

STRUCTURE-AWARE PARAMETER-EFFICIENT MACHINE UNLEARNING ON TRANSFORMER MODELS

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006 Paper under double-blind review

ABSTRACT

011 Transformer has become fundamental to a vast series of pretrained large mod-
 012 els that have achieved remarkable success across diverse applications. Machine
 013 unlearning is an emerging field focused on efficiently removing the influence of
 014 specific data from trained models, to comply with privacy regulations enforcing
 015 the right to be forgotten. The sheer size of Transformer-based models poses a
 016 significant challenge to unlearning efficiency. Existing methods find it promising
 017 to restrict unlearning updates to a small portion of influence-critical parameters.
 018 However, their parameter-efficient unlearning methods are largely devised in a
 019 structure-oblivious manner, which tends to inaccurately identify these parameters
 020 and leads to inferior unlearning performance for Transformers. In this paper, we
 021 propose SPE-Unlearn, a structure-aware parameter-efficient machine unlearn-
 022 ing approach tailored for the Transformer architecture. SPE-Unlearn intro-
 023 duces a learnable pair of masks to respectively pinpoint influence-critical parame-
 024 ters in the heads and filters of Transformers. The learning objective of these masks
 025 is derived by jointly considering both desiderata of unlearning, i.e., sufficiency in
 026 influence removal and efficiency, and optimized through an efficient algorithm fea-
 027 tured by a greedy search with a warm start. Equipped with the identified key par-
 028 ameters, SPE-Unlearn facilitates second-order unlearning, memory-free un-
 029 learning, and memory-aided unlearning scenarios. Extensive experiments on vari-
 030 ous Transformer models and datasets demonstrate the effectiveness and efficiency
 031 of SPE-Unlearn for Transformer unlearning.

1 INTRODUCTION

032 Transformer architecture ([Vaswani et al. \(2017\)](#)) has demonstrated superior performance in the field
 033 of natural language processing. Its models, e.g., BERT ([Devlin et al. \(2018\)](#)) and GPT ([Achiam et al.
 034 \(2023\)](#)), show impressive performance in a wide range of downstream tasks ([Wei et al. \(2021\)](#); [Hao et al. \(2019\)](#)). In light of privacy regulations, such as General Data Protection Regulation (GDPR)
 035 ([Hoofnagle et al. \(2019\)](#)), users are granted the right to request the removal of specific training
 036 data from models. To fulfill this requirement, machine unlearning techniques have been extensively
 037 researched ([Bourtoule et al. \(2021\)](#); [Yao et al. \(2023\)](#)). However, when applying these techniques
 038 to Transformers, which commonly involves a large number of parameters, a significant challenge
 039 lies in achieving computational efficiency while ensuring effective unlearning and preserving model
 040 fidelity ([Warnecke et al. \(2021\)](#); [Liu et al. \(2024a\)](#)).

041 Recent researches propose parameter-efficient unlearning techniques ([Liu et al. \(2024a\)](#); [Pochinkov & Schoots \(2024\)](#); [Schoepf et al. \(2024\)](#)), which identify the influence-critical parameters to govern
 042 the unlearning process. Specifically, these methods assess the importance of parameters through
 043 different evaluation strategies, allowing selective updates to reduce computational overhead and
 044 improve unlearning efficiency. However, applying parameter-efficient unlearning to address the
 045 dilemma of the unlearning tasks in Transformers faces two major limitations. First, previous eval-
 046 uation methods rely on heuristic or empirical strategies to identify parameters. For Transformer
 047 models with an immense number of parameters, identifying those specifically relevant to unlearn-
 048 ing becomes inefficient. Additionally, existing methods ([Pochinkov & Schoots \(2024\)](#); [Liu et al.
 049 \(2023b\)](#); [Shi et al. \(2023\)](#)) assess importance of parameters by comparing performance (e.g., ac-
 050 tivations) on forgetting dataset and remaining dataset may result in sub-optimal selection process
 051 for unlearning. Second, previous unlearning methods overlook the intricate interactions between
 052 053

structures in Transformers. Transformers utilize parallel attention heads and hierarchical filters to perform computation and inference [Vaswani et al. \(2017\)](#). Consequently, attempting to identify critical parameters at a fine-grained level is often inaccurate, as this approach fails to capture the broader contextual relationships inherent in Transformers.

In this paper, we propose a **Structure-aware Parameter-Efficient Unlearning** (SPE-Unlearn) approach that targets influence-critical parameters at the structural level for Transformers. Specifically, SPE-Unlearn formulates the unlearning objective through a pair of learnable masks applied to heads and filters. The derivation for this formulation ensures the effective influence removal and guides the identification of key structures. These masks are further refined by considering intra-layer interactions, and a warm-start greedy search algorithm is employed to efficiently optimize the process. Equipped with these structure-aware masks, we integrate SPE-Unlearn into second-order unlearning updates. While second-order unlearning introduces an approximation error, sparse updates using structure-aware masks can mitigate the errors, thereby preserving overall model performance. In addition, we analyze that structure-aware masks can demonstrate significant advantages in successive settings ([Hu et al. \(2023\)](#); [Liu et al. \(2023a\)](#)). In this context, we are the first to categorize second-order successive setting into two types based on whether intermediate information from previous removal requests is retained: **memory-free unlearning** ([Guo et al. \(2020\)](#); [Gu et al. \(2024\)](#)) and **memory-aided unlearning** ([Liu et al. \(2023a\)](#)). Our approach demonstrates exceptional robustness by effectively containing errors within selected structures, especially in memory-free unlearning scenarios. Our key contributions are summarized as follows:

- We introduce a new paradigm for identifying influence-critical parameters in Transformers, SPE-Unlearn, which operates at the structural level. Our approach theoretically derives importance scores for selecting key structures using a pair of learnable masks. These structure-aware masks can be seamlessly integrated into existing unlearning methods.
- We integrate SPE-Unlearn into second-order unlearning and analyze the gains with structure-aware masks. Extensive experiments across diverse datasets using three models demonstrate proposed method offers a superior trade-off among efficacy, fidelity, and efficiency.
- We categorize successive unlearning into two successive scenarios: memory-free unlearning and memory-aided unlearning. Empirical studies show that unlearning with structure-aware masks can handle a greater number of removal requests compared to standard unlearning before retraining becomes necessary, especially in memory-free scenarios.

