

Optimization Algorithms

Reinforcement Learning & Optimization

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• Reinforcement Learning is an optimization problem – how far can we get with standard optimization approaches rather than specialized RL methods?

• *The world:* An MDP $(S, A, P, R, P_0, \gamma)$ with state space S, action space A, transition probabilities $P(s_{t+1}\,|\, s_t, a_t)$, reward fct $R(s_t,a_t)$, initial state distribution $P_0(s_0)$, and discounting factor $\gamma \in [0, 1]$.

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- *The agent:* A policy $\pi(a_t|s_t)$.

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- *The agent:* A policy $\pi(a_t|s_t)$.
- Together they define the path distribution $(\xi = (s_{0:T+1}, a_{0:T}))$

T

 $t=0$

\n The path distribution
$$
(\xi = (s_{0:T+1}, a_{0:T}))
$$
\n $P_{\pi}(\xi) = P(s_0) \prod_{t=1}^{T} \pi(a_t|s_t) \cdot P(s_{t+1}|s_t, a_t)$ \n

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- *The agent:* A policy $\pi(a_t|s_t)$.
- Together they define the path distribution $(\xi = (s_0, t_{+1}, a_0, t))$

T

 $t=0$

\n The path distribution
$$
(\xi = (s_{0:T+1}, a_{0:T}))
$$
\n

\n\n The graph distribution $f(x) = P(s_0)$ is given by:\n $P_{\pi}(\xi) = P(s_0)$ \n

\n\n The graph of the graph with the following matrices:\n $P_{\pi}(\xi) = P(s_0)$ \n

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 \ddotsc

and the **expected total return**

$$
J(\pi) = \mathbb{E}_{\xi \sim P_{\pi}} \left\{ \underbrace{\sum_{t=0}^{\infty} \gamma^t R(s_t, a_t)}_{R(\xi)} \right\} = \int_{\xi} P_{\pi}(\xi) R(\xi) d\xi
$$

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- We assume the policy $\pi_\theta(a|s)$ is parameterized by some $\theta \in \mathbb{R}^n$
- The problem is

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- $-J(\theta)$ is just a function we want to optimize
- $-J(\theta)$ is a "weighted sum" over all paths (cf. additive cost function & SGD)
- We can't really compute/evaluate $f(\theta)$ exactly we can only get a sample $\xi \sim P_\theta$ and $R(\xi)$ in each iteration (cf. SGD case!)
- $-$ Different: $\sum_i f_i(x)\leftrightarrow$ fixed distribution over $i;$ $\int_\xi P_\theta(\xi)R(\xi)\leftrightarrow$ non-stationary distribution over ξ

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Can knowing about the MDP process simplify the optimization problem?

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- Yes, in at least 2 ways:
- Bellman optimality we understand sth. about the optimal policy beyond KKT
- Policy gradients we can derive gradients

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	- $-$ The value function $V^*(x)$ over state space fulfills

$$
V^*(s) = \max_{a} \left[R(s, a) + \gamma \mathbb{E}_{s'|s, a} \{ V^*(s') \} \right]
$$

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• The Bellman optimality condition truely exploits the MDP structure, and gives further conditions on the optimum beyond stationarity of $J(\theta)$.

- Learning (=optimizing) while collecting more and more data
	- Unusual from the optimization perspective \leftrightarrow instable "target" (objective, ξ -distribution)
	- Leads to breadth of RL-methodologies (model-based/model-free RL, TD-, Q-learning, etc)

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- But there are also trends to avoid this
	- "Offline RL", classical system identification, model-based RL
	- separating data collection issue from optimization issue

Stochastic Policy Gradient

• Recall

$$
J(\theta) = \int_{\xi} P_{\theta}(\xi) R(\xi) d\xi
$$

• We have

$$
\nabla_{\theta} J(\theta) = \nabla_{\theta} \int P_{\theta}(\xi) R(\xi) d\xi = \int P_{\theta}(\xi) \nabla_{\theta} \log P_{\theta}(\xi) R(\xi) d\xi
$$

\n
$$
= \mathbb{E}_{\xi|\theta} \{ \nabla_{\theta} \log P_{\theta}(\xi) R(\xi) \} = \mathbb{E}_{\xi|\theta} \{ \sum_{t=0}^{H} \nabla_{\theta} \log \pi(a_t|s_t) R(\xi) \}
$$

\n
$$
= \mathbb{E}_{\xi|\theta} \{ \sum_{t=0}^{H} \nabla_{\theta} \log \pi(a_t|s_t) \gamma^t \underbrace{\sum_{t'=t}^{H} \gamma^{t'-t} r_{t'}}_{\hat{Q}_{\xi,t}} \}
$$

Deterministic Policy Gradient

- However, in practise, policies are often not stochastic. Esp. neural networks. We have $a=\pi_\theta(s)\in\mathbb{R}^d,$ parameterized by $\theta.$ What is the correct gradient then?
- As introduced in reference [2]:

$$
\nabla_{\theta} J(\theta) = \mathbb{E}_{s \sim P_{\theta}} \left\{ \nabla_{\theta} \pi_{\theta}(s) \nabla_{a} Q^{\pi_{\theta}}(s, a) \big|_{a = \pi_{\theta}(s)} \right\}
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(NOTE: unusual convention about Jacobians... I'd write it $\partial_{\alpha}Q^{\pi_{\theta}}(s, a)\partial_{\theta}\pi(s)$) Silver et al: *Deterministic policy gradient algorithms*. 2014

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• So we in principle also have a gradient! But very noisy! Better: D4PG

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	- Bellman optimality, gradients
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- **However,** reasons to ignore structure of underlying MPD:
	- Avoid implied problems, e.g. by function approximation, value estimation, policy iteration
	- very noisy gradient estimates
	- Robustness to mis-assumptions
- → black-box or **derivative-free optimization**

References

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Evolution Strategies as a Scalable Alternative to Reinforcement Learning

Tim Salimans

Jonathan Ho.

