Multi-step learning and Value-based approximation methods

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Markov Decision Process

MDP is defined as $\mathcal{M} = (\mathcal{X}, \mathcal{A}, P, g)$, where

- \mathcal{X} is a finite state space, $\mathcal{A} = \bigcup_{x \in \mathcal{X}} \mathcal{A}(x)$ is a finite action space.
- P is a state transition probability kernel.

The system state at the next decision epoch is determined by

$$\mathbb{P}\left\{x_{k+1} = y \mid x_k = x, a_k = a\right\} = P(x, a, y)$$

for each $x_k \in \mathcal{X}$, $a_k \in A(x_k) \subset \mathcal{A}$.

Given (x_k; a_k), a new (random) state x_{k+1} is observed and a (one-step) cost g(x_k; a_k; x_{k+1}) is incurred.

• The value function of policy
$$\mu : \mathcal{X} \to \mathcal{A}$$
 is

$$J_{\mu}(x) = \mathbb{E}\Big[\sum_{k=0}^{\infty} \beta^{k} g\big(x_{k}, \mu(x_{k}), x_{k+1}\big) | x_{0} = x\Big], \text{ where } 0 < \beta < 1.$$

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Bellman operator

• For $J : \mathcal{X} \to \mathbb{R}$ and stationary policy μ

$$(T_{\mu}J)(x) = \mathbb{E}_{y \sim P(\cdot|x,\mu(x))} \Big[g(x,\mu(x),y) + \beta J(j) \Big]$$
$$= \sum_{y \in \mathcal{X}} P(x,\mu(x),y) \Big[g(x,\mu(x),y) + \beta J(y) \Big], \ x \in \mathcal{X}$$

• For $J: \mathcal{X} \to \mathbb{R}$ consider

$$(TJ)(x) = \min_{a \in \mathcal{A}(x)} \mathbb{E}_{y \sim P(\cdot|x,a)} \left[g(x,a,y) + \beta J(y) \right]$$
$$= \min_{a \in \mathcal{A}(x)} \sum_{y \in \mathcal{X}} P(x,a,y) \left[g(x,a,y) + \beta J(y) \right], \ x \in \mathcal{X}$$

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Bellman equation

Theorem (Bertsekas, Proposition 1.2.4)

(a) For every stationary policy μ , the associated value function satisfies for all $i \in \mathcal{X}$:

$$J_{\mu}(x) = \mathbb{E}_{y \sim P(\cdot \mid x, \mu(x))} \left[g(x, \mu(x), y) + \beta J_{\mu}(y) \right] \quad \text{or} \quad J_{\mu} = T_{\mu} J_{\mu}$$

(b) J_{μ} is the unique solution of equation (1).

The main purpose of this talk is to find an efficient way to estimate J_{μ} .

Policy evaluation: Dynamic programming (P_{μ} , g_{μ} are known.)

A fixed point problem $J_{\mu} = T_{\mu}J_{\mu}$ is equivalent to linear system of equations

$$J_{\mu} = g_{\mu} + \beta P_{\mu} J_{\mu},$$

where g_{μ} is a \mathcal{X} -vector with entries $g_{\mu}(x) = \sum_{y \in \mathcal{X}} P(x, \mu(x), y) g(x, \mu(x), y)$ and P_{μ} is an $\mathcal{X} \times \mathcal{X}$ matrix with entries $P_{\mu}(x, y) = P(x, \mu(x), y)$.

Monte-Carlo Simulation: P_{μ} , g_{μ} are unknown

- Let $c_m(x_0) = g(x_0, x_1) + \beta g(x_1, x_2) \dots + \beta^N g(x_N, x_{N+1})$ be the cumulative cost of the *m*th episode, s.t. x_{N+1} is a terminal state or $\frac{\beta^N}{1-\beta} \max_{x,y \in \mathcal{X}} g(x,y) \approx 0$.
- For all states $x \in \mathcal{X}$ and for all m we have

$$J_{\mu}(x) = \mathbb{E}[c_m(x)]$$

thus we can estimate $J_{\mu}(x)$ forming a sample mean:

$$J^{K}(x) \approx \frac{1}{K} \sum_{m=1}^{K} c_{m}(x) \tag{1}$$

Equation (1) can be iteratively calculated

$$J^{m}(x) = J^{m-1}(x) + \gamma_{m} \Big(c_{m}(x) - J^{m-1}(x) \Big), \quad m = 1, .., K$$

where $J^0(x) = 0$ and $\gamma_m = \frac{1}{m}$.

