

Maximum Likelihood Reinforcement Learning

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Abstract

Reinforcement learning is the method of choice to train models in *sampling-based* setups with binary outcome feedback, such as navigation, code generation, and mathematical problem solving. In such settings, models implicitly induce a likelihood over correct rollouts. However, we observe that reinforcement learning does not maximize this likelihood, and instead optimizes only a lower-order approximation. Inspired by this observation, we introduce **Maximum Likelihood Reinforcement Learning (MaxRL)**, a sampling-based framework to approximate maximum likelihood using reinforcement learning techniques. MaxRL addresses the challenges of non-differentiable sampling by defining a compute-indexed family of sample-based objectives that interpolate between standard reinforcement learning and exact maximum likelihood as additional sampling compute is allocated. The resulting objectives admit a simple, unbiased policy-gradient estimator and converge to maximum likelihood optimization in the infinite-compute limit. Empirically, we show that MaxRL Pareto-dominates existing methods in all models and tasks we tested, achieving up to **20×** test-time scaling efficiency gains compared to its GRPO-trained counterpart. We also observe MaxRL to scale better with additional data and compute. Our results suggest MaxRL is a promising framework for scaling RL training in correctness based settings.¹

1 Introduction

Maximum likelihood (ML) and reinforcement learning (RL) are two highly successful optimization paradigms that significantly shaped the landscape of modern machine learning. Maximum likelihood training is a foundational principle behind modern generative and predictive models (Bishop, 2006; Murphy, 2012); in fully differentiable settings, optimizing log-likelihood objectives has reliably translated increases in model capacity, data, and compute into consistent performance improvements (Krizhevsky et al., 2012; Radford et al., 2018). Reinforcement learning, by contrast, originated in optimal control and sequential decision-making (Bertsekas, 1995; Sutton et al., 1998), where learning proceeds through interaction with an environment and the objective is to maximize expected return. The generality of this formulation enables reinforcement learning to address problems involving *non-differentiable* intermediate sampling, and has led to models capable of superhuman performance in complex domains (Mnih et al., 2015; Silver et al., 2016).

Many modern learning problems are typically addressed with reinforcement learning even if they define an implicit likelihood over successes. Examples include navigation (Thrun et al., 2005; Anderson et al., 2018), program synthesis (Chen et al., 2018; Bunel et al., 2018), structured prediction (Smith, 2011; Mensch and Blondel, 2018), and multi-step reasoning in large language models (Wei et al., 2023). In these tasks, success is determined by an external verifier only after a stochastic generation process, yielding a binary outcome. From an end-to-end perspective, the model induces a probability of success for each input, defining an implicit likelihood over correctness. Maximizing this likelihood would be the principled approach, but non-differentiable intermediate sampling precludes direct optimization. Reinforcement learning is used instead, not because it offers a better objective, but as a workaround to this non-differentiability.

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¹Project website and code: <https://zanette-labs.github.io/MaxRL/>

Suppose that both maximum likelihood and reinforcement learning could be applied to a given task, regardless of differentiability. The two objectives induce markedly different optimization behavior as we explain below. To make this distinction precise, we compare the population-level objectives implied by each framework. Let $p_\theta(x)$ denote the probability that a model with parameters θ produces a correct output for input x . The corresponding gradients of reinforcement learning, $\nabla_\theta J_{\text{RL}}$, and of maximum likelihood, $\nabla_\theta J_{\text{ML}}$, take the form

$$\begin{aligned}\nabla_\theta J_{\text{RL}} &= \mathbb{E}_x[\nabla_\theta p_\theta(x)] \\ \nabla_\theta J_{\text{ML}} &= \mathbb{E}_x[\nabla_\theta \log p_\theta(x)] = \mathbb{E}_x\left[\frac{1}{p_\theta(x)} \nabla_\theta p_\theta(x)\right]\end{aligned}$$

The inverse-probability reweighting induced by maximum likelihood places greater emphasis on hard, low-success inputs, leading to very different optimization dynamics, as we empirically demonstrate in this paper (cf. [Section 6](#)).

When viewed end-to-end, maximum likelihood emerges as the principled objective, for the very same reasons it is the method of choice in differentiable supervised learning with binary correctness. In non-differentiable problems, however, this objective is difficult to optimize directly because correctness is observed only after a non-differentiable stochastic generation process, and the success probability $p_\theta(x)$ may be small. This computational challenge motivates a framework that leverages additional sampling compute to more faithfully approximate likelihood-based training. We call this framework **Maximum Likelihood Reinforcement Learning (MaxRL)**. At a high level, MaxRL bridges standard reinforcement learning and exact maximum likelihood by progressively incorporating higher-order correctness information as more sampling compute is utilized, ultimately recovering likelihood optimization in the infinite-compute limit. We discuss the relationship between MaxRL and recent work in this direction ([Xiong et al., 2025b](#); [Davis and Recht, 2025](#)) in [Section 7](#).

Our contributions are threefold:

1. We formalize correctness-based reinforcement learning as a latent-generation maximum likelihood problem and show that standard reinforcement learning optimizes only the first-order approximation of the maximum likelihood objective (cf. [Section 3](#)).
2. We introduce a compute-indexed family of objectives that interpolates between expected reward and exact maximum likelihood via a Maclaurin expansion in pass@k events (cf. [Section 3](#)).
3. We derive a simple on-policy estimator whose expected gradient exactly matches the compute-indexed approximation of the likelihood objective, implying that increased sampling improves the optimized objective itself rather than merely reducing variance.

Empirically, MAXRL Pareto-dominates standard RL objectives (RLOO ([Ahmadian et al., 2024](#)), GRPO ([Shao et al., 2024](#))) on all settings that we tested on (cf. [Section 6](#)). Specifically, MAXRL show better scaling trends when additional compute and data are available, and on reasoning tasks, it achieves up to 20× test-time scaling efficiency gains with a perfect verifier. Together, our results illustrate MAXRL to be a promising direction for scaling RL in correctness based domains.

2 Preliminaries

In this work, we consider reinforcement learning settings that involve *generalization*, where models learn from a set of tasks and are evaluated on a heldout task distribution. We focus on correctness-based problems that can be abstracted as a *binary success or failure* outcome for each input. Formally, let \mathcal{X} and \mathcal{Y} denote the input and output spaces, and let $x \sim \rho$ be the distribution over tasks. For each input x , we denote $y^*(x) \in \mathcal{Y}$ as the correct label or answer. Equality between outputs is defined up to a task-dependent equivalence relation, so that $y = y^*(x)$ denotes semantic correctness rather than exact output equality. Finally, we let the learner be parameterized by θ and denote the predictive distribution induced by the model as $p_\theta(y | x)$, where $p_\theta(\cdot | x) \in \Delta(\mathcal{Y})$ is a conditional probability distribution over outputs for a fixed input x . In mathematical reasoning, x is the prompt and y is the final solution produced by the model. All logarithms use base e unless stated otherwise.

Latent generation models. In many modern settings, the model does not sample outputs directly from \mathcal{Y} , but instead generates a latent variable $z \in \mathcal{Z}$ according to a conditional distribution $m_\theta(z | x)$.

The final output $y \in \mathcal{Y}$ is then obtained via a deterministic decoding function $y = f(z)$, such as parsing a generated program or extracting a boxed answer from a chain of thought. Correctness is evaluated only on the decoded output, i.e., a trajectory z is successful if $f(z) = y^*(x)$. This induces a marginal probability of correctness

$$p_\theta(y^*(x) | x) = \sum_{z \in \mathcal{Z}} m_\theta(z | x) \mathbb{I}\{f(z) = y^*(x)\}. \quad (1)$$

Throughout the paper, expectations with respect to model outputs should be understood as expectations over latent samples $z \sim m_\theta(\cdot | x)$ followed by deterministic decoding.

Pass rate. We define the *pass rate* as the probability that the model produces the correct answer for a fixed input x :

$$p_\theta^{\text{pass}}(x) := p_\theta(y^*(x) | x) = \mathbb{E}_{y \sim p_\theta(\cdot | x)}[\mathbb{I}\{y = y^*(x)\}].$$

Similarly, let $y_1, \dots, y_k \stackrel{\text{i.i.d.}}{\sim} p_\theta(\cdot | x)$. We define *pass@k* as the probability of at least one correct sample:

$$\text{pass@}k(x) := \mathbb{P}(\exists i \in [k] \text{ s.t. } y_i = y^*(x)).$$

Next, we consider two optimization frameworks for training our models: *maximum likelihood* and *reinforcement learning*.

Maximum likelihood (ML). Maximum likelihood selects parameters that maximize the log-probability of the observed data under the model. In our binary correctness setting, each input x yields a single binary observation indicating whether the model produces a correct output. Under the latent generation model $z \sim m_\theta(\cdot | x)$ with deterministic decoding $y = f(z)$, the probability of observing correctness is $p_\theta(y^*(x) | x)$. Maximizing the corresponding log-likelihood therefore yields the objective

$$J_{\text{ML}}(\theta) := \mathbb{E}_{x \sim \rho} [\log p_\theta(y^*(x) | x)] \quad \text{with} \quad p_\theta(y^*(x) | x) = \mathbb{E}_{z \sim m_\theta(\cdot | x)}[\mathbb{I}\{f(z) = y^*(x)\}], \quad (2)$$

which is directly analogous to cross-entropy training in differentiable supervised learning.

Reinforcement learning (RL). For correctness based tasks, we also define a binary reward function $r(x, y) = \mathbb{I}\{y = y^*(x)\}$, and similarly under the latent variable case define $r(x, z) = \mathbb{I}\{f(z) = y^*(x)\}$. In this binary reward setting, the RL objective becomes (using the latent version without loss of generality):

$$J_{\text{RL}}(\theta) := \mathbb{E}_{x \sim \rho} [\mathbb{E}_{z \sim m_\theta(\cdot | x)} [r(x, z)]] = \mathbb{E}_{x \sim \rho} [p_\theta^{\text{pass}}(x)]. \quad (3)$$

This gives an objective equivalent to maximizing the expected population pass rate directly.

3 Maximum Likelihood Reinforcement Learning (MAXRL)

In this section, we show that reinforcement learning on expected reward optimizes only a first-order approximation of the ML objective. Specifically, the maximum likelihood objective admits a population-level expansion in terms of *pass@k* events, with standard RL optimizing only the first-order term. This suggests a compute-indexed family of objectives that incorporate higher-order terms, converging to ML as more compute is allocated.

3.1 Maclaurin Expansion of Maximum Likelihood

For simplicity, let us consider a single task x in the development to follow, as the final objective and gradients can be obtained by taking an expectation over $x \sim \rho$. Moreover, we write $p := p_\theta^{\text{pass}}(x)$ to simplify our notation. The maximum likelihood objective admits the *Maclaurin expansion* in terms of failure events:

$$J_{\text{ML}}(x) = \log p = - \sum_{k=1}^{\infty} \frac{(1-p)^k}{k} = - \sum_{k=1}^{\infty} \frac{\text{fail@}k(x)}{k}, \quad (4)$$

where $\text{fail}@k(x) = 1 - \text{pass}@k(x)$ denotes the probability that all k i.i.d. samples from the model fail. Differentiating (4) yields the population-level gradient identity

$$\nabla_{\theta} J_{\text{ML}}(x) = \sum_{k=1}^{\infty} \frac{1}{k} \nabla_{\theta} \text{pass}@k(x). \quad (5)$$

Thus, maximum likelihood optimizes an infinite harmonic mixture of $\text{pass}@k$ gradients, with higher-order terms encoding rare success which are critical when p is small. In contrast, the classical reinforcement learning approach is to optimize only the expected $\text{pass}@1$ objective (Koenig and Simmons, 1993; Silver et al., 2016; Vecerik et al., 2018; Guo et al., 2025):

$$\nabla_{\theta} J_{\text{RL}}(x) = \nabla_{\theta} \text{pass}@1(x),$$

corresponding to retaining solely the leading term of (5). From this observation, we can claim:

Reinforcement learning optimizes a first-order approximation of the maximum likelihood objective.

3.2 MAXRL Objective Function

The maximum likelihood gradient in (5) is difficult to estimate with finite samples. In particular, estimating $\text{pass}@k$ gradients for large k requires an increasing number of samples, especially when the pass rate p is small. This finite-sample difficulty is precisely what motivates Maximum Likelihood Reinforcement Learning. We define MAXRL as the *class of reinforcement-learning methods that explicitly target the maximum likelihood objective* rather than the pass rate, while remaining implementable under finite sampling and non-differentiable generation. We consider a principled way to do so below.

Consider approximating the maximum likelihood objective by truncating the Maclaurin expansion (5) to a finite order and then estimating such an objective instead. For a truncation level $T \in \mathbb{N}$, we define the truncated maximum likelihood objective for a fixed input x as

$$J_{\text{MAXRL}}^{(T)}(x) := - \sum_{k=1}^T \frac{(1-p)^k}{k}. \quad (6)$$

Differentiating (6) yields the truncated population gradient

$$\nabla_{\theta} J_{\text{MAXRL}}^{(T)}(x) = \sum_{k=1}^T \frac{1}{k} \nabla_{\theta} \text{pass}@k(x). \quad (7)$$

This defines a family of objectives: **$T = 1$ recovers reinforcement learning**, **$T \rightarrow \infty$ recovers maximum likelihood**, and intermediate T values interpolate between them. Thus, the truncation level T directly controls the order of correctness events that contribute to learning. As we will soon see, it becomes viable to estimate higher-order $\nabla_{\theta} J_{\text{MAXRL}}^{(T)}(x)$ as more compute is expended in terms of rollouts. In other words:

MAXRL provides a principled framework for trading additional compute for higher-fidelity approximations to the maximum likelihood objective.

The remaining question is whether these truncated objectives admit simple, unbiased estimators under finite sampling, a question that we answer affirmatively in the next section.

4 Gradient Estimators for MAXRL

Eq. (7) already provides a viable approach for constructing an unbiased estimator: approximate *each* term in the finite series using a $\text{pass}@k$ gradient estimator, as provided in recent work (Walder and Karkhanis,

2025; Chen et al., 2025d). Under this strategy, any improvement in pass@k estimators directly translates into improved estimators for the truncated maximum likelihood objective in Eq. (7).

