_k)).$$

Although with biased samples, the Markov property means $\mathcal{P}(s_k = s | s_{k-T} = s', a_{k-T} = a')$ is sufficiently close to $\mu(s)$ when T is large. Such a technique then yields following lemma.

Lemma 2 *Assume* $(\boldsymbol{\theta}^k)_{k \geq 0}$ *is generated by neural adaptive TD, and condition* (9) *holds. Given an integer* $T \in \mathbb{Z}^+$, *we have*

$$\left\| \mathbb{E}(\overline{\mathbf{h}}(\boldsymbol{\theta}^{k-T}) \mid \chi^k) - \mathbf{h}(\boldsymbol{\theta}^{k-T}) \right\| \leq (2 + \gamma) C_3 C_4 \bar{\kappa} \sqrt{m \log \frac{K}{\delta}} \rho^T$$

with probability at least $1 - 2\delta - 3L^2 \exp(-C_6 m \omega^{2/3} L)$ *over the randomness of the initial point.*

Proof of Lemma 2: Given a fixed integer T, direct calculations give us

$$
\begin{aligned}
&\mathbb{E}(\overline{\mathbf{h}}(\boldsymbol{\theta}^{k-T}; s_k, a_k, s_{k+1}, a_{k+1}) \mid \sigma^{k-T}) \\
&= \sum_{s,s' \in \mathcal{S}, a, a' \in \mathcal{A}} \mathcal{P}(s_k = s \mid s_{k-T}, a_{k-T}) \mathcal{P}(a, s', a'|s) \\
&\qquad \times \nabla_{\boldsymbol{\theta}} \widehat{f}\left(\boldsymbol{\theta}^{k-T}; \phi(s, a)\right) \widehat{\Delta}\left(\boldsymbol{\theta}^{k-T}; s, a, s', a'\right) \\
&= \sum_{s,s' \in \mathcal{S}, a, a' \in \mathcal{A}} \mu(s) \mathcal{P}(a, s', a'|s) \nabla_{\boldsymbol{\theta}} \widehat{f}\left(\boldsymbol{\theta}^{k-T}; \phi(s, a)\right) \times \widehat{\Delta}\left(\boldsymbol{\theta}^{k-T}; s, a, s', a'\right) \\
&\quad + \sum_{s,s' \in \mathcal{S}, a, a' \in \mathcal{A}} \mathcal{P}(a, s', a'|s)(\mathcal{P}(s_k = s \mid s_{k-T}, a_{k-T}) - \mu(s)) \nabla_{\boldsymbol{\theta}} \widehat{f}\left(\boldsymbol{\theta}^{k-T}; \phi(s, a)\right) \\
&\qquad \times \widehat{\Delta}\left(\boldsymbol{\theta}^{k-T}; s, a, s', a'\right).
\end{aligned}
\tag{15}
$$

Notice that the following expectation

$$\sum_{s,s' \in \mathcal{S}, a, a' \in \mathcal{A}} \mu(s)) \mathcal{P}(a, s', a'|s) \nabla_{\boldsymbol{\theta}} \widehat{f}\left(\boldsymbol{\theta}^{k-T}; \phi(s, a)\right) \widehat{\Delta}\left(\boldsymbol{\theta}^{k-T}; s, a, s', a'\right) = \mathbf{h}(\boldsymbol{\theta}^{k-T}).$$

The Markovian property tells us $\sum_{s \in \mathcal{S}} |\mathcal{P}(s_k = s \mid s_{k-T}, a_{k-T}) - \mu(s)| \leq \bar{\kappa} \rho^T$. Due to that $\widehat{f} \in \mathcal{F}_{\mathbf{V}, m}$, $\nabla_{\boldsymbol{\theta}} \widehat{f}\left(\boldsymbol{\theta}^{k-T}; \phi(s, a)\right) = \nabla_{\boldsymbol{\theta}} f\left(\boldsymbol{\theta}^{\text{init}}; \phi(s, a)\right)$. With Lemma 1, $\|\nabla_{\boldsymbol{\theta}} \widehat{f}\left(\boldsymbol{\theta}^{k-T}; \phi(s, a)\right)\| \leq C_3 \sqrt{m}$ and

