Policy Gradient Methods

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Instead of predicting expected returns, we could train the method to directly predict the policy

$$\pi(a|s; \theta).$$

Obtaining the full distribution over all actions would also allow us to sample the actions according to the distribution $\pi$ instead of just $\varepsilon$-greedy sampling.

However, to train the network, we maximize the expected return $v_\pi(s)$ and to that account we need to compute its gradient $\nabla_\theta v_\pi(s)$. 
In addition to discarding $\varepsilon$-greedy action selection, policy gradient methods allow producing policies which are by nature stochastic, as in card games with imperfect information, while the action-value methods have no natural way of finding stochastic policies (distributional RL might be of some use though).
Let \( \pi(a|s; \theta) \) be a parametrized policy. We denote the initial state distribution as \( h(s) \) and the on-policy distribution under \( \pi \) as \( \mu(s) \). Let also \( J(\theta) \overset{\text{def}}{=} \mathbb{E}_{s \sim h} v_{\pi}(s) \).

Then

\[
\nabla_{\theta} v_{\pi}(s) \propto \sum_{s' \in S} P(s \to \ldots \to s'|\pi) \sum_{a \in A} q_{\pi}(s', a) \nabla_{\theta} \pi(a|s'; \theta)
\]

and

\[
\nabla_{\theta} J(\theta) \propto \sum_{s \in S} \mu(s) \sum_{a \in A} q_{\pi}(s, a) \nabla_{\theta} \pi(a|s; \theta),
\]

where \( P(s \to \ldots \to s'|\pi) \) is the probability of getting to state \( s' \) when starting from state \( s \), after any number of 0, 1, \ldots steps. The \( \gamma \) parameter should be treated as a form of termination, i.e., \( P(s \to \ldots \to s'|\pi) \propto \sum_{k=0}^{\infty} \gamma^k P(s \to s' \text{ in } k \text{ steps }|\pi) \).
Proof of Policy Gradient Theorem

\[ \nabla v_\pi(s) = \nabla \left[ \sum_a \pi(a | s; \theta) q_\pi(s, a) \right] \]

\[ = \sum_a \left[ \nabla \pi(a | s; \theta) q_\pi(s, a) + \pi(a | s; \theta) \nabla q_\pi(s, a) \right] \]

\[ = \sum_a \left[ \nabla \pi(a | s; \theta) q_\pi(s, a) + \pi(a | s; \theta) \nabla \left( \sum_{s'} p(s' | s, a) (r + \gamma v_\pi(s')) \right) \right] \]

\[ = \sum_a \left[ \nabla \pi(a | s; \theta) q_\pi(s, a) + \gamma \pi(a | s; \theta) \left( \sum_{s'} p(s' | s, a) \nabla v_\pi(s') \right) \right] \]

We now expand \( v_\pi(s') \).

\[ = \sum_a \left[ \nabla \pi(a | s; \theta) q_\pi(s, a) + \gamma \pi(a | s; \theta) \left( \sum_{s'} p(s' | s, a) \right) \right] \]

\[ \sum_{a'} \left[ \nabla \pi(a' | s'; \theta) q_\pi(s', a') + \gamma \pi(a' | s'; \theta) \left( \sum_{s''} p(s'' | s', a') \nabla v_\pi(s'') \right) \right] \]

Continuing to expand all \( v_\pi(s'') \), we obtain the following:

\[ \nabla v_\pi(s) = \sum_{s' \in S} \sum_{k=0}^{\infty} \gamma^k P(s \rightarrow s' \text{ in } k \text{ steps} | \pi) \sum_{a \in A} q_\pi(s', a) \nabla_{\theta} \pi(a | s'; \theta). \]
Proof of Policy Gradient Theorem

To finish the proof of the first part, recall that

\[ \sum_{k=0}^{\infty} \gamma^k P(s \rightarrow s' \text{ in } k \text{ steps } | \pi) \propto P(s \rightarrow \ldots \rightarrow s' | \pi). \]

