

Distributed RL

Joash Lee

Pan Liangming

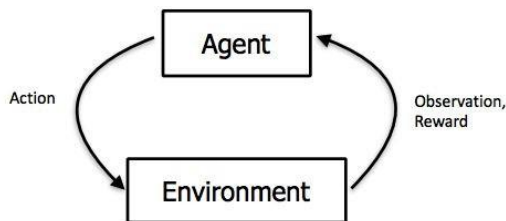
Vicky Feliren

Lesson Objectives

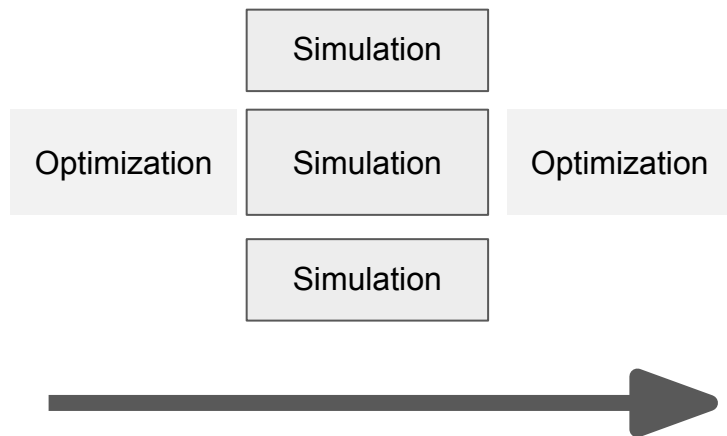
1. Why parallelise?
2. Understand how the computation of standard RL algorithms can be distributed to decrease wall-clock training time.
3. How these distributed RL algorithms can be modularised.
4. How modularised distributed RL algorithms can be implemented on real systems - case study: RLlib
5. Examples on using RLlib

Common Computational Patterns for RL

Original

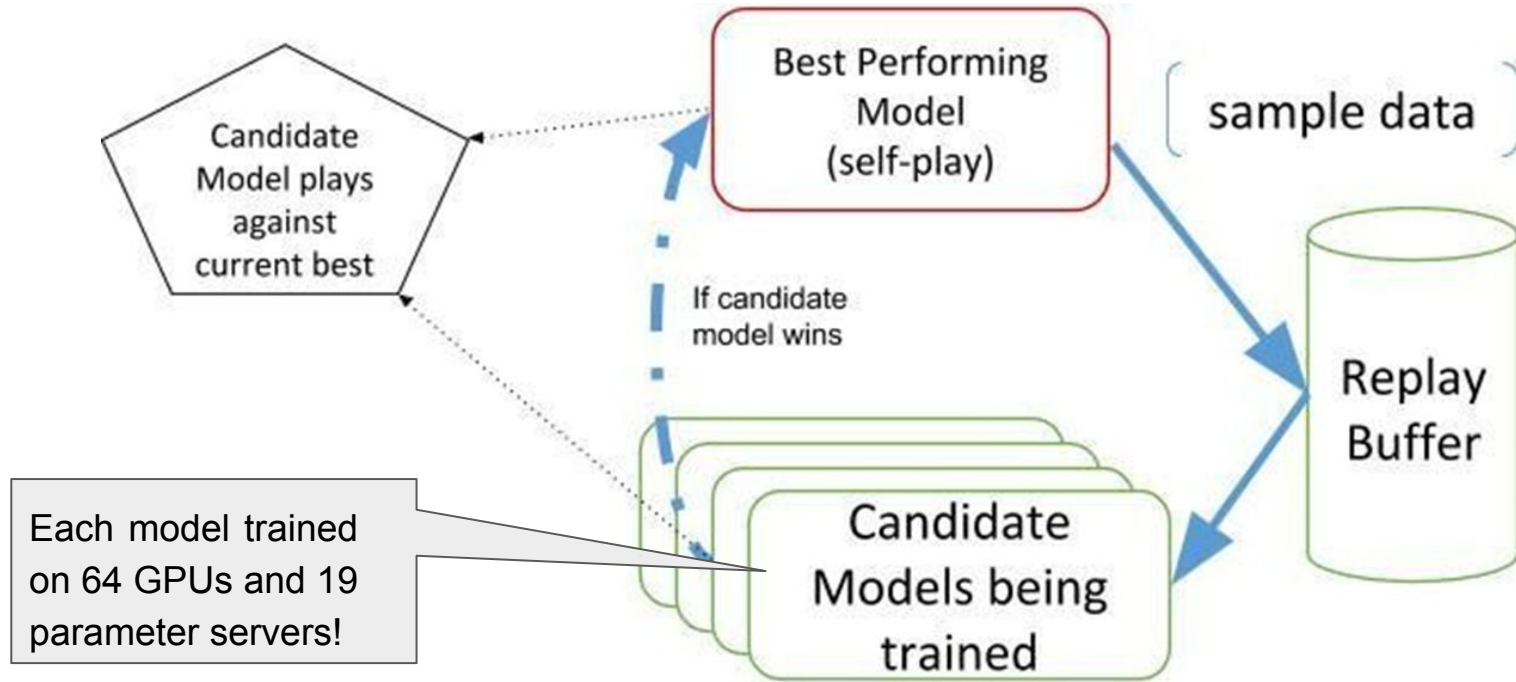


Batch Optimization



How can we **better utilize** our computational resources **to accelerate** RL progress?

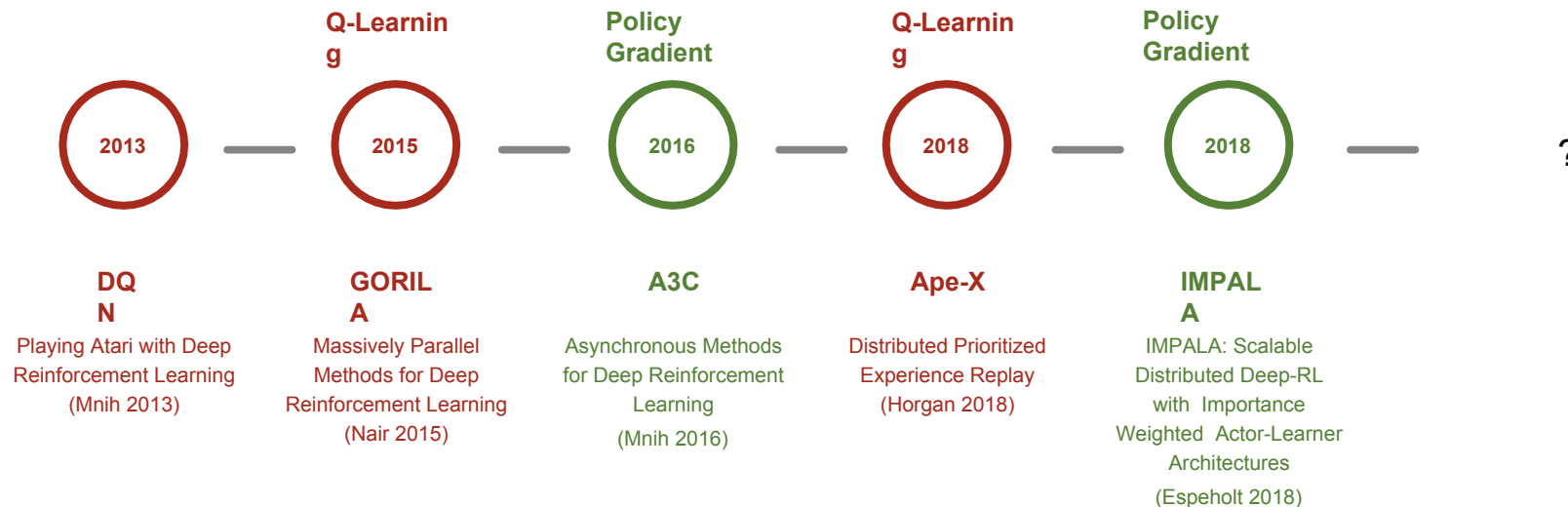
AlphaZero



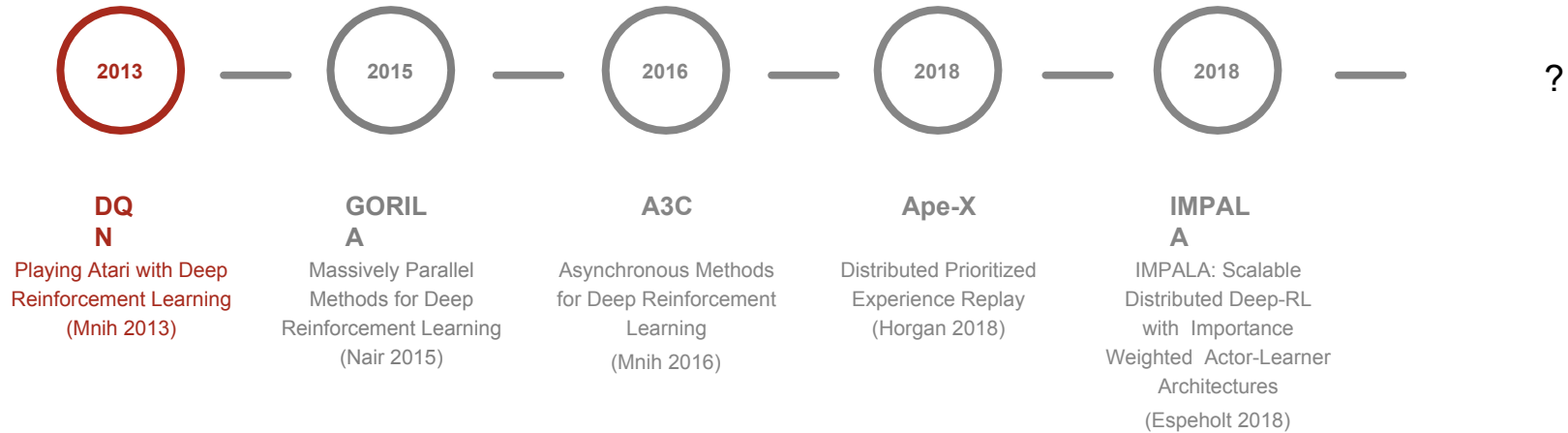
Lesson Objectives

1. Why parallelise?
2. **Understand how the computation of standard RL algorithms can be distributed to decrease wall-clock training time.**
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4. How modularised distributed RL algorithms can be implemented on real systems - case study: RLlib
5. Examples on using RLlib

