Improving Exploration in Evolution Strategies for Deep Reinforcement Learning via a Population of Novelty-Seeking Agents

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Abstract

Evolution strategies (ES) are a family of black-box optimization algorithms able to train deep neural networks roughly as well as Q-learning and policy gradient methods on challenging deep reinforcement learning (RL) problems, but are much faster (e.g. hours vs. days) because they parallelize better. However, many RL problems require directed exploration because they have reward functions that are sparse or deceptive (i.e. contain local optima), and it is unknown how to encourage such exploration with ES. Here we show that algorithms that have been invented to promote directed exploration in small-scale evolved neural networks via populations of exploring agents, specifically novelty search (NS) and quality diversity (QD) algorithms, can be hybridized with ES to improve its performance on sparse or deceptive deep RL tasks, while retaining scalability. Our experiments confirm that the resultant new algorithms, NS-ES and two QD algorithms, NSR-ES and NSRA-ES, avoid local optima encountered by ES to achieve higher performance on Atari and simulated robots learning to walk around a deceptive trap. This paper thus introduces a family of fast, scalable algorithms for reinforcement learning that are capable of directed exploration. It also adds this new family of exploration algorithms to the RL toolbox and raises the interesting possibility that analogous algorithms with multiple simultaneous paths of exploration might also combine well with existing RL algorithms outside ES.

1 Introduction

In RL, an agent tries to learn to perform a sequence of actions in an environment that maximizes some notion of cumulative reward [1]. However, reward functions are often *deceptive*, and solely optimizing for reward without some mechanism to encourage intelligent exploration can lead to getting stuck in local optima and the agent failing to properly learn [1–3]. Unlike in supervised learning with deep neural networks (DNNs), wherein local optima are not thought to be a problem [4, 5], the training data in RL is determined by the actions an agent takes. If the agent greedily takes actions that maximize reward, the training data for the algorithm will be limited and it may not discover alternate strategies with larger payoffs (i.e. it can get stuck in local optima) [1–3]. Sparse reward signals can also be a problem for algorithms that only maximize reward, because at times there may be no reward gradient to follow. The possibility of deceptiveness and/or sparsity in the reward signal motivates the need for efficient and *directed* exploration, in which an agent is motivated to visit unexplored states in order to learn to accumulate higher rewards. Although deep RL algorithms have performed amazing feats in recent years [6–8], they have mostly done so despite relying on simple, *undirected* (aka dithering) exploration strategies, in which an agent hopes to explore new areas of its environment by taking random actions (e.g. epsilon-greedy exploration) [1].

A number of methods have been proposed to promote directed exploration in RL [9, 10], including recent methods that handle high-dimensional state spaces with deep neural networks. A common idea is to encourage an agent to visit states it has rarely or never visited (or take novel actions in

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those states). Methods proposed to track state (or state-action pair) visitation frequency include (1) approximating state visitation counts based on either auto-encoded latent codes of states [11] or pseudo-counts from state-space density models [12, 13], (2) learning a dynamics model that predicts future states (assuming predictions will be worse for rarely visited states/state-action pairs) [14–16], and (3) methods based on compression (novel states should be harder to compress) [9].

Those methods all count each state separately. A different approach to is to hand-design (or learn) an abstract, holistic description of an agent's lifetime of behavior, and then encourage the agent to exhibit different behaviors from those previously performed. That is the approach of novelty search (NS) [3] and quality diversity (QD) algorithms [17–19], which are described in detail below. Such algorithms are also interestingly different, and have different capabilities, because they perform exploration with a population of agents rather than a single agent (discussed in SI Sec. 6.2). NS and QD have shown promise with smaller neural networks on problems with low-dimensional input and output spaces [17–22]. Evolution strategies (ES) [23] was recently shown to perform well on high-dimensional deep RL tasks in a short amount of wall clock time by scaling well to many distributed computers. In this paper, for the first time, we study how these two types of algorithms can be hybridized with ES to scale them to deep neural networks and thus tackle hard, high-dimensional deep reinforcement learning problems, without sacrificing the speed/scalability benefits of ES. We first study NS, which performs exploration only (ignoring the reward function) to find a set of novel solutions [3]. We then investigate algorithms that balance exploration and exploitation, specifically novel instances of QD algorithms, which seek to produce a set of solutions that are both novel and high-performing [17–20]. Both NS and QD are explained in detail in Sec. 3.

ES directly searches in the parameter space of a neural network to find an effective policy. A team from OpenAI recently showed that ES can achieve competitive performance on many reinforcement learning (RL) tasks while offering some unique benefits over traditional gradient-based RL methods [24]. Most notably, ES is highly parallelizable, which enables near linear speedups in runtime as a function of CPU/GPU workers. For example, with hundreds of parallel CPUs, ES was able to achieve roughly the same performance on Atari games with the same DNN architecture in 1 hour as A3C did in 24 hours [24]. In this paper, we investigate adding NS and QD to ES only; in future work, we will investigate how they might be hybridized with Q-learning and policy gradient methods. We start with ES because (1) its fast wall-clock time allows rapid experimental iteration, and (2) NS and QD were originally developed as neuroevolution methods, making it natural to try them first with ES, which is also an evolutionary algorithm.

Here we test whether encouraging novelty via NS and QD improves the performance of ES on sparse and/or deceptive control tasks. Our experiments confirm that NS-ES and two simple versions of QD-ES (NSR-ES and NSRA-ES) avoid local optima encountered by ES and achieve higher performance on tasks ranging from simulated robots learning to walk around a deceptive trap to the high-dimensional pixel-to-action task of playing Atari games. Our results add these new families of exploration algorithms to the RL toolbox, opening up avenues for studying how they can best be combined with RL algorithms, whether ES or others.

2 Background

2.1 Evolution Strategies

Evolution strategies (ES) are a class of black box optimization algorithms inspired by natural evolution [23]: At every iteration (generation), a population of parameter vectors (genomes) is perturbed (mutated) and, optionally, recombined (merged) via crossover. The fitness of each resultant offspring is then evaluated according to some objective function (reward) and some form of selection then ensures that individuals with higher reward tend to produce offspring for the next generation. Many algorithms in the ES class differ in their representation of the population and methods of recombination; the algorithms subsequently referred to in this work belong to the class of Natural Evolution Strategies (NES) [25, 26]. NES represents the population as a distribution of parameter vectors θ characterized by parameters ϕ : $p_{\phi}(\theta)$. Under a fitness function, $f(\theta)$, NES seeks to maximize the average fitness of the population, $\mathbb{E}_{\theta \sim p_{\phi}}[f(\theta)]$, by optimizing ϕ with stochastic gradient ascent.

