RL's Razor: Why On-Policy Reinforcement Learning Forgets

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Abstract

We compare fine-tuning models with supervised fine-tuning (SFT) and reinforcement learning (RL) and find that, even at matched new-task accuracy, RL consistently forgets less. We investigate the cause and show that the degree of forgetting is not determined by the training algorithm itself, but by the distributional shift, namely the KL divergence between the fine-tuned and base policy when evaluated on the new task distribution. RL's advantage arises because on-policy updates bias optimization toward KL-minimal solutions among the many that solve a task, whereas SFT can converge to distributions arbitrarily far from the base model. We validate this across experiments with large language models and controlled toy settings, as well as provide theory on why on-policy RL updates lead to a smaller KL change. We term this principle *RL's Razor*: among all ways to solve a new task, RL prefers those closest in KL to the original model.

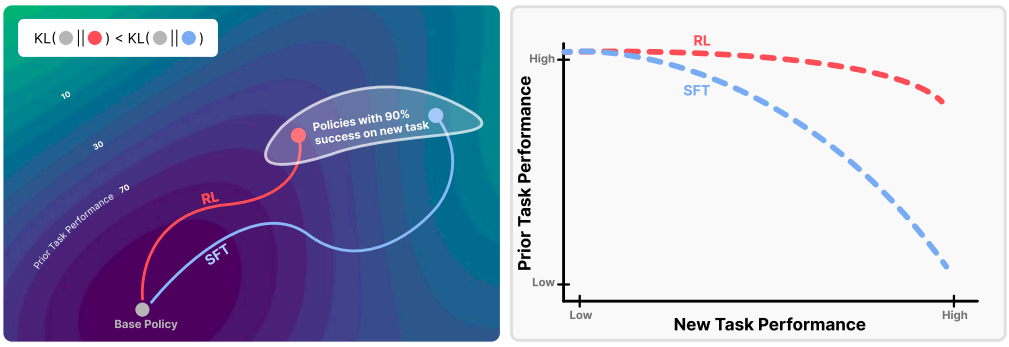

Figure 1: **RL prefers KL-minimal solutions.** Left: RL converges to policies close in KL to the base model. Right: This reduces forgetting at matched new-task accuracy compared to SFT.

1 Introduction

Foundation models have rapidly become the backbone of modern AI. Despite their remarkable capabilities, today's models are largely *static* once deployed: they are not designed to self-improve and continually acquire new capabilities. We imagine a future where deployed models are long-lived *agents* assisting humans in the long-term and continuously adapting to new needs. As such, models must improve and adapt to new data, environments, and objectives [1, 2, 3, 4, 5, 6]. A central challenge to this vision is *catastrophic forgetting*—the tendency for models to lose previously acquired capabilities when trained on new tasks [7, 8, 9, 10]. To enable foundation models to serve as long-term agents, we need to develop post-training methods that allow models to acquire new skills without erasing old ones.

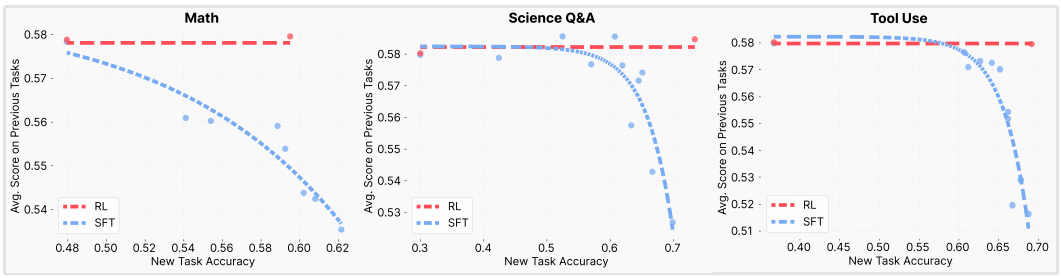

Figure 2: **Pareto frontiers of RL and SFT.** Each point represents a fine-tuned model. We sweep hyperparameters and plot only those on the Pareto-frontier. RL improves new-task performance while preserving prior knowledge, whereas SFT increases performance at the expense of forgetting.

To further this goal, we analyze two widely used post-training schemes of supervised fine-tuning (SFT) and reinforcement learning (RL). Our experiments reveal a surprising finding: even when SFT and RL achieve the same performance on the new task, we observe that **SFT often achieves new-task gains by erasing prior knowledge, while RL better preserves old skills**. This striking empirical gap raises the question: what underlying mechanism allows RL to improve on new tasks, but unlike SFT, not disturb the model's prior knowledge?

In search of this governing principle, we ablated many possible confounding variables proposed in prior work, and uncovered an empirical forgetting law: **When fine-tuning a model, π, on a new task τ, the degree of forgetting is accurately predicted by $\mathbb{E}_{x\sim\tau}\left[\mathbf{KL}(\pi_0||\pi)\right]$**, the Kullback–Leibler (KL) divergence between the finetuned and the base policy computed on the new task distribution τ. Although the underlying reason for this phenomenon remains unclear, its consistency across settings suggests it captures a fundamental property of forgetting.

This law also clarifies the surprising difference between SFT and RL. Our analysis reveals a simple but powerful principle we call ***RL's Razor*: among the many high-reward solutions for a new task, on-policy methods such as RL are inherently biased toward solutions that remain closer to the original policy in KL divergence**. Figure 1 (left) highlights this effect: among the many policies that reach a high success rate on the new task, RL is biased toward KL-minimal solutions, while SFT can converge to distant ones.