2 PRELIMINARY

2.1 PROBLEM FORMULATIONS

Machine unlearning aims to remove the influence of targeted data from a trained model. Let $\mathcal{D} = \{x_i\}_{i=1}^M$ denote a training dataset containing M data points, where each x_i corresponds to an individual data point. Starting with the original model θ^* which was trained on \mathcal{D} , the objective of unlearning is to effectively remove the sensitive or compliance-related data while maintaining overall performance. Specifically, for the unlearning task, the dataset \mathcal{D} is grouped into two subsets: **forgetting dataset** \mathcal{D}_f and **remaining dataset** \mathcal{D}_r , i.e., $\mathcal{D} = \mathcal{D}_f \cup \mathcal{D}_r$. The forgetting dataset \mathcal{D}_f consists of the targeted data we aim to remove from the model. Accordingly, the remaining dataset \mathcal{D}_r includes the data we intend to retain and potentially further optimize. Given a loss function ℓ for targeted task, the objective of unlearning can be framed as learning an optimal model θ_U^* :

$$\theta_U^* = \arg \min_{\theta} \mathcal{L}(\theta; \mathcal{D}_r) = \arg \min_{\theta} \sum_{x \in \mathcal{D}_r} \ell(\theta; x) + \lambda \Omega(\theta), \quad (1)$$

where $\mathcal{L}(\theta; \mathcal{D}_r)$ represents the total loss on the dataset \mathcal{D}_r with θ , and $\lambda \Omega(\theta)$ is a common regularization term ([Hart et al. \(2000\)](#)). The most viable solution to address this optimization problem is retraining the model from scratch. However, retraining can be costly in terms of time and computing resources. A practical alternative, known as the second-order unlearning update ([Guo et al. \(2020\)](#); [Golatkar et al. \(2020\)](#); [Izzo et al. \(2021\)](#); [Warnecke et al. \(2021\)](#); [Liu et al. \(2024b\)](#)), deduces the general close-form parameter modification from the original model θ^* :

$$\theta \approx \theta^* + \mathbf{H}_{\theta^*}^{-1} \sum_{x \in \mathcal{D}_f} \nabla_{\theta} \ell(\theta^*; x), \quad (2)$$

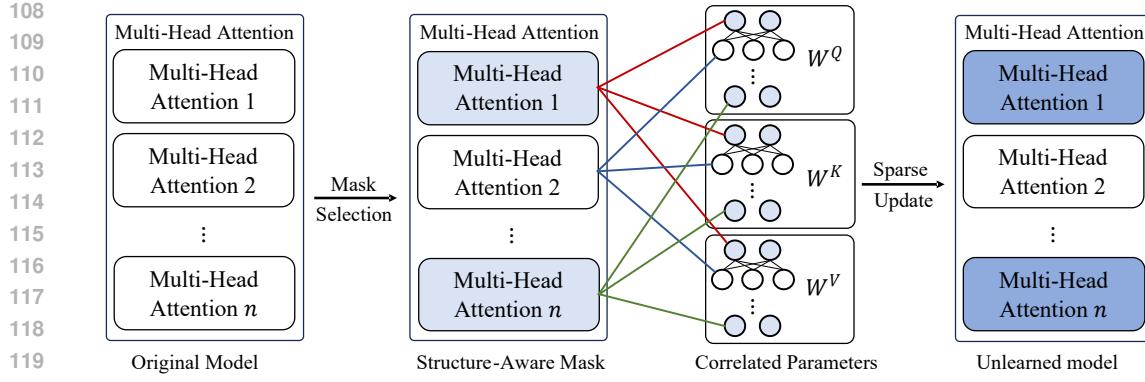


Figure 1: Illustration of our method applied to obtain important heads. Starting with the original model, key heads are identified highlighted in light blue. The different colored dashed lines (e.g., red, blue, green) represent the connections between heads and their correlated parameters. Last, we update the active parameters within heads highlighted in blue to represent unlearning process.

where $\mathbf{H}_{\theta^*}^{-1}$ is the inverse of the Hessian matrix $\nabla_{\theta}^2 \mathcal{L}(\theta^*; \mathcal{D}_r)$ evaluated at θ^* . This method is derived from influence function (Koh & Liang (2017)), which provides a bounded approximation error to facilitate effective unlearning (Guo et al. (2020)).

However, second-order unlearning involves the inverse Hessian computation, which is highly sensitive to parameters. Given the large number of parameters in large-scale models, this unlearning method cannot be applied directly. A common practice to approximate it is using the empirical FIM (Peste et al. (2021); Liu et al. (2024a); Gu et al. (2024)). Additionally, studies (Amari et al. (2019)) have shown that the off-diagonal elements of the FIM tend to be much smaller than the diagonal elements, usually by a factor $\frac{1}{\sqrt{n}}$, where n represents the dimension of the FIM. This insight highlights the effectiveness of using a diagonal approximation, particularly in large models with vast parameter counts (Hwang (2024)). As a result, we further adopt the empirical diagonal FIM $\widehat{\mathcal{I}}$ to approximate the Hessian matrix:

$$\widehat{\mathcal{I}}(\theta; \mathcal{D}) = \frac{1}{|\mathcal{D}|} \sum_{x \in \mathcal{D}} \nabla \ell(\theta; x)^2. \quad (3)$$

The storage of the diagonal FIM requires only $\mathcal{O}(d)$ space, and the inverse operation takes only $\mathcal{O}(d)$ time, where d denotes the number of model parameters. This makes second-order unlearning method straightforward and efficient to implement.