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TD(0) learning

Bellman equation

$$J_{\mu}(x) = \mathbb{E}\Big[g(x,y) + \beta J_{\mu}(y)\Big]$$

• Assume that we have an episode {*x*₀, *x*₁, ..., *x_N*} and *k_m* is a *m*th time when the state *x* is visited.

$$J_{\mu}(x) \approx \frac{1}{K} \sum_{m=1}^{K} \left[g(x_{k_m}, x_{k_m+1}) + \beta J_{\mu}(x_{k_m+1}) \right]$$

Update step for TD(0) learning

$$\begin{cases} J^{k}(x) = J^{k-1}(x_{k}) + \gamma_{m} \Big(g(x_{k}, x_{k+1}) + \beta J^{k-1}(x_{k+1}) - J^{k-1}(x_{k}) \Big), & \text{if } x = x_{k} \\ J^{k}(x) = J^{k-1}(x_{k}), & \text{if } x \neq x_{k} \end{cases}$$

N-step TD learning

N-step Bellman operator:

$$(T^{N}_{\mu}J)(x_{k}) = \mathbb{E}\Big[\sum_{m=0}^{N} \beta^{m}g(x_{k+m}, x_{k+m+1}) + \beta^{N+1}J(x_{k+N+1})\Big]$$

N-step Bellman equation:

$$(T^N_\mu J)(x) = J(x), \quad \text{for each } x \in \mathcal{X}$$

• After k + N steps

 $J_{\mu}^{k+N}(x_k) \approx g_k + \beta g_{k+1} + \dots + \beta^N g_{k+N} + \beta^{N+1} J_{\mu}^{k+N-1}(x_{k+N+1})$

Update step for N-step TD learning

$$\begin{cases} J_{\mu}^{k+N+1}(x_k) = J_{\mu}^{k+N}(x_k) - \gamma_k \Big[J_{\mu}^{k+N}(x_k) - \sum_{n=0}^N \beta^n g_{k+n} - \beta^N J_{\mu}^{k+N}(x_{k+N+1}) \Big] \\ J_{\mu}^{k+N+1}(y) = J_{\mu}^{k+N}(y), \quad y \neq x_k \end{cases}$$

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λ -weighted multistep Bellman equation

• Consider $N \sim \text{Geometric}(\lambda), 0 \leq \lambda < 1$.

Define an operator:

$$T_{\mu}^{(\lambda)} = (1 - \lambda) \sum_{n=0}^{\infty} \lambda^n T_{\mu}^{n+1}$$

The λ-weighted multistep Bellman equation:

$$J_{\mu}(x_{k}) = (1-\lambda)\mathbb{E}\Big[\sum_{n=0}^{\infty} \lambda^{n} \Big(\sum_{m=0}^{n} \beta^{m} g(x_{k+m}, x_{k+m+1}) + \beta^{n+1} J_{\mu}(x_{k+n+1})\Big)\Big]$$

$$J_{\mu}(x_{k}) = (1-\lambda)\mathbb{E}\Big[\sum_{n=0}^{\infty} \lambda^{n} \Big(\sum_{m=0}^{n} \beta^{m} g(x_{k+m}, x_{k+m+1}) + \beta^{n+1} J_{\mu}(x_{k+n+1})\Big)\Big]$$