Szymon Sidor

Ilva Sutskever

Abstract

Xi Chen

OpenAI

We explore the use of Evolution Strategies (ES), a class of black box optimization algorithms, as an alternative to popular MDP-based RL techniques such as Qlearning and Policy Gradients. Experiments on MuJoCo and Atari show that ES is a viable solution strategy that scales extremely well with the number of CPUs available: By using a novel communication strategy based on common random numbers, our ES implementation only needs to communicate scalars, making it possible to scale to over a thousand parallel workers. This allows us to solve 3D humanoid walking in 10 minutes and obtain competitive results on most Atari games after one hour of training. In addition, we highlight several advantages of ES as a black box optimization technique: it is invariant to action frequency and delayed rewards, tolerant of extremely long horizons, and does not need temporal discounting or value function approximation.

• The ES they employ:

Algorithm 1 Evolution Strategies

- 1: **Input:** Learning rate α , noise standard deviation σ , initial policy parameters θ_0
- 2: for $t = 0, 1, 2, \ldots$ do
- Sample $\epsilon_1, \ldots \epsilon_n \sim \mathcal{N}(0, I)$ $3:$
- Compute returns $F_i = F(\theta_t + \sigma \epsilon_i)$ for $i = 1, ..., n$ $4[·]$
- Set $\theta_{t+1} \leftarrow \theta_t + \alpha \frac{1}{n} \sum_{i=1}^n F_i \epsilon_i$ $5:$
- $6:$ end for
- Is an instance of our "General Stochastic Search" scheme:
	- θ is the mean; $\lambda = n$ samples $x_t \sim \mathcal{N}(\theta, \sigma^2 I)$; evaluations $f(x_i)$
- The update is more like a stochastic gradient step rather than selection In expectation, $F_i \epsilon_i \dot{=} (F_\theta + \nabla F^\top \sigma \epsilon_i) \epsilon_i \approx 0 + \sigma \nabla F$.

• Ratio of ES timesteps to TRPO timesteps needed to reach various percentages of TRPO's learning progress at 5 million timesteps:

Deep Neuroevolution: Genetic Algorithms are a Competitive Alternative for **Training Deep Neural Networks for Reinforcement Learning**

Felipe Petroski Such Vashisht Madhavan Edoardo Conti Joel Lehman Kenneth O. Stanley Jeff Clune Uber AI Labs {felipe.such, jeffclune}@uber.com

(Do you spend your time training nets, or simulating?)

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Table 1. On Atari a simple genetic algorithm is competitive with Q-learning (DQN), policy gradients (A3C), and evolution strategies (ES). Shown are game scores (higher is better). Comparing performance between algorithms is inherently challenging (see main text), but we attempt to facilitate comparisons by showing estimates for the amount of computation (operations, the sum of forward and backward neural network passes), data efficiency (the number of game frames from training episodes), and how long in wall-clock time the algorithm takes to run. The ES, DQN, A3C, and GA (1B) perform best on 3, 3, 4, and 3 games, respectively. Surprisingly, random search often finds policies superior to those of DQN, A3C, and ES (see text for discussion). Note the dramatic differences in the speeds of the algorithm, which are much faster for the GA and ES, and data efficiency, which favors DQN. The scores for DQN are from Hessel et al. (2017) while those for A3C and ES are from Salimans et al. (2017). For A3C, DQN, and ES, we cannot provide error bars because they were not reported in the original literature; GA and random search error bars are visualized in (SI Fig. 3). The wall-clock times are approximate because they depend on a variety of hard-to-control-for factors. We found the GA runs slightly faster than ES on average. The \dagger symbol indicates state of the art performance. GA 6B scores are bolded if best, but do not prevent bolding in other columns.

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Table 4. Head-to-head comparison between algorithms on the 13 Atari games. Each value represents how many games for which the algorithm listed at the top of a column produces a higher score than the algorithm listed to the left of that row (e.g. GA 6B beats DON on 7 games).

• Conclusion: It varies from problem to problem what is better. And it is suprising that "naive" black-box ES can beat elaborate RL-methods

- Overall, it is not so clear at all whether RL is actually better than black-box/derivative-free optimization
- For small problems where we have little function approximation or noisy gradient problems, yes; but for large scale problems?
- For discussion:
	- If you have a problem with dense rewards and smooth dynamics...?

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	- If you have a sparse reward problem with hard dynamics...?
- RL-methods rely on reward signals/gradients that can be "propagated" through time/steps (credit assignment, Q-learning, Bellman)
- Black-box search is ignorant to this, sometimes to the better

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