For the second part, we know that

\[ \nabla_{\theta} J(\theta) = \mathbb{E}_{s \sim h} \nabla_{\theta} v_\pi(s) \propto \mathbb{E}_{s \sim h} \sum_{s' \in S} P(s \rightarrow \ldots \rightarrow s' | \pi) \sum_{a \in A} q_\pi(s', a) \nabla_{\theta} \pi(a | s'; \theta), \]

therefore using the fact that \( \mu(s') = \mathbb{E}_{s \sim h} P(s \rightarrow \ldots \rightarrow s' | \pi) \) we get

\[ \nabla_{\theta} J(\theta) \propto \sum_{s \in S} \mu(s) \sum_{a \in A} q_\pi(s, a) \nabla_{\theta} \pi(a | s; \theta). \]
The REINFORCE algorithm (Williams, 1992) uses directly the policy gradient theorem, minimizing $-J(\theta) \overset{\text{def}}{=} -\mathbb{E}_{s \sim h} v_\pi(s)$. The loss gradient is then

$$\nabla_{\theta} - J(\theta) \propto -\sum_{s \in S} \mu(s) \sum_{a \in A} q_\pi(s, a) \nabla_{\theta} \pi(a|s; \theta) = -\mathbb{E}_{s \sim \mu} \sum_{a \in A} q_\pi(s, a) \nabla_{\theta} \pi(a|s; \theta).$$

However, the sum over all actions is problematic. Instead, we rewrite it to an expectation which we can estimate by sampling:

$$\nabla_{\theta} - J(\theta) \propto \mathbb{E}_{s \sim \mu} \mathbb{E}_{a \sim \pi} q_\pi(s, a) \nabla_{\theta} - \ln \pi(a|s; \theta),$$

where we used the fact that

$$\nabla_{\theta} \ln \pi(a|s; \theta) = \frac{1}{\pi(a|s; \theta)} \nabla_{\theta} \pi(a|s; \theta).$$
REINFORCE Algorithm

REINFORCE therefore minimizes the loss

$$\mathbb{E}_{s \sim \mu} \mathbb{E}_{a \sim \pi} q_{\pi}(s, a) \nabla_{\theta} - \ln \pi(a|s; \theta),$$

estimating the $q_{\pi}(s, a)$ by a single sample.

Note that the loss is just a weighted variant of negative log-likelihood (NLL), where the sampled actions play a role of gold labels and are weighted according to their return.

### REINFORCE: Monte-Carlo Policy-Gradient Control (episodic) for $\pi_*$

- **Input:** a differentiable policy parameterization $\pi(a|s, \theta)$
- **Algorithm parameter:** step size $\alpha > 0$
- **Initialize policy parameter** $\theta \in \mathbb{R}^{d'}$ (e.g., to 0)

Loop forever (for each episode):

1. Generate an episode $S_0, A_0, R_1, \ldots, S_{T-1}, A_{T-1}, R_T$, following $\pi(\cdot|\cdot, \theta)$
2. Loop for each step of the episode $t = 0, 1, \ldots, T - 1$:
   - $G \leftarrow \sum_{k=t+1}^{T} \gamma^{k-t-1} R_k$
   - $\theta \leftarrow \theta + \alpha G \nabla \ln \pi(A_t|S_t, \theta)$

Modified from Algorithm 13.3 of "Reinforcement Learning: An Introduction, Second Edition" by removing $\gamma^t$ from the update of $\theta$. 
On-policy Distribution in REINFORCE

In the proof, we assumed $\gamma$ is used as a form of termination in the definition of the on-policy distribution.

However, even when discounting is used during training (to guarantee convergence even for very long episodes), evaluation is often performed without discounting.

Consequently, the distribution $\mu$ used in the REINFORCE algorithm is almost always the unterminated (undiscounted) on-policy distribution (I am not aware of any implementation or paper that would use it), so that we learn even in states that are far from the beginning of an episode.

Note that this is actually true even for DQN and its variants. Therefore, the discounting parameter $\gamma$ is used mostly as a variance-reduction technique.
The returns can be arbitrary – better-than-average and worse-than-average returns cannot be recognized from the absolute value of the return.