Recent work from OpenAI outlines a version of NES applied to standard RL benchmark problems [24]. We will refer to this variant simply as ES going forward. In their work, a fitness function $f(\theta)$ represents the stochastic reward experienced over a full episode of agent interaction, where θ is the parameters of a policy π_{θ} . From the population distribution p_{ϕ_t} , parameters $\theta_t^i \sim \mathcal{N}(\theta_t, \sigma^2 I)$ are

sampled and evaluated to obtain $f(\theta_i^i)$. In a manner similar to REINFORCE [27], θ_t is updated using an estimate of approximate gradient of expected reward:

$$\nabla_{\phi} \mathbb{E}_{\theta \sim \phi}[f(\theta)] \approx \frac{1}{n} \sum_{i=1}^{n} f(\theta_{t}^{i}) \nabla_{\phi} \log p_{\phi}(\theta_{t}^{i})$$

where n is the number of samples evaluated per generation. Intuitively, NES samples parameters in the neighborhood of θ_t and determines the direction in which θ_t should move to improve expected reward. Since this gradient estimate has high variance, NES relies on a large n for variance reduction. Generally, NES also evolves the covariance of the population distribution, but for the sake of fair comparison with Salimans et al. [24] we consider only static covariance distributions, meaning σ is fixed throughout training.

To sample from the population distribution, Salimans et al. [24] apply additive Gaussian noise to the current parameter vector: $\theta_t^i = \theta_t + \sigma \epsilon_i$ where $\epsilon_i \sim \mathcal{N}(0, I)$. The gradient is then estimated by taking a sum of sampled parameter perturbations weighted by their reward:

$$\nabla_{\theta_t} \mathbb{E}_{\epsilon \sim \mathcal{N}(0, I)}[f(\theta_t + \sigma \epsilon)] \approx \frac{1}{n\sigma} \sum_{i=1}^n f(\theta_t^i) \epsilon_i$$

To ensure that the scale of reward between domains does not bias the optimization process, we follow the approach of Salimans et al. [24] and rank-normalize $f(\theta_t^i)$ before taking the weighted sum. Overall, this NES variant exhibits performance on par with contemporary, gradient-based algorithms on difficult RL domains, including simulated robot locomotion and Atari environments [28].

2.2 Novelty Search (NS)

Optimizing for reward only can often lead an agent to local optima. NS, however, avoids deception in the reward signal by ignoring reward altogether. Inspired by nature's drive towards diversity, NS encourages policies to engage in notably different behaviors than those previously seen. The algorithm encourages different behaviors by computing the *novelty* of the current policy with respect to previously generated policies and then encourages the population distribution to move towards areas of parameter space with high novelty. NS outperforms reward-based methods in maze and biped walking domains, which possess deceptive reward signals that attract agents to local optima [3]. In this work, we investigate the efficacy of NS at the scale of DNNs by combining it with ES. In NS, a policy π is assigned a domain-dependent *behavior characterization* $b(\pi)$ that describes its behavior. For example, in the case of a humanoid locomotion problem, $b(\pi)$ may be as simple as a two-dimensional vector containing the humanoid's final $\{x,y\}$ location. Throughout training, every π_{θ} evaluated adds a behavior characterization $b(\pi_{\theta})$ to an archive set A with some probability. A particular policy's novelty $N(b(\pi_{\theta}), A)$ is then computed by selecting the k-nearest neighbors of $b(\pi_{\theta})$ from A and computing the average distance between them:

$$N(\theta, A) = N(b(\pi_{\theta}), A) = \frac{1}{|S|} \sum_{j \in S} ||b(\pi_{\theta}) - b(\pi_{j})||_{2}$$
$$S = kNN(b(\pi_{\theta}), A)$$
$$= \{b(\pi_{1}), b(\pi_{2}), ..., b(\pi_{k})\}$$

Above, the distance between behavior characterizations is calculated with an L2-norm, but any distance function can be substituted. Previously, NS has been implemented with a genetic algorithm [3]. We next explain how NS can now be combined with ES, to leverage the advantages of both.

3 Methods

3.1 NS-ES

We use the ES optimization framework, described in Sec. 2.1, to compute and follow the gradient of expected novelty with respect to θ_t . Given an archive A and sampled parameters $\theta_t^i = \theta_t + \sigma \epsilon_i$, the gradient estimate can be computed:

$$\nabla_{\theta_t} \mathbb{E}_{\epsilon \sim \mathcal{N}(0, I)}[N(\theta_t + \sigma \epsilon, A)|A] \approx \frac{1}{n\sigma} \sum_{i=1}^n N(\theta_t^i, A) \epsilon_i$$

The gradient estimate obtained tells us how to change the current policy's parameters θ_t to increase the average novelty of our parameter distribution. We condition the gradient estimate on A, as the archive is fixed at the beginning of a given iteration and updated only at the end. We add only the behavior characterization corresponding to each θ_t , as adding those for each sample θ_t^i would inflate the archive and slow the nearest-neighbors computation. As more behavior characterizations are added to A, the novelty landscape changes, resulting in commonly occurring behaviors becoming "boring." Optimizing for expected novelty leads to policies that move towards unexplored areas of behavior space.

NS-ES could operate with a single agent that is rewarded for acting differently than its ancestors. However, to encourage additional diversity and get the benefits of population-based exploration described in SI Sec. 6.2, we can instead create a population of M agents, which we will refer to as the *meta-population*. Each agent, characterized by a unique θ^m , is rewarded for being different from all prior agents in the archive (ancestors, other agents, and the ancestors of other agents). In this paper, we have multiple agents in the meta-population (i.e. M > 1), but we did not conduct a thorough analysis on how varying this hyperparameter affects performance on different domains. We hypothesize that the selection of M is domain dependent and that identifying which domains favor which regime is a fruitful area for future research.

We initialize M random parameter vectors and at every iteration select one to update. For our experiments, we probabilistically select which θ^m to advance from a discrete probability distribution as a function of θ^m 's novelty. Specifically, at every iteration, for a set of agent parameter vectors $\Pi = \{\theta^1, \theta^2, ..., \theta^M\}$, we calculate each θ^m 's probability of being selected $P(\theta^m)$ as its novelty normalized by the sum of novelty across all policies:

$$P(\theta^{m}) = \frac{N(\theta^{m}, A)}{\sum_{j=1}^{M} N(\theta^{j}, A)}$$
(1)

Having multiple, separate agents represented as independent Gaussians is a simple choice for the meta-population distribution (i.e. how the meta-population distribution is represented). In future work, more complex sampling distributions that represent the multi-modal nature of meta-population parameter vectors could be tried.