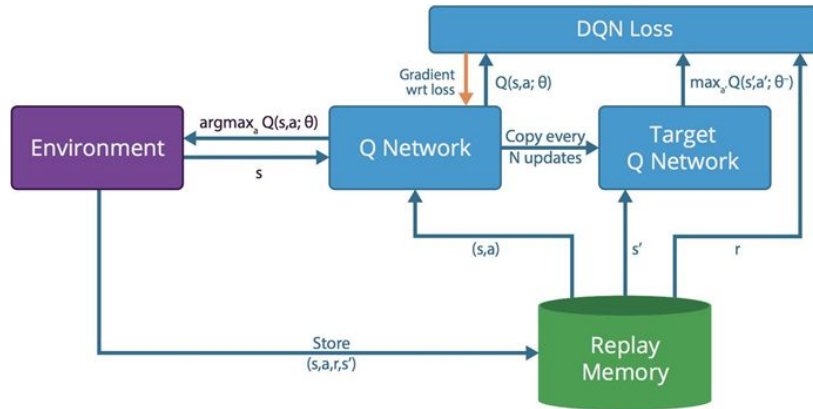
History of large scale distributed RL



History of large scale distributed RL

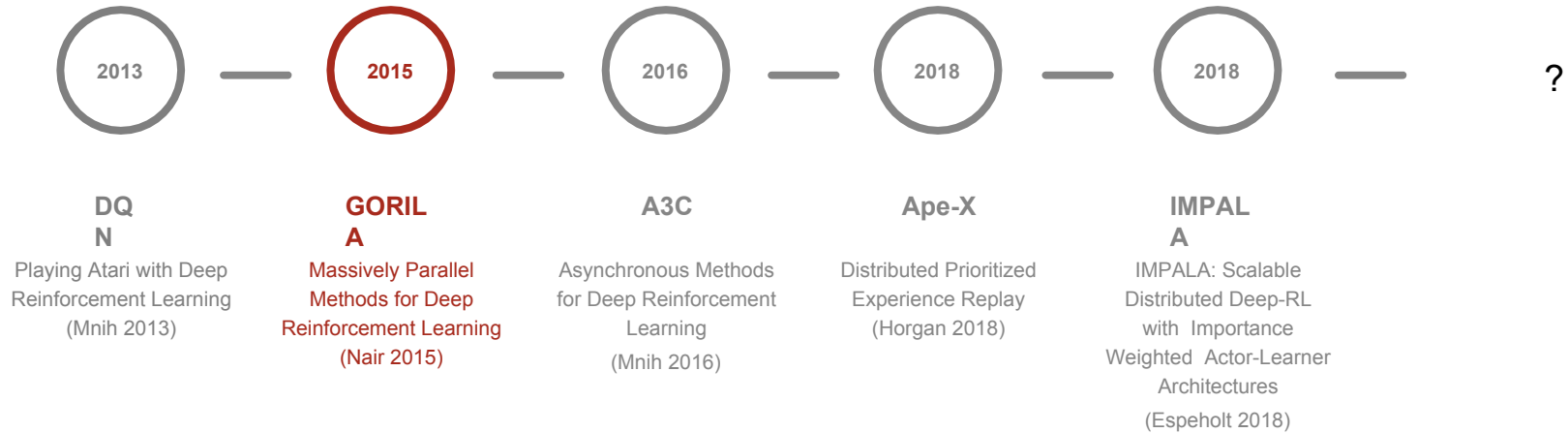


2013/2015: DQN

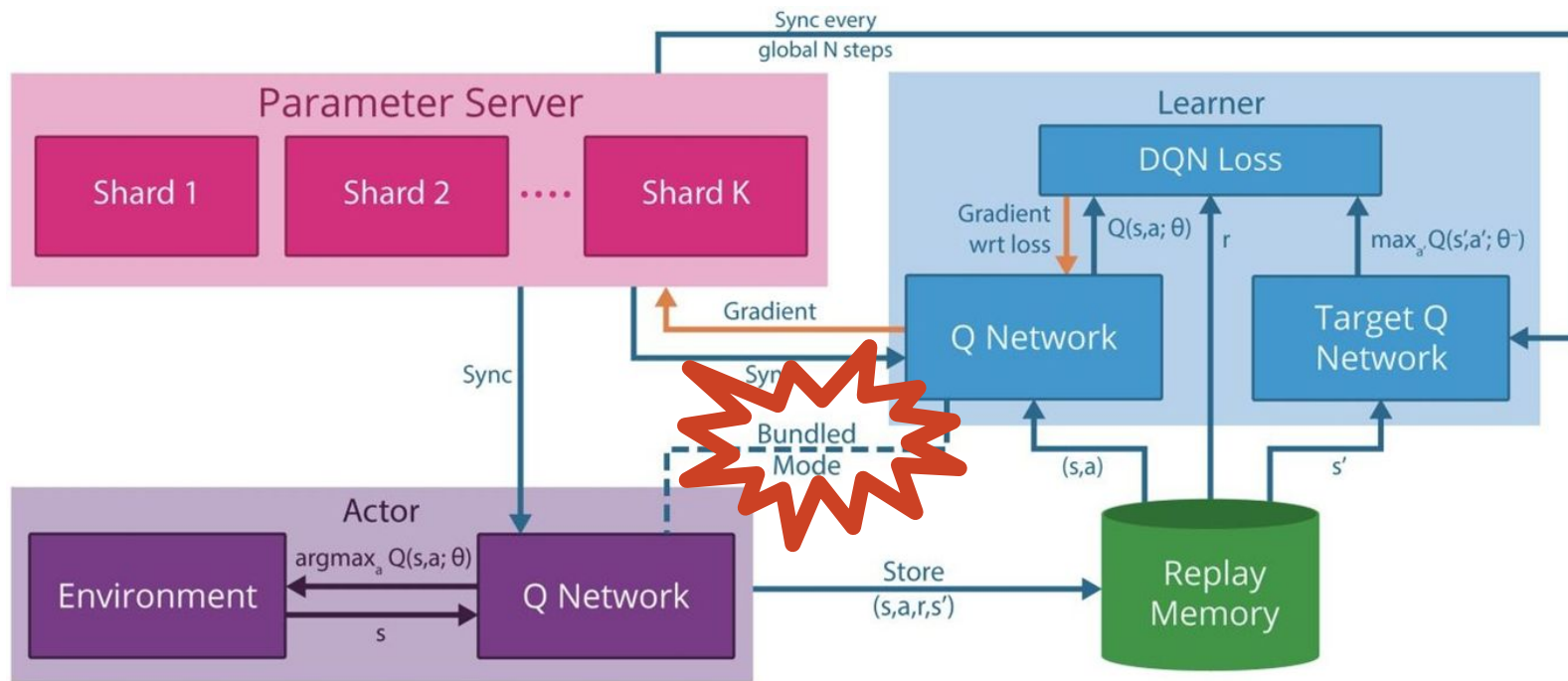


- 1. $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$ $N \times$
Store (s_i, a_i, s'_i, r_i) in \mathbf{B}
 2. Sample batch (s_j, a_j, s'_j, r_j) from \mathbf{B} $K \times$
Update Q network
 3. Update target network parameters: $\phi' \leftarrow \phi$

History of large scale distributed RL



2015: General Reinforcement Learning Architecture (GORILA)



Nair, A., Srinivasan, P., Blackwell, S., Allicek, C., Fearon, R., De Maria, A., . . . Petersen, S. (2015). Massively parallel methods for deep reinforcement learning. *arXiv preprint arXiv:1507.04296*.

2015: General Reinforcement Learning Architecture (GORILA)

-

Standard DQN

1. $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$
Store (s_i, a_i, s'_i, r_i) in \mathbf{B}
2. Sample batch (s_j, a_j, s'_j, r_j) from \mathbf{B}
Update Q network
3. Update target network parameters: $\phi' \leftarrow \phi$

-

Distributed DQN

Actor 1. $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$
Store (s_i, a_i, s'_i, r_i) in \mathbf{B}

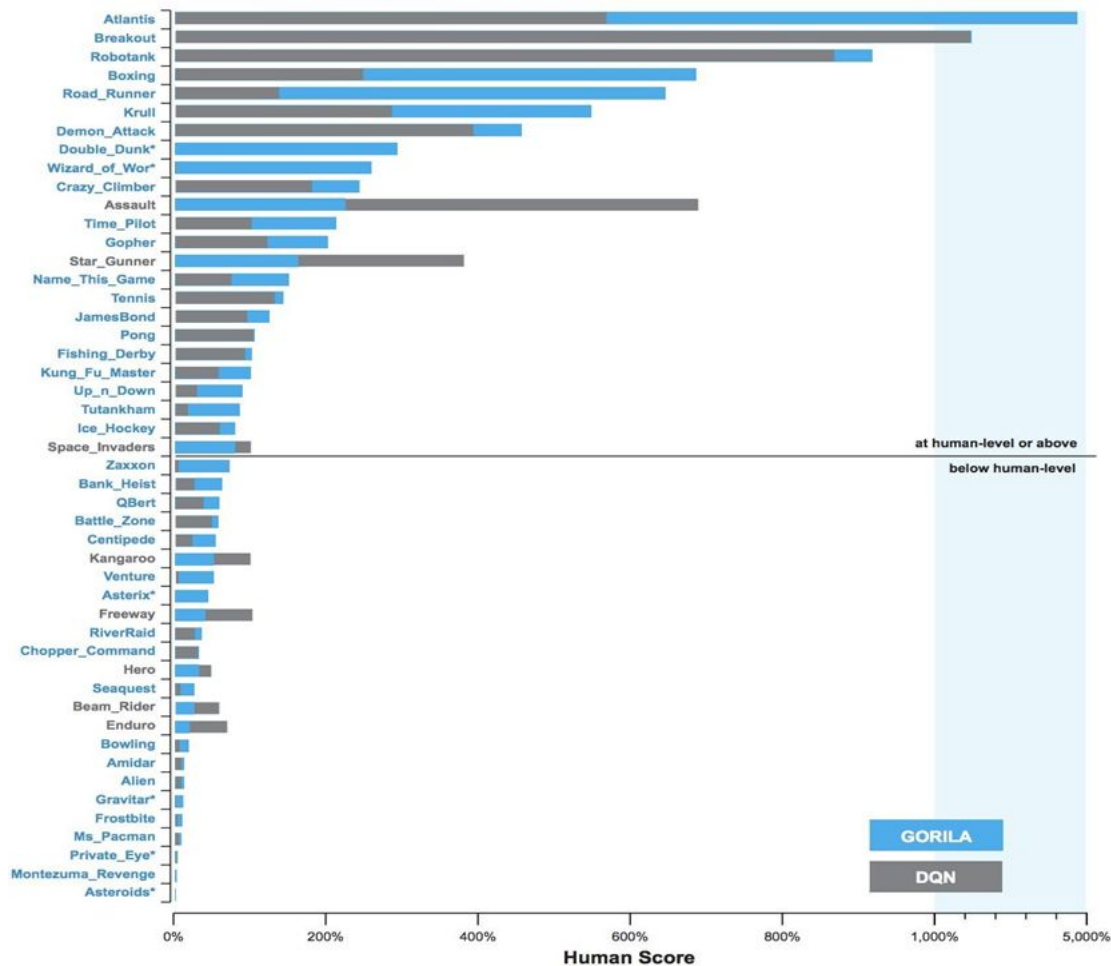
Learner 2. Sample batch (s_j, a_j, s'_j, r_j) from \mathbf{B}
Update θ with θ^+ from parameter server
Calculate gradients w.r.t. θ
Send gradients to parameter server

Parameter Server 3. Update Q network

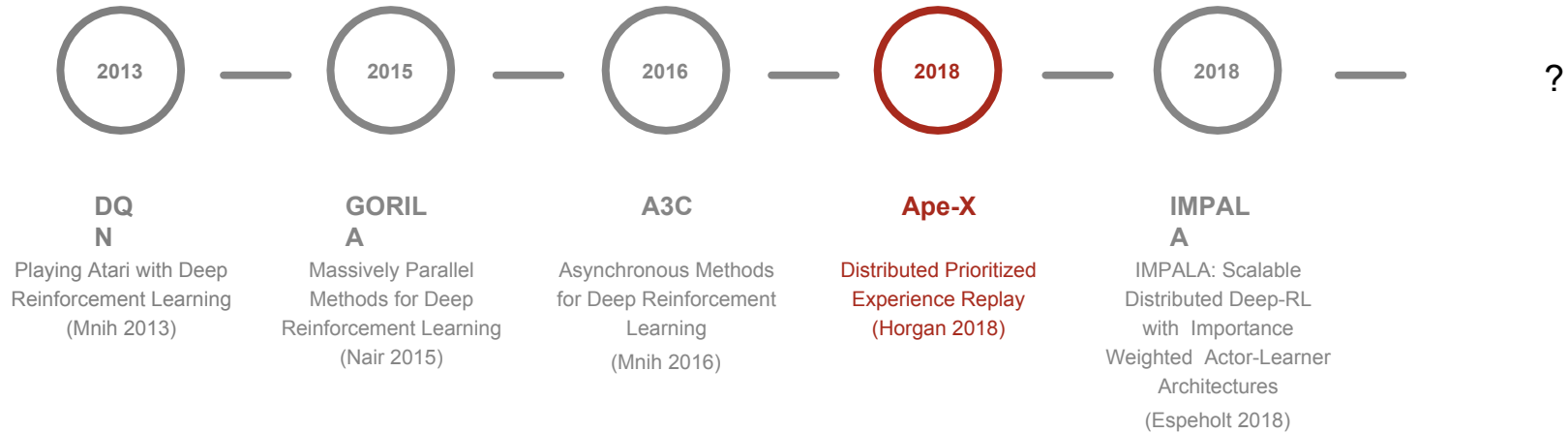
Server

Learner 4. Update target network parameters θ^-
with θ^+ from the parameter server every N steps