In this work, we take an alternate approach, one that will lead to a simpler estimator and to a new viewpoint. The key insight is that the maximum likelihood gradient can be expressed as an expectation under the *success-conditioned* distribution (Davis and Recht (2025) also recently made a similar observation), as established by the following theorem. We provide the proof in Appendix B.

Theorem 1 (Conditional Form of the Maximum Likelihood Gradient). *The gradient of the maximum likelihood objective admits the following conditional expectation representation:*

$$\nabla_{\theta} J_{\text{ML}}(x) = \mathbb{E}[\nabla_{\theta} \log m_{\theta}(z | x) | f(z) = y^*(x)]. \quad (8)$$

The theorem establishes that the maximum likelihood gradient is the average gradient from successful trajectories only. This interpretation naturally leads to a concrete gradient estimator by replacing the expectation with sample averages.

4.1 Empirical Gradient Estimator

Theorem 1 suggests drawing samples from the success-conditioned policy. Recent works have proposed rejection fine-tuning (Touvron et al., 2023; Yuan et al., 2023; Dong et al., 2023; Xiong et al., 2025a; Davis and Recht, 2025) and adaptive sampling (Xiong et al., 2025b) as mechanisms to sample from this conditional distribution. However, doing so is computationally demanding when the pass rate is small or requires a more complex implementation regarding adaptive sampling. Instead, we adopt a simpler approach: we sample from the unconditional policy $m_{\theta}(\cdot | x)$ and then *average over only the successful trajectories*. Fix an input x and draw N latent trajectories $z_1, \dots, z_N \sim m_{\theta}(\cdot | x)$. Let $r_i := \mathbb{I}\{f(z_i) = y^*(x)\}$ indicate success based reward, $S_i := \nabla_{\theta} \log m_{\theta}(z_i | x)$ denote the score function, and $K := \sum_{i=1}^N r_i$ be the number of successful samples. We average score functions *only over successful trajectories* and obtain the following REINFORCE-style estimator:

$$\hat{g}_N(x) := \begin{cases} \frac{1}{K} \sum_{i=1}^N r_i S_i, & K \geq 1, \\ 0, & K = 0. \end{cases} \quad (9)$$

The estimator constructed in this way is such that some inputs may receive zero gradient if there are no successes within N samples, making the resulting estimator no longer unbiased with respect to Eq. (9). We show that this estimator is however unbiased for the gradient of the truncated maximum likelihood objective in Eq. (7), $\nabla_{\theta} J_{\text{MAXRL}}^{(T)}(x)$, with truncation level $T = N$:

Theorem 2 (Estimator-objective equivalence). *The estimator $\hat{g}_N(x)$ is an unbiased estimator for the MAXRL gradient of order $T = N$, i.e.,*

$$\mathbb{E}[\hat{g}_N(x)] = \nabla_{\theta} J_{\text{MAXRL}}^{(N)}(x).$$

We present the proof of this result in Appendix B. Theorem 2 reveals an elegant alignment between the estimator in Eq. (9) and the gradient of the truncated Maclaurin expansion in Eq. (7). It is worth highlighting the most important property of the estimator:

Increasing compute as rollouts N leads to a better approximation of the maximum likelihood gradient.

Table 1 compares our estimator with the REINFORCE² estimator, whose expected value underlies most RL algorithms. At the estimator level, the difference is simple: both average score functions over

²Modern policy-gradient methods such as PPO (Schulman et al., 2017) introduce additional mechanisms (importance weight truncation via clipping) that trade bias for robustness. In the fully on-policy setting, these reduce to REINFORCE, our canonical baseline. GRPO (Shao et al., 2024) is a notable exception due to its division by standard deviation in the advantage calculation, which we discuss further in Section 5.

sampled trajectories, but REINFORCE normalizes by total samples N while MaxRL normalizes by successful samples K . This difference in normalization determines the objective each estimator is unbiased for.

Consequently, increasing number of samples, N , for the two estimators has different effects: REINFORCE reduces variance of a fixed objective (pass@1), while MaxRL increases the approximation order to maximum likelihood. Additional compute thus improves the *objective itself* for MAXRL, not just estimation quality.

4.2 Variance Reduction via Control Variates

Like REINFORCE, the estimator (9) can exhibit high variance when successful samples K is small. Policy-gradient baselines are typically introduced to reduce variance without changing the expected gradient (Sutton, 1988). However, standard arguments for policy-gradient baselines are not directly applicable in this setting, as the estimator normalizes by the random variable K which depends on all samples which makes it correlated with the observed rollouts.

We instead proceed from first principles and use a simple zero-mean control variate, the unconditional average score:

$$V_N := \frac{1}{N} \sum_{i=1}^N \nabla_{\theta} \log m_{\theta}(z_i | x),$$

which satisfies $\mathbb{E}[V_N] = 0$. Subtracting V_N preserves unbiasedness while reducing variance:

$$\tilde{g}_N(x) = \frac{1}{K} \sum_{i=1}^N r_i S_i - \frac{1}{N} \sum_{i=1}^N S_i = \sum_{i=1}^N \left(\frac{r_i}{K} - \frac{1}{N} \right) S_i, \quad (10)$$

with the convention that the first term, $(\sum_{i=1}^N r_i S_i) / K$, is zero when $K = 0$.

4.3 On-Policy Implementation

In Algorithm 1, we present a simple *on-policy* implementation of MAXRL that differs from standard REINFORCE-style policy gradient methods by a *single-line* modification to the advantage calculation. We adopt the variance-reduced formulation in Eq. (10), and drop both terms when $K = 0$, consistent with standard policy-gradient practice in which no gradient is computed for tasks with no successful rollouts: this choice is simpler and performs better empirically. Concretely, the advantage is normalized by the per-task mean reward, rather than left unnormalized as in RLOO (Ahmadian et al., 2024) or normalized by the reward standard deviation as in GRPO (Shao et al., 2024). The modified line is highlighted in blue.

Algorithm 1 On-Policy Implementation of MAXRL.

```

require Batch of inputs  $B$ , number of rollouts  $N$ , latent policy  $m_{\theta}(\cdot | \cdot)$ 
1: for each input  $x \in B$  do
2:   Sample  $z_1, \dots, z_N \sim m_{\theta}(\cdot | x)$ 
3:   for  $j = 1$  to  $N$  do
4:      $r_j \leftarrow \mathbb{I}\{f(z_j) = y^*(x)\}$ 
5:      $S_j \leftarrow \nabla_{\theta} \log m_{\theta}(z_j | x)$ 
6:   end for
7:    $\hat{r}(x) \leftarrow \frac{1}{N} \sum_{j=1}^N r_j$ 
8:    $\hat{g}(x) \leftarrow \begin{cases} \frac{1}{N \hat{r}(x)} \sum_{j=1}^N (r_j - \hat{r}(x)) S_j, & \hat{r}(x) > 0, \\ 0, & \text{otherwise} \end{cases}$ 
9: end for
10:  $\hat{g} \leftarrow \frac{1}{|B|} \sum_{x \in B} \hat{g}(x)$ 
11: return  $\hat{g}$ 

```

5 A Unifying Weight-Function View

Maximum likelihood, MAXRL, classical reinforcement learning, and GRPO all admit population-level gradients of the form

$$\nabla_{\theta} J = \mathbb{E}_{x \sim \rho} [w(p_{\theta}(x)) \nabla_{\theta} p_{\theta}(x)], \quad (11)$$

where $p_{\theta}(x) = p_{\theta}^{\text{pass}}(x)$ and $w(p)$ is a scalar weight that depends only on the pass rate. The function $w(p)$ determines how learning signal is allocated across inputs of varying difficulty and fully characterizes the differences between these objectives at the population level. Figure 1 plots the resulting weight functions for each method; derivations are provided in Appendix C. The key distinction among objectives is *how strongly they emphasize hard, low-pass-rate inputs*. As T increases, MAXRL uniquely approaches maximum likelihood weighting in the low pass rate regime.

This weight perspective also provides a useful reinterpretation of GRPO (Shao et al., 2024). Although GRPO is heuristically motivated by Z-normalization using the empirical standard deviation, such normalization induces a fundamentally different population-level objective from REINFORCE, a conclusion also reached by recent works (Davis and Recht, 2025; Liu et al., 2025b; Xiong et al., 2025b). Table 2 summarizes the population-level weighting functions; relative to standard expected-reward optimization, GRPO upweights low-pass-rate inputs approximately as $1/\sqrt{p}$ when p is small, placing it between classical reinforcement learning and maximum likelihood. However, increasing compute via additional sampling under GRPO does not yield a better approximation to the maximum likelihood objective, as the induced population loss is fundamentally distinct. Moreover, as shown in Figure 1, the GRPO weighting function *inverts* for sufficiently large pass rates, increasing as $p \rightarrow 1$, unlike likelihood-based objectives. Consequently, GRPO assigns increased weight to very easy inputs when they are present, in contrast to the other formulations.³

Table 2: Population-level weighting functions $w(p)$.

	RL	GRPO	MAXRL (T)	ML
$w(p)$	1	$\frac{1}{\sqrt{p(1-p)}}$	$\frac{1-(1-p)^T}{p}$	$\frac{1}{p}$

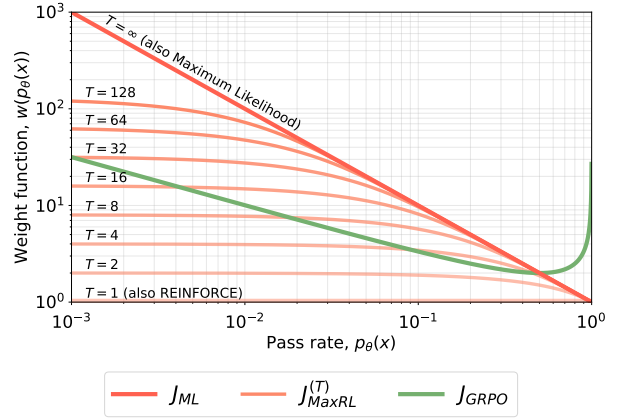


Figure 1: Population-level weighting functions $w(p)$ as a function of pass rate p . Truncated objectives $J_{\text{MAXRL}}^{(T)}$ interpolate between REINFORCE and maximum likelihood as T increases.

6 Experiments

We now turn to the empirical evaluation of MAXRL. We begin in Section 6.1 with a controlled setting where exact maximum likelihood optimization is possible, allowing direct comparison with MAXRL as compute increases. We then study non-differentiable correctness-based tasks in two regimes: (i) an *effectively infinite-data* setting with large number of novel tasks (Section 6.2), and (ii) a *data-scarce* setting with a fixed training dataset where we can scale compute nonetheless by training for many epochs over the same dataset (Section 6.3). Finally, in Section 6.4, we train and evaluate billion-parameter reasoning models on mathematical problem-solving, testing whether the benefits of MAXRL extend to larger-scale LLM training.

Because we compare training objectives rather than algorithms, all methods are trained *on-policy*. We compare against REINFORCE with a leave-one-out baseline (RLOO) (Ahmadian et al., 2024) and Group Relative Policy Optimization (GRPO) (Shao et al., 2024) as primary baselines.

6.1 Comparisons with Exact Maximum Likelihood

As a first step, we evaluate how closely MAXRL approximates *exact* maximum likelihood in a setting where the latter can be implemented exactly. We compare three objectives: (i) reinforcement learning on expected reward, (ii) MAXRL, and (iii) exact maximum likelihood training. We consider a standard image

³We conjecture that this inversion may contribute to distribution sharpening (Yue et al., 2025; Wu et al., 2026) when datasets contain a substantial fraction of overly easy inputs, and leave a detailed analysis to future work.

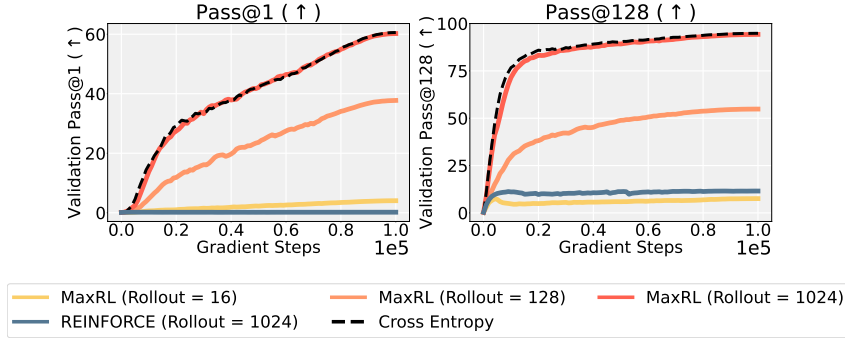


Figure 2: (ImageNet) Comparison of training dynamics under exact maximum likelihood, MAXRL, and REINFORCE in a controlled image classification setting. With sufficient rollouts, MAXRL closely matches cross-entropy training, while REINFORCE fails to make progress from low initial pass rates even with high number of rollouts.

classification task, where maximum likelihood corresponds to minimizing cross-entropy. The reinforcement learning reward is defined as 1 if the predicted class matches the ground-truth label and 0 otherwise.

We instantiate this comparison on ImageNet (Deng et al., 2009) using a ResNet-50 (He et al., 2015) trained under each objective; full experimental details are provided in Appendix D. Figure 2 summarizes the results: REINFORCE (with a standard baseline) fails to achieve meaningful improvements even with very high per-input sampling budget, whereas exact maximum likelihood training yields steady gains.

In contrast, MAXRL is trained on the same samples and observes the same sparse set of successful trajectories as REINFORCE, but makes more effective use of this limited learning signal through likelihood-inspired reweighting. As the compute increases by means of higher rollout counts, *MAXRL improves consistently and closely tracks exact maximum likelihood*. We also analyze the gradient norm resulting from different objectives in Figure 8: MAXRL and cross-entropy concentrate learning signal on harder tasks and are characteristically similar given sufficient compute for MAXRL, whereas GRPO and REINFORCE exhibit very different behavior. For additional experiments such as comparison to GRPO, we refer the reader to Appendix D.

Takeaway 1: MAXRL approaches exact maximum likelihood given infinite compute.