= $\mathbb{E}\Big[(1-\lambda)\sum_{m=0}^{\infty} \beta^{m} g(x_{k+m}, x_{k+m+1}) \sum_{n=m}^{\infty} \lambda^{n} + \sum_{n=0}^{\infty} \beta^{n+1} J_{\mu}(x_{k+n+1})(\lambda^{n} - \lambda^{n+1})\Big]$
= $\mathbb{E}\Big[\sum_{m=0}^{\infty} \beta^{m} \lambda^{m} \Big(g(x_{k+m}, x_{k+m+1}) + \beta\lambda J_{\mu}(x_{k+m+1}) - J_{\mu}(x_{k+m})\Big)\Big] + J_{\mu}(x_{k})$
= $\mathbb{E}\Big[\sum_{m=0}^{\infty} \lambda^{m} \beta^{m} d_{m+k}\Big] + J_{\mu}(x_{k})$

Temporal Difference: $d_m = g(x_m, x_{m+1}) + \beta \lambda J_\mu(x_{m+1}) - J_\mu(x_m)$

- λ -weighted Bellman equation: $J_{\mu}(x_k) = \mathbb{E}\left[J_{\mu}(x_k) + \sum_{m=0}^{\infty} (\lambda \beta)^m d_{m+k}\right]$
- Online iterations:

$$J(x_k) := J(x_k) + \gamma \Big[J(x_k) + \sum_{m=0}^{\infty} (\lambda \beta)^m d_{m+k} - J(x_k) \Big] = J(x_k) + \gamma \sum_{m=0}^{\infty} (\lambda \beta)^m d_{m+k}$$

- Assume we have a single infinitely long trajectory $(x_0, g_0, x_1, g_1, x_2, ...)$.
- Since we cannot afford to wait until the end of the trajectory we need an on-line version of the algorithm.

Let J^0 is an initial guess. The first two updated are:

• Following the transition (x_0, x_1) :

$$J^{1}(x_{0}) = J^{0}(x_{0}) + \gamma d_{0}$$

• Following the transition
$$(x_1, x_2)$$
:

$$\begin{cases}
J^2(x_0) = J^1(x_0) + \gamma \lambda \beta d_1 = J^0(x_0) + \gamma d_0 + \gamma \lambda \beta d_1 \\
J^2(x_1) = J^1(x_1) + \gamma d_1 = J^0(x_1) + \gamma d_1
\end{cases}$$
If $x_0 = x_1$ there are three variants of the TD(λ) algorithm:

- The restart variant: $J^2(x_0) = J^0(x_0) + \gamma d_0 + \gamma d_1$.
- The first-visit variant: $J^2(x_0) = J^0(x_0) + \gamma d_0 + \gamma \lambda \beta d_1$
- The every-visit variant: $J^2(x_0) = J^0(x_0) + \gamma d_0 + \gamma \lambda \beta d_1 + \gamma d_1$.

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The update equation for $TD(\lambda)$ becomes:

$$J^{k+1}(x) = J^k(x) + \gamma_k(x)z_k(x)d_k(x)$$
, for each $x \in \mathcal{X}$

where $z_{-1} = 0$ and

• The restart variant:
$$z_k(x) = \begin{cases} 1, & \text{if } x_k = x \\ \beta \lambda z_{k-1}(x), & \text{if } x_k \neq x \end{cases}$$

• The first-visit variant: $z_k(x) = \begin{cases} 1, & \text{if } x = x_k \text{ and } x_k \text{ is visited first time} \\ \beta \lambda z_{k-1}(x), & \text{otherwise} \end{cases}$
• The every-visit variant: $z_k(x) = \begin{cases} \beta \lambda z_{k-1}(x) + 1, & \text{if } x_k = x \\ \beta \lambda z_{k-1}(x), & \text{if } x_k \neq x \end{cases}$

Approximate Dynamic Programming

Approximate dynamic programming (neuro-dynamic programming, reinforcement learning):

A principle aim is to address problems with very large number of states in \mathcal{X} .

- $|\mathcal{X}|$ -dimensional inner product are time-consuming.
- It may impossible to store $|\mathcal{X}|$ -vector in a computer memory.

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Approximation is value space

 $\tilde{J}(x,r)$ is a function of some chosen form and $r = (r_1,...,r_s)$ is a parameter vector of relatively small dimension s.

Examples:

Linear form:

$$\tilde{J}(x,r) = \sum_{k=1}^{s} r_k \phi_k(x),$$

where $\phi_k : \mathcal{X} \to \mathbb{R}, \ k = 1, ..., s$ are known as feature functions.