Together, these findings suggest a new perspective on post-training: to achieve continual adaptation, algorithms should explicitly aim to minimize KL divergence from the base model. This principle opens the door to designing future training methods that combine RL's ability to preserve prior knowledge with the efficiency of SFT, enabling foundation models that can *learn for life*.

2 Reinforcement Learning Forgets Less than SFT

In this section, we compare the degree of catastrophic forgetting induced by SFT and RL.

Experimental Setup. For each new task, we trained a Qwen 2.5 3B-Instruct [11]using either SFT, with annotations from either the original dataset or GPT-4o or RL, specifically GRPO [12]. Evaluation was twofold: performance on the held-out test set of the new task measured training gains, while performance on a diverse set of unrelated benchmarks quantified catastrophic forgetting. To obtain a reliable comparison, we trained dozens of models for each method under a variety of hyperparameter choices. Importantly, all RL experiments were done *without explicit KL regularization*. We then plotted only the models lying on the Pareto frontier. For more details, see Appendix C.

Tasks and Datasets. We repeated this experiment across three distinct domains: *Math reasoning*: math questions from the Open-Reasoner-Zero dataset [13], annotated with GPT-4o [14] responses filtered for correctness. *Science Q&A*: Chemistry L-3 subset of SciKnowEval [15], also annotated with GPT-4o. *Tool use*: ToolAlpaca dataset [16], using available annotations. For the evaluation of catastrophic forgetting, we used established benchmarks: Hellaswag [17], TruthfulQA [18], MMLU [19], IFEval [20], Winogrande [21], and HumanEval [22]. These serve as proxies for diverse prior abilities that the model should ideally retain.

Results. Figure 2 illustrates the learning-forgetting trade-offs for all tasks. Across all of them, RL training produces nearly horizontal Pareto frontiers as gains on the new task are achieved without loss on previous tasks. In contrast, SFT exhibits a steep downward slope—as new task accuracy rises, average performance on prior tasks consistently degrades.

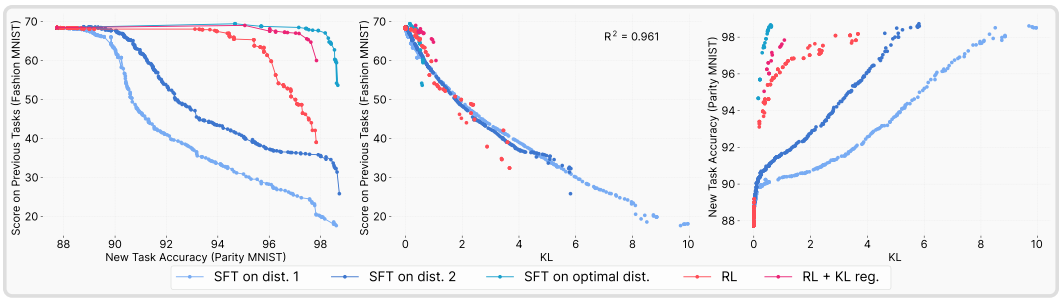

Figure 3: **KL divergence predicts catastrophic forgetting.** (Left) Learning-Forgetting Trade-offs. SFT outperforms RL only when an oracle distribution is used as a source of annotation. (Middle) Forgetting aligns to a single curve when plotted against KL divergence. (Right) RL improves new-task accuracy with smaller KL shifts than SFT, highlighting the conservativeness of on-policy updates.

Takeaway 1

While SFT boosts new-task performance by sacrificing prior knowledge, RL achieves comparable improvements with substantially less forgetting.

3 Smaller KL divergences lead to less forgetting

A central question raised by our initial results is why RL fine-tuning tends to forget less than SFT. To address this, we sought a confounding variable that could robustly explain the behavior of both methods. We systematically tested several candidates, including weight change under different norms, sparsity, and gradient rank. None of these explained the observed differences (see Appendix D.2). What ultimately emerged is that the *KL divergence between the trained model and the base model on the new task distribution* is an excellent predictor of catastrophic forgetting.

To test this hypothesis in a setting where we could run large numbers of experiments and push RL until it starts to forget, we developed a toy problem we call ParityMNIST which mimic realistic setting by allowing multiple different distributions to be all correct. We pretrained a 3-layer MLP on a subset of ParityMNIST and FashionMNIST [23], then fine-tuned on ParityMNIST. For more details, see Appendix C.2. This design allowed us to replicate the phenomenon where RL reached high accuracy on the new task with substantially slower degradation of prior knowledge, while SFT exhibited a steeper trade-off (Figure 3, left). Importantly, *reproducing the effect in this simple MLP setting shows that it is not specific to transformers or language modeling, but a more general property of fine-tuning deep generative models*.

KL as Predictor. When we plot forgetting directly against KL divergence from the base model on the ParityMNIST distribution, results across both RL and SFT collapse onto the same curve (Figure 3, middle). This shows that KL, not the training algorithm itself, predicts forgetting. A second-order polynomial fit achieves $R^2 = 0.961$ in this toy setting. The same correlation is observed in our LLM experiments with $R^2 = 0.61$. Although the statistic is lower, the residuals are mean-zero and not explained by predictors, consistent with random noise possibly due to approximated measures of KL and task accuracy.