When direct maximum-likelihood optimization is available, MAXRL converges to it as sampling compute increases.

6.2 Infinite Data Regime

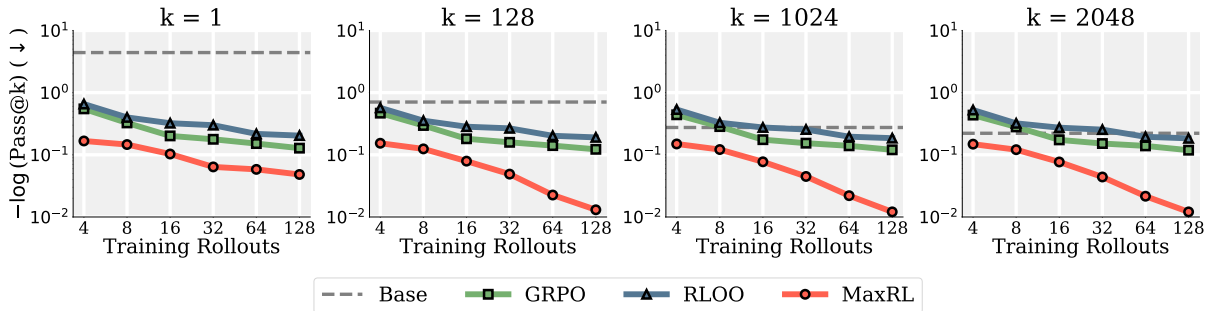


Figure 3: (Maze) Motivated by Schaeffer et al. (2025), we record $-\log(\text{Pass}@k)$ (lower is better) as a function of training rollouts for different objectives. We see that for across different all inference rollout budgets (k), MAXRL exhibit better scaling compared to GRPO and RLOO as we increase number of training rollouts.

Next, we study MAXRL in non-differentiable settings. For the first experiment, we study how MAXRL behave in data-rich domains. To simulate training with infinite data, we construct a procedurally generated maze-navigation environment with 1 million unique 17×17 mazes for training where multiple valid solution paths might exist for a given task. We reserve a held-out set of 256 mazes for evaluation and apply a brief supervised pretraining phase to ensure a non-zero initial pass rate. The complete details of the task are provided in the Appendix E.1.

We train a lightweight transformer model (Vaswani et al., 2017) with approximately 3M parameters

and simulate extended training by running 9K RL steps with up to 128 rollouts per prompt, varying the number of rollouts to control compute. We report performance after 9K steps in Figure 3 as a function of training compute, implemented as different number of rollouts per prompt from 4 to 128. Notice that this is a substantial amount of compute for the size of the model.

All three objectives (RLOO, GRPO, and MAXRL) improve upon the base model, but the magnitude of improvement differs markedly across methods. Even at the highest compute budget (128 rollouts per prompt), RLOO fails to match the performance achieved by MAXRL at the lowest budget (4 rollouts) across all pass@k metrics. GRPO-trained models at the highest compute similarly trail MAXRL trained with only 16 rollouts per prompt at every pass@k evaluation. These results highlight that MAXRL *scales with compute* far more effectively than competing frameworks when large amounts of unique training data is available. Comparisons with additional baselines are provided in Table 3 and Appendix L.

Table 3: Performance comparison in maze with additional baselines.

Method	Pass@1	Pass@128	Pass@256
GRPO (Shao et al., 2024)	43.6	49.0	49.6
RLOO (Ahmadian et al., 2024)	25.2	28.7	29.0
GRPO (with entropy bonus)	47.2	53.4	54.0
PKPO (T = 16) (Walder and Karkhanis, 2025)	74.5	77.6	77.9
Differential Smoothing (Gai et al., 2025)	50.0	57.8	58.7
MaxRL	84.4	92.0	94.3

Takeaway 2: MAXRL scales better with additional compute in the infinite data regime.

In a data-rich training regime, MAXRL scales more favorably with additional compute compared to existing methods.

6.3 Data-Scarce Regime

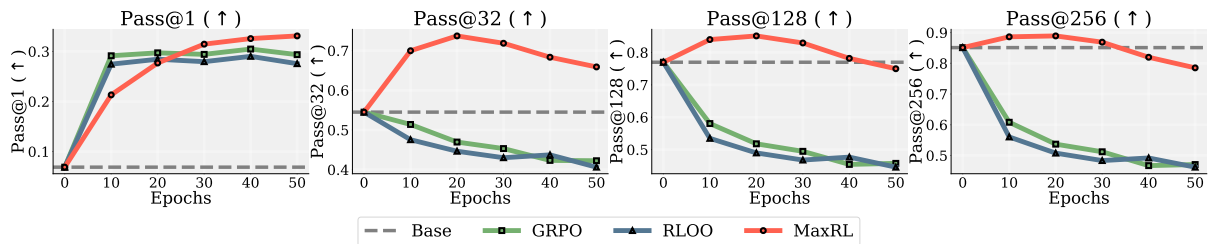


Figure 4: (GSM8K) Training dynamics on GSM8K with a fixed dataset and increasing training compute in terms of RL steps. MAXRL shows slower initial gains but ultimately achieves higher performance with substantially less pass@k degradation compared to GRPO and REINFORCE.

We next consider a data-scarce regime where models are trained for many epochs on a fixed dataset until peak performance is reached, with the goal of identifying which method extracts the highest attainable performance. Unlike the infinite-data setting in Section 6.2, longer training in this regime does not necessarily translate into improved performance due to the increased risk of overfitting. Specifically, we train a SmoLM2-360M-Instruct model (Allal et al., 2025) on GSM8K (Cobbe et al., 2021) for up to 50 epochs. Training dynamics are reported in Figure 4, with additional experimental details provided in Appendix E.2.

All methods improve upon the base model in terms of pass@1 performance; however, only MAXRL consistently exceeds the base model in pass@k metrics. In contrast, RLOO and GRPO exhibit massive pass@k degradation with extended training, mirroring the behavior observed by Yue et al. (2025) and exacerbated here by prolonged training on a fixed dataset. RLOO and GRPO reach their peak pass@1 performance faster than MAXRL (around 10 epochs), but MAXRL overtakes competing methods at approximately 30 epochs and continues to increase through the end of training, reaching a higher peak. At the same time, pass@k remains substantially healthier for MAXRL and exceeds the base model for a large portion of training. This behavior suggests MAXRL *is more resistant to overfitting*, particularly with regards to output diversity. Additional comparisons with baseline methods are reported in Table 4.

Table 4: Performance comparison across methods on GSM8K.

Method	Pass@1	Pass@128	Pass@1024
GRPO (Shao et al., 2024)	29.3	45.8	48.8
RLOO (Ahmadian et al., 2024)	27.5	44.6	48.5
GRPO (with entropy bonus)	31.1	48.1	51.6
PKPO (T = 16) (Walder and Karkhanis, 2025)	30.7	67.2	75.9
Differential Smoothing (Gai et al., 2025)	31.4	48.5	52.3
MaxRL	33.2	75.0	83.4

Takeaway 3: MAXRL is more resistant to overfitting.

In a data-scarce regime, MAXRL can sustain improvement over a large number of epochs, demonstrating less pass@k degradation (overfitting) and converging to a higher average performance.

6.4 Large Reasoning Model Training

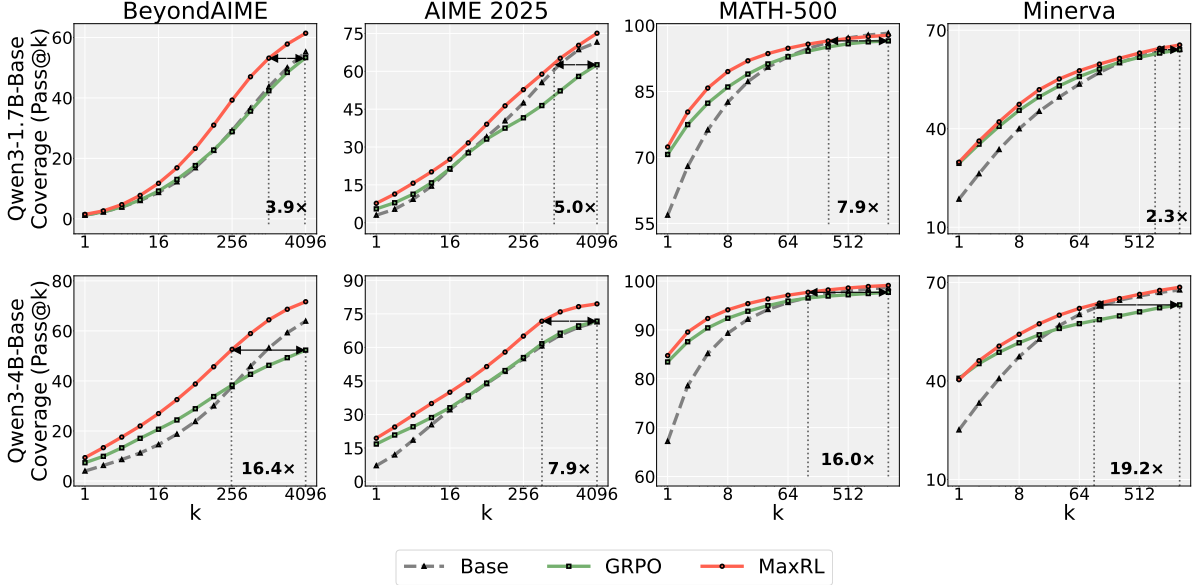


Figure 5: (Qwen3 training results) Evaluation of final checkpoints from training Qwen3-1.7B-Base and Qwen3-4B-Base models, on 4 benchmarks: AIME 2025, BeyondAIME, MATH-500 and Minerva. MAXRL match or outperform GRPO in all 4 evaluation datasets and shows little to no degradation at coverage (pass@k) for very high k values. We also note the increase in inference efficiency: MAXRL can provide 2.3x - 19.2x speedup compared to GRPO while generating multiple samples with a perfect verifier and maintains similar or better pass@1 performance.

We next demonstrate that the benefits of MAXRL extend to larger-scale LLM reasoning training. We train Qwen3-1.7B-Base and Qwen3-4B-Base models on POLARIS-53K (An et al., 2025), a dataset of approximately 50K mathematical reasoning prompts, using 256 prompts per batch, 16 rollouts per prompt, and 1000 RL steps. Notice that this setup utilizes lower rollout counts and RL steps than prior experiments to allow training larger models within our compute budget. We evaluate on four standard math benchmarks: AIME 2025, BeyondAIME (ByteDance-Seed, 2025), MATH-500 (Hendrycks et al., 2021; Lightman et al., 2023), and Minerva (Lewkowycz et al., 2022). We compare against GRPO (Shao et al., 2024), a widely used baseline for large-scale reasoning (Guo et al., 2025; Yang et al., 2025). Additional details are provided in Appendix E.3.

Figure 5 summarizes our main results. Across both model sizes, MAXRL consistently Pareto dominates GRPO, achieving higher pass@1 while simultaneously improving pass@k. Consistent with prior work (Yue et al., 2025; Wu et al., 2026), GRPO exhibits pronounced pass@k degradation at larger k. In contrast, MAXRL improves pass@k relative to both the pretrained base model and the GRPO-trained checkpoint in 7 out of 8 evaluation settings. Improved pass@k directly translates into inference efficiency under repeated sampling. As shown in Figure 5, MAXRL achieves up to 20x test-time scaling efficiency gains when using a perfect verifier to filter wrong solutions, yielding substantial practical savings at inference time. We note that many settings admit a strong verifier, such as programming (Chen et al., 2021) or Lean (de Moura et al., 2015), where MAXRL can show strong benefits over other RL objectives. We provide additional results, such as evaluation on 4 additional benchmarks, and training dynamics statistics such as mean generated response length, entropy and gradient norm, in Appendix H, K, and J.

Takeaway 4: MAXRL’s benefits transfer to larger scale mathematical reasoning

On larger scale mathematical reasoning, MAXRL Pareto-dominates GRPO, shows little to no diversity degradation with respect to the base model, and leads to strong (up to 20x) test-time scaling efficiency gains.

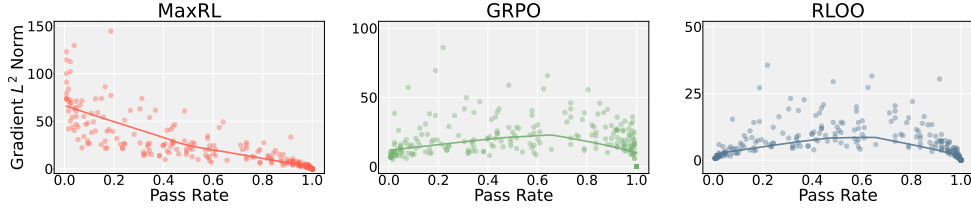


Figure 6: (Gradient norm analysis) To compare different objectives qualitatively, we show a scatter plot of gradient L^2 norm vs pass rate over individual prompts. We use Qwen2.5-1.5B-Instruct on MATH-500 dataset for this analysis. MAXRL generate larger gradient norms over prompts with close to 0 pass rates.

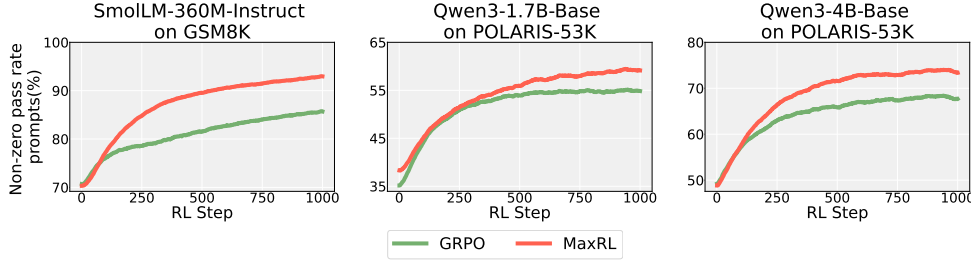


Figure 7: (Training dynamics comparison) Fraction of prompts where the model generates at least one correct rollout (out of 128, 16, and 16 rollouts for SmolLM-360M-Instruct, Qwen3-1.7B-Base, and Qwen3-4B-Base, respectively) during training. MAXRL consistently produces at least one correct rollout for more prompts across all settings, demonstrating its effectiveness at extracting more learning signal from the training dataset.