Hopefully, we can generalize the policy gradient theorem using a baseline $b(s)$ to

$$
\nabla_\theta J(\theta) \propto \sum_{s \in S} \mu(s) \sum_{a \in A} (q_{\pi}(s, a) - b(s)) \nabla_\theta \pi(a|s; \theta).
$$

The baseline $b(s)$ can be a function or even a random variable, as long as it does not depend on $a$, because

$$
\sum_{a} b(s) \nabla_\theta \pi(a|s; \theta) = b(s) \sum_{a} \nabla_\theta \pi(a|s; \theta) = b(s) \nabla_\theta \sum_{a} \pi(a|s; \theta) = b(s) \nabla 1 = 0.
$$
A good choice for $b(s)$ is $v_\pi(s)$, which can be shown to minimize the variance of the gradient estimate (in limit $\gamma \rightarrow 1$; see L. Weaver and N. Tao, *The Optimal Reward Baseline for Gradient-Based Reinforcement Learning* for the proof). Such baseline reminds centering of returns, given that

$$v_\pi(s) = \mathbb{E}_{a \sim \pi} q_\pi(s, a).$$

Then, better-than-average returns are positive and worse-than-average returns are negative. The resulting $q_\pi(s, a) - v_\pi(s)$ function is also called the **advantage** function

$$a_\pi(s, a) \overset{\text{def}}{=} q_\pi(s, a) - v_\pi(s).$$

Of course, the $v_\pi(s)$ baseline can be only approximated. If neural networks are used to estimate $\pi(a|s; \theta)$, then some part of the network is usually shared between the policy and value function estimation, which is trained using mean square error of the predicted and observed return.
REINFORCE with Baseline (episodic), for estimating $\pi(\theta) \approx \pi_*$

Input: a differentiable policy parameterization $\pi(a|s, \theta)$
Input: a differentiable state-value function parameterization $\hat{v}(s, w)$
Algorithm parameters: step sizes $\alpha^\theta > 0$, $\alpha^w > 0$
Initialize policy parameter $\theta \in \mathbb{R}^{d'}$ and state-value weights $w \in \mathbb{R}^d$ (e.g., to 0)

Loop forever (for each episode):
  Generate an episode $S_0, A_0, R_1, \ldots, S_{T-1}, A_{T-1}, R_T$, following $\pi(\cdot|\cdot, \theta)$
  Loop for each step of the episode $t = 0, 1, \ldots, T - 1$:
  \[
  G \leftarrow \sum_{k=t+1}^{T} \gamma^{k-t-1} R_k \\
  \delta \leftarrow G - \hat{v}(S_t, w) \\
  w \leftarrow w + \alpha^w \delta \nabla \hat{v}(S_t, w) \\
  \theta \leftarrow \theta + \alpha^\theta \delta \nabla \ln \pi(A_t|S_t, \theta)
  \]

Modified from Algorithm 13.4 of "Reinforcement Learning: An Introduction, Second Edition" by removing $\gamma^t$ from the update of $\theta$. 
REINFORCE with Baseline

$G_0$
Total reward on episode
averaged over 100 runs

REINFORCE
$\alpha = 2^{-13}$

REINFORCE with baseline
$\alpha^0 = 2^{-9}$, $\alpha^w = 2^{-6}$

Figure 13.2 of "Reinforcement Learning: An Introduction, Second Edition".
Actor-Critic

It is possible to combine the policy gradient methods and temporal difference methods, creating a family of algorithms usually called the **actor-critic** methods.

The idea is straightforward – instead of estimating the episode return using the whole episode rewards, we can use $n$-step temporal difference estimation.
One-step Actor–Critic (episodic), for estimating $\pi_\theta \approx \pi_*$

Input: a differentiable policy parameterization $\pi(a|s, \theta)$
Input: a differentiable state-value function parameterization $\hat{v}(s, w)$
Parameters: step sizes $\alpha^\theta > 0$, $\alpha^w > 0$
Initialize policy parameter $\theta \in \mathbb{R}^{d'}$ and state-value weights $w \in \mathbb{R}^d$ (e.g., to 0)
Loop forever (for each episode):
   Initialize $S$ (first state of episode)

Loop while $S$ is not terminal (for each time step):
   $A \sim \pi(\cdot|S, \theta)$
   Take action $A$, observe $S', R$
   $\delta \leftarrow R + \gamma \hat{v}(S', w) - \hat{v}(S, w)$  
      (if $S'$ is terminal, then $\hat{v}(S', w) \equiv 0$)
   $w \leftarrow w + \alpha^w \delta \nabla \hat{v}(S, w)$
   $\theta \leftarrow \theta + \alpha^\theta \delta \nabla \ln \pi(A|S, \theta)$
   $S \leftarrow S'$
The A3C was introduced in a 2016 paper from Volodymyr Mnih et al. (the same group as DQN) *Asynchronous Methods for Deep Reinforcement Learning*. The authors propose an asynchronous framework, where multiple workers share one neural network, each training using either an off-line or on-line RL algorithm.