After selecting an individual m from the meta-population, we compute the gradient of expected novelty with respect to m's current parameter vector, θ_t^m , and perform an update step accordingly:

$$\theta_{t+1}^m \leftarrow \theta_t^m + \alpha \frac{1}{n\sigma} \sum_{i=1}^n N(\theta_t^{i,m}, A) \epsilon_i$$

Where n is the number of sampled perturbations to θ^m_t , α is the stepsize, and $\theta^{i,m}_i = \theta^m_t + \sigma \epsilon_i$, where $\epsilon_i \sim \mathcal{N}(0, I)$. Once the current parameter vector is updated, $b(\pi_{\theta^m_{t+1}})$ is computed and added to the shared archive A with probability 1. The whole process is repeated for a pre-specified number of iterations, as there is no true convergence point of NS. In this work, the algorithm simply returns the highest-performing parameter vector found. Algorithm 1 in SI Sec. 6.4 outlines a simple, parallel implementation of NS-ES. It is important to note that the addition of the archive and the replacement of the fitness function with novelty does not damage the scalability of the ES optimization procedure (SI Sec. 6.3).

3.2 QD-ES Algorithms: NSR-ES and NSRA-ES

NS-ES alone can enable agents to avoid deceptive local optima in the reward function. Reward signals, however, are still very informative and discarding them completely may cause performance to suffer. Consequently, we train a variant of NS-ES, which we call NSR-ES, that combines the reward ("fitness") and novelty calculated for a given set of policy parameters θ . Similar to NS-ES and ES, NSR-ES operates on entire episodes and can thus evaluate reward and novelty simultaneously for any sampled parameter vector: $\theta_t^{i,m} = \theta_t^m + \epsilon_i$. Specifically, we compute $f(\theta_t^{i,m})$ and $N(\theta_t^{i,m}, A)$, average the two values, and set the average as the weight for the corresponding ϵ_i . The averaging process is integrated into the parameter update rule as: $\theta^m_{t+1} \leftarrow \theta^m_t + \alpha \frac{1}{n\sigma} \sum_{i=1}^n \frac{f(\theta^{i,m}_t) + N(\theta^{i,m}_t,A)}{2} \epsilon_i$

$$\theta_{t+1}^m \leftarrow \theta_t^m + \alpha \frac{1}{n\sigma} \sum_{i=1}^n \frac{f(\theta_t^{i,m}) + N(\theta_t^{i,m}, A)}{2} \epsilon_i$$

Intuitively, the algorithm follows the approximated gradient in parameter-space towards policies that both exhibit novel behaviors and achieve high rewards. Often, however, the scales of $f(\theta)$ and $N(\theta, A)$ differ. To combine the two signals effectively, we rank-normalize $f(\theta_t^{i,m})$ and $N(\theta_t^{i,m}, A)$ independently before computing the average. The algorithm is a QD algorithm because it has a set of M agents being optimized to be both high-performing, yet different from each other.

NSR-ES has an equal weighting of the performance and novelty gradients that is static across training. We explore a further extension of NSR-ES called NSR Adapt-ES (NSRA-ES), which takes advantage of the opportunity to dynamically weight the priority given to the performance gradient $f(\theta_t^{i,m})$ vs. the novelty gradient $N(\theta_t^{i,m}, A)$ by intelligently adapting a weighting parameter w during training. By doing so, the algorithm can follow the performance gradient when it is making progress, increasingly try different things if stuck in a local optimum, and switch back to following the performance gradient once unstuck. For a specific w at a given generation, the parameter update rule for NSRA-ES is expressed as follows:

$$\theta_{t+1}^m \leftarrow \theta_t^m + \alpha \frac{1}{n\sigma} \sum_{i=1}^n w f(\theta_t^{i,m}) \epsilon_i + (1-w) N(\theta_t^{i,m}, A) \epsilon_i$$

We set w=1.0 initially and decrease it if performance stagnates across a fixed number of generations. We continue decreasing w until performance increases, at which point we increase w. SI Sec. 6.4 provides a more detailed description of how we adapt w as well as pseudocode for NSR-ES and NSRA-ES. Source code and hyperparameter configurations for all of our experiments are in the SI.

4 Experiments

4.1 Simulated Humanoid Locomotion problem

We first tested our implementation of NS-ES, NSR-ES, and NSRA-ES on the problem of having a simulated humanoid learn to walk. We chose this problem because it is a challenging continuous control benchmark where most would presume a reward function is necessary to solve the problem. With NS-ES, we test whether searching through novelty alone can find solutions to the problem. A similar result has been shown for much smaller neural networks (~50-100 parameters) on a more simple simulated biped [20], but here we test whether NS-ES can enable locomotion at the scale of deep neural networks on a much more sophisticated environment. NSR-ES and NSRA-ES experiments then test the effectiveness of combining exploration and reward pressures on this difficult continuous control problem. SI Sec. 6.6 has complete experimental details.

The first experiment is in a slightly modified version of OpenAI Gym's Humanoid-v1 environment. Because the heart of this challenge is to learn to walk efficiently, not to walk in a particular direction, we modified the environment reward to be isotropic (i.e. indifferent to the direction the humanoid traveled) by setting the velocity component of reward to distance traveled from the origin as opposed to distance traveled in the positive x direction.

As described in section 2.2, novelty search requires a domain-specific behavior characterization (BC) for each policy, which we denote as $b(\pi_{\theta_i})$. For the Humanoid Locomotion problem the BC is the agent's final $\{x,y\}$ location, as it was in Lehman and Stanley [20]. NS also requires a distance function between two BCs. Following Lehman & Stanley 2011 [20], the distance function is the square of the Euclidean distance:

$$dist(b(\pi_{\theta_i}), b(\pi_{\theta_i})) = ||b(\pi_{\theta_i}) - b(\pi_{\theta_i})||_2^2$$

The first result is that ES obtains a higher final reward than NS-ES (p < 0.05) and NSR-ES (p < 0.05); these and all future p values are via a Mann-Whitney U test. The performance gap is more pronounced for smaller amounts of computation (Fig. 1 (b)). However, many will be surprised that NS-ES is still able to consistently solve the problem despite ignoring the environment's multi-part reward function. While the BC is aligned [29] with the problem in that reaching new $\{x,y\}$ positions tends to also encourage walking, there are many parts of the reward function that the BC ignores (e.g. energy-efficiency, impact costs).