GORILA Performance



History of large scale distributed RL




Prioritised Experience Replay

Standard DQN

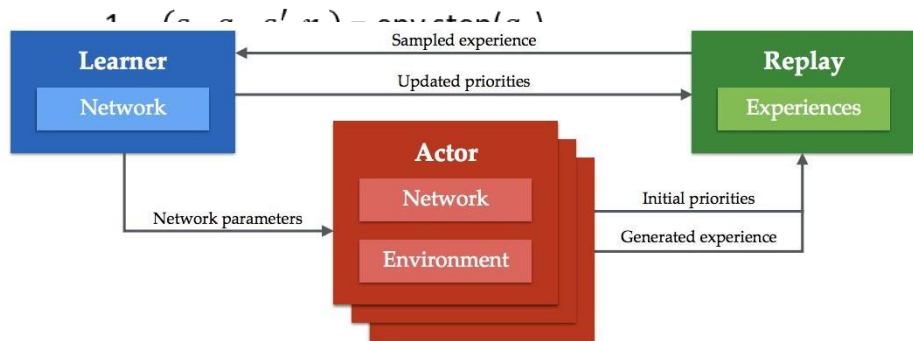
1. $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$
Store (s_i, a_i, s'_i, r_i) in \mathbf{B}
2. Sample K transitions (s_j, a_j, s'_j, r_j) uniformly
from \mathbf{B}
Update Q network
3. Update target network parameters: $\phi' \leftarrow \phi$

Prioritised Experience Replay

1. $(s_t, a_t, s'_t, r_t) = \text{env.step}(a_t)$
Store (s_t, a_t, s'_t, r_t) in \mathbf{B} with **max priority** $p_t = \max_{i < t} p_i$
2. Sample transition $j \sim P(j) = p_j^\alpha / \sum_i p_i^\alpha$ from \mathbf{B}
Compute TD error δ_j
Update transition probability $p_j \leftarrow |\delta_j|$  $K \times$
3. Update Q network
4. Update target network parameters: $\phi' \leftarrow \phi$

Distributed Prioritized Experience Replay (Ape-X)

Prioritised Experience Replay



3. Update Q network

4. Update target network parameters: $\phi' \leftarrow \phi$

Ape-X

Actors

- $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$
- Compute TD error δ_j
- Update transition probability $p_j \leftarrow |\delta_j|$
- Store (s_i, a_i, s'_i, r_i) in \mathbf{B}

Learners

- Update θ with θ^+ from parameter server
- Sample transition (s_j, a_j, s'_j, r_j) from \mathbf{B}
- Calculate gradients w.r.t. θ
- Update parameters θ of Q -network
- Compute TD error δ_j
- Update transition probability $p_j \leftarrow |\delta_j|$
- Update target network parameters θ^- with θ^+ from the parameter server every N steps

Distributed Prioritized Experience Replay (Ape-X)

Gorila

Actors

- $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$
Store (s_i, a_i, s'_i, r_i) in \mathbf{B}

Learners

- Update θ with θ^+ from parameter server
- Sample batch (s_j, a_j, s'_j, r_j) from \mathbf{B}
- Calculate gradients w.r.t. θ
- **Send gradients to parameter server**
- Update target network parameters θ^-
with θ^+ from the parameter server every N steps

Parameter Server

- Update parameters θ of Q network

Ape-X

Actors

- $(s_i, a_i, s'_i, r_i) = \text{env.step}(a_i)$
- **Compute TD error δ_j**
Update transition probability $p_j \leftarrow |\delta_j|$
- Store (s_i, a_i, s'_i, r_i) in \mathbf{B}

Learners

- Update θ with θ^+ from parameter server
- Sample transition (s_j, a_j, s'_j, r_j) from \mathbf{B}
- Calculate gradients w.r.t. θ
- **Update parameters θ of Q -network**
- **Compute TD error δ_j**
Update transition probability $p_j \leftarrow |\delta_j|$
- Update target network parameters θ^-
with θ^+ from the parameter server every N steps

Ape-X Performance

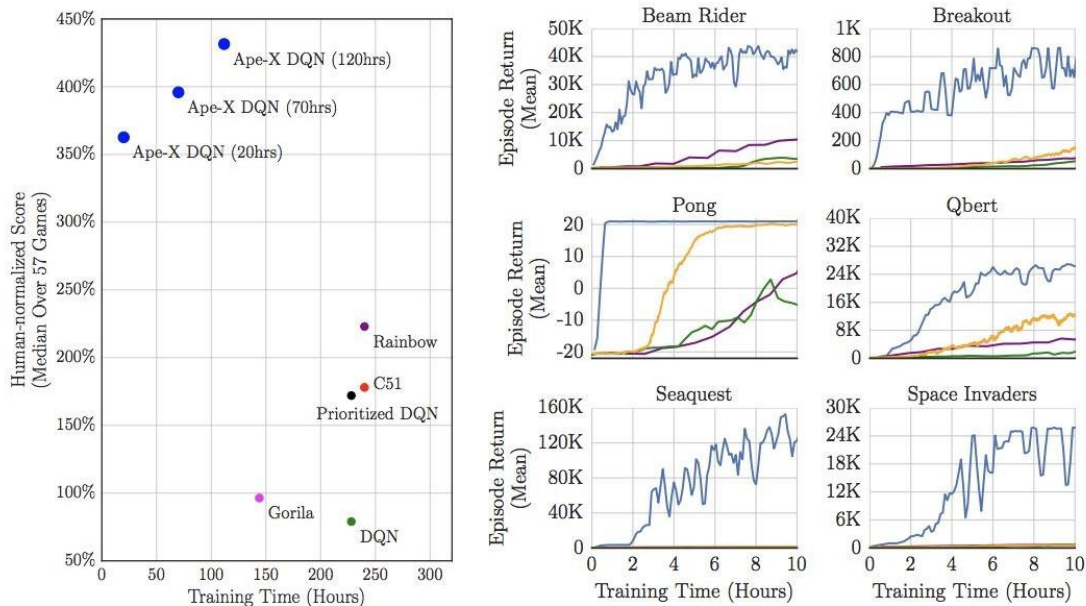
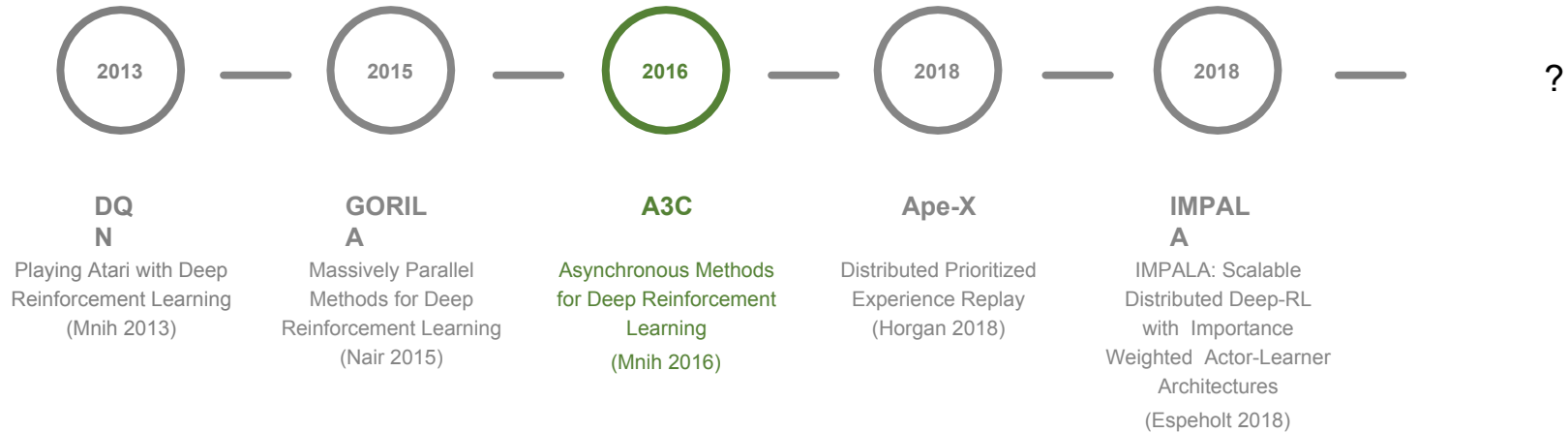


Figure 2: Left: Atari results aggregated across 57 games, evaluated from random no-op starts. Right: Atari training curves for selected games, against baselines. Blue: Ape-X DQN with 360 actors; Orange: A3C; Purple: Rainbow; Green: DQN. See appendix for longer runs over all games.

History of large scale distributed RL



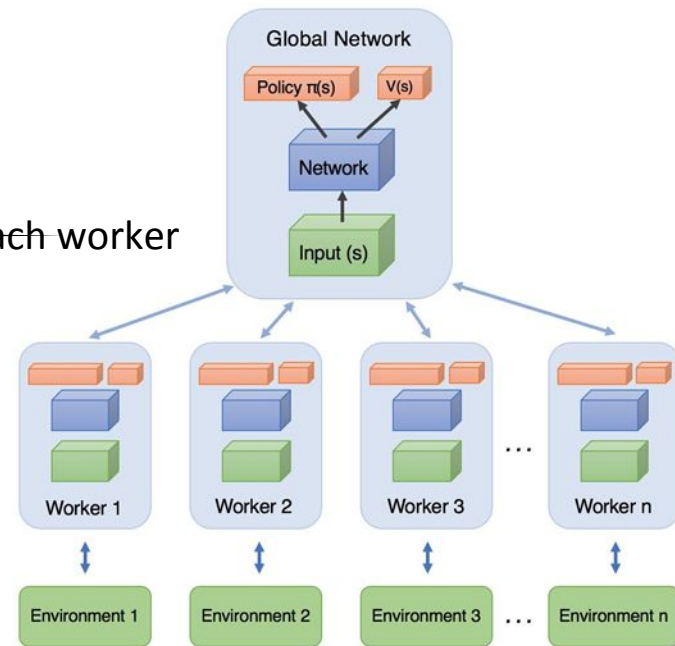
Recap: Online actor-critic

- 1. Take action $a \sim \pi_\theta(a|s)$, get (s, a, s', r)
 2. Update \hat{V}_ϕ^π using target $r + \hat{V}_\phi^\pi(s')$
 3. Evaluate $\hat{A}^\pi(s, a) = r(s, a) + \gamma \hat{V}_\phi^\pi(s') - \hat{V}_\phi^\pi(s)$
 4. $\nabla_\theta J(\theta) \approx \nabla_\theta \log \pi_\theta(a|s) \hat{A}^\pi(s, a)$
 5. Update policy
 $\theta \leftarrow \theta + \alpha \nabla_\theta J(\theta)$

Asynchronous advantage actor-critic (A3C)

1. Sync weights θ and ϕ from master
2. Take action $a \sim \pi_\theta(a|s)$, get (s, a, s', r)
3. Compute gradient of \hat{V}_ϕ^π using target $r + \hat{V}_\phi^\pi(s')$
4. Evaluate $\hat{A}^\pi(s, a) = r(s, a) + \gamma \hat{V}_\phi^\pi(s') - \hat{V}_\phi^\pi(s)$
5. $\nabla_\theta J(\theta) \approx \nabla_\theta \log \pi_\theta(a|s) \hat{A}^\pi(s, a)$

Each worker

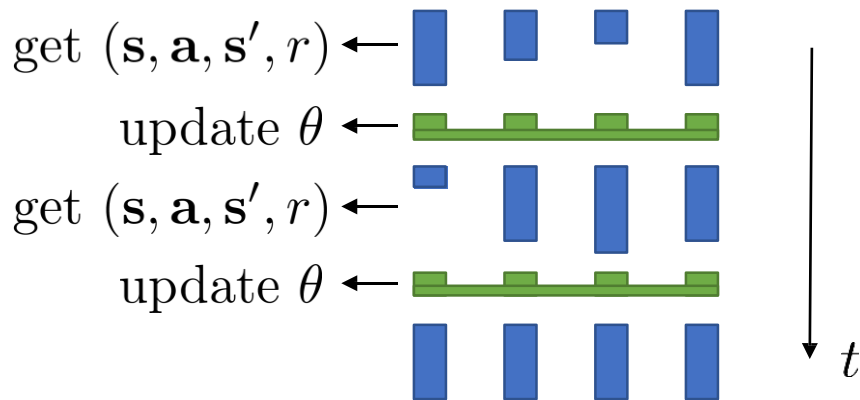


Each has different exploration -> more diverse samples!