3 STRUCTURE-AWARE PARAMETER-EFFICIENT MACHINE UNLEARNING

Inspired by the lottery hypothesis (Frankle & Carbin (2018)), recent research suggests that localizing functional regions within neural networks can make the model more effective for specific tasks (Zhang et al. (2024b)). Given the high dimension for large models, empirically identifying influence-critical parameters from a too fine-grained perspective is both inefficient and often sub-optimal. To this end, we propose SPE-Unlearn, which derived a pair of masks to pinpoint influence-critical parameters within heads and filters in Section 3.1. By selectively targeting the most influence-critical parameters, SPE-Unlearn is integrated into second-order unlearning in 3.2, enabling more efficient and effective unlearning processes. At last, we extend SPE-Unlearn to support successive unlearning, demonstrating its robustness in Section 3.3.

3.1 STRUCTURE-AWARE PARAMETER LOCALIZATION

While parameter-efficient methods involve identifying critical parameters, this process can be framed as finding an optimal binary mask. In this context, a mask value of 1 indicates that the corresponding parameter should be updated, while a value of 0 represents that the corresponding parameter should remain frozen. Given that the number of structures is significantly smaller than the number of parameters (e.g., 37K vs. 110M in case of BERT-base), SPE-Unlearn adapt a coarse-grained method to pinpoint influence-critical parameters in heads and filters. Thus, we formulate

the unlearning objective (1) with a learnable pair of masks for the heads and filters as a constrained optimization problem. To streamline the problem, we provide a general expression for the heads and filters by introducing the mask variables m :

$$m^* = \arg \min_m \mathcal{L}(m; \theta^*, \mathcal{D}_r) \quad \text{s.t. } \frac{\sum_{i=1}^{|m|} m_i}{|m|} < 1 - S, \quad (4)$$

where $|m|$ is the number of mask variables, θ^* represents the original model, and S denotes the sparsity (e.g., 90%) which determines the proportion of frozen structures. Since we focus exclusively on the mask variables, we henceforth regard the parameters θ^* as constants. Thus, the total loss $\mathcal{L}(\theta; \mathcal{D}_r)$ can be mapped to $\mathcal{L}(m; \theta^*, \mathcal{D}_r)$. If \mathcal{L} is differentiable with respect to m , we then approximate $\mathcal{L}(m; \theta^*, \mathcal{D}_r)$ using the second-order Taylor series around the mask variables $\mathbb{1}$:

$$\mathcal{L}(m; \theta^*, \mathcal{D}_r) \approx \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r) - (\mathbb{1} - m) \nabla_m \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r) + \frac{1}{2} (\mathbb{1} - m)^T \nabla_m^2 \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r) (\mathbb{1} - m). \quad (5)$$

As the original model θ^* has converged to a local minimum of $\nabla_m \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D})$, we can assume that $\nabla_m \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}) = 0$ (LeCun et al. (1989)). Incorporating this assumption, we simplify gradient term in the Taylor series approximation, i.e., $\nabla_m \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r) = \nabla_m \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}) - \sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x) = -\sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x)$. As $\mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r)$ is a constant, we can adjust the unlearning objective with mask variables:

$$m^* \approx \arg \min_m (\mathbb{1} - m) \sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x) + \frac{1}{2} (\mathbb{1} - m)^T \nabla_m^2 \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r) (\mathbb{1} - m). \quad (6)$$

Thus, the optimization problem depends on the two factors: the gradient with respect to the forgetting dataset \mathcal{D}_f (i.e., $\sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x)$) and the Hessian matrix with respect to the remaining dataset \mathcal{D}_r (i.e., $\nabla_m^2 \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_r)$). These components together reflect the effectiveness of influence removal. Since forming the Hessian matrix directly is computationally prohibitive, we approximate it using the empirical **diagonal** FIM of the mask variables with Equation (3). This leads to a simplified form of the optimization objective in Equation (6):

$$m^* \approx \arg \min_m (\mathbb{1} - m) \sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x) + \frac{1}{2} (\mathbb{1} - m)^2 \widehat{\mathcal{I}}(\mathbb{1}; \theta^*, \mathcal{D}_r). \quad (7)$$

Given that the mask variable can only be set to 0 or 1, we transform the optimization problem into a mask selection problem with heads and filters:

$$m^* \approx \arg \min_m \sum_i \left[(1 - m_i) \left[\sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x) \right]_i + \frac{1}{2} (1 - m_i)^2 [\widehat{\mathcal{I}}(\mathbb{1}; \theta^*, \mathcal{D}_r)]_i \right]. \quad (8)$$

Therefore, we propose importance scores to identify influence-critical heads and filters. Each head or filter can be assessed based on the sum of its corresponding gradient and half of the diagonal FIM element. Heads or filters with higher scores will be prioritized for selection. Additionally, to better understand the influence of off-diagonal elements on mask selection for each layer, we replace the diagonal FIM with the block diagonal FIM, where each block is associated with a layer. Thus, Equation (7) decomposes into *layer-wise* optimization problems:

$$m_l^* \approx \arg \min_{m_l} (\mathbb{1} - m_l) \left[\sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; \theta^*, x) \right]_l + \frac{1}{2} (\mathbb{1} - m_l)^2 [\widehat{\mathcal{I}}(\mathbb{1}; \theta^*, \mathcal{D}_r)]_l, \quad (9)$$

where l represents the layer being optimized. This optimization problem can be efficiently solved using a greedy search with warm start (Kwon et al. (2022)), i.e., initializing the mask variables m_l derived from Equation (8). In this process, we iteratively swap unselected each head (or filter) with the highest importance score for selected one in the current mask to further optimize Equation (9), yielding an approximate solution after one round of swapping. Consequently, the rearranged mask variables captures the impact of intra-layer interactions, enabling precise localization of the parameters within the model structures. Additionally, our approach can be integrated with other methods for identifying influence-critical parameters, offering enhanced flexibility. Detailed information about these techniques can be found in Appendix A.4. In practice, our derivation can also be applied to other unlearning objectives, such as maximizing the loss on the forgetting dataset (Jia et al. (2024)). Detailed information is presented in Appendix B.