We hypothesize that with a sophisticated BC that encourages diversity in all of the behaviors the multi-part reward function cares about, there would be no performance gap. However, such a BC may be difficult to construct and would likely further exaggerate the amount of computation required for NS to match ES. NSR-ES demonstrates faster learning than NS-ES due to the addition of reward pressure, but ultimately results in similar final performance after 600 generations (p > 0.05, Fig. 1 (b)). Promisingly, on this non-deceptive problem, NSRA-ES does not pay a cost for its latent exploration capabilities and performs similarly to ES (p > 0.05).

The Humanoid Locomotion problem does not appear to be a deceptive problem, at least for ES. To test whether NS-ES, NSR-ES, and NSR-ES specifically help with deception, we also compare ES to these algorithms on a variant of this environment we created that adds a deceptive trap (a local optimum) that must be avoided for maximum performance (Fig. 1, (c)). In this new environment, a small three-sided enclosure is placed at a short distance in front of the starting position of the humanoid and the reward function is simply distance traveled in the positive \boldsymbol{x} direction.

Fig. 1 (d) and SI Sec. 6.7 show the reward received by each algorithm and Fig. 2 shows how the algorithms differ qualitatively during search on this problem. In every run, ES gets stuck in the local optimum due to following reward into the deceptive trap. NS-ES is able to avoid the local optimum as it ignores reward completely and instead seeks to thoroughly explore the environment, but doing so also means it makes slow progress according to the reward function. NSR-ES demonstrates superior performance to NS-ES (p < 0.01) and ES (p < 0.01) as it benefits from both optimizing for reward and escaping the trap via the pressure for novelty. Like ES, NSRA-ES learns to walk into the

deceptive trap initially, as it initially is optimizing for reward only. Once stuck in the local optimum, the algorithm continually increases its pressure for novelty, allowing it to escape the deceptive trap and ultimately achieve much higher rewards than NS-ES (p < 0.01) and NSR-ES (p < 0.01). Based just on these two domains, NSRA-ES seems to be the best algorithm across the board because it can exploit well when there is no deception, add exploration dynamically when there is, and return to exploiting once unstuck. The latter is likely why NSRA-ES outperforms even NSR-ES on the deceptive humanoid locomotion problem.

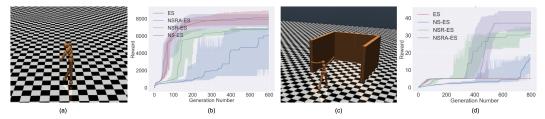


Figure 1: **Humanoid Locomotion Experiment.** The humanoid locomotion task without a deceptive trap (a) and with one (c), and results on them in (b) and (d), respectively. Here and in similar figures below, the median reward (of the best seen policy so far) per generation across 10 runs is plotted as the bold line with 95% bootstrapped confidence intervals of the median (shaded). Following [24], policy performance is measured as the average performance over \sim 30 stochastic evaluations.

Fig. 2 also shows the benefit of maintaining a meta-population (M=5) in the NS-ES, NSR-ES, and NSRA-ES algorithms. Some lineages get stuck in the deceptive trap, incentivizing other policies to explore around the trap. At that point, all three algorithms begin to allocate more computational resources to this newly discovered, more promising strategy via the probabilistic selection method outlined in Sec. 3.1. Both the novelty pressure and having a meta-population thus appear to be useful, but in future work we look to disambiguate the relative contribution made by each.



Figure 2: ES gets stuck in the deceptive local optimum while NS-ES, NSR-ES & NSRA-ES explore to find better solutions. An overhead view of a representative run is shown for each algorithm on the Humanoid Locomotion with Deceptive Trap problem. The black star represents the humanoid's starting point. Each diamond represents the final location of a generation's policy, i.e. $\pi(\theta_t)$, with darker shading for later generations. For NS-ES, NSR-ES, & NSRA-ES plots, each of the M=5 agents in the meta-population and its descendants are represented by different colors. Similar plots for all 10 runs of each algorithm are provided in SI 6.9.

4.2 Atari

We also tested NS-ES, NSR-ES, and NSRA-ES on numerous games from the Atari 2600 environment in OpenAI Gym [30]. Atari games serve as an informative benchmark due to their high-dimensional pixel input and complex control dynamics; each game also requires different levels of exploration to solve. To demonstrate the effectiveness of NS-ES, NSR-ES, and NSRA-ES for local optima avoidance and directed exploration, we tested on 12 different games with varying levels of complexity, as defined by the taxonomy in [12]. Primarily, we focused on games in which, during preliminary experiments, we observed ES prematurely converging to local optima (Seaquest, Q*Bert, Freeway, Frostbite, and Beam Rider). However, we also included a few other games where ES did not converge to local optima to understand the performance of our algorithm in less-deceptive domains (Alien, Amidar, Bank Heist, Gravitar, Zaxxon, and Montezuma's Revenge). SI Sec. 6.5 describes additional experimental details. We report the median reward across r independent runs of the best policy found in each run (see Fig. 3, left).

For the behavior characterization, we follow an idea from Naddaf [31] and concatenate Atari game RAM states for each timestep in an episode. RAM states in Atari 2600 games are integer-valued vectors of length 128 in the range [0, 255] that describe all the state variables in a game (e.g. the location of the agent and enemies). Ultimately, we want to automatically learn behavior characterizations directly from pixels. A plethora of recent research suggests that this is a viable approach [12, 32, 33]. For example, low-dimensional, latent representations of the state space could be extracted from auto-encoders [11, 34] or networks trained to predict future states [14, 16]. In this work, however, we focus on learning with a pre-defined, informative behavior characterization and leave the task of jointly learning a policy and latent representation of states for future work. In effect, basing novelty on RAM states provides a confirmation of what is possible in principle with a sufficiently informed behavior characterization. We also emphasize that, while during training NS-ES, NSR-ES, and NSRA-ES use RAM states to guide novelty search, the policy itself, π_{θ_t} , operates only on image input and can be evaluated without any RAM state information. The distance between behavior characterizations is the sum of L2-distances at each timestep t:

$$dist(b(\pi_{\theta_i}), b(\pi_{\theta_j})) = \sum_{t=1}^{T} ||(b_t(\pi_{\theta_i})) - b_t(\pi_{\theta_j}))||_2$$

Fig. 3, left compares the performance of each algorithm discussed above to each other and with those from two previously works that promote exploration in RL, namely Noisy DQN [35] and A3C+ [12]. Although Noisy DQN and A3C+ are popular methods for exploration in RL, they only outperform all the ES variants considered in this paper on 3/12 games and 2/12 games respectively. NSRA-ES, however, outperforms the other algorithms on 5/12 games, suggesting that NS and QD are viable alternatives to contemporary exploration methods.