Optimal SFT Distribution. Finally, the simplicity of ParityMNIST allowed us to analytically compute the optimal SFT distribution, that is, the optimal distribution with minimal KL divergence to the base model, see C.2 for details. Training SFT directly on this distribution outperformed RL, showing that the key factor is the KL distance, with RL being naturally biased toward such solutions. In addition, we also trained the SFT model on data generated by an RL trained model and found that it can reach the same level of performance and forgetting as the RL model, see Figure 7.

Takeaway 2

Across both SFT and RL, the amount of catastrophic forgetting is determined by how far the model moves from its base distribution on the new task, as measured by KL divergence.

4 On-policy methods leads to smaller KL divergence

Having established that the KL divergence between the trained model and its base distribution on the new task predicts catastrophic forgetting, we now ask: why are RL fine-tuned models able to achieve strong task performance while moving less in KL than SFT models?

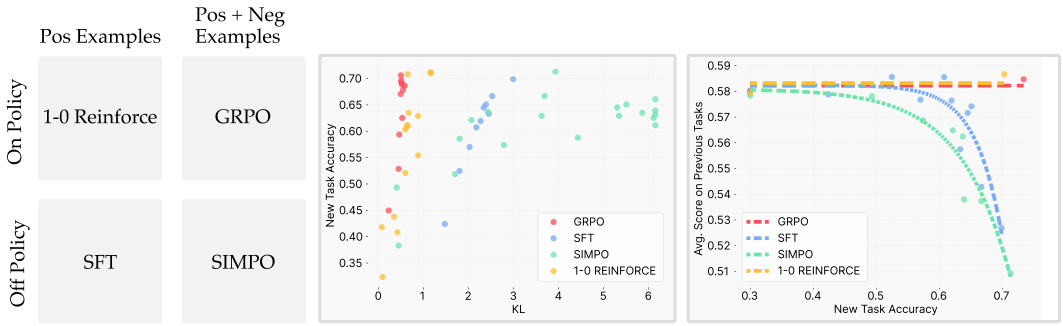

Figure 4: **Comparison of algorithm classes.** (Left) The four quadrants illustrate algorithm types, defined by whether they are on- or off-policy and whether they incorporate negative gradients. (Middle) On-policy methods retain prior knowledge more effectively. (Right) The on-policy methods achieve higher new-task accuracy while incurring smaller KL shifts from the base model.

Experimental Evidence To answer this, we examine the differences in training objectives. For discrete output spaces, SFT and policy gradient RL use the following objectives[1]:

$$\mathcal{L}_{\text{SFT}}(\pi) = -\mathbb{E}_{x\sim\mathcal{D},y\sim\pi_\beta}[\log \pi(y|x)] \quad \mathcal{L}_{\text{RL}}(\pi) = -\mathbb{E}_{x\sim\mathcal{D},y\sim\pi}\left[A(x,y)\log \pi(y|x)\right]$$

where π_β is the source of annotations, usually humans, and $A(x, y)$ is an Advantage function (the reward of y normalized with respect to other rewards for the same x). Two features distinguish the objectives: *Sampling Distribution*—RL trains on outputs drawn from the model's own distribution, whereas SFT relies on fixed external annotations; and *Negative Examples*—in RL, some sampled responses receive negative coefficients $A(x, y)$, pushing probability mass away from poor outputs, a mechanism absent in SFT.

Our hypothesis is that one of these two differences drives RL's slower forgetting. To test this, we compare four objectives: **GRPO** is an on-policy objective with negative examples, where $A(x, y)$ is the normalized reward. **1–0 Reinforce** is on-policy without negative examples, setting $A(x, y) = 1$ for correct responses and 0 otherwise, equivalent to sampling from the model and applying SFT only on correct answers. **SFT** is offline and does not use negative examples. **SimPO** [25] is offline with negative examples (see Appendix C.4).

We compared the four objectives on the Science Q&A task, measuring their learning–forgetting trade-offs as in Section 4. The results, shown in Figure 4, reveal that 1–0 Reinforce behaves similarly to GRPO, while SimPO resembles SFT. Thus, the critical factor is not the presence of negative gradients but the use of on-policy data. Plotting KL divergence confirms this conclusion: on-policy methods reach the same task performance with significantly smaller KL divergence from the base model than offline methods.

Theoretical Perspective To support our empirical findings, we also analyzed the policy gradient objective and found that, in a simplified setting, it has an implicit bias toward the minimum KL solution. for formal theorem and details, see Appendix B.

Takeaway 3

On-policy training explains why RL maintains smaller KL divergence than SFT. Sampling from the model's own distribution keeps it close to the base model, while SFT pushes it toward arbitrary external distributions.

5 Discussion and Conclusion

We show that forgetting is governed by KL divergence from the base policy rather than the training algorithm itself. RL forgets less than SFT because on-policy updates naturally bias toward KL-minimal solutions, preserving prior knowledge. Open questions remain about why large KL shifts disrupt old skills. More broadly, this work highlights a new direction for optimization design: developing methods that minimize KL divergence from first principles.

[1] In practice, the expectation is taken over outputs sampled from the current policy, and the policy gradient trick [24] ensures gradients flow only through the log-probability term, not through the sampling distribution.