$$
\begin{aligned}
|\widehat{\Delta}\left(\boldsymbol{\theta}^{k-T}; s, a, s', a'\right)| &= \left| \widehat{f}(\boldsymbol{\theta}^{k-T}; \phi(s, a)) - r(s, s') - \gamma \widehat{f}(\boldsymbol{\theta}^{k-T}; \phi(s', a')) \right| \\
&\leq (2 + \gamma) C_4 \sqrt{\log \frac{K}{\delta}},
\end{aligned}
$$

with probability at least $1 - 2\delta - 3L^2 \exp(-C_6 m \omega^{2/3} L)$. $\qquad \square$

The factor ρ^T in the upper bound in Lemma 2 comes from the Markovian noise. The lemma tells that when T is large enough, the difference between $\mathbb{E}\overline{\mathbf{h}}(\boldsymbol{\theta}^{k-T})$ and $\mathbf{h}(\boldsymbol{\theta}^{k-T})$ will be very small.

The lower bound directly holds based on the definitions above.

Lemma 3 *For* $\forall \boldsymbol{\theta} \in \mathbf{V}$, *it follows*

$$\langle \mathbf{h}(\boldsymbol{\theta}), \boldsymbol{\theta} - \boldsymbol{\theta}^* \rangle \geq (1 - \gamma) \mathbb{E}(\widehat{f}(\boldsymbol{\theta}; s, a) - \widehat{f}(\boldsymbol{\theta}^*; s, a))^2. \tag{16}$$

3 Main Results

This section contains the finite-time analysis of adaptive TD using deep neural network approximations. The descriptions of convergence results are dependent on the approximate stationary point $\boldsymbol{\theta}^*$.

Figure 1: The roadmap of the proof.

3.1 Roadmap

We first provide a high level explanation of the proofs and how the technical lemmas work in the proofs. Our proofs begin with a traditional idea of algorithmic convergence analysis, i.e., bounding the gap $(\|\boldsymbol{\theta}^* - \boldsymbol{\theta}^k\|^2)_{k\geq 0}$ versus the iteration. It is easy to have the following estimate

$$2\eta\langle\mathbf{m}^k, \boldsymbol{\theta}^* - \boldsymbol{\theta}^k\rangle/(v^k)^{\frac{1}{2}} \leq \|\boldsymbol{\theta}^* - \boldsymbol{\theta}^k\|^2 - \|\boldsymbol{\theta}^* - \boldsymbol{\theta}^{k+1}\|^2 + \eta^2\|\mathbf{m}^k\|^2/v^k.$$

Summing this inequality from $k = 1$ to K and taking expectation, we are led to

$$2\eta\sum_{k=1}^{K}\mathbb{E}(\langle\mathbf{m}^k, \boldsymbol{\theta}^* - \boldsymbol{\theta}^k\rangle/(v^k)^{\frac{1}{2}}) \leq \mathbb{E}\|\boldsymbol{\theta}^* - \boldsymbol{\theta}^1\|^2 + \eta^2\sum_{k=1}^{K}\mathbb{E}(\|\mathbf{m}^k\|^2/v^k). \tag{17}$$

Notice that (17) is very similar to the recursion of the well-known SGD. Recalling the finite-time convergence analysis of SGD, we then get the big picture for the proof of our algorithms, which contains two major steps and can be presented by Figure 1.

- **Step 1**. The first step is quite straightforward, i.e., bounding the summation $\sum_{k=1}^{K}\mathbb{E}(\|\mathbf{m}^k\|^2/v^k)$ with high probability. This step is proved by Lemma 4 in the proofs part.