6.5 MAXRL Behaves Characteristically Different from Other RL Objectives

Finally, we study whether MAXRL shows different characteristics compared to commonly used RL objectives besides performance metrics. We first study the gradient norms produced by different objectives: Figure 6 illustrates that MAXRL generates higher gradient norms on harder prompts and lower gradient norms on easier prompts. This behavior matches that of cross-entropy in fully differentiable image classification setting (Figure 8): showing that MAXRL concentrate learning signal on harder problems unlike GRPO and RLOO. This larger gradient norms on more difficult prompts then translates into the model’s capability to generate correct solutions for a larger fraction of problems during training: Figure 7 demonstrates this for 3 different base models where MAXRL consistently generate at least one correct rollout for a larger fraction of training prompts, and the gap between MAXRL and GRPO persists as we train longer. We have also run this analysis for our maze and GSM8K training settings: Section I shows similar behavior in these setups as well. Overall, we demonstrate that MAXRL exhibit interesting differences compared to other RL objectives, and we leave their further study to future work.

Takeaway 5: MAXRL shows characteristically different optimization dynamics.

Besides performance metrics, MAXRL also exhibits different optimization dynamics. Most notably, it produces stronger gradients on harder prompts, and also leads to a larger fraction of prompts with at least one correct rollout during training.

7 Related Works

Supervised training vs reinforcement learning. Supervised learning and reinforcement learning (RL) are complementary but fundamentally different paradigms. Supervised training is stable, sample-efficient, and well-calibrated within the training distribution (Ng and Jordan, 2001), but it is limited by the quality and scope of available data and cannot directly optimize non-differentiable objectives such as correctness or preferences. In contrast, RL can optimize such objectives directly, typically via policy gradients (Williams, 1992; Sutton et al., 1999, 1998; Schulman et al., 2017; Guo et al., 2017), and improve performance beyond available demonstrations by having access to interactions with an environment and the resulting reward-based feedback. Although recent work has reframed RL objectives as supervised ones (Rafailov et al., 2023), on-policy learning, characteristic of online RL algorithms, appears crucial for optimal performance (Tajwar et al., 2024; Xu et al., 2024). Modern foundation model training often combines supervised learning on human data with subsequent RL (Ouyang et al., 2022). Unlike these

approaches, we assume no access to high-quality demonstrations or a stronger model, and instead study a purely interactive RL setting that nonetheless optimizes an objective that mimics cross-entropy. We discuss additional related works in Appendix A.

Training LLMs for strong reasoning abilities. Reinforcement learning from verifiable rewards (RLVR), where LLMs receive reward from a ground truth verifier instead of using a trained reward model, has emerged as the dominant paradigm for instilling strong reasoning capabilities into LLMs (OpenAI et al., 2024; Guo et al., 2025; Team et al., 2025; Lambert et al., 2025; Yang et al., 2025). Whereas supervised training learns better behavior from fixed static datasets, reinforcement learning uses policy gradient algorithms (e.g., PPO (Schulman et al., 2017), GRPO (Shao et al., 2024), RLOO (Ahmadian et al., 2024)) to learn from self-generated responses and non-differentiable rewards. However, these algorithms and their variants (Zheng et al., 2025; Liu et al., 2025b; MiniMax et al., 2025) optimize expected reward or pass rate and only differs in how the advantage or off-policy updates are calculated. In contrast, the goal of our work is to propose a fundamentally different objective for RL training.

RL training causes distribution sharpening. Despite its usefulness, questions remain on whether RLVR teaches LLMs fundamentally new behavior/skills, or simply sharpens existing good behavior from the pretrained model. Prior works (Liu et al., 2025b; Zhao et al., 2025; AI et al., 2025) demonstrated that certain reasoning skills like reflection already exist in the pretrained model, and Gandhi et al. (2025) shows that good reasoning behaviors learned from pretraining is crucial for the success of RLVR in the post-training phase. More recently, studies (Yue et al., 2025; Dang et al., 2025; Wu et al., 2026) found that RLVR decreases the model’s diversity by reducing pass@k. In our paper, we confirm these findings and attribute this to the RL objective itself as optimizing expected reward tends to marginalize learning signal from harder prompts, which results in distribution sharpening.

Learning to solve hard problems. Due to RLVR’s shrinking of model coverage, significant attention has been drawn to new RL algorithms mitigating pass@k collapse. Approaches range from directly optimizing for pass@k during training (Walder and Karkhanis, 2025; Tang et al., 2025) to employing exploration bonuses in RL (Song et al., 2025; Tuyls et al., 2025). We show in our work that pass@k optimization objectives are a special case of our objective, since it optimizes an infinite harmonic series of pass@k objectives. On the other hand, the other works carry the fundamental limitation of RLVR of maximizing expected reward or pass rate over a batch of prompts, which we demonstrate to have vanishing gradient for prompts with low pass rate. This issue is also recognized by Nguyen et al. (2025b), which introduces selective learning only on prompts where greedy response fails, but unlike us, does not weigh prompts differently based on their pass rate.

Closely related works. One closely related line of work is that of Xiong et al. (2025b), which also considers non-linear functions of the pass rate in reinforcement learning. Both works are motivated by the observation that expected-reward objectives can underweight low-pass-rate prompts, and maximizing a log-likelihood like objective can mitigate this issue. However, the focus of Xiong et al. (2025b) is on adaptive rollout budget allocation and sampling strategies, treating the choice of non-linear weighting as part of an algorithmic design space. In contrast, we do not employ adaptive sampling, and instead derive a sampling-based on-policy estimator that approaches maximum likelihood as sampling compute budget is increased. We also empirically focus on demonstrating better data and compute scaling with our framework, whereas Xiong et al. (2025b) focuses on comparing against other adaptive sampling frameworks (Yu et al., 2025a). Furthermore, Davis and Recht (2025) provides a theoretical argument characterizing the population-level objectives approached by specific binary-reward reinforcement-learning algorithms in asymptotic limits, showing that certain procedures induce log-like weighting of the pass rate. Our work addresses a complementary question: how finite sampling defines explicit population-level objectives. We establish an exact estimator-objective equivalence at each finite rollout count and empirically evaluate how this objective-level interpolation manifests as compute is increased in both controlled settings and large-scale LLM post-training experiments. We discuss additional related works in Appendix A.

8 Conclusion

In this work, we establish Maximum Likelihood Reinforcement Learning as a principled optimization framework for non-differentiable binary reward settings. We showed that MAXRL approaches maximum

likelihood in differentiable settings as compute increases and that in non-differentiable settings it offers key advantages over traditional RL, scaling more effectively with additional compute and data. More broadly, our results suggest that some limitations attributed to reinforcement learning with foundation models arise from objective choice rather than optimization or sampling. Our work currently assumes a binary reward setting and does not extend directly to continuous or arbitrarily valued rewards. Moreover, one can also consider objectives other than maximum likelihood to optimize following our framework. Generalizing MAXRL to continuous rewards, multi-turn reinforcement learning, and off-policy settings such as PPO-style training are promising directions for future work.

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A Extended Related Works

Cross-entropy objective. Maximizing log-likelihood via optimizing cross-entropy is a widely used framework in machine learning due to its simplicity and favorable theoretical properties. In particular, cross-entropy is a strictly proper scoring rule, meaning that the expected loss is uniquely minimized by the true probability distribution, which encourages statistically calibrated predictions (Good, 1992; Savage, 1971; Gneiting and Raftery, 2007; Waghmare and Ziegel, 2025). Moreover, it yields statistically efficient estimators under standard assumptions (Vaart, 1998; Casella and Berger, 2002; Lehmann and Casella, 2006), and induces gradients that concentrate learning signal on low-probability or uncertain outcomes through logarithmic weighting (Wang et al., 2020). As a result, cross-entropy (or log-loss) often yields consistent estimators for classification and tends to generalize well in practice (Ng and Jordan, 2001; Zhang, 2004). However, more recent work has shown that models trained to maximize log-likelihood can still overfit and exhibit miscalibration, motivating post-hoc techniques such as temperature scaling (Niculescu-Mizil and Caruana, 2005; Guo et al., 2017). Moreover, the unbounded nature of cross-entropy and its sensitivity to small perturbations in the predicted distribution suggest that alternative strictly proper scoring rules may be more suitable in certain settings (Kornblith et al., 2021). Because cross-entropy is extensively studied, we refer interested readers to Mao et al. (2023); Li et al. (2025); Terven et al. (2025) for a comprehensive review.

Supervised training vs reinforcement learning. Supervised learning has been the go-to training paradigm in machine learning, beginning with early “learning with a teacher” neural network based systems such as the perceptron (Rosenblatt, 1958) and later becoming practical with backpropagation-based neural network training (Rumelhart et al., 1986). It has been used to tackle a broad range of problems, from financial fraud detection (Afriyie et al., 2023; Editya et al., 2025), to sentiment analysis (Zhang et al., 2018c) and spam detection (Jain et al., 2022; Jamil et al., 2025). More recently, supervised training has been used in modern image classification systems (Lecun et al., 1998; Krizhevsky et al., 2012) to achieve strong performance. In modern foundation models, “pretraining” is typically done by supervised learning via minimizing cross-entropy over next-token prediction (Radford et al., 2018, 2019) on large corpus of text, followed by additional supervised training on high quality human written demonstrations (Ouyang et al., 2022; Zhang et al., 2023) to teach models how to respond to prompts, often known as instruction-tuning. In sequential decision making domains, supervised training appears as behavior cloning from expert demonstration, used in early autonomous driving and robotic systems (Pomerleau, 1988; Bojarski et al., 2016; Codevilla et al., 2018). However, supervised learning in sequential decision making breaks the i.i.d. assumption since the learned policy’s actions affect which states it visits during execution, causing compounding errors over time (Ross and Bagnell, 2010; Belkhale et al., 2023). The classic no-regret reductions line (DAgger) makes this explicit and addresses drift by iteratively querying the expert on states visited by the learner, turning the sequential problem into an online supervised learning loop with improved guarantees (Ross et al., 2010).

In such scenarios, reinforcement learning is an attractive alternative paradigm that formalizes learning under delayed, sparse, and evaluative feedback in Markov decision processes (MDPs), with foundational roots in dynamic programming and MDP theory (Bellman, 1957). Mechanistically, the key contrast with supervised learning / behavior cloning is that supervised learning assumes (or benefits strongly from) an i.i.d. dataset of correct targets under a fixed data distribution, whereas RL’s data distribution is policy-induced and nonstationary, and gradients arise from credit assignment through rewards rather than direct target labels. This mismatch shows up starkly in sequential prediction/imitation: naive behavior cloning trains on expert state distributions but at test time visits states induced by its own errors, causing compounding error (a form of distribution shift / “state drift”). Classical RL algorithms include temporal-difference learning (Sutton, 1988; Sutton et al., 1998) and value-based control such as Q-learning (Watkins and Dayan, 1992), while policy-gradient methods (Williams, 1992; Sutton et al., 1999) directly optimize expected return via likelihood-ratio gradients (REINFORCE) and later stabilized large-scale learning through variants like trust-region policy optimization (TRPO) (Schulman et al., 2015) and off-policy actor-critic methods for continuous control (e.g., DDPG (Lillicrap et al., 2015)) and maximum-entropy actor-critic (e.g., SAC (Haarnoja et al., 2018)). Empirically, deep RL’s modern resurgence is often associated with representation learning + RL on high-dimensional inputs (e.g., DQN (Mnih et al., 2013)).

A large body of work blends supervised and RL to get the best of both: (i) Imitation + online correction methods like DAgger (Ross et al., 2010) explicitly combine supervised learning with interactive data collection to mitigate distribution shift ; (ii) Inverse RL / MaxEnt IRL reframes imitation as learning a

reward/cost model that explains expert behavior, with maximum-entropy formulations giving a principled probabilistic objective (Ziebart et al., 2008); (iii) Adversarial imitation (GAIL) (Ho and Ermon, 2016) avoids explicit reward learning by matching occupancy measures via a GAN-like discriminator, typically trained with policy optimization; (iv) Learning from demonstrations in deep RL injects supervised losses and/or demonstration replay into RL to improve exploration and sample efficiency—e.g., DQfD (Hester et al., 2017) combines TD learning with supervised large-margin imitation terms, and demonstration-augmented continuous-control methods address sparse-reward exploration failures (Nair et al., 2018); and (v) Trajectory-optimization-guided policy learning (guided policy search (Levine and Koltun, 2013)) explicitly produces supervised targets for a policy network from trajectory optimization / local controllers, bridging optimal control, RL, and supervised regression. In modern LLM alignment, the same hybrid template appears as “Supervised fine-tuning + preference-based RL”: InstructGPT (Ouyang et al., 2022) first performs supervised fine-tuning on demonstrations and then applies RL from human feedback (RLHF), while more recent approaches like Direct Preference Optimization (DPO) (Rafailov et al., 2023) recast parts of RLHF into a supervised-style classification objective — illustrating an active trend of recovering supervised-like training signals even when the underlying goal is preference/reward optimization.

Finally, control as inference is also a closely related topic (Millidge et al., 2020; O’Donoghue et al., 2020; Rawlik et al., 2012; Ito and Kashima, 2024; Tarbouriech et al., 2023), and we point the reader to Levine (2018) for more details.

Training LLMs for strong reasoning abilities. A few different approaches for post-training have demonstrated success, including supervised fine-tuning on human-crafted high quality demonstrations (Wang et al., 2023b), iterative supervised training on self-generated good quality responses (Zelikman et al., 2022; Gulcehre et al., 2023), reinforcement learning from a learned reward model on human preferences (Ouyang et al., 2022), and more recently preference-based contrastive learning (Rafailov et al., 2023; Pang et al., 2024). In our work, we focus on recovering the cross-entropy based classification objective in an RL training pipeline, fundamentally differing from the prior works. Since the recent advent of RLVR, multiple followup works have studied the RLVR pipeline (Zeng et al., 2025; Liu et al., 2025b; Khatri et al., 2025) and proposed alternative algorithms such as Dr.GRPO (Liu et al., 2025b), DAPO (Yu et al., 2025a), GSPO (Zheng et al., 2025) and CISPO (MiniMax et al., 2025), RAFT (Xiong et al., 2025a). The idea of normalizing advantages by mean reward, similar to ours, has been explored in (Huang et al., 2025), but whereas we normalize advantage by group mean reward (mean reward over the rollouts associated with a particular prompt), Huang et al. (2025) normalizes by the batch mean reward (mean reward over all prompts in a batch of policy gradient updates). Finally, recent work such as Zhang et al. (2025) has also studied how RL training is influenced by pretraining and midtraining in toy didactic settings, establishing the importance of good pretraining/midtraining for the success of RL, similar to Gandhi et al. (2025).