They compare 1-step Q-learning, 1-step Sarsa, \(n\)-step Q-learning and A3C (an *asynchronous advantage actor-critic* method). For A3C, they compare a version with and without LSTM. The authors also introduce *entropy regularization term* \(-\beta H(\pi(s; \theta))\) to the loss to support exploration and discourage premature convergence (they use \(\beta = 0.01\)).
Algorithm 1 Asynchronous one-step Q-learning - pseudocode for each actor-learner thread.

// Assume global shared $\theta$, $\theta^-$, and counter $T = 0$.
Initialize thread step counter $t \leftarrow 0$
Initialize target network weights $\theta^- \leftarrow \theta$
Initialize network gradients $d\theta \leftarrow 0$
Get initial state $s$

repeat
  Take action $a$ with $\epsilon$-greedy policy based on $Q(s, a; \theta)$
  Receive new state $s'$ and reward $r$
  $y = \begin{cases} 
  r & \text{for terminal } s' \\
  r + \gamma \max_{a'} Q(s', a'; \theta^-) & \text{for non-terminal } s'
  \end{cases}$
  Accumulate gradients wrt $\theta$: $d\theta \leftarrow d\theta + \frac{\partial(y - Q(s, a; \theta))^2}{\partial \theta}$
  $s = s'$
  $T \leftarrow T + 1$ and $t \leftarrow t + 1$
  if $T \mod I_{\text{target}} == 0$ then
    Update the target network $\theta^- \leftarrow \theta$
  end if
  if $t \mod I_{\text{AsyncUpdate}} == 0$ or $s$ is terminal then
    Perform asynchronous update of $\theta$ using $d\theta$.
    Clear gradients $d\theta \leftarrow 0$.
  end if
  until $T > T_{\text{max}}$

Algorithm 1 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.
**Algorithm S2** Asynchronous n-step Q-learning - pseudocode for each actor-learner thread.

// Assume global shared parameter vector $\theta$.
// Assume global shared target parameter vector $\theta^-$.  
// Assume global shared counter $T = 0$.  
Initialize thread step counter $t \leftarrow 1$
Initialize target network parameters $\theta^- \leftarrow \theta$
Initialize thread-specific parameters $\theta' = \theta$
Initialize network gradients $d\theta \leftarrow 0$

repeat

Clear gradients $d\theta \leftarrow 0$
Synchronize thread-specific parameters $\theta' = \theta$
$t_{start} = t$
Get state $s_t$

repeat

Take action $a_t$ according to the $\epsilon$-greedy policy based on $Q(s_t, a; \theta')$
Receive reward $r_t$ and new state $s_{t+1}$
$t \leftarrow t + 1$
$T \leftarrow T + 1$

until terminal $s_t$ or $t - t_{start} = t_{\text{max}}$

$R = \begin{cases} 0 & \text{for terminal } s_t \\ \max_a Q(s_t, a; \theta^-) & \text{for non-terminal } s_t \end{cases}$

for $i \in \{t - 1, \ldots, t_{\text{start}}\}$ do

$R \leftarrow r_t + \gamma R$

Accumulate gradients wrt $\theta'$: $d\theta \leftarrow d\theta + \frac{\partial(R - Q(s_t, a_t; \theta'))^2}{\partial \theta'}$

end for

Perform asynchronous update of $\theta$ using $d\theta$.
if $T \mod I_{\text{target}} = 0$ then

$\theta^- \leftarrow \theta$

end if

until $T > T_{\text{max}}$

Algorithm S2 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.
**Algorithm S3** Asynchronous advantage actor-critic - pseudocode for each actor-learner thread.