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A Related work

Foundation Models and Post-training In modern deep learning, large-scale models trained on broad, diverse datasets (usually termed Foundation models) serve as general-purpose backbones across domains such as language, vision, robotics, and multimodal reasoning [26, 27, 28, 29, 30]. Pre-training often relies on self-supervised or weakly supervised objectives, producing models with broad domain knowledge and some zero-shot capabilities [31, 32]. However, raw pre-trained models may not directly meet the requirements of specific applications or align with domain-specific constraints. Post-training methods address this gap by adapting foundation models to downstream tasks through supervised fine-tuning on curated datasets [33, 34, 35, 36], reinforcement learning from human or automated feedback [37, 38, 39, 40], and other techniques [41].

Catastrophic Forgetting. While fine-tuning primarily aims to improve performance on a new specific task, preserving the model's pre-existing general capabilities is equally critical. Unfortunately, fine-tuning often leads to catastrophic forgetting—a phenomenon where learning new information significantly deteriorates previously acquired knowledge [7, 8, 9]. Many works have sought to reduce forgetting by constraining updates, for example through weight penalties, feature preservation, or output matching [42]. These approaches highlight important factors, but what remains missing is a unifying principle that predicts forgetting across different algorithms and settings.

In contrast, our work does not propose a new method but instead identifies a simple law: the KL divergence between the fine-tuned and base policy, measured on the new task distribution, reliably predicts the degree of forgetting. This explains the success of some popular forgetting mitigation methods like Elastic Weight Update [9], which can be seen as approximation of KL minimization [43]. Interestingly, practitioners have also observed that KL regularization, originally introduced in RL fine-tuning of LLMs to stabilize optimization or prevent reward hacking [44, 45], helps reduce catastrophic forgetting [38].

SFT versus RL. Most prior comparisons between SFT and RL have focused on performance on the new task being learned. A seminal result by [46] showed that in sequential decision making, on-policy learning can achieve stronger performance even when the learning signal is identical. Building on this, recent studies have found that RL fine-tuned models often exhibit superior generalization beyond the training distribution [47, 48, 49] and transfer more effectively to related tasks [50] compared to SFT trained models. However, none of this work has examined their relative susceptibility to catastrophic forgetting, which is the focus of our study.

Concurrently, [51] report that RL forgets less than SFT, but ascribe RL's advantage to negative examples and ignore sampling-distribution effects. Section 4 shows that this assumption is inconsistent with our results.

B Theory

Policy gradient methods can be understood as a form of conservative projection. At each step, the policy samples outputs it already finds likely, then re-weights those samples according to reward, shifting probability mass toward higher-reward outcomes while suppressing lower-reward ones. Crucially, because updates are defined relative to the model's own distribution, they nudge the policy toward a nearby re-weighted distribution, rather than pulling it toward a potentially distant external distribution (as in SFT). This explains why policy gradient methods tend to remain close to the base model in KL divergence.

This perspective can be formalized by observing that, in the binary-reward case, the re-weighted distribution targeted by policy gradient is exactly the minimum-KL projection of the current policy onto the set of optimal ones.

Lemma B.1 (Rejection sampling as an I-projection)**.** *Let p be a distribution over a finite set Y, and let $R : Y \to \{0, 1\}$ be a reward function. Rejection sampling from p with acceptance condition $R(y) = 1$ yields a distribution q_{RS}. This distribution can be equivalently characterized as the solution to:*

$$q_{RS} =_q D_{KL}(q||p) \quad s.t \quad \mathbb{E}_{y\sim q}[R(y)] = 1$$

Equivalently, q_{RS} is the I-projection of p onto the set $\{q : \mathbb{E}_q[R] = 1\}$

Proof. Let $S = \{y \in Y : R(y) = 1\}$. Rejection sampling produces the conditional distribution

$$q_{\mathrm{RS}}(y) = \begin{cases} \frac{p(y)}{p(S)} & y \in S, \\ 0 & y \notin S, \end{cases}$$

where $p(S) = \sum_{y\in S} p(y)$ and we assume $P(S) > 0$.

Now consider the optimization problem. The constraint $\mathbb{E}_q[R] = 1$ means

$$\sum_{y\in Y} q(y)R(y) = \sum_{y\in S} q(y) = 1$$

so q must put all of its mass on S. Thus the feasible set is exactly all distributions supported on S.

For any q supported on S, we can write $p(y) = p(S)\, p(y|S)$ for $y \in S$, and then

$$\begin{aligned} D_{\mathrm{KL}}(q\|p) &= \sum_{y\in S} q(y) \log \frac{q(y)}{p(y)} = \sum_{y\in S} q(y) \log \frac{q(y)}{p(y \mid S)} - \log p(S) \sum_{y\in S} q(y) \\ &= D_{\mathrm{KL}}\big(q\|p(\cdot \mid S)\big) - \log p(S) \end{aligned}$$

where we used $\sum_{y\in S} q(y) = 1$ in the last step. The second term is constant in q, so minimizing $D_{\mathrm{KL}}(q||p)$ is the same as minimizing $D_{\mathrm{KL}}(q||p(\cdot|S))$. By strict convexity of $D_{\mathrm{KL}}(\cdot\|\cdot)$ in its first argument, the unique minimizer is $q = p(\cdot \mid S) = q_{\mathrm{RS}}$. □