Asynchronous advantage actor-critic (A3C)

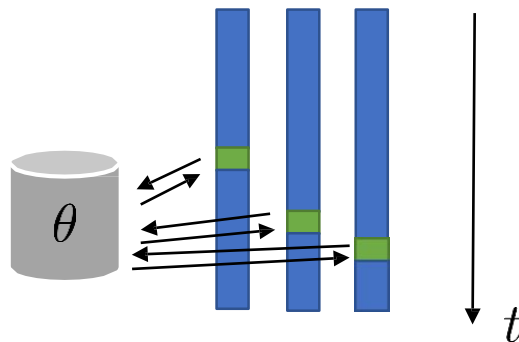
A2C

- can lead to low GPU utilisation due to rendering time variance within a batch



A3C

- decouples acting from learning



Asynchronous advantage actor-critic (A3C)

Some extra features:

- n-step estimation: $\hat{A}^\pi(s, a) = \sum_{i=0}^{k-1} \gamma^i r(s_t, a_t) + \gamma^k \hat{V}_\phi^\pi(s_{t+k}) - \hat{V}_\phi^\pi(s_t)$
- Entropy of the policy π_θ was added to the objective function to improve exploration:

$$\nabla_\theta J(\theta) \approx \nabla_\theta \log \pi_\theta(a|s) \hat{A}^\pi(s, a) + \beta \nabla_\theta H(\pi_\theta(s))$$

A3C Performance

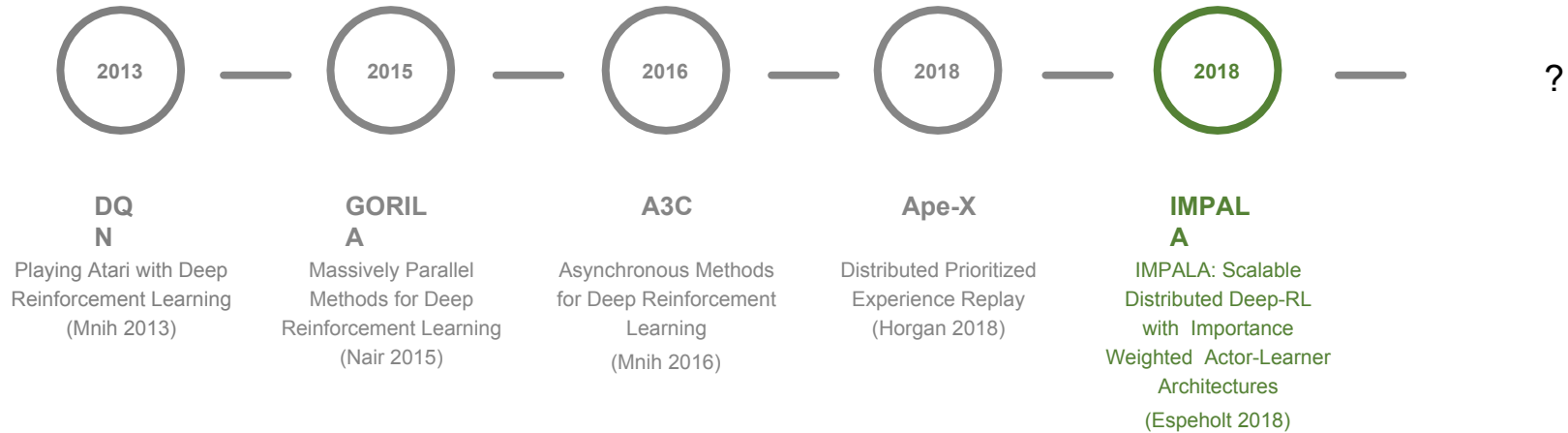
Changes to GORILA:

1. **Faster updates**
2. **No replay buffer**
3. **Actor-critic**

| Method | Training Time | Mean | Median |
|-----------------|----------------------|--------|--------|
| DQN | 8 days on GPU | 121.9% | 47.5% |
| Gorila | 4 days, 100 machines | 215.2% | 71.3% |
| D-DQN | 8 days on GPU | 332.9% | 110.9% |
| Dueling D-DQN | 8 days on GPU | 343.8% | 117.1% |
| Prioritized DQN | 8 days on GPU | 463.6% | 127.6% |
| A3C, FF | 1 day on CPU | 344.1% | 68.2% |
| A3C, FF | 4 days on CPU | 496.8% | 116.6% |
| A3C, LSTM | 4 days on CPU | 623.0% | 112.6% |

Table 1. Mean and median human-normalized scores on 57 Atari games using the human starts evaluation metric. Supplementary

History of large scale distributed RL



Importance Weighted Actor-Learner Architectures (IMPALA)

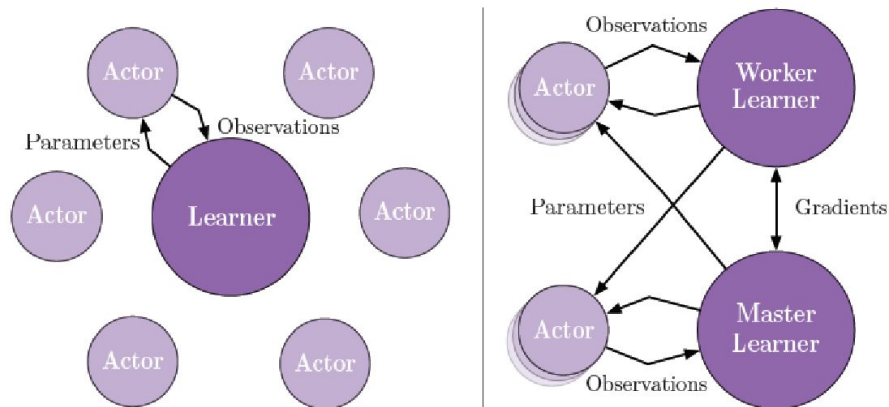
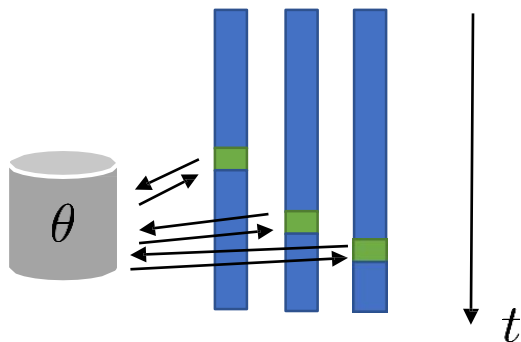


Figure 1. Left: Single Learner. Each *actor* generates trajectories and sends them via a queue to the *learner*. Before starting the next trajectory, *actor* retrieves the latest policy parameters from *learner*. **Right: Multiple Synchronous Learners.** Policy parameters are distributed across multiple *learners* that work synchronously.

How to correct for Policy Lag? Importance Sampling!

Shortcoming of A3C:

- *Policy-lag*



Apply importance sampling:

1. to policy gradient

$$\mathbb{E}_{a_s \sim \mu(\cdot|x_s)} \left[\frac{\pi_{\bar{\rho}}(a_s|x_s)}{\mu(a_s|x_s)} \nabla \log \pi_{\bar{\rho}}(a_s|x_s) q_s | x_s \right]$$

2. to critic update

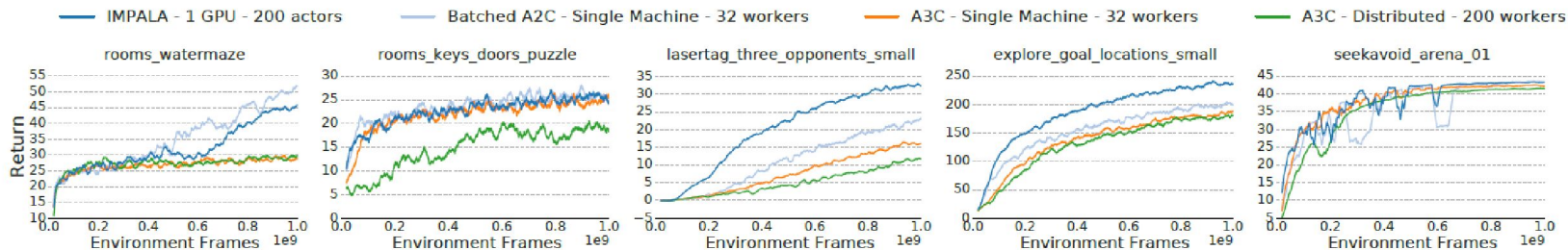
4.1. V-trace target

Consider a trajectory $(x_t, a_t, r_t)_{t=s}^{t=s+n}$ generated by the actor following some policy μ . We define the n -steps V-trace target for $V(x_s)$, our value approximation at state x_s , as:

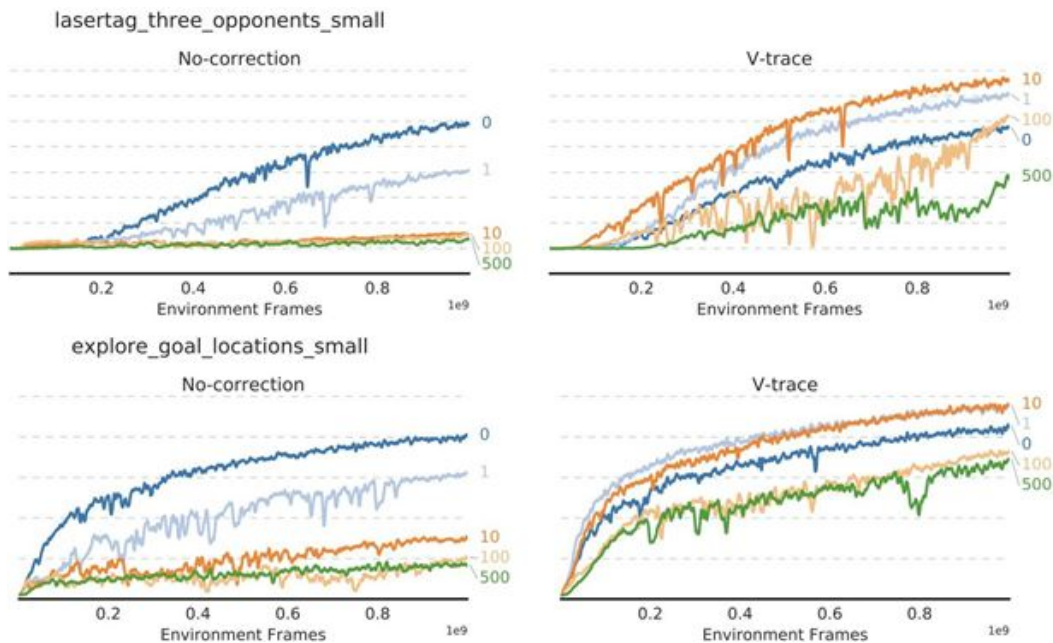
$$v_s \stackrel{\text{def}}{=} V(x_s) + \sum_{t=s}^{s+n-1} \gamma^{t-s} \left(\prod_{i=s}^{t-1} c_i \right) \delta_t V, \quad (1)$$