// Assume global shared parameter vectors $\theta$ and $\theta_v$ and global shared counter $T = 0$
// Assume thread-specific parameter vectors $\theta'$ and $\theta'_v$
Initialize thread step counter $t \leftarrow 1$

repeat
  Reset gradients: $d\theta \leftarrow 0$ and $d\theta_v \leftarrow 0$.
  Synchronize thread-specific parameters $\theta' = \theta$ and $\theta'_v = \theta_v$
  $t_{\text{start}} = t$
  Get state $s_t$
  repeat
    Perform $a_t$ according to policy $\pi(a_t|s_t; \theta')$
    Receive reward $r_t$ and new state $s_{t+1}$
    $t \leftarrow t + 1$
    $T \leftarrow T + 1$
  until terminal $s_t$ or $t - t_{\text{start}} = t_{\text{max}}$
  $R = \begin{cases} 0 & \text{for terminal } s_t \\ V(s_t, \theta'_v) & \text{for non-terminal } s_t \end{cases}$
  // Bootstrap from last state
  for $i \in \{t - 1, \ldots, t_{\text{start}}\}$ do
    $R \leftarrow r_i + \gamma R$
    Accumulate gradients wrt $\theta'$: $d\theta \leftarrow d\theta + \nabla_{\theta'} \log \pi(a_i|s_i; \theta')(R - V(s_i; \theta'_v))$
    Accumulate gradients wrt $\theta'_v$: $d\theta_v \leftarrow d\theta_v + \partial (R - V(s_i; \theta'_v))^2 / \partial \theta'_v$
  end for
  Perform asynchronous update of $\theta$ using $d\theta$ and of $\theta_v$ using $d\theta_v$.
until $T > T_{\text{max}}$

*Algorithm S3 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.*
Asynchronous Methods for Deep RL

All methods performed updates every 5 actions \( t_{\text{max}} = I_{\text{AsyncUpdate}} = 5 \), updating the target network each 40,000 frames.

The Atari inputs were processed as in DQN, using also action repeat 4.

The network architecture is: 16 filters \( 8 \times 8 \) stride 4, 32 filters \( 4 \times 4 \) stride 2, followed by a fully connected layer with 256 units. All hidden layers apply a ReLU non-linearity. Values and/or action values were then generated from the (same) last hidden layer.

The LSTM methods utilized a 256-unit LSTM cell after the dense hidden layer.

All experiments used a discount factor of \( \gamma = 0.99 \) and used RMSProp with momentum decay factor of 0.99.
Asynchronous Methods for Deep RL

![Figure 1 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.](image)

<table>
<thead>
<tr>
<th>Method</th>
<th>Training Time</th>
<th>Mean</th>
<th>Median</th>
</tr>
</thead>
<tbody>
<tr>
<td>DQN</td>
<td>8 days on GPU</td>
<td>121.9%</td>
<td>47.5%</td>
</tr>
<tr>
<td>Gorila</td>
<td>4 days, 100 machines</td>
<td>215.2%</td>
<td>71.3%</td>
</tr>
<tr>
<td>D-DQN</td>
<td>8 days on GPU</td>
<td>332.9%</td>
<td>110.9%</td>
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<tr>
<td>Dueling D-DQN</td>
<td>8 days on GPU</td>
<td>343.8%</td>
<td>117.1%</td>
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<td>Prioritized DQN</td>
<td>8 days on GPU</td>
<td>463.6%</td>
<td>127.6%</td>
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<td>A3C, FF</td>
<td>1 day on CPU</td>
<td>344.1%</td>
<td>68.2%</td>
</tr>
<tr>
<td>A3C, LSTM</td>
<td>4 days on CPU</td>
<td>496.8%</td>
<td>116.6%</td>
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<tr>
<td>A3C</td>
<td>4 days on CPU</td>
<td>623.0%</td>
<td>112.6%</td>
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<table>
<thead>
<tr>
<th>Method</th>
<th>Number of threads</th>
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<tbody>
<tr>
<td>1</td>
<td>1</td>
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<tr>
<td>1-step Q</td>
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</tr>
<tr>
<td>1-step SARSA</td>
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<td>n-step Q</td>
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<tr>
<td>A3C</td>
<td>16</td>
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Table 1 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.
Figure 3 of the paper “Asynchronous Methods for Deep Reinforcement Learning” by Volodymyr Mnih et al.
Asynchronous Methods for Deep RL

Figure 4 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.