216 **Algorithm 1** Structure-aware Parameter-Efficient Second-Order Unlearning

217 **Input:** remaining dataset \mathcal{D}_r , forgetting dataset \mathcal{D}_f , Transformer model T , loss function ℓ , model
218 parameter θ , sparsity S , unlearning rate η

219 **Output:** Updated model parameter θ

220 1: Initialize mask $m \leftarrow \mathbb{1}$, parameter FIM $\widehat{\mathcal{I}} \leftarrow 0$, parameter gradients $g_\theta \leftarrow 0$, mask gradients
221 $g_m^r \leftarrow [], g_m^f \leftarrow []$

222 2: **for** each x in \mathcal{D}_r **do** ▷ Iterate data points in \mathcal{D}_r

223 3: $\nabla_m \ell(\theta, x), \nabla_\theta \ell(\theta, x) \leftarrow T(\theta, m, \ell, x)$

224 4: $\widehat{\mathcal{I}} += \frac{1}{|\mathcal{D}_r|} \nabla_\theta \ell(\theta, x)^2$ ▷ Obtain the parameter diagonal FIM in \mathcal{D}_r

225 5: Append $\nabla_m \ell(\theta, x)$ to g_m^r ▷ Gather the mask gradients in \mathcal{D}_r

226 6: **end for**

227 7: **for** each x in \mathcal{D}_f **do** ▷ Iterate data points in \mathcal{D}_f

228 8: $\nabla_m \ell(\theta, x_i), \nabla_\theta \ell(\theta, x_i) \leftarrow T(\theta, m, \ell, x_i)$

229 9: $g_\theta += \nabla_\theta \ell(\theta, x_i)$ ▷ Obtain the parameter gradient in \mathcal{D}_f

230 10: Append $\nabla_m \ell(\theta, x_i)$ to g_m^f ▷ Gather the mask gradients in \mathcal{D}_f

231 11: **end for**

232 12: $SC \leftarrow \frac{1}{2}(g_m^r)^2 + g_m^f$ ▷ Compute importance scores of structures

233 13: $IN \leftarrow$ indices of unimportant heads ▷ Find the optimal mask indices

234 14: $IN^* \leftarrow$ rearrange the mask indices with warm start

235 15: $m[IN^*] = 0$ ▷ Set unimportant indices to 0

236 16: $\theta += \eta * m \circ \widehat{\mathcal{I}}^{-1} g_\theta$ ▷ Sparse Second-Order unlearning update

237 17: **return** θ

238

239 3.2 STRUCTURE-AWARE PARAMETER-EFFICIENT SECOND-ORDER UNLEARNING

240

241 By pinpointing influence-critical parameters within the heads and filters, SPE-Unlearn enables
242 efficient integration with widely-adopted unlearning methods, e.g., fine-tuning ([Golatkar et al.
243 \(2020\)](#)) and gradient difference ([Liu et al. \(2022\); Jia et al. \(2024\)](#)). A key observation is that
244 both SPE-Unlearn and second-order unlearning share the computational need for gradients and
245 FIM. Therefore, we leverage second-order unlearning as a representative case study to showcase the
246 efficacy of our approach.

247 Following the insights of SPE-Unlearn, we formalize **Structure-aware Parameter-Efficient
248 Second-Order unlearning** (SPE-SO) by introducing sparse mask variables linked to the outputs of
249 heads and filters:

250
$$\theta \approx \theta^* + m \circ \left[[\widehat{\mathcal{I}}(\theta^*; \mathcal{D}_r)]^{-1} \sum_{x \in \mathcal{D}_f} \nabla_\theta \ell(\theta^*; x) \right], \quad (10)$$

251

252 where m are the binary mask variables, and \circ denotes the Hadamard product. Note that Equation 2
253 can be represented by setting all mask variables to 1. This method introduces several key advantages
254 over standard unlearning techniques. First, by incorporating sparsity through structure-aware masks,
255 SPE-SO significantly reduces the number of parameters required for the expensive computation
256 of the Hessian matrix. This leads to lower computational complexity, making the method more
257 scalable and efficient when applied to large-scale models. Second, SPE-SO offers a more tightly
258 bounded approximation error compared to standard methods. The approximation error is reduced
259 by a factor that is directly proportional to the sparsity introduced by the mask variables. This ensures
260 that the unlearning process remains highly accurate while avoiding unnecessary parameter updates.
261 Furthermore, by restricting the influence-critical parameters within the heads and filters, SPE-SO
262 provides fine-grained control over the error bounds.

263 Algorithm 1 presents the workflow of SPE-SO, which handles removal requests by accumulating
264 and processing them collectively. The algorithm can be adapted to various constraints, such as
265 time or memory. For scenarios where computational efficiency is the primary concern, SPE-SO
266 allows for pre-computation of the gradient and diagonal FIM for the entire training dataset. Upon
267 receiving removal requests, we can compute the data information about forgetting dataset to obtain
268 the required data, i.e., the gradient of forgetting dataset and diagonal FIM of remaining dataset.
269 Alternatively, to reduce memory consumption, SPE-SO can retrieve only the necessary information
by utilizing selected structures tied to specific parameters.