• Feedforward neural network with a single hidden layer with K neurons:

$$\tilde{J}(x,r) = \sum_{k=1}^{K} r(k) \sigma \left(\sum_{l=1}^{L} r(k,l) v_l(x) \right),$$

where state x is encoded as a L -dimensional vector v(x), $\sigma(\cdot)$ is an activation function.

Linear Approximation

$$J_{\mu}(x) = \mathbb{E}\Big[\sum_{k=0}^{\infty} \beta^k g(x_k, \mu(x_k), x_{k+1}) | x_0 = x\Big]$$

• We approximate $J_{\mu}(x)$ with a linear architecture

$$\tilde{J}(x,r) = \sum_{k=1}^{s} r_k \phi_k(x) = \phi(x)^T r, \ x \in \mathcal{X},$$

where $\phi(x)$ is an *s*-dimensional feature vector.

$$\tilde{J}_r = \begin{pmatrix} \phi(x_1)^T r \\ \vdots \\ \phi(x_{|X|})^T r \end{pmatrix} = \Phi r,$$

where Φ is $|\mathcal{X}| \times s$ matrix that has as rows the feature vectors $\phi(x)^T$.

The Projected Equation

- We want to approximate J_{μ} within space $S = \{\Phi r \mid r \in \mathbb{R}^s\}$.
- The goal is to find \tilde{J}^* w.r.t. ξ -weighted norm:

$$\tilde{J}^* = \underset{\tilde{J} \in S}{\arg\min} \sum_{x \in \mathcal{X}} |J_{\mu}(x) - \tilde{J}(x, r)|^2 \xi(x) = \underset{\tilde{J} \in S}{\arg\min} ||J_{\mu} - \tilde{J}_r||_{\xi}^2$$

• The problem is equivalent to finding $r^* \in \mathbb{R}^s$ s.t.

$$r^* = \underset{r \in \mathbb{R}^s}{\operatorname{arg\,min}} ||J_{\mu} - \Phi r||_{\xi}^2$$

Let Π denote the projection operation onto S w.r.t the ξ-weighted norm.
 Then

$$\Pi J_{\mu} = \Phi r^*. \tag{2}$$

The Projected Bellman Equation

We assume

$$J_{\mu} \approx T_{\mu}(\Phi r^*) \tag{3}$$

Combing (3) and (2) we get

 $\Pi T_{\mu}(\Phi r^*) \approx \Phi r^*$

The Projected Bellman equation:

$$\Pi T_{\mu}(\Phi r) = \Phi r \tag{4}$$

One can show that ΠT_μ is a contraction operator w.r.t. || · ||_ξ norm when ξ is a stationary distribution of the DTMC with transition matrix P_μ.

The Projected Bellman Equation

By the definition of projection the unique solution of (4) satisfies

$$r^* = \underset{r \in \mathbb{R}^s}{\operatorname{arg\,min}} ||\Phi r - (g_\mu + \beta P_\mu \Phi r^*)||_{\xi}^2$$

By setting to 0 the gradient w.r.t. r we obtain

$$\Phi^T D(\Phi r^* - (g_\mu + \beta P_\mu \Phi r^*)) = 0,$$
(5)

where D is the diagonal matrix with ξ along the diagonal.

Equation (5) can be compactly written as

$$Cr^* = d_i$$

where $C = \Phi^T D (I - \beta P_\mu) \Phi$ and $d = \Phi^T D g_\mu$.