On exploration for reinforcement learning for LLMs. Exploration, or taking actions to discover new information, is a widely studied topic in reinforcement learning. A closely related topic is *curiosity*, where an agent seeks new information about its environment via interactions. *Intrinsic motivation* is a popular notion for curiosity, where the agent is driven by an exploration bonus that is not necessarily related to the task to be achieved (Schmidhuber, 1991, 2007). Followup works have built on this notion to mitigate problems of sparse reward (reward is observed at a very belated phase of interactions) or no reward at all (Pathak et al., 2017, 2019; Eysenbach et al., 2018; Burda et al., 2018; Sharma et al., 2019; Yang et al., 2024c; Houthoofd et al., 2016). Count-based bonuses have also been introduced as a way of computing intrinsic motivation (Bellemare et al., 2016). Prompt-level reweighting of gradients has also been studied (Yu et al., 2025b), though under a different context (self-training) and a different weighting mechanism. Finally, adding noise to network parameters or optimization has been another line of work to improve exploration during RL training (Fortunato et al., 2017; Ishfaq et al., 2024b,a, 2025). Maximum entropy RL, the principle where one attempts to recover an agent that achieves high reward but is as stochastic as possible, can be seen as another attempt at solving exploration for classical RL (Haarnoja et al., 2018; Boucher et al., 2025; Eysenbach and Levine, 2022; Dong et al., 2025). In summary, exploration-exploitation tradeoff (Sutton, 1988; Auer et al., 2002; Thompson, 1933) has been a crucial topic for ensuring RL agents’ success.

More recently, exploration has emerged as an important topic for building modern LLM based systems. There are two types of exploration to consider. The first is *inference-time exploration*, where an agent has to efficiently gather information during deployment by strategically choosing its interactions with its environment, Tajwar et al. (2025) is an important work in this line of research. More importantly,

pass@k degradation (mode collapse) during RLVR (Yue et al., 2025; Wu et al., 2026; GX-Chen et al., 2025) has prompted research into *train-time exploration*, where the challenge is to go beyond the pretrained model’s capabilities and discover new knowledge. Primary approaches include directly optimizing for pass@k (Walder and Karkhanis, 2025; Tang et al., 2025; Chen et al., 2025d), curriculum learning (Tajwar et al., 2025; Chen et al., 2025c; Setlur et al., 2025; Motwani et al., 2025), learning from additional hints or abstractions (Qu et al., 2025; Chen et al., 2025a; Anonymous, 2025b), increasing number of rollouts to prevent RL gains from saturating (Hu et al., 2025), employing data curation algorithm to redirect effort to problems with low success rate (Nguyen et al., 2025b), leveraging expert guidance (Chang et al., 2024; Qu et al., 2026), or differential smoothing by penalizing entropy on low reward trajectories and encouraging entropy on high reward trajectories. Entropy based bonuses to encourage exploration during RL training (Hao et al., 2025; Chen et al., 2025b; Cheng et al., 2025a; Wang et al., 2025; Anonymous, 2025a; Ged and Veiga, 2024) is another popular line of work for improving exploration. A few modern approaches for exploration bonus utilized for LLM training are Song et al. (2025); Tuyls et al. (2025). The idea of curiosity-driven exploration from classical RL discussed above has also been adopted for LLMs (Dai et al., 2025). Although some works have reported pass@k degradation during RL training, others have found the opposite results. For example, ProRL (Liu et al., 2025a) has shown that RL training on a mixture of reasoning puzzles (Stojanovski et al., 2025) can improve pass@k on a heldout reasoning task. Similarly, Yuan et al. (2025) has shown that LLMs can learn new skills via RL by composing old ones, showing the promise of going beyond pre-training knowledge, and Cheng et al. (2025b) also found pass@k to improve, particularly on tasks less likely to appear during the pre-training stage. Ray interference (Schaul et al., 2019) has been proposed as an explanation for the observed pass@k degradation. Overall, this line of research remains important as focus moves to LLMs discovering new information during RL training and it is therefore an ongoing field of research.

Generalization in reinforcement learning. Research on generalization in reinforcement learning asks a core question: does an agent learn principles that transfer beyond the exact environments it trained in, or does it just memorize experience (Zhang et al., 2018b,a; Schaul et al., 2019; Bengio et al., 2020)? Empirical work shows deep RL agents often overfit to training seeds, visuals, or dynamics, performing poorly on new levels, layouts, or slightly shifted physics. To study this, researchers built procedural benchmarks (like CoinRun/Procgen (Cobbe et al., 2020)) and multi-task suites (e.g., robotics task collections (Yu et al., 2021; Atamuradov, 2025) or LLM sequential decision-making task suites (Tajwar et al., 2025)) that separate train and test environments. A major line of work improves generalization through regularization (Kostrikov et al., 2021) and invariances (Zhang et al., 2021) — especially data augmentation (Laskin et al., 2020; Raileanu et al., 2021), mixup-style methods (WANG et al., 2020), and representation learning tricks (Higgins et al., 2018; Srinivas et al., 2020; Wang et al., 2021) that make policies rely less on superficial visual details. Another branch focuses on task and domain shift, using meta-RL (Duan et al., 2016; Finn et al., 2017; Liang et al., 2024), multi-task learning (Brunskill and Li, 2013), domain randomization, and distributionally robust RL (Clavier et al., 2022; Lu et al., 2024; Shi et al., 2025) to handle new tasks or uncertain dynamics. Recent lines of work has also directly studied exploration as a mean to achieve generalization in RL (Jiang et al., 2023), and have shown that simple architectural changes and scale can often improve generalization in the ProcGen benchmark (Jesson and Jiang, 2024). On the theory side, classical PAC-MDP (Strehl et al., 2009) and robust MDP frameworks (Nilim and Ghaoui, 2004; Iyengar, 2005) formalize when policies learned from limited samples can be expected to work in new situations. In the LLM settings, analogous questions appear in RLHF (Ouyang et al., 2022), where reinforcement learning is used to align models with human preferences, and researchers now study how this training affects generalization to unseen prompts, behaviors, and user distributions (Kirk et al., 2024; Lin et al., 2024; Lambert et al., 2024; Jia, 2024; Li et al., 2026). Overall, the field has moved from “can RL learn?” to “what exactly does it learn, and when does that knowledge transfer?”

B Theoretical Results

Here we present the proofs of theorems mentioned in the main paper. First we restate and prove [Theorem 1](#).

Theorem 3 (Restatement of Theorem 1). *The gradient of the maximum likelihood objective admits the following conditional expectation representation:*

$$\nabla_{\theta} J_{\text{ML}}(x) = \mathbb{E}[\nabla_{\theta} \log m_{\theta}(z | x) | f(z) = y^*(x)].$$

Proof. Recall the standard REINFORCE identity for the gradient of the pass rate:

$$\nabla_{\theta} p_{\theta}^{\text{pass}}(x) = \nabla_{\theta} \mathbb{E}_{z \sim m_{\theta}(\cdot | x)}[\mathbb{I}\{f(z) = y^*(x)\}] = \mathbb{E}_{z \sim m_{\theta}(\cdot | x)}[\mathbb{I}\{f(z) = y^*(x)\} \nabla_{\theta} \log m_{\theta}(z | x)].$$

The gradient of the maximum likelihood objective is:

$$\nabla_{\theta} J_{\text{ML}}(x) = \nabla_{\theta} \log p_{\theta}^{\text{pass}}(x) = \frac{\nabla_{\theta} p_{\theta}^{\text{pass}}(x)}{p_{\theta}^{\text{pass}}(x)} = \frac{\mathbb{E}_{z \sim m_{\theta}(\cdot | x)}[\mathbb{I}\{f(z) = y^*(x)\} \nabla_{\theta} \log m_{\theta}(z | x)]}{\mathbb{E}_{z \sim m_{\theta}(\cdot | x)}[\mathbb{I}\{f(z) = y^*(x)\}]}.$$

By the definition of conditional expectation for an event A with $\mathbb{P}(A) > 0$:

$$\mathbb{E}[X | A] = \frac{\mathbb{E}[X \cdot \mathbb{I}_A]}{\mathbb{P}(A)}.$$

Letting $X = \nabla_{\theta} \log m_{\theta}(z | x)$ and $A = \{z : f(z) = y^*(x)\}$, and noting that $p_{\theta}^{\text{pass}}(x) = \mathbb{P}(A)$, we obtain:

$$\nabla_{\theta} J_{\text{ML}}(x) = \mathbb{E}[\nabla_{\theta} \log m_{\theta}(z | x) | f(z) = y^*(x)].$$

□

Next, we restate and prove [Theorem 2](#).

Theorem 4 (Restatement of Theorem 2). *The estimator $\hat{g}_N(x)$ is an unbiased estimator for the MAXRL gradient of order $T = N$, i.e.,*

$$\mathbb{E}[\hat{g}_N(x)] = \nabla_{\theta} J_{\text{MAXRL}}^{(N)}(x).$$

Proof. Conditioned on $K \geq 1$, the successful samples are i.i.d. draws from the success-conditioned distribution, so by Theorem 1:

$$\mathbb{E}[\hat{g}_N(x) | K \geq 1] = \nabla_{\theta} \log p_{\theta}^{\text{pass}}(x).$$

Since $\hat{g}_N(x) = 0$ when $K = 0$:

$$\mathbb{E}[\hat{g}_N(x)] = \nabla_{\theta} \log p_{\theta}^{\text{pass}}(x) \cdot \mathbb{P}(K \geq 1) = \nabla_{\theta} \log p_{\theta}^{\text{pass}}(x) \cdot \text{pass}@N(x).$$

Writing $p = p_{\theta}^{\text{pass}}(x)$ and using $\text{pass}@k(x) = 1 - (1 - p)^k$:

$$\frac{\nabla_{\theta} p}{p} \cdot (1 - (1 - p)^N) = \nabla_{\theta} p \sum_{k=1}^N (1 - p)^{k-1} = \sum_{k=1}^N \frac{1}{k} \nabla_{\theta} \text{pass}@k(x) = \nabla_{\theta} J_{\text{MAXRL}}^{(N)}(x),$$

where the second equality uses $\nabla_{\theta} \text{pass}@k(x) = k(1 - p)^{k-1} \nabla_{\theta} p$.

□

C More on Unifying Weight-Function View on RL Objectives

Here we provide full derivations following [Section 5](#). Recall that we want to express population-level gradients of different objectives in the following form:

$$\nabla_{\theta} J = \mathbb{E}_{x \sim \rho} [w(p_{\theta}(x)) \nabla_{\theta} p_{\theta}(x)]$$

where $p_{\theta}(x) = p_{\theta}^{\text{pass}}(x)$ and $w(p_{\theta}(x))$ is the weighting function. In this section, we show that all objectives of our consideration can be written in this form. Furthermore, we will derive the weighting function $w(p_{\theta}(x))$ for each of them.

Classical RL (REINFORCE). For classical reinforcement learning, i.e., the REINFORCE objective, we have:

$$\begin{aligned} J_{\text{RL}} &= \mathbb{E}_{x \sim \rho} [\mathbb{E}_{z \sim m_{\theta}(\cdot|x)} [r(x, z)]] \\ &= \mathbb{E}_{x \sim \rho} [\mathbb{E}_{z \sim m_{\theta}(\cdot|x)} [\mathbb{I} \{f(z) = y^*(x)\}]] \\ &= \mathbb{E}_{x \sim \rho} [p_{\theta}^{\text{pass}}(x)] \end{aligned}$$

Therefore, its gradient is:

$$\nabla_{\theta} J_{\text{RL}} = \mathbb{E}_{x \sim \rho} [\nabla_{\theta} p_{\theta}^{\text{pass}}(x)]$$

giving the corresponding w_{RL} to be 1.

GRPO. Our analysis is similar to that of [Davis and Recht \(2025\)](#). The gradient of the population level GRPO objective gradient can be written as:

$$\nabla_{\theta} J_{\text{GRPO}} = \mathbb{E}_{x \sim \rho} \left[\mathbb{E}_{z \sim m_{\theta}(\cdot|x)} \left[\left(\frac{r(x, z) - \mathbb{E}_{z \sim m_{\theta}(\cdot|x)} [r(x, z)]}{\sqrt{\text{Var}_{z \sim m_{\theta}(\cdot|x)} [r(x, z)]}} \right) \nabla_{\theta} \log m_{\theta}(z|x) \right] \right]$$

Since we consider a binary reward setting, we have $\mathbb{E}_{z \sim m_{\theta}(\cdot|x)} [r(x, z)] = p_{\theta}^{\text{pass}}(x)$. Similarly, considering the variance of a Bernoulli random variable, we get:

$$\text{Var}_{z \sim m_{\theta}(\cdot|x)} = p_{\theta}^{\text{pass}}(x) (1 - p_{\theta}^{\text{pass}}(x))$$

Therefore, the objective becomes:

$$\nabla_{\theta} J_{\text{GRPO}} = \mathbb{E}_{x \sim \rho} \left[\frac{1}{\sqrt{p_{\theta}^{\text{pass}}(x) (1 - p_{\theta}^{\text{pass}}(x))}} \nabla_{\theta} p_{\theta}^{\text{pass}}(x) \right]$$

which thereby gives us the weighting function to be $1/\sqrt{p_{\theta}(x)(1 - p_{\theta}(x))}$, as desired.