IMPALA - Performance

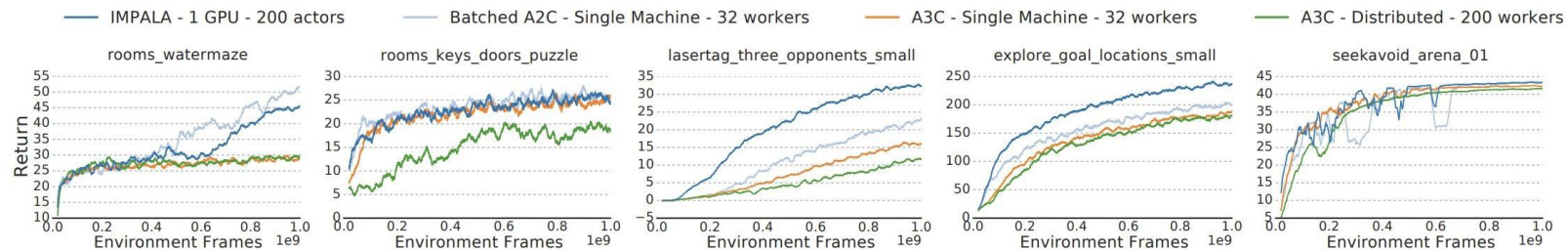
A comparison between IMPALA, A3C and batched A2C



IMPALA - Performance



IMPALA Performance



Evolution Strategies

Evolution Strategies as a Scalable Alternative to Reinforcement Learning

Tim Salimans

Jonathan Ho

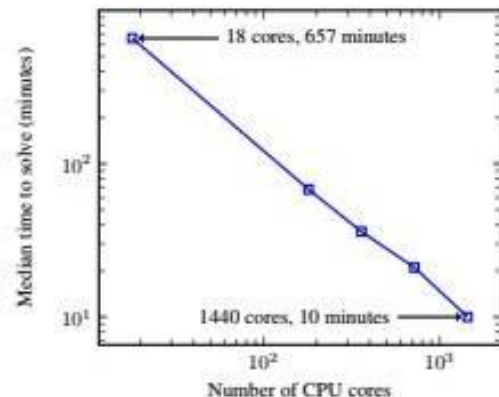
Xi Chen
OpenAI

Szymon Sidor

Ilya Sutskever

Algorithm 2 Parallelized Evolution Strategies

```
1: Input: Learning rate  $\alpha$ , noise standard deviation  $\sigma$ , initial policy parameters  $\theta_0$ 
2: Initialize:  $n$  workers with known random seeds, and initial parameters  $\theta_0$ 
3: for  $t = 0, 1, 2, \dots$  do
4:   for each worker  $i = 1, \dots, n$  do
5:     Sample  $\epsilon_i \sim \mathcal{N}(0, I)$ 
6:     Compute returns  $F_i = F(\theta_t + \sigma \epsilon_i)$ 
7:   end for
8:   Send all scalar returns  $F_i$  from each worker to every other worker
9:   for each worker  $i = 1, \dots, n$  do
10:    Reconstruct all perturbations  $\epsilon_j$  for  $j = 1, \dots, n$  using known random seeds
11:    Set  $\theta_{t+1} \leftarrow \theta_t + \alpha \frac{1}{n\sigma} \sum_{j=1}^n F_j \epsilon_j$ 
12:   end for
13: end for
```



Summary

| Algorithm | Policy Evaluation | Gradient- based optimizer | CPU | GPU | Replay Buffer | Prioritised Replay | Parameter Server | Importance Sampling |
|-----------|----------------------|---------------------------------|------|-----|------------------|-----------------------|---------------------|------------------------|
| DQN | X | X | 1 | 1 | X | | | |
| Gorila | X | X | | | X | | X | |
| Ape-X | X | X | | | X | X | | |
| A3C | X | X | many | 0 | | | | |
| Impala | X | X | many | | | | | X |

Lesson Objectives

1. Why parallelise?
2. Understand how the computation of standard RL algorithms can be distributed to decrease wall-clock training time.
3. **How these distributed RL algorithms can be modularised.**
4. **How modularised distributed RL algorithms can be implemented on real systems - case study: RLlib**
5. Examples on using RLlib

RLlib: Abstractions for Distributed Reinforcement Learning (ICML'18)

Eric Liang^{*}, Richard Liaw^{*}, Philipp Moritz, Robert Nishihara, Roy Fox, Ken Goldberg, Joseph E. Gonzalez, Michael I. Jordan, Ion Stoica

RL research scales with compute

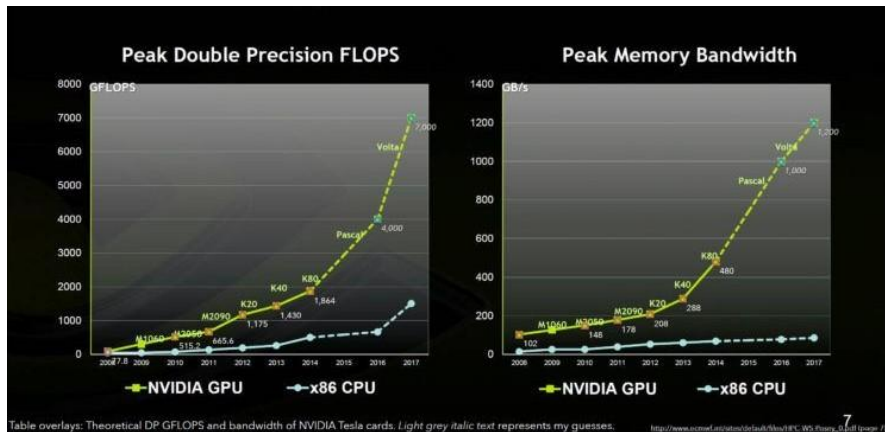


Fig. courtesy NVIDIA Inc.



CPU



GPU



TPU

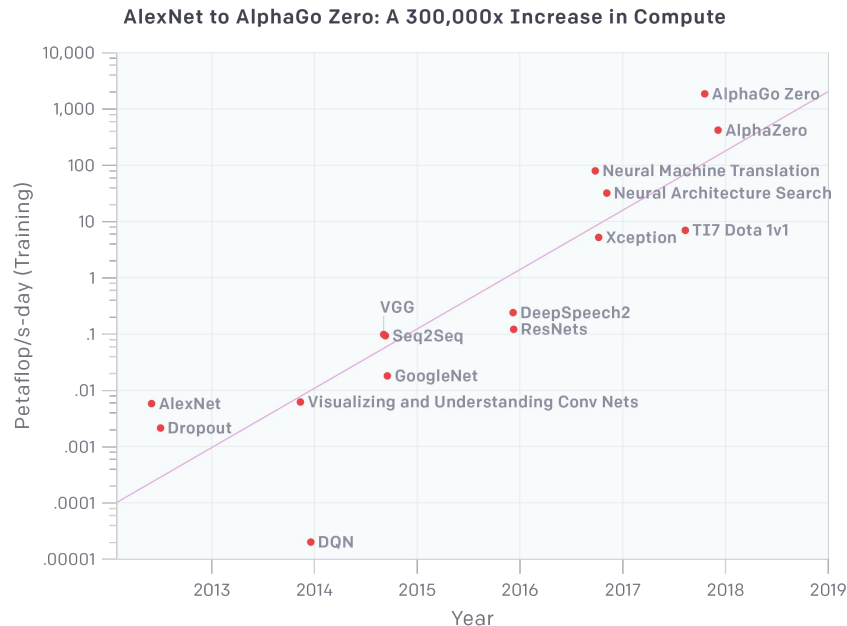
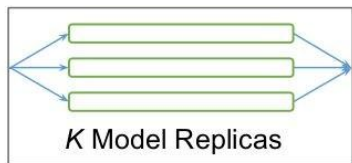
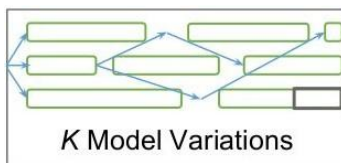


Fig. courtesy OpenAI

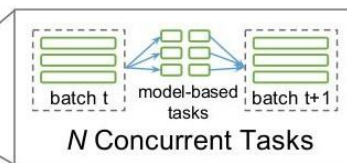
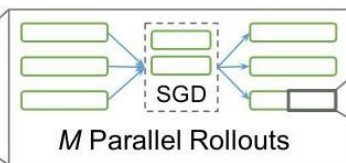
How do we leverage this hardware?



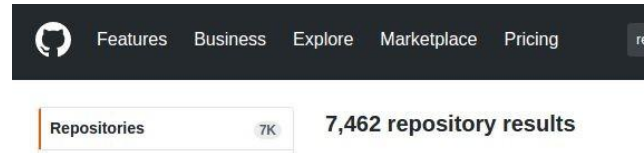
(a) Supervised Learning



(b) Reinforcement Learning



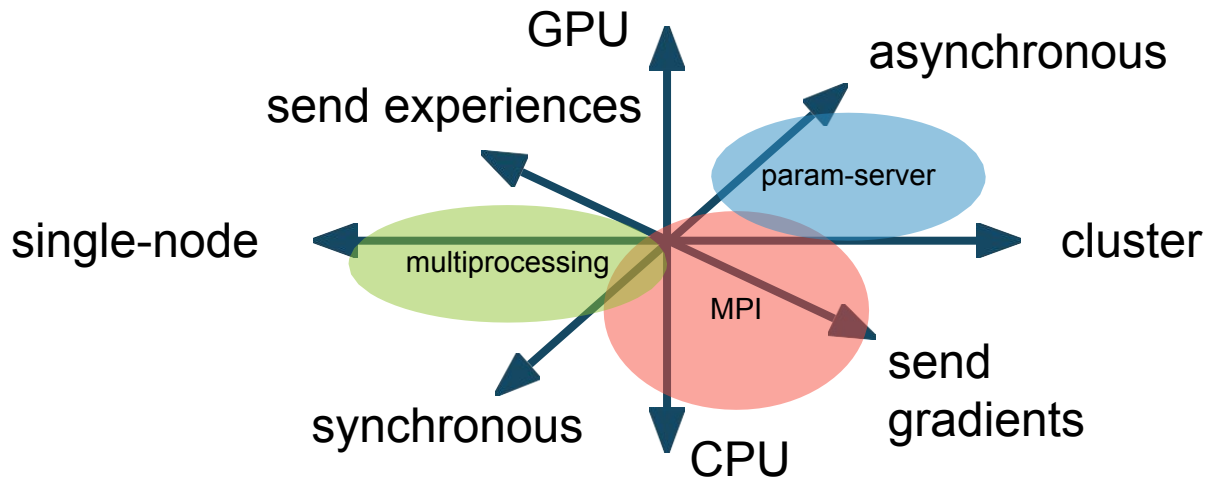
Systems for RL today



- Many implementations (7000+ repos on GitHub!)
 - how general are they (and do they scale)?
 - PPO: multiprocessing, MPI
 - AlphaZero: custom systems
 - Evolution Strategies: Redis
 - IMPALA: Distributed TensorFlow
 - A3C: shared memory, multiprocessing, TF
- Huge variety of algorithms and distributed systems used to implement, but little reuse of components

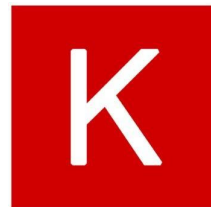
Challenges to reuse

1. Wide range of physical execution strategies for one "algorithm"



Challenges to reuse

2. Tight coupling with deep learning frameworks



Different parallelism paradigms:

- Distributed TensorFlow vs TensorFlow + MPI?