Lemma B.2 (Policy gradient as an M-projection)**.** *Let Y be a finite set and let $\Pi \subseteq \Delta(Y)$ be a set of admissible policies (distributions over Y). Consider the single-step reinforcement learning objective*

$$\max_{\pi} \mathbb{E}_{y\sim\pi}[R(y)]$$

where $R : Y \to \mathbb{R}_{\geq 0}$ is a reward function. By the policy gradient theorem, this objective is equivalently optimized by

$$\max_{\pi} \mathbb{E}_{y\sim\bar{\pi}}\big[R(y) \log \pi(y)\big]$$

where $\bar{\pi}$ indicates that gradients are not propagated through the sampling distribution. Define the distribution

$$q(y) = \frac{\pi(y)R(y)}{Z}, \qquad Z = \sum_{y\in Y} \pi(y)R(y)$$

Then taking a policy gradient step is equivalent to taking a gradient step on the following objective:

$$\min_{\pi} -\mathbb{E}_{y\sim q}[\log \pi(y)]$$

In other words, optimizing the RL objective using policy gradient is equivalent to finding the M-projection of q onto the set of feasible policies π using gradient descent.

Proof. Expanding the policy gradient objective gives

$$\mathbb{E}_{y\sim\bar{\pi}}[R(y) \log \pi(y)] = \sum_{y\in Y} \pi(y)R(y) \log \pi(y)$$

Let $Z = \sum_{y\in Y} \pi(y)R(y)$. Define $q(y) = \pi(y)R(y)/Z$. Then the above becomes

$$\sum_{y\in Y} \pi(y)R(y) \log \pi(y) = Z \sum_{y\in Y} q(y) \log \pi(y) = Z\, \mathbb{E}_{y\sim q}[\log \pi(y)]$$

Since Z does not depend on π in the gradient computation (it is treated as a constant in the $\bar{\pi}$ sense), maximizing the original objective is equivalent to maximizing $\mathbb{E}_{y\sim q}[\log \pi(y)]$.

Finally, recall that the M-projection of a distribution q onto a set of distributions Π is given by

$$\min_{\pi\in\Pi} \mathrm{KL}(q\|\pi) = \mathbb{E}_q[\log \frac{q}{\pi}] \; = \mathbb{E}_q[\log q] \; - \; \mathbb{E}_q[\log \pi]$$

since $\mathbb{E}_q[\log q]$ does not depend on π, the maximizer of $\mathbb{E}_{\bar{\pi}}[R \log \pi]$ over Π coincides with $\arg\min_{\pi\in\Pi} \mathrm{KL}(q\|\pi)$. Thus, the policy gradient update corresponds to the M-projection of q onto the policy class. □

Theorem B.3 (RL with binary reward as an EM algorithm)**.** *Let Y be a finite set and let $\Pi \subseteq \Delta(Y)$ be a set of feasible policies. Let $R : Y \to \{0, 1\}$ be a binary reward function and P^* the set of all optimal policies $P^* = \{q : \mathbb{E}_q[R] = 1\}$. Then, solving the Single-step reinforcement learning objective using policy gradients is equivalent to performing the following optimization procedure:*

$$q_t = \arg\min_{q\in P^*} \mathrm{KL}(q\|\pi_t), \qquad \pi_{t+1} = \arg\min_{\pi\in\Pi} \mathrm{KL}(q_t\|\pi)$$

This procedure is also known as EM with information projection.

Proof. Sampling $y \sim \pi$ and accepting iff $R(y) = 1$ is exactly rejection sampling onto the event $S = \{y \in Y : R(y) = 1\}$. The resulting distribution is $\pi(\cdot|S)$. By Lemma A.1 with $p \leftarrow \pi$, this $\pi(\cdot|S)$ solves

$$\min_q D_{\mathrm{KL}}(q\|\pi) \quad \text{s.t.} \quad \mathbb{E}_q[R] = 1$$

establishing the I-projection. Applying Lemma A.2 on the RL objective gives us the M-projection. □

Proposition B.4 (Convergence to minimum KL solution)**.** *Under the setting appear in theorem B.3 and assume* Π *is an e-flat (exponential-family) model with full support, the optimal set* P^* *is nonempty and realizable (i.e.,* $\Pi \cap P^* \neq \varnothing$*). Then:*

(1) *If the M-projection is exact at every step, then* (π_t) *converges to*

$$\pi^\dagger = \arg \min_{\pi \in P^* \cap \Pi} D_{\mathrm{KL}}(\pi \,\|\, \pi_0)$$

(2) *If the M-projection is inexact but, for some errors* $\varepsilon_t \geq 0$*, it holds that*

$$D_{\mathrm{KL}}(q_t\|\pi_{t+1}) \;\leq\; \min_{\pi \in \Pi} D_{\mathrm{KL}}(q_t\|\pi) \;+\; \varepsilon_t \quad \textit{with} \quad \sum_{t=0}^{\infty} \varepsilon_t < \infty$$

then π_t *also converges to the same limit* $\pi^\dagger$*.*

Proof. The I-step is always an exact I-projection (Lemma A.1). In the case of an exact M-step, the iterative process is EM with information projections. The e-/m-flat geometry yields the Pythagorean identities implying convergence to $\pi^\dagger$ [52, 53, 54]. When the M-step only ensures a (near-)minimization up to summable errors, the iteration is GEM: monotone improvement and convergence follow from the GEM theory of [55] together with generalized alternating minimization for Bregman divergences [56], which, under the same e-/m-flat assumptions, selects the same minimum-KL limit $\pi^\dagger$. □

Practical considerations. Our theoretical equivalence should be interpreted with the following caveats:

- Beyond REINFORCE. In practice, many policy gradient algorithms such as GRPO and PPO replace the raw reward $R(y)$ with an advantage estimate $A(y)$. Since this substitution is a control variate technique, it leaves the expected gradient direction unchanged while reducing its variance. Thus, our projection-based interpretation continues to hold.
- The optimal policy set P^* defined by the linear constraint $\mathbb{E}_q[R] = 1$ is an m-flat family, but the representable policy set Π induced by a neural network parametrization is not in general e-flat. This may prevent exact convergence to the minimum-KL solution described above. Nevertheless, our theorem provides a principled explanation for the bias observed in practical RL algorithms.