While the novelty pressure in NS-ES does help it avoid local optima in some cases (discussed below), optimizing for novelty alone does not result in higher reward in most games (although it does in some). However, it is surprising how well NS-ES does in many tasks given that it is not explicitly attempting to increase reward. Because NSR-ES combines exploration with reward maximization, it is able to avoid local optima encountered by ES while also learning to play the game well. In each of the 5 games in which we observed ES converge to premature local optima (i.e. Seaquest, Q*Bert, Freeway, Beam Rider, Frostbite), NSR-ES achieves a higher median reward. In the other games, ES does not benefit from adding an exploration pressure and NSR-ES performs worse. It is expected that if there are no local optima and reward maximization is sufficient to perform well, the extra cost of encouraging exploration will hurt performance. Mitigating such costs, NSRA-ES optimizes solely for reward until a performance plateau is reached. After that, the algorithm will assign more weight to novelty and thus encourage exploration. We found this to be beneficial, as NSRA-ES achieves higher median rewards than ES on 8/12 games and NSR-ES on 9/12 games. It's superior performance validates NSRA-ES as the best among the evolutionary algorithms considered and suggests that using an *adaptive* weighting between novelty and reward is a promising direction for future research.

In the game Seaquest, the avoidance of local optima is particularly important (Fig. 3, right). ES performance flatlines early at a median reward of 960, which corresponds to a behavior of the agent descending to the bottom, shooting fish, and never coming up for air. This strategy represents a classic local optima, as coming up for air requires temporarily foregoing reward, but enables far higher rewards to be earned in the long run (Salimans et al. [24] did not encounter this particular local optimum with their hyperparameters, but the point is that ES without exploration can get stuck indefinitely on whichever major local optima it encounters). NS-ES learns to come up for air in all 5 runs and achieves a slightly higher median reward of 1044.5 (p < 0.05). NSR-ES also avoids this local optimum, but its additional reward signal helps it play the game better (e.g. it is better at shooting enemies), resulting in a significantly higher median reward of 2329.7 (p < 0.01). Because NSRA-ES takes reward steps initially, it falls into the same local optimum as ES. Because we chose (without performing a hyperparameter search) to change the weighting w between performance and novelty infrequently (only every 50 generations), and to change it by a small amount (only 0.05), 200 generations was not long enough to emphasize novelty enough to escape this local optimum. We found that by changing w every 10 generations, this problem is remedied and the performance of NSRA-ES equals that of NSR-ES (p > 0.05, Fig. 3, right). These results motivate future research into better hyperparameters for changing w, and into more complex, intelligent methods of dynamically adjusting w, including with a population of agents with different dynamic w strategies.

The Atari results illustrate that NS is an effective mechanism for encouraging directed exploration, given an appropriate behavior characterization, for complex, high-dimensional control tasks. A novelty pressure alone produces impressive performance on many games, sometimes even beating

ES. Combining novelty and reward performs far better, and improves ES performance on tasks where it appears to get stuck on local optima.

| Game | ES | NS-ES | NSR-ES | NSRA-ES | DQN | Noisy DQN | A3C+ | 2500 | — ES |
|-------------------------|--------|--------|--------|---------|-------|-----------|---------|--------|---|
| Alien | 3283.8 | 1124.5 | 2186.2 | 4846.4 | 2404 | 2403 | 1848.3 | 2500 | — NS-ES |
| Amidar | 322.2 | 134.7 | 255.8 | 305.0 | 924 | 1610 | 964.7 | | - NSR-ES |
| Bank Heist | 140.0 | 50.0 | 130.0 | 152.9 | 455 | 1068 | 991.9 | 2000 | - NSRA-ES 50 gens |
| Beam Rider [†] | 871.7 | 805.5 | 876.9 | 906.4 | 10564 | 20793 | 5992.1 | | — NSRA-ES 10 gens |
| Freeway [†] | 31.1 | 22.8 | 32.3 | 32.9 | 31 | 32 | 27.3 | 1500 g | |
| Frostbite [†] | 367.4 | 250.0 | 2978.6 | 3785.4 | 1000 | 753 | 506.6 | ewa | |
| Gravitar | 1129.4 | 527.5 | 732.9 | 1140.9 | 366 | 447 | 246.02 | 1000 | |
| Montezuma's Revenge | 0.0 | 0.0 | 0.0 | 0.0 | 2 | 3 | 142.5 | 1000 | |
| Ms. Pacman | 4998.0 | 2252.2 | 3495.2 | 5171.0 | 2674 | 2722 | 2380.6 | | |
| Q*Bert [†] | 1075.0 | 1234.1 | 1400.0 | 1350.0 | 11241 | 15545 | 15804.7 | 500 | |
| Seaquest [†] | 960.0 | 1044.5 | 2329.7 | 960.0 | 4163 | 2282 | 2274.1 | | |
| Zaxxon | 9885.0 | 1761.9 | 6723.3 | 7303.3 | 4806 | 6920 | 7956.1 | 0 | 7 |
| | | | | | | | | | 0 25 50 75 100 125 150 175 200 Generation Number |

Figure 3: Atari Results:Left: The scores are the median, across 5 runs, of the mean reward (over $\sim\!\!30$ stochastic evaluations) of each run's final best policy. SI Sec. 6.8 plots performance over time, along with bootstrapped confidence intervals of the median, for each ES algorithm for each game. In some cases rewards reported here for ES are lower than those in Salimans et al. [24], which could be due to differing hyperparameters (SI Sec. 6.5). Games with a † are those in which we observed ES to converge prematurely, presumably due to it encountering local optima. The DQN and A3C results are reported after 200M frames of, and one to many days of, training. All evolutionary algorithm results are reported after \sim 2B frames of , and $\sim\!\!2$ hours of, training

5 Discussion and Conclusion

NS and QD are classes of evolutionary algorithms designed to avoid local optima and promote exploration in RL environments, but have only been previously shown to work with small neural networks (on the order of hundreds of connections). ES was recently shown to be capable of training deep neural networks that can solve challenging, high-dimensional RL tasks [24]. It also is much faster when many parallel computers are available. Here we demonstrate that, when hybridized with ES, NS and QD not only preserve the attractive scalability properties of ES, but also help ES explore and avoid local optima in domains with deceptive reward functions. To the best of our knowledge, this paper reports the first attempt at augmenting ES to perform directed exploration in high-dimensional environments. We thus provide an option for those interested in taking advantage of the scalability of ES, but who also want higher performance on domains that have reward functions that are sparse or have local optima. The latter scenario will likely hold for most challenging, real-world domains that machine learning practitioners will wish to tackle in the future.