Challenges to reuse

3. Large variety of algorithms with different structures

| Algorithm Family | Policy Evaluation | Replay Buffer | Gradient-Based Optimizer | Other Distributed Components |
|----------------------|-------------------|---------------|--------------------------|------------------------------------|
| DQNs | X | X | X | |
| Policy Gradient | X | | X | |
| Off-policy PG | X | X | X | |
| Model-Based/Hybrid | X | | X | Model-Based Planning |
| Multi-Agent | X | X | X | |
| Evolutionary Methods | X | | | Derivative-Free Optimization |
| AlphaGo | X | X | X | MCTS, Derivative-Free Optimization |

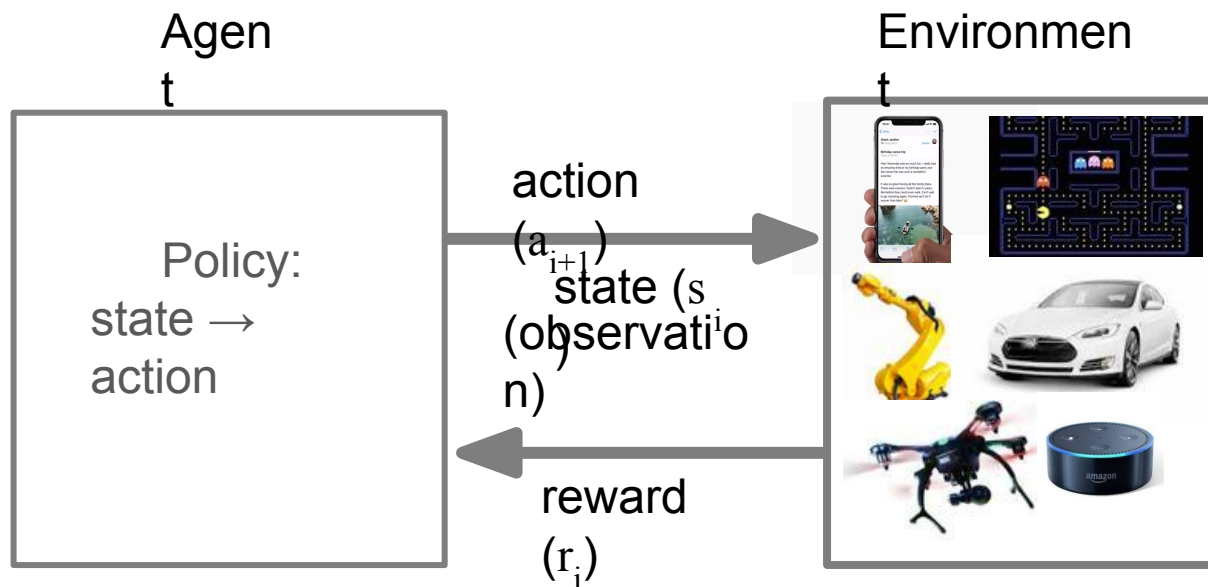
We need abstractions for RL

Good abstractions decompose RL algorithms into reusable components.

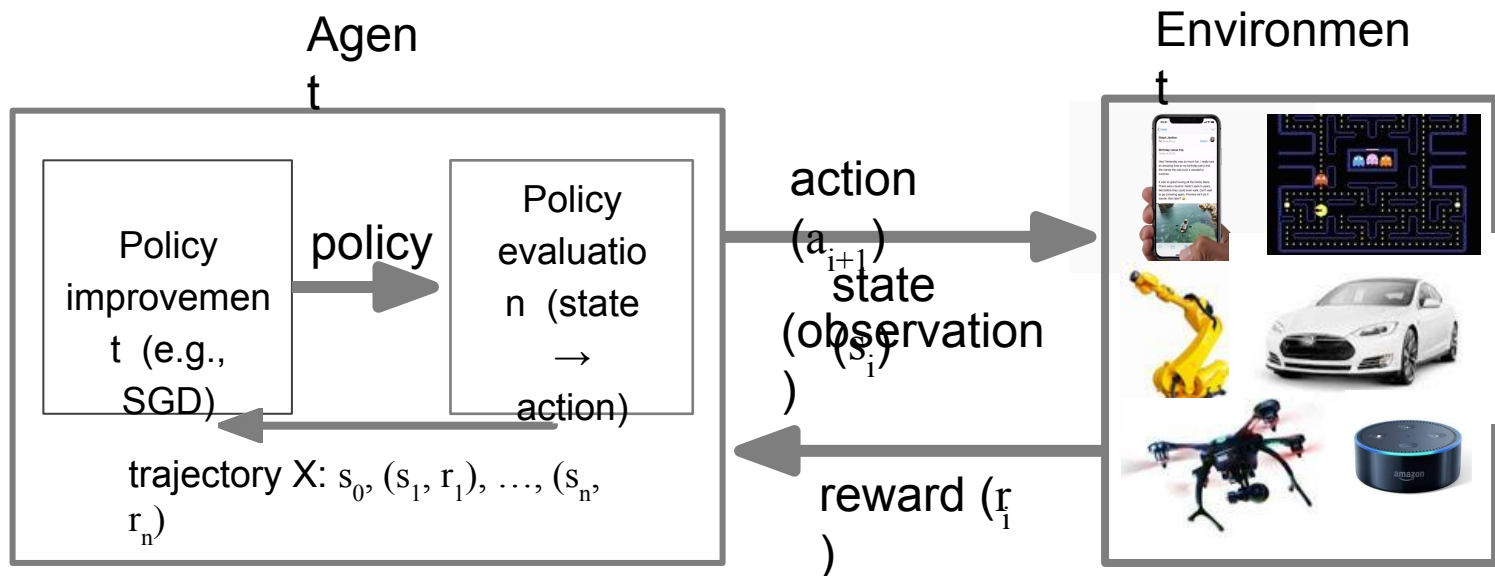
Goals:

- Code reuse across deep learning frameworks
- Scalable execution of algorithms
- Easily compare and reproduce algorithms

Structure of RL computations

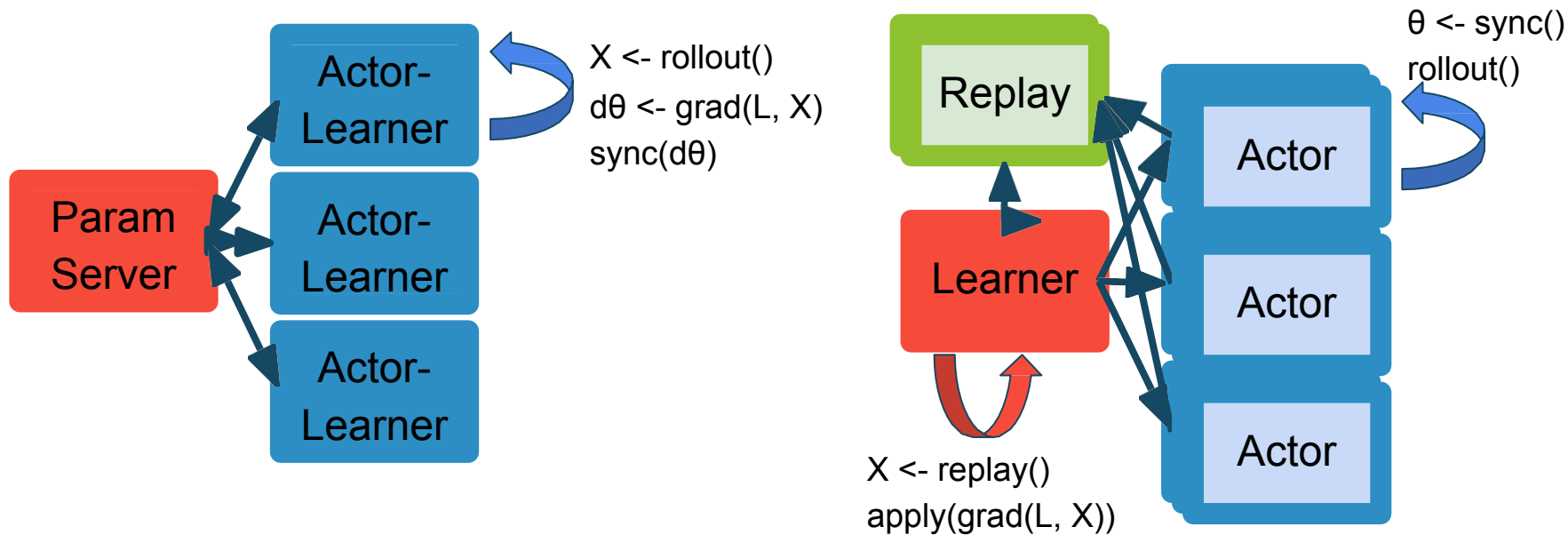


Structure of RL computations



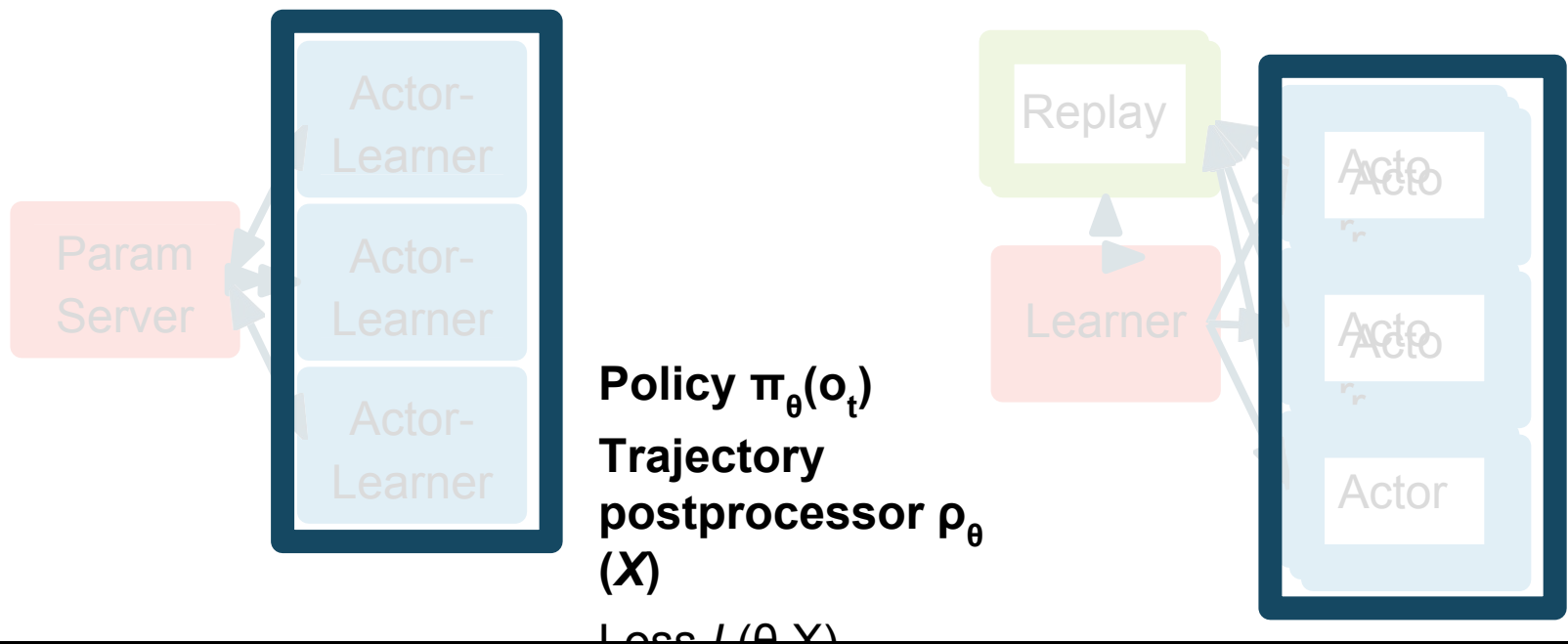
Many RL loop decompositions

Async DQN (Mnih et al; 2016) Ape-X DQN (Horgan et al; 2018)