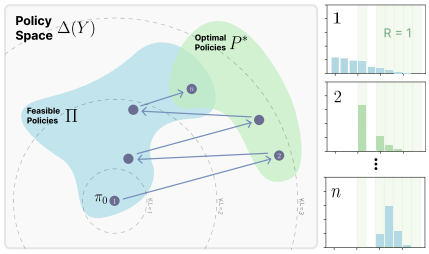

Figure 5: **KL-minimal path to optimality.** Alternating I-projection into the set of optimal policies and M-projection into Π carries π_0 into P^* while preferring the closest solution in KL.

C Training and Evaluation Details

C.1 LLM experiments

Unless otherwise stated, all reinforcement learning experiments were conducted using GRPO [12].

For the *Math* reasoning task, the training set provided final answers but lacked reasoning chains required for SFT training. To obtain these, we queried DeepSeek R1 [39], sampling up to 16 responses per prompt and retaining a single response that matched the correct final answer. This yielded valid annotations for 96% of the dataset. For the *Science Q&A* task, we applied the same procedure with GPT-4o, obtaining correct annotations for the entire dataset.

To construct the learning–forgetting trade-off curves (e.g., Figure 2), we followed the protocol below:

1. Hyperparameter sweep. We trained multiple models under a broad sweep of hyperparameters (see Table 1).
2. New-task evaluation. For *Math* and *Science Q&A*, accuracy was measured by comparing the model's final answer to the ground truth, ignoring intermediate reasoning chains. For Tool Use, we extracted API calls from the output and matched them against ground-truth calls via regular expressions.
3. Previous-task evaluation. We assessed performance on unrelated benchmarks as described in Section 2, using the `Language Model Evaluation Harness` [57].
4. Pareto filtering. From the trained models, we retained only those lying within 2 accuracy points of the Pareto frontier.
5. Curve fitting. An exponential function was fit to the filtered points to produce the trade-off curves.

Hyperparameter	**SFT / SIMPO**	**RL**
Base Model	Qwen2.5 3B-Instruct	Qwen2.5 3B-Instruct
Learning Rate	{1e-5, 3e-5, 5e-5, 7e-5, 9e-5}	{1e-5, 2e-5, 3e-5, 4e-5, 5e-5}
Optimizer	adamw	adamw
LR Scheduler	{constant w. warmup, cosine w. warmup}	constant w. warmup
Warmup steps	50	50
Epochs	{1,2}	1
Batch Size	{16,32,64,128}	See Below
Max Grad Norm	1	1
bfloat16	True	True
Weight Decay	0	0
GRPO-only hyperparameters		
KL reg.		0
Group Size		64
Prompts per generation		8
num iterations (μ)		{1,2}
Loss type		Dr. GRPO [58]

Table 1: Hyperparameters used for the LLM experiments. Curly braces {} indicate a sweep over the specified values. Additional parameters such as weight decay and max gradient norm were manually ablated; since they showed no significant effect on results, they were not included in the final sweep.]

C.2 MNIST Experiments

All MNIST experiments were conducted using a 3-layer MLP with input dimension 785, hidden layers of sizes 512 and 256, and output dimension 10. The input consisted of a flattened 28×28 image concatenated with a binary indicator: $+1$ for ParityMNIST and -1 for FashionMNIST.

Pretraining. We pretrained the network jointly on ParityMNIST and FashionMNIST using small subsets of the original datasets (500 images from each). For ParityMNIST, the label was chosen uniformly at random among all digit labels with the correct parity.

Fine-tuning methods. We evaluated five fine-tuning strategies:

1. **GRPO**.
2. **GRPO + KL regularization** with coefficient 0.1.
3. **SFT 1**: all even digits mapped to label 0, all odd digits to label 1.
4. **SFT 2**: even digits randomly mapped to $\{0, 4\}$, odd digits to $\{1, 5\}$.
5. **SFT with oracle distribution**: annotations drawn from the minimum-KL distribution consistent with task correctness.

Oracle distribution. Motivated by the KL–forgetting connection, we define the oracle distribution as the one that achieves perfect task accuracy while remaining closest (in KL divergence) to the pretraining distribution π_0. Concretely, for an input image x we compute $\pi_0(\cdot|x) \in \mathbb{R}^{10}$ and the binary indicator vector $R \in \{0,1\}^{10}$ encoding which labels are correct given the digit's parity. The oracle distribution q^* is the solution to:

$$q^* = \arg\min_q D_{\mathrm{KL}}(\pi_0 \| q) \quad \text{s.t.} \quad q^\top R = 1.$$

Since KL is convex and the constraint is linear, we can calculate a closed-form solution for every image. We then sample from q^* to produce SFT annotations.