- **Step 2**. The second step is to establish a lower bound of $\sum_{k=1}^{K}\mathbb{E}(\langle\mathbf{m}^k, \boldsymbol{\theta}^* - \boldsymbol{\theta}^k\rangle/(v^k)^{\frac{1}{2}})$. However, it is much more complicated than the first step. To this end, we need to exploit the relation between $\mathbb{E}\langle\mathbf{m}^k, \boldsymbol{\theta}^* - \boldsymbol{\theta}^k\rangle/(v^k)^{\frac{1}{2}}$ and $\mathbb{E}\langle\mathbf{h}(\boldsymbol{\theta}^k), \boldsymbol{\theta}^* - \boldsymbol{\theta}^k\rangle/(v^{k-1})^{\frac{1}{2}}$ (proved by Lemma 6 in the supplementary materials). Nevertheless, we cannot directly get Lemma 6; we have to develop a technical lemma that connects $\mathbb{E}\langle\boldsymbol{\theta}^k - \boldsymbol{\theta}^*, \mathbf{h}(\boldsymbol{\theta}^k)\rangle/(v^{k-1})^{\frac{1}{2}}$ with $\mathbb{E}\langle\boldsymbol{\theta}^* - \boldsymbol{\theta}^k, \mathbf{g}^k\rangle/(v^{k-1})^{\frac{1}{2}}$ (proved by Lemma 5 in the supplementary materials). Further with the connection between $\mathbb{E}\langle\mathbf{h}(\boldsymbol{\theta}^k), \boldsymbol{\theta}^k - \boldsymbol{\theta}^*\rangle$ and $\mathbb{E}\big(\hat{f}(\boldsymbol{\theta}^k; s, a) - \hat{f}(\boldsymbol{\theta}^*; s, a)\big)^2$ (Lemma 3), we then prove the lower bound.

3.2 Finite-Time Analysis of Neural Adaptive TD

Now, we are prepared to present the convergence of neural adaptive TD.

Theorem 1 *Suppose* $(\boldsymbol{\theta}^k)_{k\geq 0}$ *is generated by neural adaptive TD under the Markovian observation, and condition* (9) *holds, and Assumptions 1, 2, and condition* (14) *hold. Given the integer* $T \in \mathbb{Z}^+$, $\eta > 0, v_0 \geq \varpi > 0, 0 \leq \beta < 1$, *for* $K \geq 2^{\frac{2}{2-\alpha}}T$, *we have*

$$\min_{1\leq k\leq K}\mathbb{E}_\pi\big(\hat{f}(\boldsymbol{\theta}^k; s, a) - \hat{f}(\boldsymbol{\theta}^*; s, a)\big)^2 \leq \frac{c_1(m, \eta, \alpha, T)}{K^{1-\alpha/2}} + \frac{c_2(m, \eta, \omega, \alpha, T, K)}{K^{1-\alpha/2}}$$
$$+ c_3(m, \omega, \alpha, K)\rho^T K^{\frac{\alpha}{2}} + c_4(m, \omega, K) \tag{18}$$

with probability at least $1 - 2\delta - 3L^2\exp(-C_6 m\omega^{2/3}L)$ *over the randomness of the initial point, where*

$$\begin{aligned}
c_1(m, \eta, \alpha, T) &= \widetilde{\mathcal{O}}([m^{\frac{\alpha}{2}}T^2\eta^2 + m^{\frac{\alpha}{2}}\eta]\log K),\\
c_2(m, \eta, \omega, \alpha, T, K) &= \mathcal{O}(\omega m^{\frac{1+\alpha}{2}}\log K + m^{\frac{1+\alpha}{2}}\omega^2 T + \omega\sqrt{m}/(K-T)),\\
c_3(m, \omega, \alpha, K) &= \mathcal{O}([\omega m^{\frac{\alpha+1}{2}}]\log K),\\
c_4(m, \omega, K) &= \widetilde{\mathcal{O}}([\omega^{\frac{4}{3}}\sqrt{m} + \omega^{\frac{7}{3}}\sqrt{m} + \omega^3 m]\log K),
\end{aligned} \tag{19}$$

and their details are given by (30) *in supplementary materials.*

If we set the radius as $\omega = \Theta(m^{-1/2})$ [3], and $T = \frac{\ln(mK)}{\ln \frac{1}{\rho}}$, and $\eta = m^{-1/2}$, with (19), the right-hand side of (18) is in the order of

$$\widetilde{\mathcal{O}}\Big(1/(K^{1-\alpha/2}\ln^2\frac{1}{\rho}) + 1/(K^{1-\alpha/2}\ln\frac{1}{\rho}) + m^{-\frac{1}{6}}\Big).$$