270 3.3 STRUCTURE-AWARE PARAMETER-EFFICIENT SUCCESSIVE UNLEARNING
 271

272 Successive unlearning presents a practical scenario where data owners request the removal of data
 273 points from the model at intervals, necessitating prompt deletion (e.g., machine learning as a ser-
 274 vice (MLaaS) (Hu et al. (2023))). While prior work has proposed different approaches to suc-
 275 ccessive unlearning, we introduce the classification to better differentiate how unlearning algorithm is
 276 used. Specifically, we categorize second-order successive unlearning into two distinct types based
 277 on whether or not the algorithm retains information from removed data: memory-free (Guo et al.
 278 (2020); Gu et al. (2024)) and memory-aided (Liu et al. (2023a)).
 279

279 Memory-free unlearning iteratively update the **latest model** following each removal request without
 280 retaining any information from the removed data. However, this method increases the unlearning
 281 approximation error, as the updates are based solely on the latest model, which can be more severe
 282 for Transformers. In contrast, memory-aided unlearning retains data information (i.e., gradients and
 283 FIM) to efficient unlearn on the **original model**. In what follows, we apply structure-aware masks
 284 into these successive unlearning scenarios and discuss the advantages of these masks.
 285

285 3.3.1 MEMORY-FREE UNLEARNING
 286

287 The way to apply SPE-Unlearn into the memory-free unlearning is straightforward. Upon each
 288 data removal request, we can directly compute the structure-aware mask and apply second-order
 289 unlearning. Specifically, the model is progressively updated based on the state from the previous un-
 290 learning cycle. At timestamp t (i.e., the t -th unlearning request), structure-aware parameter-efficient
 291 memory-free unlearning can be formalized:

$$m^t \circ \left[[\widehat{\mathcal{I}}(\theta^{t-1}; \mathcal{D}_r^t)]^{-1} \nabla_{\theta} \ell(\theta^{t-1}; x^t) \right], \quad (11)$$

294 where θ^{t-1} represents the unlearned model parameters at timestamp $t - 1$, \mathcal{D}_r^t and x^t denote the
 295 remaining dataset and the data point to be removed at timestamp t . Additionally, m^t is the structure-
 296 aware mask corresponding to the t -th removal request.

297 Although memory-free unlearning is simple and easy to implement, it suffers a major drawbacks.
 298 This method inherently diverges from the Taylor series approximation, which tends to introduce
 299 small errors during each approximation. As these errors accumulate with each successive update,
 300 the model is continually adjusted based on its latest state rather than retaining the original form.
 301 Consequently, with an increasing number of removal requests, the disparity between the original
 302 and updated models widens, resulting in a gradual decline in model performance.
 303

304 Once the number of unlearning requests surpasses a certain threshold, the model needs to
 305 be retrained from scratch to recover its per-
 306 formance (detailed in Table 1). Fortunately,
 307 structure-aware masks allows for more removal
 308 requests before retraining becomes necessary
 309 (as shown in Figure 4). This improvement is
 310 likely due to selectively adjust only the structures directly related to the data being removed. By
 311 confining the cumulative errors to a minimal subset of parameters, the overall impact on the model
 312 performance is reduced. Consequently, the model remains robust even after multiple unlearning
 313 operations, delaying the need for costly retraining.
 314

315 3.3.2 MEMORY-AIDED UNLEARNING
 316

316 Compared to memory-free unlearning, memory-aided unlearning approximates directly through a
 317 Taylor expansion at original model parameters. In contrast, memory-aided unlearning (Liu et al.
 318 (2023a)) accumulates the gradients on forgotten data and FIM on remaining dataset to achieve
 319 unlearning. Specifically, upon receiving the t -th unlearning request, structure-aware parameter-
 320 efficient memory-aided unlearning at timestamp t can be expressed as follows:
 321

$$m^t \circ \left\{ \left[\frac{|\mathcal{D}_r^{t-1}| \cdot \widehat{\mathcal{I}}(\theta^*; \mathcal{D}_r^{t-1}) - \widehat{\mathcal{I}}(\theta^*; x^t)}{|\mathcal{D}_r^{t-1} - 1|} \right]^{-1} \left[\sum_{x \in \mathcal{D}_f^{t-1}} \nabla_{\theta} \ell(\theta^*; x) + \nabla_{\theta} \ell(\theta^*; x^t) \right] \right\}, \quad (12)$$

324 where \mathcal{D}_f^{t-1} represents the data points that have already been removed at timestamp $t - 1$, \mathcal{D}_r^{t-1}
 325 denotes the remaining dataset at timestamp $t - 1$. In practice, rather than storing these data points
 326 directly, we retain the gradients or FIM associated with the data in memory. With each new unlearn-
 327 ing request, these data information are updated accordingly. Furthermore, considering the proportion
 328 of the forgetting dataset is negligible, the mask selection process can be accelerated. As a result, in
 329 the mask selection Equation (8), the term $\sum_{x \in \mathcal{D}_f} \nabla_m \ell(\mathbb{1}; x)$ can be omitted, and the term $\widehat{\mathcal{I}}(\mathbb{1}; \mathcal{D}_r)$
 330 can be approximated by $\widehat{\mathcal{I}}(\mathbb{1}; \mathcal{D})$, resulting in the following simplification:
 331

$$m^* \approx \arg \min_m \sum_i (\mathbb{1} - m_i)^2 \widehat{\mathcal{I}}(\mathbb{1}; \mathcal{D})_i, \quad (13)$$

334 Since the Equation (13) is derived based solely on the entire dataset, the corresponding mask can
 335 be pre-computed during the pre-unlearning phase. Although this simplification enhances efficiency,
 336 it does not fully account for the influence of the data points slated for deletion. Thus, we finally
 337 rearrange the mask variables using Equation (9), which allows for a more targeted adjustment. In
 338 memory-aided scenario, unlearning is achieved through a single-step second-order update on the
 339 original model. Therefore, the key strength of structure-aware masks stems from the superiority of
 340 SPE-Unlearn in handling general second-order unlearning, which offers a tighter approximation
 341 error bound to facilitate more effective and accurate data removal.
 342

4 EXPERIMENTS

4.1 EXPERIMENT SETUPS

346 **Models and Datasets.** We conduct comprehensive experiments on three pretrained Transformer
 347 models: BERT-base ([Devlin et al. \(2018\)](#)), DistilBERT ([Sanh et al. \(2019\)](#)), and RoBERTa-large
 348 ([Liu et al. \(2019\)](#)). These models are accessed through the HuggingFace Transformers library ([Wolf
 349 et al. \(2020\)](#)). Our evaluation spans four GLUE tasks (MNLI, QQP, SST-2, and STS-B) ([Wang et al.
 350 \(2018\)](#)) and two SQuAD tasks (SQuAD v1.1 and SQuAD v2.0) ([Rajpurkar \(2016\)](#)). Consistent with
 351 the configurations outlined in prior works [Devlin et al. \(2018\)](#); [Sanh et al. \(2019\)](#); [Liu et al. \(2019\)](#),
 352 we fine-tune these models, treating them as the original models for our experiments.
 353