Hyperparameter sweep. For each method we trained models across a sweep of 15 learning rates logarithmically spaced between $3e-6$ and $1e-3$, using either a constant-with-warmup or cosine-with-warmup scheduler, and training for 1 or 2 epochs. Including mid-training checkpoints, this produced approximately 500 runs per method.

C.3 Centered Kernel Alignmen

Centered Kernel Alignment (CKA) [59] Given representations $X, Y \in \mathbb{R}^{n \times d}$, define kernels $K = XX^\top$, $L = YY^\top$. Let $H = I - \frac{1}{n}\mathbf{1}\mathbf{1}^\top$ be the centering matrix. The centered kernels are

$$\bar{K} = HKH, \quad \bar{L} = HLH.$$

CKA is then computed as

$$\mathrm{CKA}(K, L) \;=\; \frac{\langle \bar{K}, \bar{L} \rangle_F}{\|\bar{K}\|_F \, \|\bar{L}\|_F},$$

where $\langle A, B \rangle_F = \mathrm{tr}(A^\top B)$.

CKA with k-NN Alignment (CKNNA) [60] Let $\alpha(i,j) \in \{0,1\}$ indicate whether i, j are mutual k-nearest neighbors in both X and Y. Define the masked inner product

$$\langle A, B \rangle_\alpha = \sum_{i=1}^{n} \sum_{j=1}^{n} \alpha(i,j)\, A_{ij} B_{ij}.$$

CKNNA is then given by

$$\mathrm{CKNNA}(K, L) \;=\; \frac{\langle \bar{K}, \bar{L} \rangle_\alpha}{\sqrt{\langle \bar{K}, \bar{K} \rangle_\alpha \, \langle \bar{L}, \bar{L} \rangle_\alpha}}.$$

When $\alpha(i,j) = 1$ for all $i \neq j$, CKNNA reduces to standard CKA.

C.4 SIMPO

SimPO is an offline objective that utilizes negative examples. We create negative examples by sampling incorrect responses from an external model, and use the SFT data for positive examples. The SimPO [25] loss compares correct and incorrect outputs via a logistic term:

$$\mathcal{L}_{\text{SIMPO}}(\pi) = -\mathbb{E}_{x \sim \mathcal{D}, y_w \sim \pi_{\beta+}, y_l \sim \pi_{\beta-}} \left[\log \sigma \left(\log \pi(y_w|x) - \log \pi(y_l|x) - 1 \right) \right]$$

where $\pi_{\beta+}$ and $\pi_{\beta-}$ denote distributions for correct and incorrect responses, respectively. We used SimPO rather than naïve likelihood/negative likelihood because the latter was unstable to train.

D Additional Results

D.1 Representation Preservation

While benchmark performance provides an external view of forgetting, it can also be sensitive to superficial factors, such as formatting mismatches with the previous tasks. To probe whether the training altered the model's internal capacity more fundamentally, we analyzed changes to the model's representations.

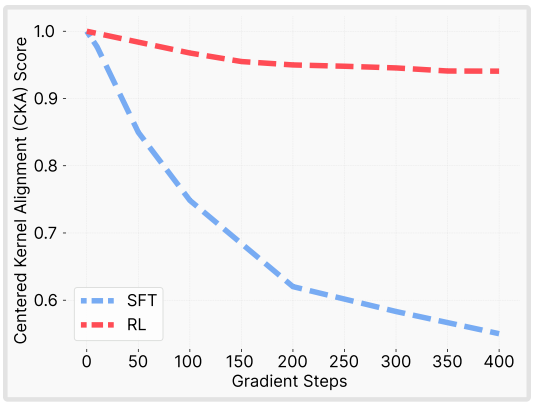

Figure 6: **CKA similarity to the base model during training.** Although SFT and RL achieve comparable task performance, SFT models diverge substantially in their representations, whereas RL models remain more closely aligned with the base model.

Experimental Setup. To study how representations change between models, we compare their embeddings on a shared dataset. Directly comparing raw embedding values is not meaningful, since these can vary arbitrarily during training. Instead, it is common to compare the relative geometry of the embeddings—that is, how different inputs relate to each other. This geometry can be summarized by a kernel (similarity) matrix, which encodes pairwise relationships among inputs. Centered Kernel Alignment (CKA) [59] is a standard measure for comparing such kernels, providing a way to quantify representational similarity between models.

For this analysis, we constructed kernels from random Wikipedia paragraphs, ensuring that the comparison probes representations of content unrelated to the new training tasks. We then measured the similarity of the kernel between the base model and its fine-tuned variant using a version of CKA called CKNNA ([60], see Appendix C.3 for more details). Specifically, we compare models that achieved similar final accuracy on the new task.

Results. Figure D.1 shows that RL-trained models maintain almost perfect representational alignment with the base model, even after learning the new task. By contrast, SFT-trained models diverge substantially, indicating that the training reshapes their internal space in ways that distort previously encoded knowledge. Together with the benchmark results, this suggests that RL is able to integrate new abilities without disturbing the broader conceptual structure, while SFT incurs representational shifts that manifest as catastrophic forgetting.