Thus, with high probability to achieve the ϵ-accuracy for $\min_{1\le k\le K}\mathbb{E}_\pi\big(\hat{f}(\boldsymbol{\theta}^k; s, a) - \hat{f}(\boldsymbol{\theta}^*; s, a)\big)^2$, we need that

$$m = \tilde{\Theta}(1/\epsilon^6), \; 1/(K^{1-\alpha/2}\ln\frac{1}{\rho}) = \widetilde{\mathcal{O}}(\epsilon).$$

Thus, we get the following worst iteration complexity of the K as

$$\widetilde{\mathcal{O}}\left(\frac{1}{\epsilon^{\frac{2}{2-\alpha}}\cdot\ln^{\frac{2}{2-\alpha}}\left(\frac{1}{\rho}\right)}\right). \tag{20}$$

We can see that when $0 \le \alpha < 1$, the worst iteration complexity of the K is smaller than $\widetilde{\mathcal{O}}\left(\frac{1}{\epsilon^2\ln^2(\frac{1}{\rho})}\right)$ when $\epsilon \ll 1/[\ln(\frac{1}{\rho})]$, which indicates a *faster speed than the neural TD*. Due to that ϵ is the desired error and thus can be very small, the acceleration always happens for a small α. Note that $\alpha = 1$ directly holds with any extra assumption. In this case, the iteration complexity of the adaptive neural TD matches existing neural TD [48] in the DNN approximation case.

It is worth mentioning that compared with the convergence rate of neural TD with ReLU DNN, we used the same searching radius and network width as presented by [48]. In applications, the sparse stochastic semi-gradients always yield the fast decaying condition. In other words, with sparse stochastic semi-gradients, the neural adaptive TD also accelerates the vanilla TD even in the ReLU DNN approximation case. Such a phenomenon resonates with the existing acceleration results of adaptive methods for stochastic optimization: "for sparse data, the adaptive methods are likely to perform better than non-adaptive methods" [17].

The scheme of neural TD uses a projection to set $\mathbf{V}$ such that all iterates are constrained in a small neighborhood around the initialization (also called neural tangent kernel (NTK) regime). Such a procedure is to use the property of DNN ReLU [7, 48]. Thus, the searching radius is small, which is reasonable because in DNN training, the parameters usually only change in a very small range in the overparameterized regime.

Proposition 1 *Assume that conditions of Theorem 1 hold, and* $0 \le \alpha < 1, \omega = \Theta(m^{-1/2}), \eta = m^{-\frac{1}{2}}, m = \tilde{\Theta}(1/\epsilon^6), \epsilon \ll 1/[\ln(\frac{1}{\rho})]$, *the neural adaptive TD enjoys a faster finite-time convergence rate than the neural TD in the ReLU DNN approximation of function values with probability at least* $1 - 2\delta - 3L^2\exp(-C_6 m\omega^{2/3}L)$ *over the randomness of the initial point.*

4 Conclusions

This paper studied the finite-time convergence analyses of TD learning with adaptive learning rates and momentum approximated by deep ReLU neural networks using the Markovian samplings. Our established theoretical results show that the neural adaptive temporal difference learning is convergent when the neural network is sufficiently wide. Our work shows convergence results and establishes the theoretical advantages of the adaptive algorithms for neural networks approximation cases, i.e., adaptive schemes achieve better rates than neural TD when the stochastic semi-gradients decay fast. There are numerous avenues for future works, including 1) Can we extend the multiple ReLU active functions to other ones, e.g., sigmoid function or more general functions? 2) Can we establish the finite-time convergence of adaptive TD with other kinds of neural network approximation, e.g., recurrent neural networks and graph neural networks. 3) Can we get a relaxed bound for the radius of the searching area?

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[3]Such a choice of the radius in neural TD is also used in [48].

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