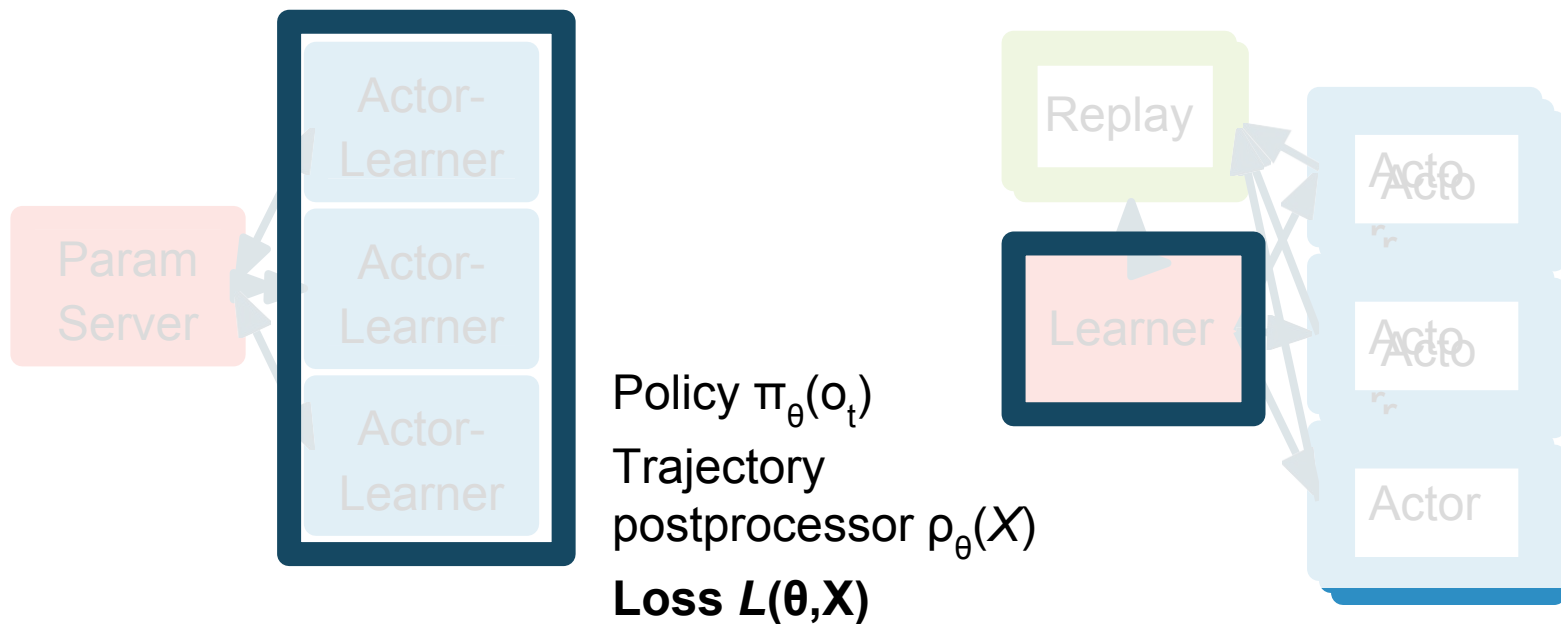
Common components

Async DQN (Mnih et al; 2016) Ape-X DQN (Horgan et al; 2018)



Common components

Async DQN (Mnih et al; 2016) Ape-X DQN (Horgan et al; 2018)



Structural differences

Async DQN (Mnih et al; 2016)

- Asynchronous optimization
- Replicated workers
- Single machine

...and this is just one family!

→ No existing system can effectively meet all the varied demands of RL workloads.

Ape-X DQN (Horgan et al; 2018)

- Central learner
- Data queues between components
- Large replay buffers
- Scales to clusters

+ Population-Based Training
(Jaderberg et al; 2017)

- Nested parallel computations
- Control decisions based on intermediate results

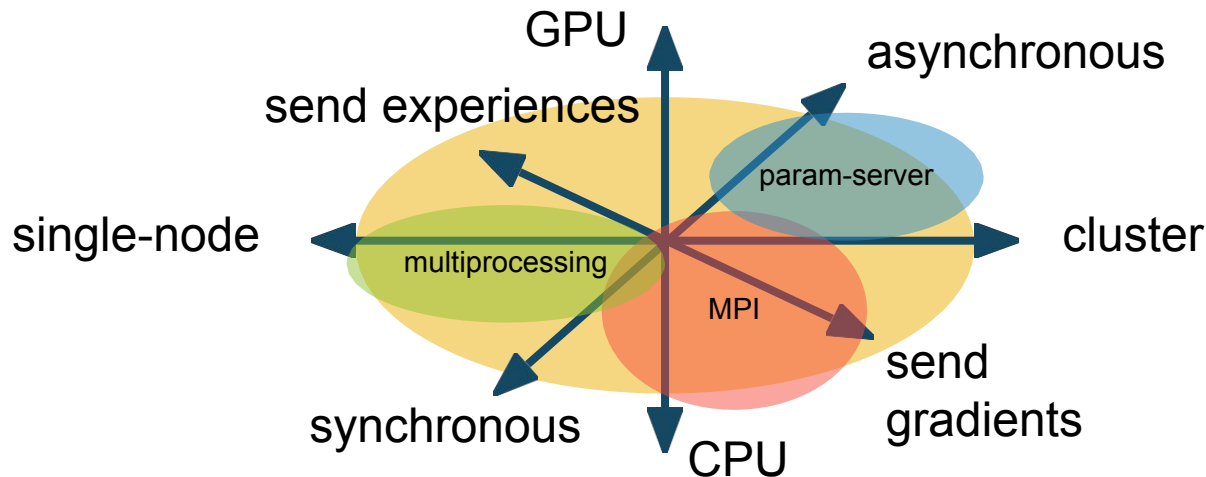
Requirements for a new system

Goal: Capture a broad range of RL workloads with high performance and substantial code reuse

1. Support stateful computations
 - e.g., simulators, neural nets, replay buffers
 - big data frameworks, e.g., Spark, are typically stateless
2. Support asynchrony
 - difficult to express in MPI, esp. nested parallelism
3. Allow easy composition of (distributed) components

Ray System Substrate

- RLlib builds on Ray to provide higher-level RL abstractions
- Hierarchical parallel task model with stateful workers
 - flexible enough to capture a broad range of RL workloads (vs specialized sys.)

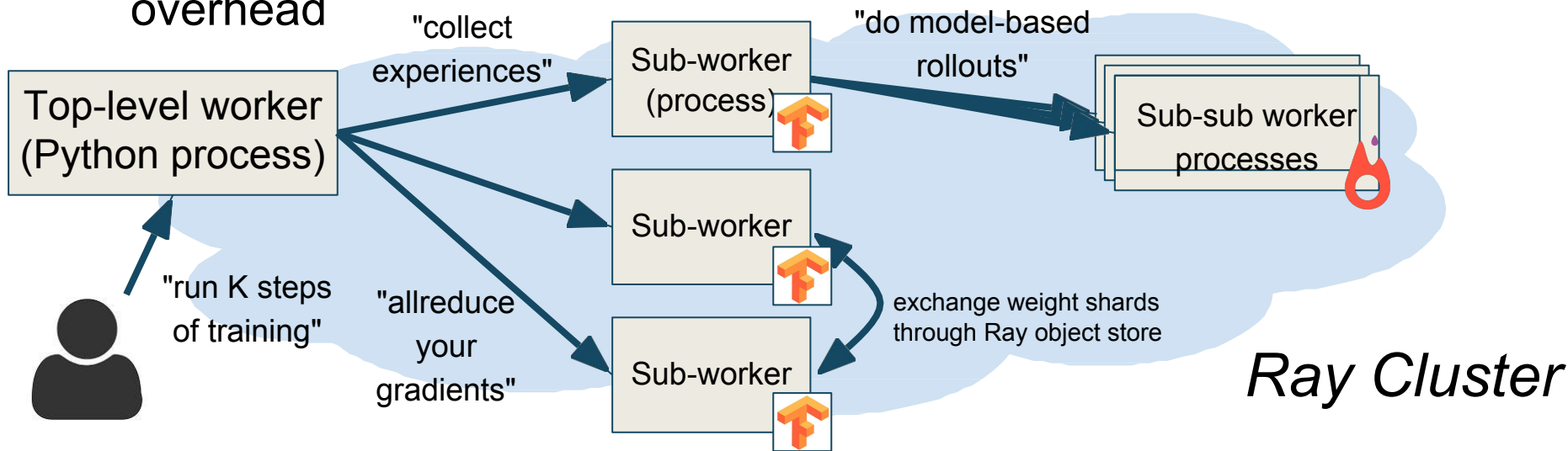


 **Hierarchical Task Model**

Hierarchical Parallel Task

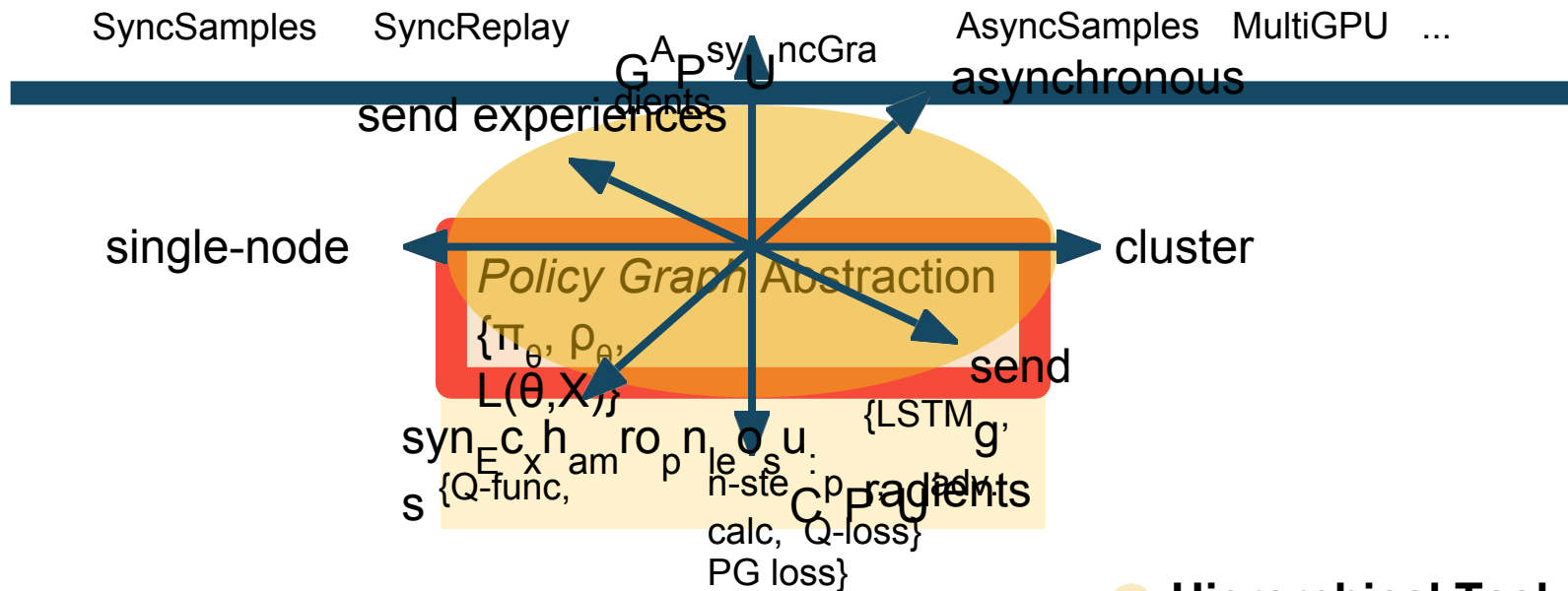
Model

1. Create Python class instances in the cluster (stateful workers)
2. Schedule short-running tasks onto workers
 - Challenge: High performance: $1e6+$ tasks/s, $\sim 200\mu s$ task overhead