Maximum Likelihood (ML). The maximum likelihood objective is given by

$$J_{\text{ML}} = \mathbb{E}_{x \sim \rho} [\log p_{\theta}^{\text{pass}}(x)].$$

Taking its gradient with respect to θ and applying the chain rule, we obtain

$$\begin{aligned} \nabla_{\theta} J_{\text{ML}} &= \mathbb{E}_{x \sim \rho} [\nabla_{\theta} \log p_{\theta}^{\text{pass}}(x)] \\ &= \mathbb{E}_{x \sim \rho} \left[\frac{1}{p_{\theta}^{\text{pass}}(x)} \nabla_{\theta} p_{\theta}^{\text{pass}}(x) \right] \end{aligned}$$

This shows that the weighting function for the maximum likelihood objective is $1/p_{\theta}^{\text{pass}}(x)$, as we claimed in [Table 2](#).

MAXRL. Finally, we consider the objective $J_{\text{MAXRL}}^{(T)}$.

Proposition 5. *For MAXRL with order T , we can rewrite it as*

$$\nabla_{\theta} J_{\text{MAXRL}}^{(T)} = \mathbb{E}_{x \sim \rho} [w(p_{\theta}(x)) \nabla_{\theta} p_{\theta}(x)],$$

where

$$w_T(p) = \sum_{k=1}^T (1-p)^{k-1} = \frac{1 - (1-p)^T}{p}.$$

Proof. From Equation (7), we have:

$$\nabla_{\theta} J_{\text{MAXRL}}^{(T)}(x) = \sum_{k=1}^T \frac{1}{k} \nabla_{\theta} \text{pass}@k(x).$$

Using $\text{pass}@k(x) = 1 - (1-p)^k$ where $p = p_{\theta}^{\text{pass}}(x)$:

$$\nabla_{\theta} \text{pass}@k(x) = k(1-p)^{k-1} \nabla_{\theta} p.$$

Substituting:

$$\nabla_{\theta} J_{\text{MAXRL}}^{(T)}(x) = \sum_{k=1}^T \frac{1}{k} \cdot k(1-p)^{k-1} \nabla_{\theta} p = \left(\sum_{k=1}^T (1-p)^{k-1} \right) \nabla_{\theta} p = w_T(p) \nabla_{\theta} p_{\theta}^{\text{pass}}(x).$$

Taking the expectation over $x \sim \rho$ completes the proof. □

D Additional Details on ImageNet Experiments

D.1 Training Procedure

Let \mathcal{X} be the input space and \mathcal{Y} be the label space. Let π_θ denote our model: given an input image $x \in \mathcal{X}$, $\pi_\theta(y|x)$ is model’s predicted probability of image x belonging to class $y \in \mathcal{Y}$. For an input image and label pair $(x, y^*(x))$, the cross-entropy loss is:

$$\mathcal{L}_{\text{CE}}(x, y^*(x); \pi_\theta) = -\log \pi_\theta(y^*(x)|x)$$

On the other hand, the corresponding RL objective for the same pair is:

$$\mathcal{L}_{\text{RL}}(x, y^*; \pi_\theta) = -\mathbb{E}_{y \sim \pi_\theta(\cdot|x)}[-\log \pi_\theta(y|x) \cdot \hat{A}(y|x)]$$

where the expectation is computed using Monte-Carlo sampling K rollouts of y from $\pi_\theta(\cdot|x)$. GRPO, REINFORCE and MAXRL vary only in the calculation of the advantage $A(y|x)$. Concretely, let $y^{(1)}, \dots, y^{(K)}$ be our K rollouts, sampled from the conditional probability distribution $\pi_\theta(\cdot|x)$. We operate under a binary reward setting, meaning the reward function $r(x, y)$ is:

$$r(x, y) = \mathbb{I}[y = y^*(x)] = \begin{cases} 1, & \text{if } y = y^*(x) \\ 0, & \text{otherwise} \end{cases}$$

Given this reward, we calculate advantage under GRPO, REINFORCE and MAXRL as follows:

$$\hat{A}_{\text{GRPO}}(x, y) = \frac{r(x, y) - \hat{\mu}}{\hat{\sigma}}$$

$$\hat{A}_{\text{REINFORCE}}(x, y) = r(x, y) - \hat{\mu}$$

$$\hat{A}_{\text{MAXRL}}(x, y) = \frac{r(x, y) - \hat{\mu}}{\hat{\mu}}$$

where $\hat{\mu} = \frac{\sum_{i=1}^K r(x, y^{(i)})}{K}$, $\hat{\sigma} = \sqrt{\frac{\sum_{i=1}^K (r(x, y^{(i)}) - \hat{\mu})^2}{K}}$ is the mean and standard deviation of rewards of the sampled rollouts.

Finally, at each training step, a batch of (input image, label) pairs are collected from the training dataset. The above computation gives us per (input image, label) loss, we average them over all the pairs in a given batch to calculate the final loss which is then used to update the model via gradient descent.

D.2 Training Hyperparameters

We use the following set of hyperparameters in all ImageNet experiments:

- **Batch size:** 256
- **Number of epochs:** 20
- **Optimizer:** SGD with momentum 0.9, no Nesterov momentum, initial learning rate 0.1. We run a sweep over the learning rate over 0.001, 0.003, 0.01, 0.03, 0.1, 0.3, 0.7, and 1.0. We find the standard learning rate, 0.1, generally works well for all objectives, and report that in our experiments.
- **Learning rate scheduler:** Cosine scheduler (Loshchilov and Hutter, 2017), with linear warmup for the first epoch.
- **Image augmentations:** No augmentations are used for evaluation, we only resize each image to 224x224 and normalize the images by mean and standard deviation of pixel values. For training, in addition to the same resizing and normalizing steps, we also add a random horizontal flip (with probability 0.5) and a random resized crop to 224 (with scale (0.08, 1.0)).
- **Number of rollouts, K:** This is usually varied for different experiments.

All training is done on single L40S GPUs for 15 hours.

D.3 Equivalence of Validation Top-1 Accuracy and Majority Voting Accuracy

In this section, we discuss the validation top-1 accuracy metric, which is the traditional metric used in image classification. Formally, validation accuracy for a single image and label pair $(x, y^*(x))$ is defined as:

$$\text{Accuracy}(x, y^*(x); \pi_\theta) = \mathbb{I} \left[\arg \max_{y \in \mathcal{Y}} \pi_\theta(y|x) = y^*(x) \right] = \begin{cases} 1, & \text{if } \arg \max_{y \in \mathcal{Y}} \pi_\theta(y|x) = y^*(x) \\ 0, & \text{otherwise} \end{cases}$$

which is then averaged over all validation examples for the final metric. In other words, validation accuracy is the same as majority voting accuracy (Wang et al., 2023a) in traditional LLM chain-of-thought reasoning tasks.

D.4 Pass@k Calculation

To calculate pass@k from a generative model, one usually samples $T \geq k$ rollouts from the model, calculate success or failure from each of them, and then uses an appropriate statistical estimator for pass@k (Chen et al., 2021; Yue et al., 2025). However, since there is no latent reasoning process involved in our didactic ImageNet experiments and since we can directly calculate the model likelihood of label $y \in \mathcal{Y}$ for an input image $x \in \mathcal{X}$, namely $\pi_\theta(y|x)$, we can also analytically compute pass@k without sampling as well. Formally, in all ImageNet experiments, we calculate pass@k for an example (image, label) pair $(x, y^*(x))$ as follows:

$$\text{Pass@k}(x, y^*(x); \pi_\theta) = 1 - (1 - \pi_\theta(y^*(x)|x))^k$$

The average pass@k is then obtained by averaging the above quantity over all example pairs in the validation dataset.

D.5 Gradient Norm Analysis

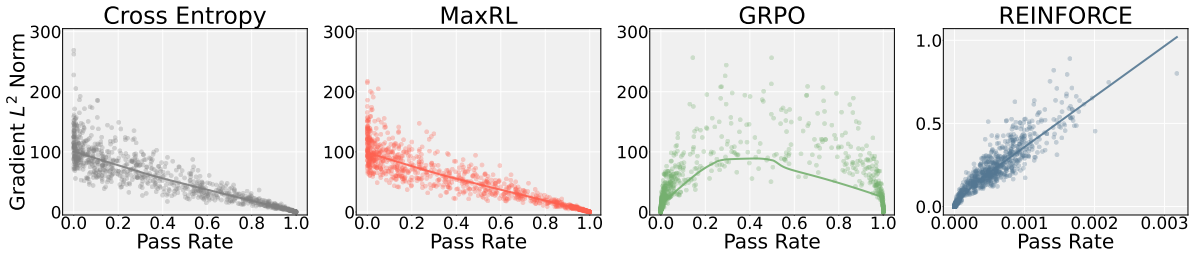


Figure 8: (ImageNet Gradient Norm Analysis) Scatter plot, where each point has the pass rate (model’s predicted probability of the correct class) of a particular image in the x-axis, and gradient L^2 norm for that image in the y-axis, for 1000 randomly selected images from the ImageNet validation dataset after 1500 steps of training on a ResNet-50 model. Sampling based algorithms’ (MAXRL, GRPO and REINFORCE) gradients are calculated using 131,072 rollouts per example to reduce sampling error and estimate the population level gradient. Cross Entropy and MAXRL have similar scatter plot: with high gradient norm for hard inputs (pass rate close to 0) and lower gradient norm for the easier ones (pass rate close to 1). In contrast, highest gradient norm for GRPO is on medium difficulty (pass rate close to 0.5) inputs, with hard inputs having very low gradient norm. Finally, REINFORCE fails to produce any significant gradient norm and its pass rate is confined below 0.003 after 1500 steps, demonstrating its difficulty to learn in this setting.

Figure 8 shows the correlation between gradient norm and pass rate (model’s predicted probability of the correct class) for a particular image on different objectives. We see that cross-entropy and MAXRL have similar scatter plot: with high gradient norm for hard inputs (pass rate close to 0) and lower gradient norm for the easier ones (pass rate close to 1). In contrast, highest gradient norm for GRPO is on medium difficulty (pass rate close to 0.5) inputs, with hard inputs having very low gradient norm. Finally, REINFORCE fails to produce any significant gradient norm compared to the other objectives and its pass rate is confined below 0.003 after 1500 steps, demonstrating its difficulty to learn in this setting. This is also reflected in our other results, where REINFORCE does not show any signs of learning. We attribute this to the very low gradient norm: since the randomly initialized model has pass rate 0.001 in expectation over all inputs, REINFORCE fails to produce sufficiently large gradients during training and therefore stalls in model improvement. One caveat: REINFORCE’s failure maybe due to us training the model from scratch — on a pretrained model, it indeed produces gradients but still shows poor gradient norm on hard inputs (see Figure 6).

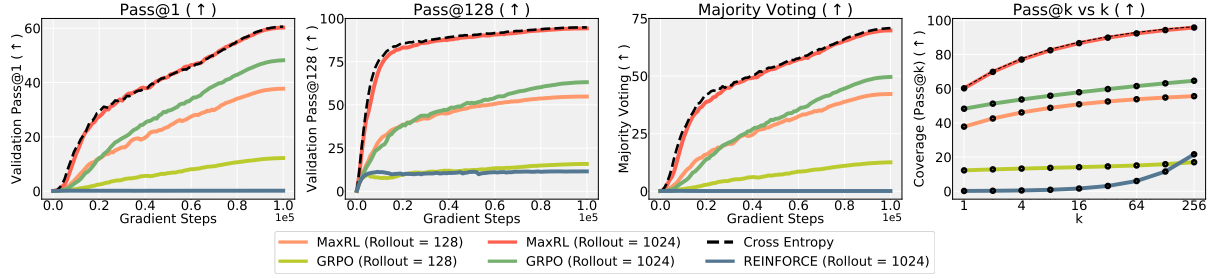


Figure 9: (Additional ImageNet results) On the didactic image classification setting, MAXRL outperforms and scales better than GRPO with additional compute, and approaches the same performance as maximum likelihood training via cross-entropy given sufficient number of rollouts (≥ 1024). Note that REINFORCE remains flat, since the initial model’s pass rate is low ($\sim 0.1\%$) and REINFORCE fails to generate significant gradient signal (Figure 8). From left, the plots show Pass@1, Pass@128, Majority Voting Accuracy (equivalent to traditional validation top-1 accuracy in image classification, see Appendix D.3), and coverage of the final checkpoint, respectively.

D.6 More Experimental Results

Here we present additional experimental results. In particular, **(1)** we compare against GRPO with varying number of rollouts, **(2)** record additional metrics such as majority voting accuracy (i.e., validation top-1 accuracy), and **(3)** show the resulting coverage (pass@k vs k) from different objectives. Figure 9 records our findings: MAXRL outperforms and scales better than GRPO with additional compute. While GRPO improves performance if given more compute unlike REINFORCE, it remains suboptimal compared to MAXRL and supervised cross-entropy training. Moreover, both GRPO and REINFORCE exhibit worse coverage as their pass@k values are significantly lower compared to MAXRL, corroborating our experiments from other sections.

E Details on Other Training Settings

E.1 Maze

E.1.1 Model Architecture

We adopt a lightweight decoder-only Transformer model following the Qwen2 architecture (Yang et al., 2024a), with a total of approximately $3M$ parameters. The model consists of 4 Transformer layers, each using full self-attention. The hidden size is set to 256, with an intermediate (feed-forward) dimension of 1024, and 4 attention heads per layer. We use grouped query attention with 2 key-value heads. The model employs RMSNorm with $\sigma = 1 \times 10^{-6}$ and uses the SiLU activation function in the feed-forward networks. Rotary positional embeddings (RoPE) (Su et al., 2021) are applied with $\theta = 1,000,000$, and the maximum sequence length is 512 tokens. The vocabulary size is 32 tokens, and input and output embeddings are tied. The model is trained and evaluated using bfloat16 precision, with attention dropout set to 0. The architecture follows a standard causal language modeling setup with autoregressive decoding.

E.1.2 Task Description

Mazes are procedurally generated using Prim’s algorithm (Prim, 1957), and task difficulty is controlled by the grid size. We use a symbolic tokenization to represent both the maze layout and the navigation policy, with tokens drawn from a small, discrete vocabulary.