Additionally, this work highlights alternate options for exploration in RL domains. The first is to holistically describe the behavior of an agent instead of defining a per-state exploration bonus. The second is to encourage a population of agents to simultaneously explore different aspects of an environment. These new options thereby open new research areas into (1) comparing holistic vs. state-based exploration, and population-based vs. single-agent exploration, more systematically and on more domains, (2) investigating the best way to combine the merits of all of these options, and (3) hybridizing holistic and/or population-based exploration with other algorithms that work well on deep RL problems, such as policy gradients and DQN. It should be relatively straightforward to combine NS with policy gradients (NS-PG). It is less obvious how to combine it with Q-learning (NS-Q), but may be possible.

As with any exploration method, encouraging novelty can come at a cost if such an exploration pressure is not necessary. In Atari games such as Alien and Gravitar, and in the Humanoid Locomotion problem without a deceptive trap, both NS-ES and NSR-ES perform worse than ES. To avoid this cost, we introduce the NSRA-ES algorithm, which attempts to invest in exploration only when necessary. NSRA-ES tends to produce better results across many different domains than ES, NS-ES, and NSR-ES, making it an attractive new algorithm for deep RL tasks. Similar strategies for adapting the amount of exploration online may also be advantageous for other deep RL algorithms. How best to dynamically balance between exploitation and exploration in deep RL remains an open, critical research challenge, and our work underscores the importance of, and motivates further, such work. Overall, our work shows that ES is a rich and unexploited parallel path for deep RL research. It is worthy of exploring not only because it is an alternative algorithm for RL problems, but also because innovations created in the ES family of algorithms could be ported to improve other deep RL algorithm families like policy gradients and Q learning, or through hybrids thereof.

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6 Supplementary Information

6.1 Videos of agent behavior

Videos of example agent behavior in all the environments can be viewed here: https://goo.gl/cVUG2U.

6.2 Population-based exploration vs. single-agent exploration

As mentioned in the introduction, aside from being holistic vs. state-based, there is another interesting difference between most exploration methods for RL [12–16] and the NS/QD family of algorithms. We do not investigate the benefits of this difference experimentally in this paper, but they are one reason we are interested in NS/QD as an exploration method for RL. One commonality among the previous methods is that the exploration is performed by a *single* agent, a choice that has interesting consequences for learning. To illustrate these consequences, we borrow an example from Stanton and Clune [36]. Imagine a cross-shaped maze (Fig. 4) where to go in each cardinal direction an agent must master a different skill (e.g. going north requires learning to swim, west requires climbing mountains, east requires walking on sand, and south requires walking on ice). Assume rewards may or may not exist at the end of each corridor, so all corridors need to be explored. A single agent has two extreme options, either a depth-first search that serially learns to go to the end of each corridor, or a breadth-first search that goes a bit further in one direction, then comes back to the center and goes a bit further in another direction, etc. Either way, to get to the end of each hallway, the agent will have to at least have traversed each hallway once and thus will have to learn all four sets of skills. With the breadth-first option, all four skillsets must be mastered, but a much longer total distance is traveled.

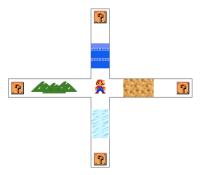


Figure 4: **Hypothetical Hard Exploration Maze.** In this maze, the agent needs to traverse 4 different terrains to obtain rewards associated with the "?" boxes. Traversing each terrain requires learning a certain skill (i.e. climbing, swimming, etc.). The sprites are from a Super Mario Bros. environment introduced by [37].

In both cases, another problem arises because, despite recent progress [38, 39], neural networks still suffer from catastrophic forgetting, meaning that as they learn new skills they rapidly lose the ability to perform previously learned ones [40]. Due to catastrophic forgetting, at the end of learning there will be an agent specialized in one of the skills (e.g. swimming), but all of the other skills will have been lost. Furthermore, if any amount of the breadth-first search strategy is employed, exploring each branch a bit further will require relearning that skill mostly from scratch each iteration, significantly slowing exploration. Even if catastrophic forgetting could be solved, there may be limits on the cognitive capacity of single agents (as occurs in humans), preventing one agent from mastering all possible skills.

A different approach is to explore with a *population* of agents. In that case, separate agents could become experts in the separate tasks required to explore in each direction. That may speed learning because each agent can, in parallel, learn only the skills required for its corridor. Additionally, at the end of exploration a specialist will exist with each distinct skill (versus only one skill remaining in the single-agent case). The resulting population of specialists, each with a different skill or way of solving a problem, can then be harnessed by other machine learning algorithms that efficiently search through the repertoire of specialists to find the skill or behavior needed in a particular situation [17, 41]. The

skills of each specialist (in any combination or number) could also then be combined into a single generalist via policy distillation [42]. A further benefit of the population-based approach is, when combining exploration with some notion of quality (e.g. maximizing reward), a population can try out many different strategies/directions and, once one or a few promising strategies are found, the algorithm can reallocate resources to pursue them. The point is not that population-based exploration methods are better or worse than single-agent exploration methods (when holding computation constant), but instead that they are a different option with different capabilities, pros, and cons, and are thus worth investigating [36]. Supporting this view, recent work has demonstrated the benefits of populations for deep learning [43, 44].

6.3 Preserving scalability

As shown in Salimans et al. [24], ES scales well with the amount of computation available. Specifically, as more CPUs are used, training times reduce almost linearly, whereas DQN and A3C are not amenable to massive parallelization. NS-ES, NSR-ES and NSRA-ES, however, enjoy the same parallelization benefits as ES because they use an almost identical optimization process. The addition of an archive between agents in the meta-population does not hurt scalability because A is only updated after θ_t^m has been updated. Since A is kept fixed during the calculation of $N(\theta_t^{i,m},A)$ and $f(\theta_t^{i,m})$ for all i=1...n perturbations, the coordinator only needs to broadcast A once at the beginning of each generation. In all algorithms, the parameter vector θ_t^i must be broadcast at the beginning of each generation and since A generally takes up much less memory than the parameter vector, broadcasting both would incur effectively zero extra network overhead. NS-ES, NSR-ES, and NSRA-ES do however introduce an additional computation conducted on the coordinator node. At the start of every generation we must compute the novelty of each candidate $\theta_t^m; m \in \{1,...,M\}$. For an archive of length n this operation is O(Mn), but since M is small and fixed throughout training this cost is not significant in practice. Additionally, there are methods for keeping the archive small if this computation becomes an issue [45].