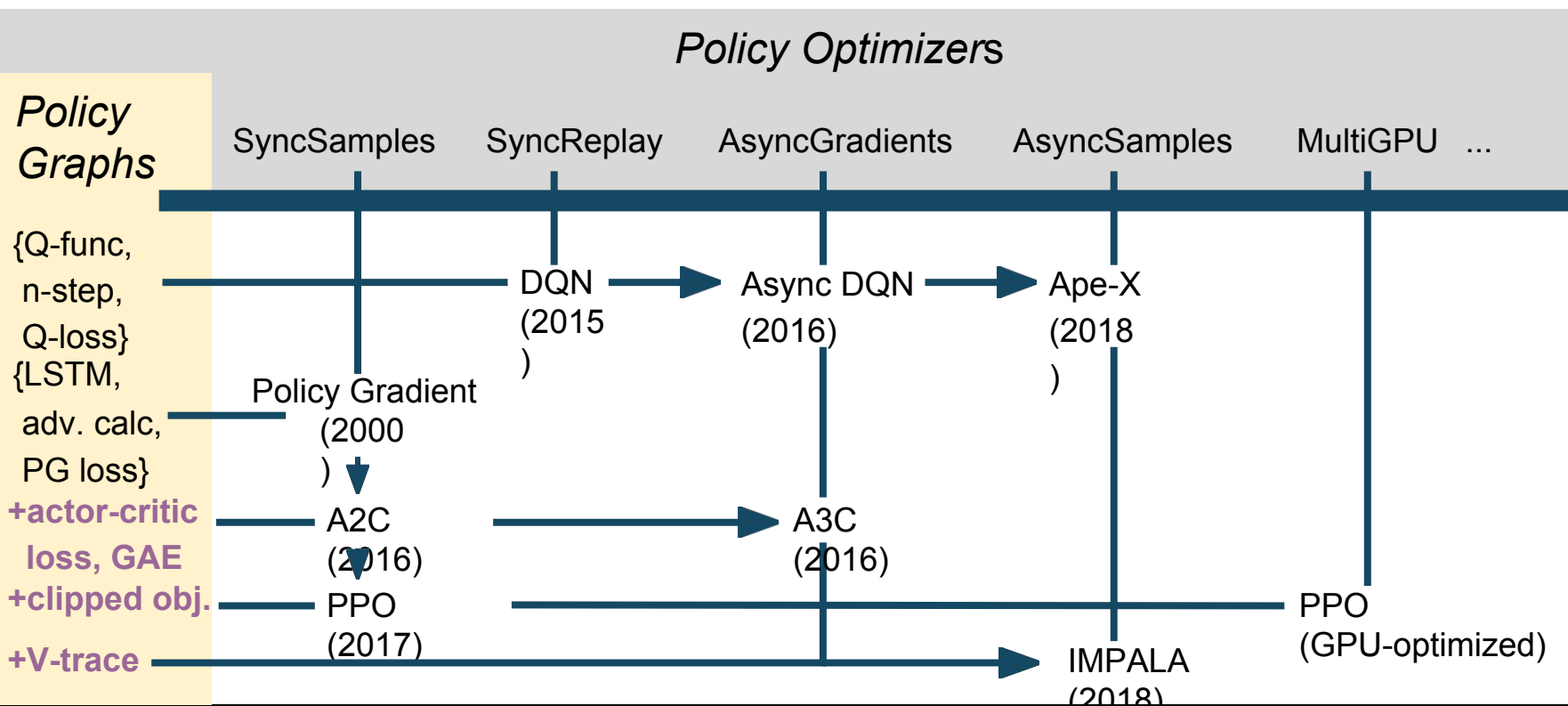
Unifying system enables RL Abstractions

Policy Optimizer Abstraction



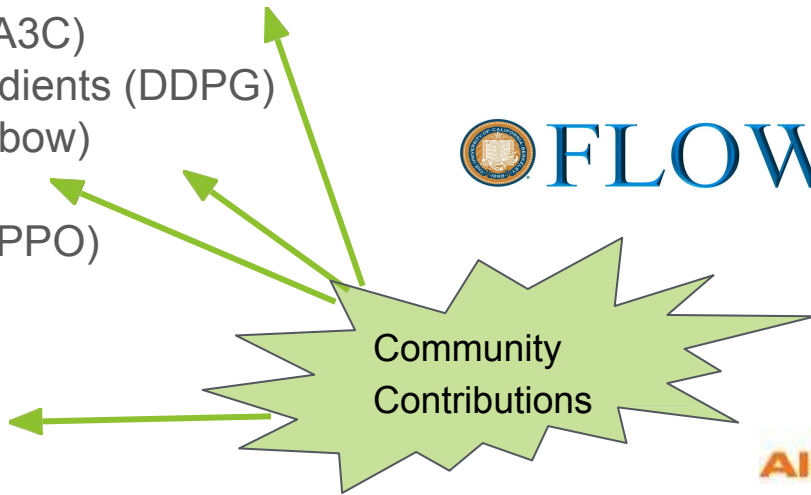
 **Hierarchical Task Model**

RLlib Abstractions in Action



RLlib Reference Algorithms

- **High-throughput architectures**
 - Distributed Prioritized Experience Replay (Ape-X)
 - Importance Weighted Actor-Learner Architecture (IMPALA)
- **Gradient-based**
 - Advantage Actor-Critic (A2C, A3C)
 - Deep Deterministic Policy Gradients (DDPG)
 - Deep Q Networks (DQN, Rainbow)
 - Policy Gradients
 - Proximal Policy Optimization (PPO)
- **Derivative-free**
 - Augmented Random Search (ARS)
 - Evolution Strategies



FLOW Lab



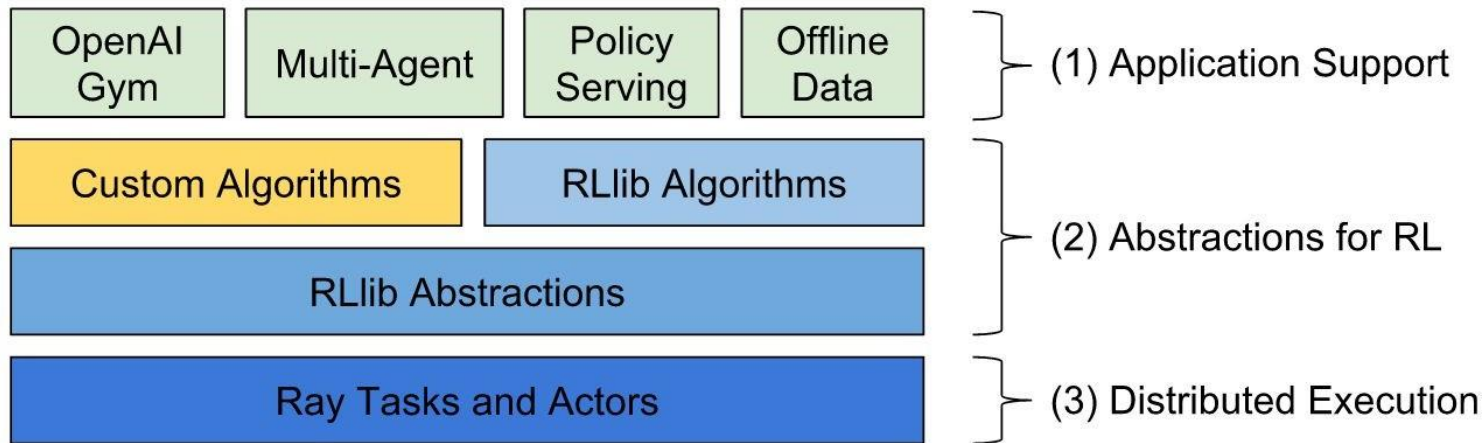
RLlib Reference Algorithms

| Atari env | RLlib IMPALA 32-workers @1 hour | Mnih et al A3C 16-workers @1 hour |
|---------------|---------------------------------|-----------------------------------|
| BeamRider | 3181 | ~1000 |
| Breakout | 538 | ~10 |
| Qbert | 10850 | ~500 |
| SpaceInvaders | 843 | ~300 |

1 GPU + 64 vCPUs (large single machine)

Scale your algorithms with RLlib

- Beyond a "collection of algorithms",
- RLlib's abstractions let you easily implement and scale new algorithms (multi-agent, novel losses, architectures, etc)



Code example: training PPO

Tutorial on google Colab:

<https://drive.google.com/open?id=1pvE7KvnhYR0Ynqt0J0fzYSmkjLg64Qq>

```
import ray
import ray.rllib.agents.ppo as ppo
from ray.tune.logger import pretty_print

ray.init()
config = ppo.DEFAULT_CONFIG.copy()
config["num_gpus"] = 0
config["num_workers"] = 1
agent = ppo.PPOAgent(config=config, env="CartPole-v0")

# Can optionally call agent.restore(path) to load a checkpoint.

for i in range(1000):
    # Perform one iteration of training the policy with PPO
    result = agent.train()
    print(pretty_print(result))

    if i % 100 == 0:
        checkpoint = agent.save()
        print("checkpoint saved at", checkpoint)
```

Code example: multi-agent RL

```
trainer = pg.PGAgent(env="my_multiagent_env", config={
    "multiagent": {
        "policy_graphs": {
            "car1": (PGPolicyGraph, car_obs_space, car_act_space, {"gamma": 0.85}),
            "car2": (PGPolicyGraph, car_obs_space, car_act_space, {"gamma": 0.99}),
            "traffic_light": (PGPolicyGraph, tl_obs_space, tl_act_space, {}),
        },
        "policy_mapping_fn":
            lambda agent_id:
                "traffic_light" # Traffic lights are always controlled by this policy
                if agent_id.startswith("traffic_light_")
                else random.choice(["car1", "car2"]) # Randomly choose from car policies
    },
},
})

while True:
    print(trainer.train())
```


Code example: hyperparam tuning

```
import ray
import ray.tune as tune

ray.init()
tune.run_experiments({
    "my_experiment": {
        "run": "PPO",
        "env": "CartPole-v0",
        "stop": {"episode_reward_mean": 200},
        "config": {
            "num_gpus": 0,
            "num_workers": 1,
            "sgd_stepsize": tune.grid_search([0.01, 0.001, 0.0001]),
        },
    },
})
```

Code example: hyperparam tuning

== Status ==

Using FIFO scheduling algorithm.

Resources requested: 4/4 CPUs, 0/0 GPUs

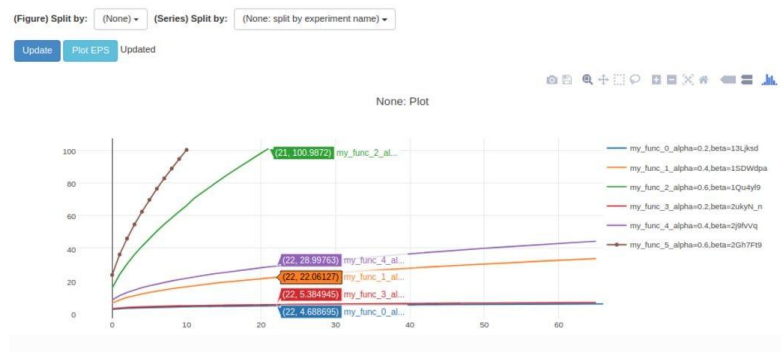
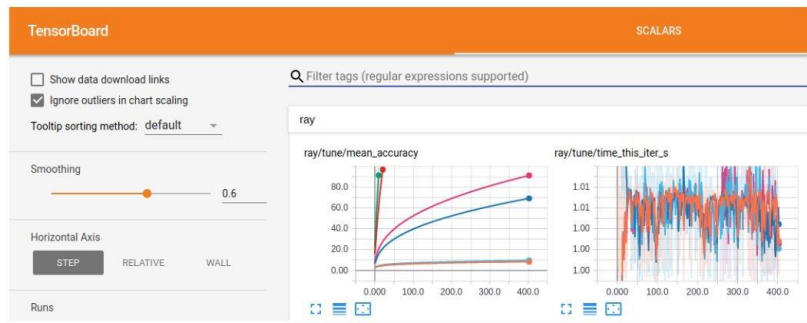
Result logdir: ~/ray_results/my_experiment

PENDING trials:

- PPO_CartPole-v0_2_sgd_stepsize=0.0001: PENDING

RUNNING trials:

- PPO_CartPole-v0_0_sgd_stepsize=0.01: RUNNING [pid=21940], 16 s, 4013 ts, 22 rew
- PPO_CartPole-v0_1_sgd_stepsize=0.001: RUNNING [pid=21942], 27 s, 8111 ts, 54.7 rew



Summary: Ray and RLlib addresses challenges in providing scalable abstractions for reinforcement learning.

RLlib is open source and available at <http://rlib.io>

Thanks!