354 **Unlearning methods.** Our experiments focus on comparing the proposed method SPE-SO with
 355 several other unlearning methods. These methods include Fine-Tuning (FT), Gradient Difference
 356 (GD) ([Liu et al. \(2022\)](#), [Jia et al. \(2024\)](#)), Sparsity-Aware unlearning (SA) ([Liu et al. \(2024a\)](#)). For
 357 FT, we continue training the original model on the remaining dataset for 3 epochs. For GD, the
 358 model is fine-tuned on entire dataset for 3 epochs, with the gradient direction reversed for the data
 359 that needs to be forgotten. For SA, fine-tuning is performed on the remaining dataset with a sparsity
 360 penalty ($\gamma = 5e - 5$) applied to the parameters for 3 epochs. Additionally, we also include the
 361 standard Second-Order unlearning (SO) method, which excludes structure-aware masks, to evaluate
 362 the effectiveness of SPE-Unlearn. Meanwhile, Retraining from scratch (RT) serves as the gold
 363 standard, where the model is fine-tuned on the remaining dataset following the configurations from
 364 [Devlin et al. \(2018\)](#); [Sanh et al. \(2019\)](#); [Liu et al. \(2019\)](#). Detailed hyperparameters are presented in
 365 Appendix A.1.
 366

367 **Evaluation metrics.** We analyze the unlearning methods from three aspects ([Warnecke et al. \(2021\)](#);
 368 [Gu et al. \(2024\)](#)): 1) **Efficacy** in removing the targeted data. We evaluate this using unlearning
 369 accuracy and membership inference attacks (MIA) on \mathcal{D}_f . Unlearning accuracy directly reflects
 370 the effectiveness of the unlearning algorithm, while MIA assesses the vulnerability of the model
 371 to attacks after unlearning. In practice, we use a confidence-based MIA predictor to gauge the
 372 likelihood of a successful attack ([Liu et al. \(2024a\)](#); [Song et al. \(2019\)](#)); 2) **Fidelity** of model utility.
 373 We measure this by examining both the remaining accuracy and the testing accuracy to assess the
 374 preservation of model performance and its generalization ability after unlearning; 3) **Efficiency** of
 375 executing the unlearning methods. We report the time required to perform unlearning as a measure
 376 of speed and computational efficiency.
 377

4.2 EXPERIMENT RESULTS

378 We present the experimental results using the MNLI dataset as a case study. Detailed results for
 379 additional datasets are provided in Appendix A. Unless otherwise specified, the default number of
 380

378
 379 Table 2: Overall results of unlearning performance using different unlearning methods under three
 380 fine-tuned models. We focus on 90% sparsity SPE-SO as our approaches.
 381

Model	Method	Efficacy		Fidelity		Efficiency
		Unlearning Accuracy ↓	MIA ↓	Remaining Accuracy ↑	Testing Accuracy ↑	
BERT-base	RT	85.16%	0.7500	97.95%	84.78%	8880s
	FT	92.19%	0.8594	99.16%	84.63%	5651s
	GD	90.62%	0.8437	99.13%	84.20%	5690s
	SA	89.84%	0.8437	92.77%	82.05%	4845s
	SO	85.94%	0.8047	94.07%	84.60%	1160s
	SPE-SO	85.94%	0.7969	94.15%	84.62%	1274s
DistilBERT	RT	82.81%	0.7266	96.61%	82.47%	4989s
	FT	94.53%	0.8906	98.94%	81.63%	2434s
	GD	91.41%	0.8750	98.72%	81.37%	2498s
	SA	90.62%	0.8750	96.49%	81.23%	2399s
	SO	89.06%	0.8516	96.37%	81.29%	587s
	SPE-SO	88.28%	0.8359	96.47%	81.62%	643s
RoBERTa-large	RT	90.62%	0.8125	98.79%	90.02%	62068s
	FT	97.66%	0.9766	99.50%	90.02%	18004s
	GD	95.31%	0.8906	99.64%	89.57%	18176s
	SA	92.97%	0.8906	96.86%	87.08%	14634s
	SO	92.97%	0.8906	94.32%	88.99%	3575s
	SPE-SO	92.19%	0.8906	95.75%	89.52%	3642s

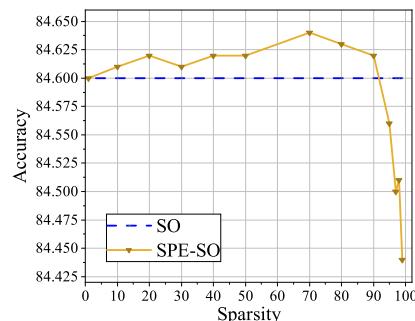
397
 398 unlearned samples is 128. We randomly select 128 samples as the forgetting dataset \mathcal{D}_f and use all
 399 orthogonal samples as the remaining dataset \mathcal{D}_r . In what follows, we compare different unlearning
 400 methods and conduct an in-depth analysis of our approach.
 401