The input sequence describes a two-dimensional grid in row-major order. Each cell is represented by a single token indicating its type (e.g., WALL, PATH, START, or GOAL). Rows are separated by a dedicated NEWLINE token, and the entire grid is delimited by special boundary tokens marking the beginning (GRID_START) and end (GRID_END) of the grid description. Following the maze specification, the model autoregressively generates a sequence of navigation actions drawn from a fixed action vocabulary (e.g., directional moves) and terminates by a DONE token.

Below, we provide an example data instance following this format.

7*7 Maze Example Model Input and Output Format

Input:

```
<bos> GRID_START WALL WALL WALL WALL WALL WALL WALL NEWLINE WALL START WALL PATH PATH PATH WALL  
NEWLINE WALL PATH WALL PATH WALL WALL WALL NEWLINE WALL PATH PATH PATH PATH PATH WALL NEWLINE WALL  
PATH WALL WALL WALL PATH WALL NEWLINE WALL PATH WALL PATH PATH GOAL WALL NEWLINE WALL WALL WALL WALL  
WALL WALL WALL NEWLINE GRID_END PATH_START
```

Output:

```
RIGHT RIGHT RIGHT RIGHT DOWN DOWN DOWN DOWN DONE <eos>
```

For reference, we also visualize one typical successful trajectory and one representative failed prediction in Figure 10.

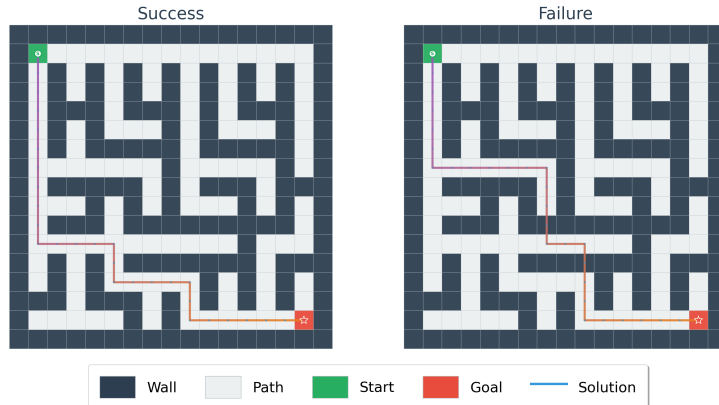


Figure 10: (Maze Data Visualization) The left plot shows a successful navigation trajectory, while the right plot illustrates a failure case produced by the trained model, where the generated action sequence deviates from the correct path before reaching the goal.

E.1.3 Training Setups

To ensure sufficient task complexity and rigorous evaluation, we construct a training set of 1 million distinct 17×17 mazes and a test set of 256 non-overlapping samples. We first pretrain the model from scratch, where it is trained to follow a provided ground-truth trajectory for each maze. During SFT, we use a learning rate of 5×10^{-4} with the AdamW optimizer (Kingma and Ba, 2017; Loshchilov and Hutter, 2019) and train for 1,500 steps with a batch size of 32. This pretraining stage initializes the model with the basic output format required for representing maze-solving trajectories. Subsequently, we perform reinforcement learning (RL) training. By default, we use a data batch size of 32, a rollout number of 128, a learning rate of 1×10^{-4} . We update the model parameters only once per RL step (fully on-policy setting (Tajwar et al., 2024)) to ensure all trajectories are on-policy, with 20,000 training steps. On this basis, we compare the performance of different RL algorithms.

In addition, we fix the data batch size to 256 and vary the rollout number from 4, 8, 16, 32, 64 to 128 when illustrating the scaling trends of GRPO and our MaxRL algorithm. Each training run uses 4 RTX 4090 GPUs and is conducted for at least 9,000 steps. Given the small number of model parameters, the model is no longer memory-bound, so we modified the rollout engine (instead of using the default vLLM (Kwon et al., 2023) engine) in the `verl` (Zhang et al., 2024; Sheng et al., 2024) codebase to increase sampling parallelism and speed up training.

E.2 GSM8K

E.2.1 Prompt template

We conduct reinforcement finetuning for the SmoLM2-360M-Instruct model (Allal et al., 2025) on GSM8K (Cobbe et al., 2021) training set for 1500 steps. We use GSM8K-Platinum (Vendrow et al., 2025), a revised version of the original GSM8K test that is non-overlapping with the GSM8K train dataset, for evaluation. An example of prompt and the model’s response are shown below. In particular, we take each individual problem, append it with “ Let’s think step by step and output the final answer within `\boxed{}`.” and process the string through the SmoLM2-360M-Instruct model’s tokenizer chat-template to obtain the final prompt for the model.

SmoLM2-360M-Instruct

```
<|im_start|>system
You are a helpful AI assistant named SmoLM, trained by Hugging Face<|im_end|>
<|im_start|>user
Jeannie hikes the 12 miles to Mount Overlook at a pace of 4 miles per hour, and then returns at
a pace of 6 miles per hour. How long did her hike take, in hours? Let’s think step by step and
output the final answer within \boxed{}. <|im_end|>
<|im_start|>assistant
To find the time it takes for Jeannie to hike the entire trip, we need to calculate her total
travel time and then divide it by the rate at which she travels.
Jeannie hikes the 12 miles at a pace of 4 miles per hour. So, the time taken is 12 miles / 4
miles per hour = 3 hours.
Next, she returns at a pace of 6 miles per hour. So, the time taken to return is 12 miles / 6
miles per hour = 2 hours.
Therefore, the total time taken is 3 hours + 2 hours = 5 hours.
So, the final answer is 5 hours. \boxed{5} <|im_end|>
```

E.2.2 Hyperparameters

Detailed hyperparameter setup is provided in Table 5. For other baselines, we use $T = 16$ for PKPO, which means we optimize Pass@16 as the objective, and $\gamma = 0.05$ for differential smoothing.

Table 5: Training hyperparameters for SmoLM2-360M-Instruct on GSM8K.

Training hyperparameters for SmoLM2-360M-Instruct

Parameter	Value	Parameter	Value
Base model	SmoLM2-360M-Instruct	Training set	GSM8K
Test set	GSM8K	Prompts per batch	256
Generations per prompt	128	Grad update per RL step	1
Max prompt length	512	Max response len	2048
Learning rate	1×10^{-5}	Training Steps	1500
KL coeff	0.0	Entropy coeff	0.0
Rollout temp	1.0	Validation top_p	0.95
Validation temp	0.6	Device	$8 \times$ Nvidia GH200

E.3 Qwen3 Training

E.3.1 Prompt template

We use the Qwen-math template (Yang et al., 2024a; Qwen et al., 2025; Yang et al., 2024b) for formatting our prompts. We show an example prompt (Yu et al., 2025a) after formatting through our template below. In particular, we take each individual problem, append it with “\nPlease reason step by step, and put your final answer within `\\boxed{}`.” and process the string through the SmolLM2-360M-Instruct model’s tokenizer chat-template to obtain the final prompt for the model.

Qwen Math Prompt Template

```
<|im_start|>system
Please reason step by step and put the final answer in \\boxed{}. <|im_end|>
<|im_start|>user
Denote by  $S(n)$  the sum of the digits of the positive integer  $n$ . Find all the solutions of the
equation  $n(S(n) - 1) = 2010$ . Let’s think step by step and output the final answer within \\boxed{}.
<|im_end|>
<|im_start|>assistant
```

E.3.2 Hyperparameters

Next, we describe the default hyperparameters for our training setup. Since there are many possible alternatives to handle off-policy updates and corresponding importance ratio (Schulman et al., 2017; Shao et al., 2024; Zheng et al., 2025; MiniMax et al., 2025; Yu et al., 2025a), to keep things simple, we choose to train in the fully on-policy setup, meaning we have no importance ratio or associated clipping. Similarly, to avoid tuning additional hyperparameters for each algorithm, following Olmo et al. (2025), we remove KL penalty and also entropy bonus in our default training comparison. Note: we train with GRPO and entropy bonus as a baseline in our SmolLM2-360M-Instruct training on GSM8K, results are recorded in Table 4: MAXRL outperform this variant, showing that entropy bonus does not fully mitigate issues resulting from GRPO though it can slightly mitigate it, as also observed by Yue et al. (2025).

We generate all training rollouts using temperature 1.0, and do not use special sampling techniques. Similarly, we also do not use any adaptive sampling (Yu et al., 2025a) or fixes for inference-training logit mismatch (He and Lab, 2025; Khatri et al., 2025) Finally, for evaluation, we follow the same protocol as Yue et al. (2025), and we run inference with temperature 0.6, top-p sampling parameter 0.95, no top-k or min-p sampling (Nguyen et al., 2025a).

Table 6 shows our default hyperparameter setting.

Table 6: Training hyperparameters for Qwen3-1.7B-Base and Qwen3-4B-Base training.

Training hyperparameters for Qwen3-1.7B-Base and Qwen3-4B-Base

Parameter	Value	Parameter	Value
Base model	Qwen3-1.7B-Base, Qwen3-4B-Base	Prompts per batch	256
Generations per prompt	16	Grad update per RL step	1
Max prompt length	1024	Max response len	4096
Learning rate	1×10^{-6}	Training Steps	1000
KL coeff	0.0	Entropy coeff	0.0
Rollout temp	1.0	Validation top-p	0.95
Validation temp	0.6	Device	$32 \times$ Nvidia H200

F Implementation Details on RL Training for LLM Experiments

Our discussion here follows that of (Shafayat et al., 2025). For continuity with existing literature, we use slightly different notations from the rest of the paper for this section. Let x represent a prompt, and let $y \sim \pi(\cdot|x)$ represent sequence of tokens autoregressively sampled from the language model π conditioned on the prompt x . Let π_θ be the current policy, and $\pi_{\theta_{\text{old}}}$ be an older policy (from earlier iterations in training) used for data generation. In our implementation (based on ver1 (Zhang et al., 2024; Sheng et al., 2024)), we use the following general RL objective:

$$\mathcal{J}(\theta) = \mathbb{E}_{x \sim \mathcal{D}, \{y_i\}_{i=1}^G \sim \pi_{\theta_{\text{old}}}(\cdot|x)} \left[\frac{1}{T} \sum_{i=1}^G \sum_{t=1}^{|y_i|} \min \left(w_{i,t}(\theta) \hat{A}_{i,t}, \text{clip}(w_{i,t}(\theta), 1 - \epsilon, 1 + \epsilon) \hat{A}_{i,t} \right) \right]$$

where T is the total number of tokens in the mini-batch (excluding tokens in the prompt etc., since we only compute loss on the model generated tokens), π_θ represents the current LLMs autoregressive probability distribution, $\pi_{\theta_{\text{old}}}$ denote the behavior policy/data generation policy’s probability distribution, $w_{i,t}(\theta)$ is the importance ratio, defined as:

$$w_{i,t}(\theta) = \frac{\pi_\theta(y_{i,t} | x, y_{i,<t})}{\pi_{\theta_{\text{old}}}(y_{i,t} | x, y_{i,<t})}$$

Since we operate fully on-policy, i.e., one RL step per one batch of generated rollouts, this is always one in our experiments, and the clipping parameter ϵ has no effect on our training. $\hat{A}_{i,t}$ represents the advantage for the t -th token in the sequence y_i . The same advantage defined at a sequence level is applied to each token in the sequence, so henceforth we will drop the t from the notation as well.

The main difference between GRPO (Shao et al., 2024), RLOO (Ahmadian et al., 2024) and MAXRL come from their use of different advantage functions. RLOO objective uses the following advantage function:

$$\frac{1}{G} \sum_{i=1}^G [R(y_{(i)}, x) - \frac{1}{G-1} \sum_{j \neq i} R(y_{(j)}, x)]$$

whereas GRPO uses the following advantage function:

$$\hat{A}_i = \frac{r(x, y_i) - \text{mean}(\{r(x, y_i)\}_{i=1}^G)}{\text{std}(\{r(x, y_i)\}_{i=1}^G) + \epsilon}$$

where ϵ is a small number (1×10^{-6}) added to avoid division by zero. Finally, the advantage for MAXRL is follows:

$$\hat{A}_i = \frac{r(x, y_i) - \text{mean}(\{r(x, y_i)\}_{i=1}^G)}{\text{mean}(\{r(x, y_i)\}_{i=1}^G) + \epsilon}$$

Here G is the number of online samples generated. RLOO, GRPO and MAXRL create a dynamic baseline for each sample without needing a separate value function (unlike PPO (Schulman et al., 2017)), effectively estimating the expected return on-the-fly during training. Not having a value networks makes the training much simpler for all three algorithms.