6.4 NS-ES, NSR-ES, and NSRA-ES Algorithms

Algorithm 1 NS-ES

```
1: Input: learning rate \alpha, noise standard deviation \sigma, number of policies to maintain M, iterations T, behavior characterization b(\pi_{\theta})
 2: Initialize: M randomly initialized policy parameter vectors \{\theta_0^1, \theta_0^2, ..., \theta_0^M\}, archive A, number of workers n
        Compute b(\pi_{\theta_0^j})
         Add b(\pi_{\theta_0^j}) to A
 5:
    for t=0 to T-1 do
          Sample \theta_t^m from \{\theta_t^1, \theta_t^2, \dots, \theta_t^M\} via eq.1
          for i = 1 to n do
              Sample \epsilon_i \sim \mathcal{N}(0, \sigma^2 I)
Compute \theta_t^{i,m} = \theta_t^m + \epsilon_i
10:
11:
               Compute b(\pi_{\theta_{+}^{i},m})
12:
               Compute N_i = N(\theta_t^{i,m}, A)
13:
               Send N_i from each worker to coordinator
14:
15:
          Set \theta_{t+1}^m = \theta_t^m + \alpha \frac{1}{n\sigma} \sum_{i=1}^n N_i \epsilon_i
16:
         Compute b(\pi_{\theta_{t+1}^m})
17:
18:
          Add b(\pi_{\theta_{t+1}^m}) to A
19: end for
```

For the NSRA-ES algorithm, we introduce 3 new hyperparameters: w, t_w , δ_w . These quantities determine the agent's preference for novelty or reward at various phases in training. For all of our experiments, w=1.0 initially, meaning that NSRA-ES initially follows the gradient of reward only. If the best episodic reward seen, f_{best} , does not increase in $t_w=50$ generations, w is decreased by $\delta_w=0.05$ and gradients with respect to the new weighted average of novelty and reward are followed. The process is repeated until a new, higher f_{best} is found, at which point w is increased by δ_w . Intuitively the algorithm follows reward gradients until the increases in episodic reward plateau, at which point the agent is increasingly encouraged to explore. The agent will continue to explore until a promising behavior (i.e. one with higher reward than seen so far) is found. NSRA-ES then increases w to pivot back towards exploitation instead of exploration.

Algorithm 2 NSR-ES

```
1: Input: learning rate \alpha, noise standard deviation \sigma, number of policies to maintain M, iterations T, behavior characterization b(\pi_{\theta})
 2: Initialize: M sets of randomly initialized policy parameters \{\theta_0^1, \theta_0^2, ..., \theta_0^M\}, archive A, number of workers n
            Compute b(\pi_{\theta_0^j})
            Add b(\pi_{\theta_0^j}) to A
 5:
 6: end for
 7: for t=0 to T-1 do
            Sample \theta_t^m from \{\theta_t^0, \theta_t^1, \dots, \theta_t^M\} via eq. 1 for i=1 to n do
                  \begin{array}{l} \mathbf{a} \cdot \mathbf{b} = \mathbf{b} \cdot \mathbf{h} \cdot \mathbf{d} \mathbf{0} \\ \text{Sample } \epsilon_i \sim \mathcal{N}(0, \sigma^2 I) \\ \text{Compute } \theta_t^{i,m} = \theta_t^m + \epsilon_i \\ \text{Compute } b(\pi_{\theta_t^{i,m}}) \end{array}
10:
11:
12:
                   Compute N_i = N(\theta_t^{i,m}, A)
13:
                   Compute F_i = f(\hat{\theta}_t^{i,m})
14.
                   Send N_i and F_i from each worker to coordinator
15:
16:
             Set \theta_{t+1}^m = \theta_t^m + \alpha \frac{1}{n\sigma} \sum_{i=1}^n \frac{N_i + F_i}{2} \epsilon_i
17:
18:
            Compute b(\pi_{\theta_{t+1}^m})
            \operatorname{Add} b(\pi_{\theta_{t+1}^{\,m}}) \text{ to } A
19:
20: end for
```

Algorithm 3 NSRA-ES

```
1: Input: learning rate \alpha, noise standard deviation \sigma, number of policies to maintain M, iterations T, behavior characterization b(\pi_{\theta})
2: Initialize: M sets of randomly initialized policy parameters \{\theta_0^1, \theta_0^2, ..., \theta_0^M\}, archive A, number of workers n, initial weight w, weight
       tune frequency t_w, weight delta \delta_w
 3: for j = 1 to M do
            Compute b(\pi_{\theta_0^j})
             Add b(\pi_{\theta_0^j}) to A
 6: end for
 7: f_{best} = -\infty
 8: t_{best} = 0
9: for t = 0 to T - 1 do
            Sample \theta_t^m from \{\theta_t^0, \theta_t^1, \dots, \theta_t^M\} via eq. 1 for i=1 to n do
10:
11:
                  Frig. 1 to n do Sample \epsilon_i \sim \mathcal{N}(0, \sigma^2 I) Compute \theta_t^{i,m} = \theta_t^m + \epsilon_i Compute b(\pi_{\theta_t^{i,m}})
12:
13:
14:
                  Compute N_i = N(\theta_i^{i,m}, A)

Compute F_i = f(\theta_i^{i,m})

Send N_i and F_i from each worker to coordinator
15:
16:
17:
18:
             Set \theta_{t+1}^m = \theta_t^m + \alpha \frac{1}{n\sigma} \sum_{i=1}^n w * N_i \epsilon_i + (1-w) * F_i \epsilon_i
19:
             Compute b(\pi_{\theta_{t+1}^m})
20:
             Compute f(\theta_{t+1}^{\tilde{m}})
21.
             if f(\theta_{t+1}^m) > f_{best} then
22:
                   w = min(1, w + \delta_w)
23:
                   t_{best} = 0
24:
25:
                   f_{best} = f(\theta_{t+1}^m)
             else
26:
             t_{best} = t_{best} + 1 end if
27:
28:
29.
             if t_{best} \geq t_w then
                   w = max(0, w - \delta_w)
30:
31.
             \begin{array}{c} t_{best} = 0 \\ \mathbf{end} \ \mathbf{if} \end{array}
32:
33:
             Add b(\pi_{\theta_{t+1}^m}) to A
34: end for
```

6.5 Atari training details

Following Salimans et al. [24], the network architecture for the Atari experiments consists of 2 convolutional layers (16 filters of size 8x8 with stride 4 and 32 filters of size 4x4 with stride 2) followed by 1 fully-connected layer with 256 hidden units, followed by a linear output layer with one neuron per action. The action space dimensionality can range from 3 to 18 for different games. ReLU activations are placed between all layers, right after virtual batch normalization units [46]. Virtual batch normalization is equivalent to batch normalization [47], except that the layer normalization statistics are computed from a reference batch chosen at the start of training. In our experiments, we collected a reference batch of size 128 at the start of training, generated by random agent gameplay. Without virtual batch normalization, Gaussian perturbations to the network parameters tend to lead to single-action policies. The lack of action diversity in perturbed policies cripples learning and leads to poor results [24].