402 **Structure-Aware sparse unlearning is effective.** Table 2 presents the unlearning performance
 403 of various unlearning methods across three mod-
 404 els. As subsequent experiments show that 90%
 405 sparsity is sufficient for effective unlearning, we
 406 focus on the SPE-SO with 90% sparsity regime
 407 for comparison with other methods. Our ex-
 408 periments reveal that FT is inefficient for unlearning in Transformers, while SA demonstrates strong
 409 unlearning efficacy but at the cost of significantly compromising model fidelity. GD generally strikes
 410 a reasonable balance between efficacy and efficiency. However, these methods demands consider-
 411 able time due to the lengthy fine-tuning process. In contrast, both SO and SPE-SO achieve effective
 412 unlearning with just a single epoch over the dataset, which provide robust efficacy guarantees with
 413 minimal impact on fidelity. As shown in Table 3, we further compare memory usage during model
 414 updates for SO and SPE-SO. Although SPE-SO takes more time to identify structure-aware mask,
 415 it has lower storage overhead and delivers superior performance compared to SO. Thus, we con-
 416 clude that SPE-SO with 90% sparsity is sufficient to strike a favorable “efficacy-fidelity-efficiency”
 417 trade-off.
 418

419 **A sparsity of 90% is sufficient for effective unlearning.**
 420 We explore the effectiveness of various sparsity strategies
 421 in facilitating unlearning. Figure 2 shows the relationship
 422 between testing accuracy and sparsity while maintaining
 423 comparable unlearning efficacy. As sparsity increases up
 424 to 90%, the model retains high utility. However, when
 425 sparsity surpasses 90%, a sharp decline in model accu-
 426 racy occurs, indicating that updating fewer than 10% of
 427 parameters may be insufficient to preserve utility. Similar
 428 effects of sparsity strategies on unlearning performance
 429 are observed across other datasets (detailed in Appendix
 430 A.3). We also delve into the functional regions respon-
 431 sible for unlearning within models, but find no single net-
 432 work layer that stands out as particularly crucial for un-
 433 learning. This suggests that the effectiveness of unlearn-
 434 ing may be task-specific, resisting any fixed structural or parametric approach. Overall, our findings

398 Table 3: Memory consumption with three mod-
 399 els. SPE-SO takes 90% sparsity.
 400

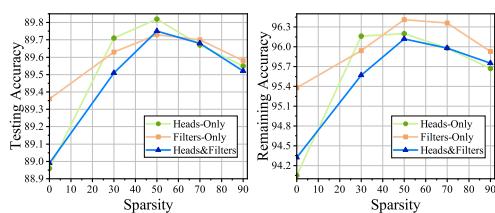
Memory (MB)	BERT-base	DistilBERT	RoBERTa-large
SO	995.7	544.0	3371.6
SPE-SO	663.8	377.4	2174.9



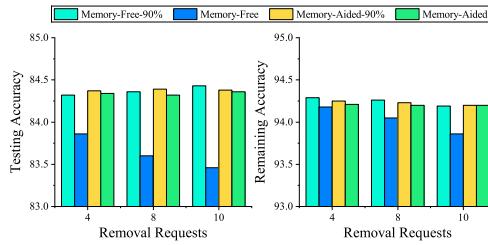
401 Figure 2: Testing accuracy of SO and
 402 SPE-SO applied to BERT-base across
 403 varying sparsity.
 404

432 emphasize that a 90% sparsity strategy strikes the sufficient balance between efficiency and effectiveness
 433 in unlearning tasks, offering a practical approach without compromising much utility.
 434

435 **Selective parameter updates only in filters can effectively accomplish unlearning.** (Pochinkov
 436 & Schoots (2024)) argued that pruning filters is more effective than heads. To further investigate this
 437 claim, we conducted a comparative analysis of three selective parameter update strategies: heads-
 438 only, filters-only, and heads&filters in Figure 3. All the experiments are designed to provide compa-
 439 rable unlearning guarantees varying sparsity. While the heads-only approach demonstrated superior
 440 testing accuracy at moderate sparsity levels (30% to 70%), it falls behind in terms of remaining accu-
 441 racy. In contrast, the filters-only strategy not only maintained stability at lower sparsity but also
 442 delivered consistently strong unlearning performance at higher sparsity. Notably, we observed that
 443 compared to updating the parameter both in heads and filters, updating only the parameters in either
 444 heads or filters can achieve better unlearning performance. This underscores that more focused up-
 445 dates may mitigate unnecessary overhead, without sacrificing performance. Among the approaches,
 446 filters-only updates consistently proved to be the most stable and effective, making it a more optimal
 447 choice for robust unlearning.



448 Figure 3: Testing accuracy and remaining accu-
 449 racy for various sparsity applied to RoBERTa-
 450 large after unlearning different structures.
 451



452 Figure 4: Results using memory-free unlearn-
 453 ing and memory-aided unlearning on BERT-
 454 base under varying removal requests.
 455

456 **Structure-aware masks benefit robust unlearning.** Our method highlights that structure-aware
 457 masks serve as an effective mechanism for guiding data removal, enabling models to meet strict un-
 458 learning guarantees while preserving model performance. Motivated by this observation, we further
 459 explore the potential of structure-aware masks in successive unlearning scenarios, focusing on both
 460 memory-free and memory-aided unlearning, as depicted in Figure 4. Our results show that sparse
 461 updates with structure-aware offer marginal improvements over full updates in memory-aided un-
 462 learning. This is likely because memory-aided unlearning operates by updating the model directly
 463 from its original state in a single step, minimizing the relative advantage of sparse updates. In con-
 464 trast, sparse updates offer significant benefits in memory-free unlearning. When all parameters are
 465 updated in memory-free unlearning, model fidelity is overly impacted, consistent with the analysis
 466 in Section 3.3.1. However, applying sparse updates with 90% sparsity in memory-free unlearning
 467 preserves high model utility, even after 10 removal requests. This suggests that structure-aware
 468 masks can support a higher volume of removal requests before retraining becomes necessary. These
 469 results highlight the potential of structure-aware masks to enhance the robustness of unlearning.
 470