TD(0) with linear approximation

•
$$Cr - d = \Phi^T D(I - \beta P_\mu)r - \Phi^T Dg_\mu = \mathbb{E}\Big[\phi(x)\Big(\phi(x)^T r - \beta\phi(y)^T r - g_\mu(x,y)\Big)\Big]$$

• Equation Cr - d = 0 is equivalent to

$$r = r - \gamma (Cr - d) = r - \gamma \mathbb{E} \Big[\phi(x) \Big(\phi(x)^T r - \beta \phi(y)^T r - g(x, y) \Big) \Big]$$

• Given an episode $(x_0, g_0, x_1, g_1, ..., x_N, g_N)$, the TD(0) iteration is

$$r_{k+1} = r_k - \gamma_k \phi(x_k) \Big(\phi(x_k)^T r_k - \beta \phi(x_{k+1})^T r_k - g_k \Big)$$

Convergence of TD(0) with linear approximation

Theorem (Tsitsiklis, Van Roy, 1997)

Assume that

- the Markov chain assosiated with policy μ is irreducible and aperiodic
- the steady-state variance of transition costs is finite $\mathbb{E}[g^2(x_k, x_{k+1})] < \infty$.
- the learning rate is s.t.

$$\sum\limits_{k=0}^{\infty}\gamma_k=\infty$$
 and $\sum\limits_{k=0}^{\infty}\gamma_k^2<\infty$

Then

- TD(0) converges to r^* that is a unique solution of $\Pi T(\Phi r) = \Phi r$.
- r* satisfies

$$||\Phi r^* - J_{\mu}||_{\xi} \le \frac{1}{1-\beta} ||\Pi J_{\mu} - J_{\mu}||_{\xi}$$

Finite Sample Analyses for TD(0) with Function Approximation

- Gal Dalal et. al., 2017 found convergence rate under assumption that one can generate iid samples from a steady-state distribution, using recently developed stochastic approximation techniques.
- Bhandari, Russo, Singal, 2018 found convergence rate for projected TD(0) algorithm (*r_k* is assumed to be ||*r_k*|| < *R*) using information theoretic techniques.
- Lei Ying, 2018 found convergence rate for TD(0) algorithm using Stein's Method.

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LSTD(0)

•
$$C = \mathbb{E}\Big[\phi(x)\Big(\phi(x) - \beta\phi(y)\Big)^T\Big]$$
 and $d = \mathbb{E}[\phi(x)g(x,y)]$

Based on simulation

$$C_N = \frac{1}{N+1} \sum_{k=0}^N \phi(x_k) \Big(\phi(x_k) - \beta \phi(x_{k+1}) \Big)^T \text{ and } d_N = \frac{1}{N+1} \sum_{k=0}^N \phi(x_k) g(x_k, x_{k+1})$$

• LSTD(0) algorithm: simulate N time-steps according to a policy μ

$$\begin{cases} C_{k+1} = C_k - \frac{1}{k} \left(\phi(x_k) \left(\phi(x_k) - \beta \phi(x_{k+1}) \right)^T - C_k \right) \\ d_{k+1} = d_k - \frac{1}{k} \left(\phi(x_k) g(x_k, x_{k+1}) - d_k \right) \end{cases}$$

After the end of simulation:

$$\tilde{r} = C_N^{-1} d_N \quad J_\mu(x) \approx \phi^T(x) \tilde{r}$$

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LSPE(0)

• LSPE method is based on an idea of Projected Value Iteration

$$\Phi r_{k+1} = \Pi T_{\mu}(\Phi r_k) \tag{6}$$

- Equation (6) is equivalent to $r_{k+1} = r_k (\Phi^T D \Phi)^{-1} (Cr_k d)$
- a simulation-based implementation:

$$r_{k+1} = r_k - G_k (C_k r_k - d_k),$$

where

$$\begin{cases} C_{k+1} = C_k - \frac{1}{k} \Big(\phi(x_k) \Big(\phi(x_k) - \beta \phi(x_{k+1}) \Big)^T - C_k \Big) \\ d_{k+1} = d_k - \frac{1}{k} \Big(\phi(x_k) g(x_k, x_{k+1}) - d_k \Big) \\ G_{k+1} = \Big(\frac{1}{k+1} \sum_{t=0}^k \phi(x_t) \phi(x_t)^T \Big)^{-1} = G_k - \frac{G_k \phi(x_k) \phi^T(x_k) G_k}{1 + \phi^T(x_k) G_k \phi(x_k)} \end{cases}$$

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Limitations

Limitations:

- No guaranty that TD(λ) with non-linear approximation will converge.
 Counterexample in Tsitsiklis, Van Roy, 1997.

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