D.2 Incorrect Hypothesis

Science advances not only through identifying the right explanations, but also by systematically ruling out incorrect ones. To this end, we tested a broad set of candidate variables as potential predictors of catastrophic forgetting. These variables fell into four categories:

- **Weight-level changes.** A natural hypothesis is that forgetting is tied to how much the parameters themselves move. We measured parameter changes under L_1, Fisher-weighted L_2, and spectral norm metrics. The Fisher matrix was computed on the basis of the model parameters, with expectation over inputs from the previous task. These metrics correlated only weakly with forgetting: large parameter shifts could occur without forgetting, and conversely, forgetting sometimes occurred despite small parameter movement.
- **Sparsity and rank of updates.** Motivated by [61], who argue that RL updates are sparse while SFT weight updates are dense, we explicitly tested this hypothesis. In our setting, however, we found that the reason for the observed sparse updates where the use of `bfloat16` for model training. Since `bfloat16` has a limited mantissa, small parameter updates (such as those produced by RL) can fail to cross the representational threshold, effectively causing no update at all. Performing the same training with `float32` resulted in models with identical performance but without any sparsity

in their weight updates. Checking the rank of the weight changes, we found that all algorithms lead to full rank weight updates.

- **Representation-level changes.** Following the CKNNAA results from subsection D.1, we examined hidden activation shifts (L1 and L2 distances) as proxies for changes in internal representations. These variables show some correlation, but the curves were distinct between training objectives.
- **Distributional distances.** We considered multiple measures of output distribution change, all measured over inputs from the new task τ: Forward KL ($\mathbb{E}_{x\sim\tau}\left[\mathrm{KL}(\pi_0||\pi)\right]$), Reverse KL ($\mathbb{E}_{x\sim\tau}\left[\mathrm{KL}(\pi||\pi_0)\right]$), Total Variation, and L_2 distance between distributions. While reverse KL showed a good signal, and TV moderately correlated with forgetting, none approached the predictive power of forward KL.

Table 2 summarizes these results. Across all candidates, forward KL divergence between the fine-tuned and base model on the new task distribution emerges as the only consistent and high-fidelity predictor of catastrophic forgetting.

Variable	R^2 (2nd deg. polynomial)
KL, forward	$\mathbf{0.96 \pm 0.01}$
KL, reverse	0.93 ± 0.01
TV	0.80 ± 0.01
Distribution change, L2	0.56 ± 0.02
Weight change, L1	0.34 ± 0.02
Weight change, Fisher Weighted L2	0.58 ± 0.02
Weight change, spectral norm	0.58 ± 0.02
Sparsity of weight change	N/A
Rank of weight change	N/A
Activation change, L1	0.52 ± 0.02
Activation change, L2	0.55 ± 0.02

Table 2: Predictive power of alternative variables compared to forward KL. None approaches the explanatory strength of forward KL divergence.

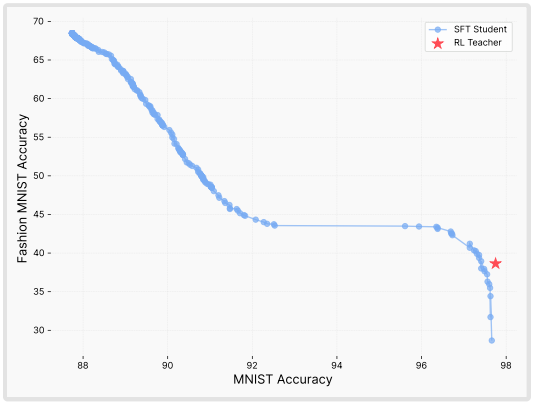

Figure 7: **SFT distillation from an RL teacher.** Accuracy trade-off between the new task (MNIST) and the prior task (FashionMNIST). Sweeping student hyperparameters shows that SFT can match the teacher within noise on both tasks. This suggests that what matters is not the optimization path, but the distribution of the final model.

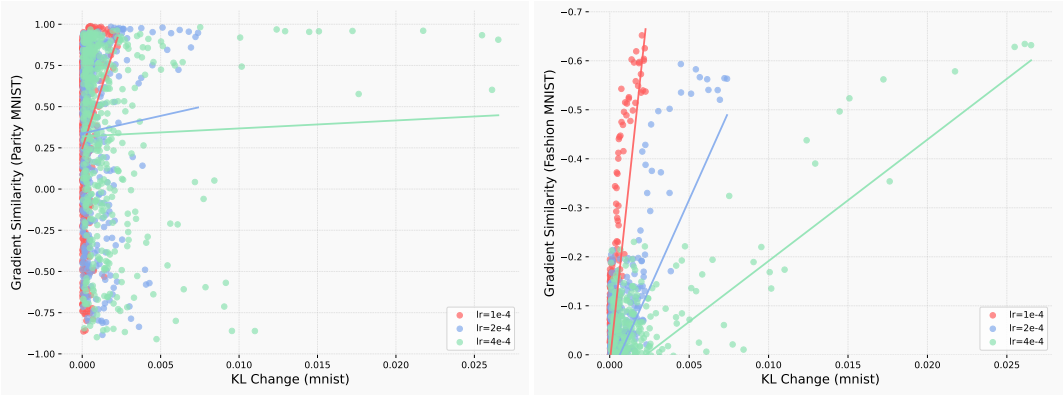

Figure 8: **Gradient similarity versus KL change.** (Left) On the new training task (ParityMNIST), gradient cosine similarity and KL change per step remain anti-correlated. (Right) On the prior task (FashionMNIST), the gradient similarity is more correlated with the KL change per step on the training task (ParityMNIST). Together, these plots show that taking a larger step on the current task induces gradients that are more similar in direction to the