NPFL122, Lecture 6
Policy Gradient Methods  REINFORCE  Baseline  Actor-Critic  A3C  PAAC
Asynchronous Methods for Deep RL

Figure 2 of the paper "Asynchronous Methods for Deep Reinforcement Learning" by Volodymyr Mnih et al.
Parallel Advantage Actor Critic

An alternative to independent workers is to train in a synchronous and centralized way by having the workers to only generate episodes. Such approach was described in May 2017 by Clemente et al., who named their agent parallel advantage actor-critic (PAAC).

Figure 1 of the paper “Efficient Parallel Methods for Deep Reinforcement Learning” by Alfredo V. Clemente et al.
Algorithm 1 Parallel advantage actor-critic

1: Initialize timestep counter $N = 0$ and network weights $\theta, \theta_v$
2: Instantiate set $e$ of $n_e$ environments
3: repeat
4:   for $t = 1$ to $t_{max}$ do
5:     Sample $a_t$ from $\pi(a_t|s_t; \theta)$
6:     Calculate $v_t$ from $V(s_t; \theta_v)$
7:     parallel for $i = 1$ to $n_e$ do
8:       Perform action $a_{t,i}$ in environment $e_i$
9:       Observe new state $s_{t+1,i}$ and reward $r_{t+1,i}$
10:     end parallel for
11:   end for
12: $R_{t_{max}+1} = \begin{cases} 
0 & \text{for terminal } s_t \\
V(s_{t_{max}+1}; \theta) & \text{for non-terminal } s_t 
\end{cases}$
13: for $t = t_{max}$ down to $1$ do
14:   $R_t = r_t + \gamma R_{t+1}$
15: end for
16: $d\theta = \frac{1}{n_e \cdot t_{max}} \sum_{i=1}^{n_e} \sum_{t=1}^{t_{max}} (R_{t,i} - v_{t,i}) \nabla_{\theta} \log \pi(a_{t,i} | s_{t,i}; \theta) + \beta \nabla_{\theta} H(\pi(s_{e,t}; \theta))$
17: $d\theta_v = \frac{1}{n_e \cdot t_{max}} \sum_{i=1}^{n_e} \sum_{t=1}^{t_{max}} \nabla_{\theta_v} (R_{t,i} - V(s_{t,i}; \theta_v))^2$
18: Update $\theta$ using $d\theta$ and $\theta_v$ using $d\theta_v$.
19: $N \leftarrow N + n_e \cdot t_{max}$
20: until $N \geq N_{max}$

Algorithm 1 of the paper "Efficient Parallel Methods for Deep Reinforcement Learning" by Alfredo V. Clemente et al.
The authors use 8 workers, $n_e = 32$ parallel environments, 5-step returns, $\gamma = 0.99$, $\varepsilon = 0.1$, $\beta = 0.01$ and a learning rate of $\alpha = 0.0007 \cdot n_e = 0.0224$.

The $\text{arch}_{\text{nips}}$ is from A3C: 16 filters $8 \times 8$ stride 4, 32 filters $4 \times 4$ stride 2, a dense layer with 256 units. The $\text{arch}_{\text{nature}}$ is from DQN: 32 filters $8 \times 8$ stride 4, 64 filters $4 \times 4$ stride 2, 64 filters $3 \times 3$ stride 1 and 512-unit fully connected layer. All nonlinearities are ReLU.

### Table 1

<table>
<thead>
<tr>
<th>Game</th>
<th>Gorila</th>
<th>A3C FF</th>
<th>GA3C</th>
<th>PAAC $\text{arch}_{\text{nips}}$</th>
<th>PAAC $\text{arch}_{\text{nature}}$</th>
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<td>22707.9</td>
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<td>Pong</td>
<td>18.3</td>
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<td>100523.3</td>
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</table>

Table 1 of the paper "Efficient Parallel Methods for Deep Reinforcement Learning" by Alfredo V. Clemente et al.
Figure 3 of the paper "Efficient Parallel Methods for Deep Reinforcement Learning" by Alfredo V. Clemente et al.
Figure 4 of the paper "Efficient Parallel Methods for Deep Reinforcement Learning" by Alfredo V. Clemente et al.
Parallel Advantage Actor Critic

Figure 2 of the paper “Efficient Parallel Methods for Deep Reinforcement Learning” by Alfredo V. Clemente et al.