G Pass@k Calculation for Tasks with Sampling

Unlike the ImageNet setting, we can’t usually directly calculate pass@k via accessing the true probability of the correct action. Therefore, we use the default pass@k calculation mechanism in ver1 (Sheng et al., 2024; Zhang et al., 2024), using the bootstrapping low variance unbiased estimator introduced by Chen et al. (2021). This employs generating $n \geq k$ samples per task, counting the number of correct samples $c(x)$ among the n samples, and estimate pass@k as:

$$\text{Pass@k} = \mathbb{E}_{x \sim \rho} \left[1 - \frac{\binom{n-c(x)}{k}}{\binom{n}{k}} \right]$$

H Qwen3 Model Evaluation on Additional Benchmarks

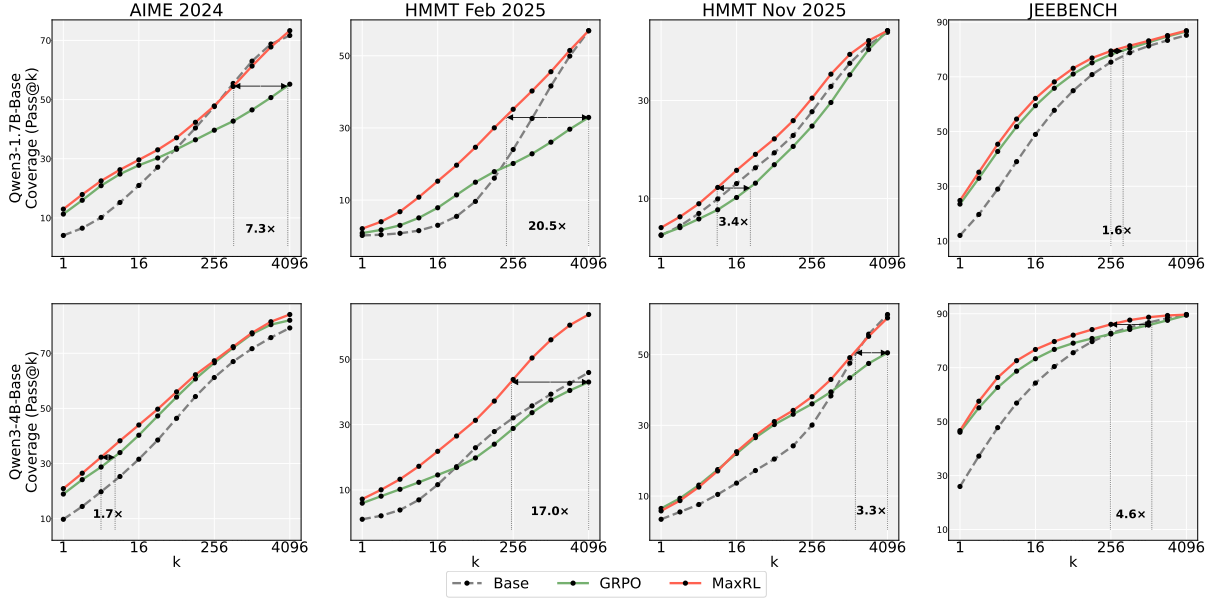


Figure 11: (Evaluation of Qwen3 model training on additional benchmarks) Here we report coverage on 4 additional benchmarks, namely AIME 2024, HMMT Feb 2025 (Balunović et al., 2025a), HMMT Nov 2025 (Balunović et al., 2025b), and JEEBENCH (Arora et al., 2023). MAXRL match or outperform both base model and GRPO, leading up to 20.5 \times speedup compared to GRPO while generating multiple samples with a perfect verifier and maintains similar or better pass@1 performance.

I More on MAXRL Extracting Better Learning Signal During Training

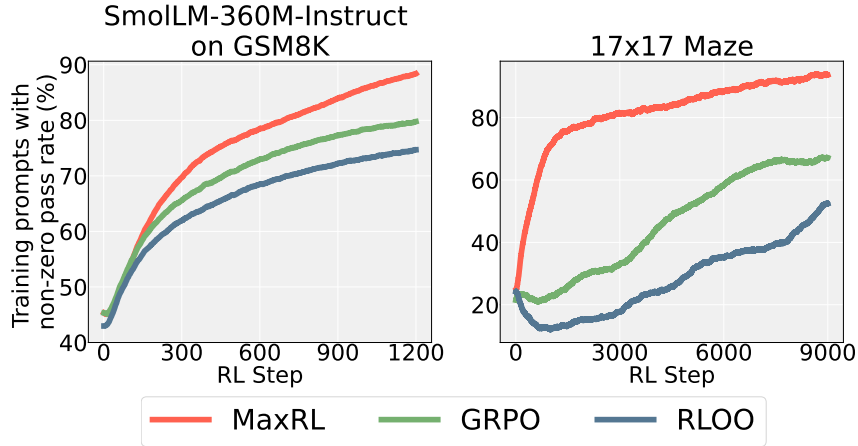


Figure 12: (Fraction of training tasks with non-zero pass rate) Similar to Figure 6, we also record the fraction of training tasks where the model generates at least one correct rollout on maze and SmolLM2-360M-Instruct training on GSM8K. We see the same trends as Figure 7, and MAXRL consistently outperforms both GRPO and RLOO, demonstrating MAXRL ability to generate better learning signal during training, as tasks with zero pass-rate contributes no gradients.

J Additional Results on Qwen3-4B-Base

J.1 Majority Voting Performance

Table 7: (Majority Voting Performance Comparison on Qwen3-4B-Base) We compare the performance of MAXRL in terms of majority voting against the pretrained base model and GRPO.

	AIME 2024 (majority@4096)	AIME 2025 (majority@4096)	BeyondAIME (majority@4096)	MATH-500 (majority@2048)	Minerva (majority@2048)
Base	23.3	23.3	7.0	69.8	18.8
GRPO	23.3	23.3	7.0	72.4	27.2
MaxRL	26.7	26.7	14.0	74.0	28.7

Here we present comparisons across one other metric, majority voting (Wang et al., 2023a), a commonly used verifier free method for scaling test-time compute, where we generate N i.i.d. rollouts from the model for a single task x , group the responses by the final answer, and take the most frequent answer as our outcome. Table 7 shows our results across all five benchmarks, we outperform both the pre-trained base model and GRPO trained model on majority voting across all benchmarks.

J.2 Training Dynamics

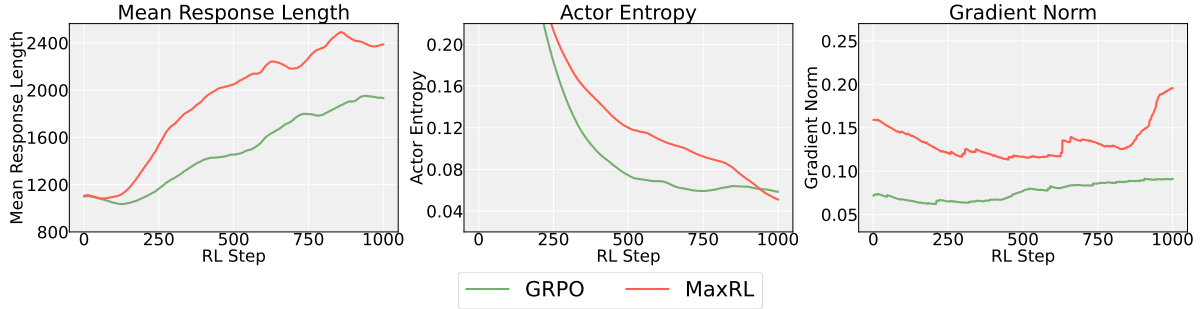


Figure 13: (Additional training dynamics metrics for Qwen3-4B-Base) We show comparison between GRPO and MAXRL in terms of mean response length, entropy of the actor, and gradient norm during training for the Qwen3-4B-Base model. MAXRL generally produces longer chains-of-thought, and also retains higher actor entropy during training. MAXRL also produce larger gradient norm during training.

J.3 Validation Accuracy During Training

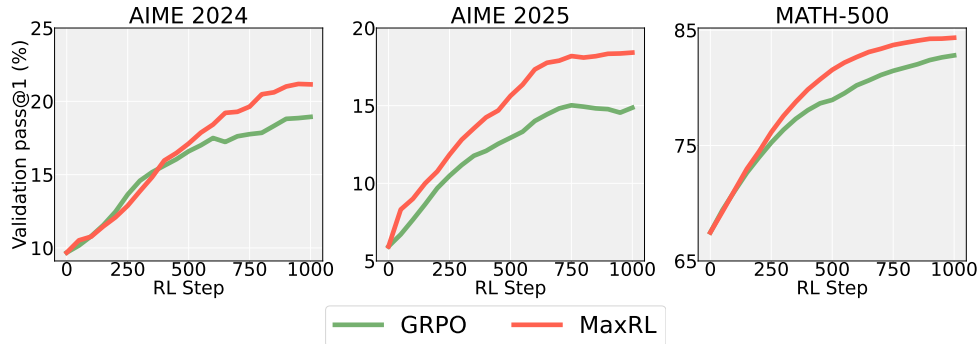


Figure 14: (Qwen3-4B-Base validation pass@1 during training) Pass@1 (estimated using 32 samples) during training of Qwen3-4B-Base, on 3 different evaluation dataset. MAXRL consistently outperform GRPO during training.

K Additional Results on Qwen3-1.7B-Base

K.1 Validation Accuracy During Training

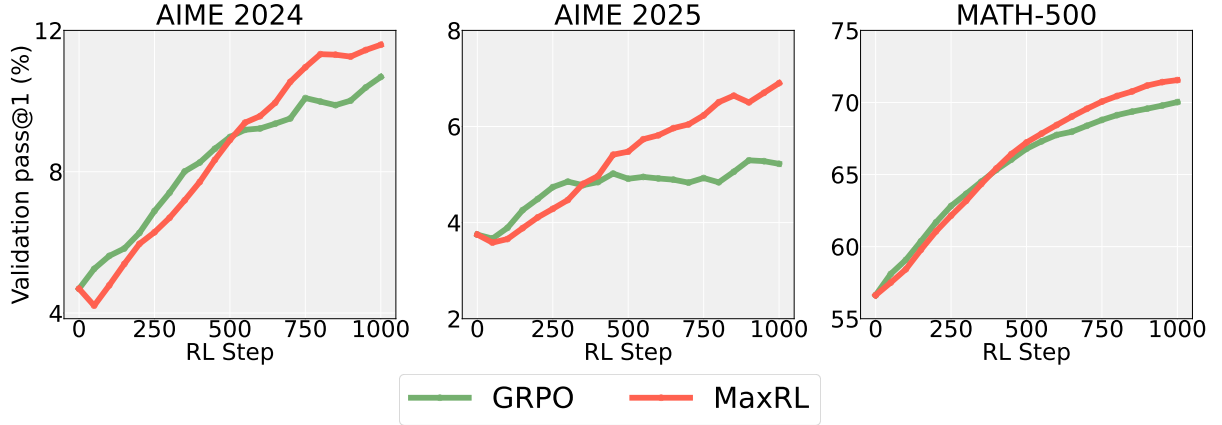


Figure 15: (Qwen3-1.7B-Base validation accuracy during intermediate training) We record validation pass@1 (using mean over 32 rollouts per prompt) over AIME 2024, AIME 2025 and MATH-500 during Qwen3-1.7B-Base model training. Similar to Figure 4, we observe that MAXRL initially trail behind GRPO at pass@1, but catches up with extended training and then converges to a higher value.

K.2 Training Dynamics

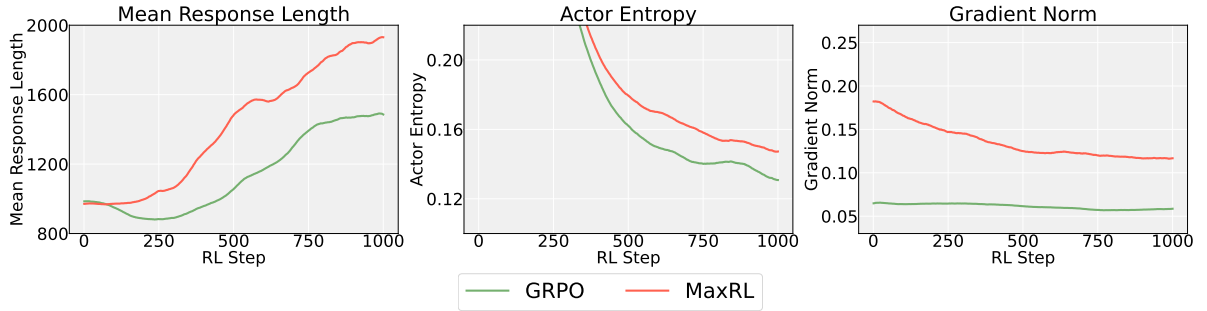


Figure 16: (Additional training dynamics metrics for Qwen3-1.7B-Base) We show comparison between GRPO and MAXRL in terms of mean response length, entropy of the actor, and gradient norm during training for the Qwen3-1.7B-Base model. MAXRL generally produces longer chains-of-thought, and also retains higher actor entropy during training. MAXRL also produce larger gradient norm during training.

L Additional Experimental Results on Maze

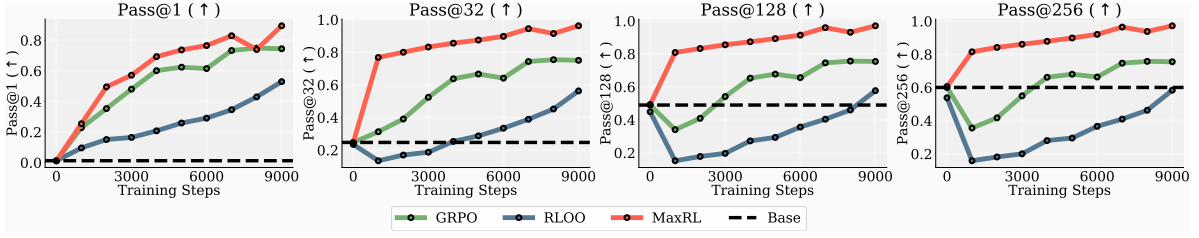


Figure 17: (Infinite training compute in maze experiment) We investigate how different objectives perform when we train a 3M model to solve 17x17 maze puzzles. MAXRL performs significantly better compared with GRPO and REINFORCE in Pass@1, Pass@32, Pass@128 and Pass@256. These results signify MAXRL’s effectiveness in computation scaling during RL.

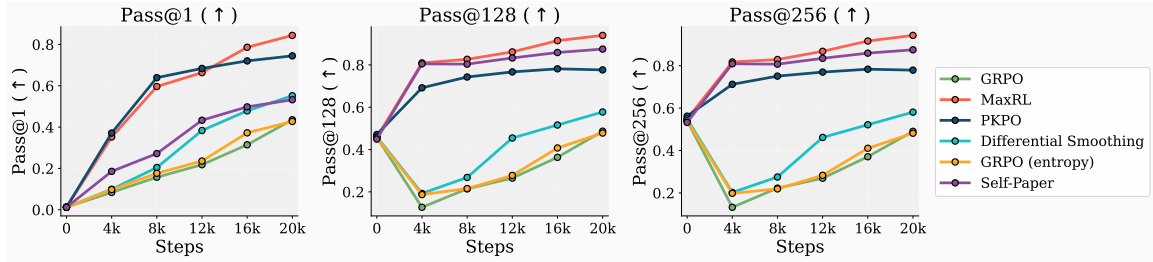


Figure 18: (Training curves compared with other baselines.) We compare MAXRL with other RL algorithms, including entropy regularization, PKPO (Waldner and Karkhanis, 2025), Differential Smoothing (Gai et al., 2025) and SELF (Nguyen et al., 2025b). MAXRL significantly outperforms other methods in all metrics, and is the only method to maintain both good average performance (pass@1) and coverage (pass@k).