The preprocessing is identical to that in Mnih et al. [7]. Each frame is downsampled to 84x84 pixels, after which it is converted to grayscale. The actual observation to the network is a concatenation of 4 subsequent frames. There is a *frameskip* of 4. Each episode of training starts with up to 30 random, no-operation actions.

For all experiments, we fixed the training hyperparameters for fair comparison. Each network is trained with the Adam optimizer [48] with a learning rate of $\eta=10^{-2}$ and a noise standard deviation of $\sigma=0.02$. The number of samples drawn from the population distribution each generation was W=5000. For NS-ES, NSR-ES, and NSRA-ES we set M=3 as the meta-population size and k=10 for the nearest-neighbor computation, values that were both chosen through an informal hyperparameter search. We lowered M because the Atari network is much larger and thus each generation is more computationally expensive. A lower M enables more generations to occur in training. We trained ES, NS-ES, NSR-ES, and NSRA-ES for a the same number of generations T for each game. The value of T varies between 150 and 300 depending on the number of timesteps per episode of gameplay (i.e. games with longer episodes are trained for 150 generations and vice versa). The figures in SI Sec. 6.8 show how many generations of training occurred for each game.

6.6 Humanoid Locomotion problem training details

The domain is the MuJoCo Humanoid-v1 environment in OpenAI Gym [30]. In it, a humanoid robot receives a scalar reward composed of four components per timestep. The robot gets positive reward for standing and velocity in the positive x direction, and negative reward for ground impact energy and energy expended. These four components are summed across every timestep in an episode to get the total reward. Following the neural network architecture outlined by Salimans et al. [24], the neural network is a multilayer perceptron with two hidden layers containing 256 neurons each, resulting in a network with 166.7K parameters. While small (especially in the number of layers) compared to many deep RL architectures, this network is still orders of magnitude larger than what NS has been tried with before. The input to the network is the observation space from the environment, which is a vector $\in \mathbb{R}^{376}$ representing the state of the humanoid (e.g. joint angles, velocities) and the output of the network is a vector of motor commands $\in \mathbb{R}^{17}$ [30].

The network architecture for the Humanoid Locomotion experiments is a multilayer perceptron with two hidden layers containing 256 neurons (with tanh activations) resulting in a network with 166,700 parameters. This architecture is the one in the configuration file included in the source code released by Salimans et al. [24]. The architecture described in their paper is similar, but smaller, having 64 neurons per layer Salimans et al. [24].

For all experiments, we fixed the training hyperparameters for fair comparison. Each network was trained with the Adam optimizer [48] with a learning rate of $\eta=10^{-2}$ and a noise standard deviation of $\sigma=0.02$. The number of samples drawn from the population distribution each generation was W=10000. For NS-ES, NSR-ES, and NSRA-ES we set M=5 as the meta-population size and k=10 for the nearest-neighbor computation, values that were both chosen through an informal hyperparameter search. We trained ES, NS-ES, NSR-ES, and NSRA-ES for the same number of generations T for each game. The value of T is 600 for the Humanoid Locomotion problem and 800 for the Humanoid Locomotion with Deceptive Trap problem.

6.7 Humanoid Locomotion problem tabular results

| ENVIRONMENT | ES | NS-ES | NSR-ES | NSRA-ES |
|-------------|--------|--------|--------|---------|
| ISOTROPIC | 8098.5 | 6010.5 | 6756.9 | 7923.1 |
| DECEPTIVE | 5.3 | 16.5 | 31.2 | 37.1 |

Table 1: **Final results for the Humanoid Locomotion problem.** The reported scores are computed by taking the median over 10 independent runs of the rewards of the highest scoring policy per run (each of which is the mean over \sim 30 evaluations).

6.8 Plots of Atari learning across training (generations)

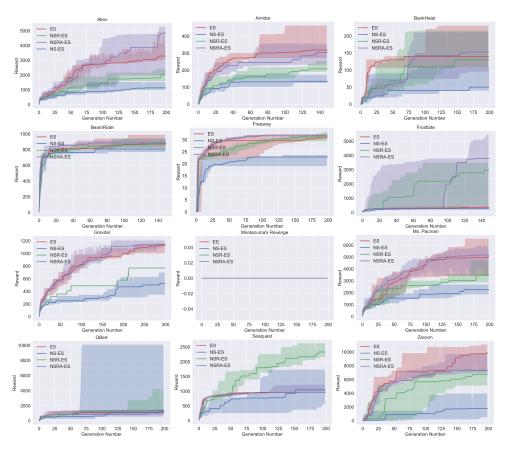


Figure 5: Learning comparison plots for the 12 Atari games we tested on

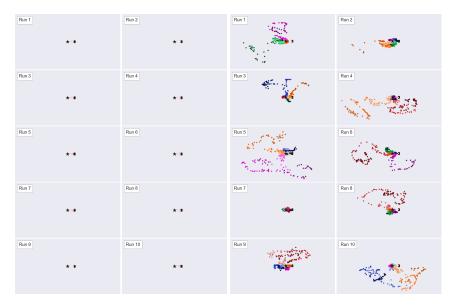


Figure 6: Overhead plot of ES (left) and NS-ES (right) across 10 independent runs on the Humanoid Locomotion with Deceptive Trap problem.

6.9 Overhead plots of agent behavior on the Humanoid Locomotion with Deceptive Trap Problem.

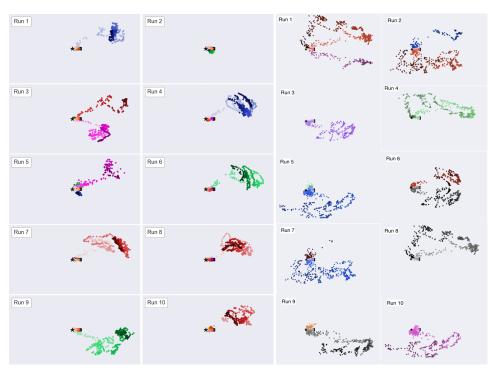


Figure 7: Overhead plot of NSR-ES (left) and NSRA-ES (right) across 10 independent runs on the Humanoid Locomotion with Deceptive Trap problem.