5 RELATED WORK

471 **Transformer Unlearning.** The concept of machine unlearning was first introduced by (Cao & Yang
 472 (2015)). Initially applied to simple model, machine unlearning has since been extended to Trans-
 473 former models (Jang et al. (2022); Eldan & Russinovich (2023); Yao et al. (2023; 2024); Chen et al.
 474 (2024); Jia et al. (2024); Gu et al. (2024)). (Jang et al. (2022)) proposed inverting the training objec-
 475 tive on forgetting sequences and utilize straightforward gradient ascent. As gradient ascent signifi-
 476 cantly degrades performance, (Yao et al. (2024)) refined the objective function by employing gradi-
 477 ent descent on in-distribution data to enhance robustness. Subsequently, (Jia et al. (2024)) provided
 478 a comprehensive overview of unlearning objectives and developed a second-order optimization un-
 479 learning approach. (Gu et al. (2024)) further investigated the effectiveness of second-order updates
 480 in Transformers. However, these methods primarily focus on updating all model parameters, which
 481 is computationally expensive and time-consuming. In our work, we study the parameter-efficient
 482 methods to achieve effective unlearning in Transformers.
 483

Parameter-efficient Unlearning. Parameter-efficient unlearning methods focus on identifying influence-critical parameters and updating only those to accelerate the unlearning process. Several strategies (Ma et al. (2022); Pochinkov & Schoots (2024); Shi et al. (2023); Liu et al. (2023b); Wu & Harandi (2024); Foster et al. (2024); Schoepf et al. (2024)) have been proposed to assess parameter importance. Although these approaches may be applicable to Transformers, they are largely heuristic or empirical, which can result in sub-optimal outcomes for unlearning tasks. Recently, (Liu et al. (2024a)) highlighted that unlearning can be effective when performed on a pruned model with a theoretical foundation. However, pruning primarily focuses on identifying parameters critical to maintain model performance, which does not align with the desiderata of unlearning. Additionally, the focus on parameter ignore the complex intra-structural interactions within Transformers, which results in inaccurate identification of the parameters. Therefore, we specifically target at heads and filters within Transformers and derive an efficient strategy to identify influence-critical parameters.

6 CONCLUDING REMARKS

In this work, we propose structure-aware parameter-efficient unlearning (SPE-Unlearn), a novel approach tailored for Transformers. SPE-Unlearn derives an optimal masking strategy to identify influence-critical parameters within heads and filters. By selectively targeting these key parameters, SPE-Unlearn infuses into second-order unlearning update to demonstrate its efficacy and strengths. We further analyze the advantages of our method across both memory-free and memory-aided successive unlearning scenarios. Empirical study demonstrate that our method accommodates more removal requests than standard second-order unlearning in memory-free unlearning scenarios. Additionally, comprehensive experiments conducted on various Transformer models and datasets reveal that our method with 90% sparsity outperforms existing approaches.

For future work, we suggest extending to other existing unlearning methods to demonstrate the effectiveness of SPE-Unlearn in Transformers. While our experiments focus on small-scale Transformers, we plan to explore larger-scale models (e.g, OPT-13b and LLaMA2-13b) to better understand the behavior of structure-aware masks. Furthermore, our study concentrates on fine-tuned models, which limits the ability to unlearn deeply ingrained undesired information from pre-trained models. To address this, we aim to identify structure-aware masks directly in pre-trained models.

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666 A ADDITIONAL EXPERIMENTAL DETAILS

668 A.1 HYPERPARAMETERS

669 We fine-tune BERT-base, DistilBERT, and RoBERTa-large on different datasets using AdamW with
 670 weight decay of 0 as vanilla models. The learning rate is selected from 10^{-5} , $2 \cdot 10^{-5}$, $3 \cdot 10^{-5}$ and
 671 $5 \cdot 10^{-5}$. Such learning rate is also applied to unlearning methods such as Retraining, Fine-Tuning
 672 (FT), Structure-Aware unlearning (SA), and Gradient Difference (GD). The number of epochs is set
 673 to 5. For unlearning, the number of epochs is fixed at 3 for FT, SA, and GD, while it is fixed at 5 for
 674 retraining. The unlearning rate for Second-Order unlearning (SO) are chosen through grid search in
 675 the range $[10^{-6}, 10^{-7}]$. For Structure-aware Parameter-Efficient Second-Order SO (SPE-SO), the
 676 unlearning rate increases proportionally with the fraction of updated parameters relative to the total
 677 parameters compared to SO.

679 A.2 COMPARE TO OTHER UNLEARNING METHODS

680 We compare ours to other unlearning methods in three GLUE tasks (QQP, SST-2 and STS-B) and
 681 two SQuAD tasks (SQuAD v1.1 and SQuAD v2.0) under three models (detailed in Table 4 to Table
 682 8). The evaluation metrics vary depending on the task. For example, we use Spearman correlations
 683 to assess STS-B, while F1 scores are reported for both SQuAD v1.1 and SQuAD v2.0. Higher values
 684 for these metrics indicate better model performance.

687 A.3 FIND THE APPROPRIATE SPARSITY

688 We aim to determine the level of sparsity that can ensure adequate model performance while pro-
 689 viding sufficient unlearning guarantees. Therefore, we conducted a detailed sparsity analysis on
 690 additional datasets, as shown in Figure 5. Our results indicate that updating all parameters is not
 691 the most effective strategy for unlearning, as it can lead to excessive forgetting, causing a rapid de-
 692 cline in model performance. In contrast, we found that a sparsity of 50% offers the most efficient
 693 improvement in unlearning. Moreover, sparsity levels between 80% and 90% perform on par with,
 694 and sometimes even surpass, the performance of other methods.

696 A.4 IDENTIFY INFLUENCE-CRITICAL PARAMETERS IN STRUCTURES

697 In our approach, the mask is applied to specific heads and filters, resulting in a relatively coarse
 698 granularity for unlearning. To achieve a more refined and precise method, we further investigate the
 699 importance of individual parameters within these selected heads and filters. Our hypothesis is that
 700 by focusing on individual parameters, we can identify more fine-grained regions that are critical for
 701 effective unlearning.

Table 4: Overall results of unlearning performance using different unlearning methods under three fine-tuned models on QQP dataset.

Model	Method	Efficacy		Fidelity		Efficiency
		Unlearning Accuracy ↓	MIA ↓	Remaining Accuracy ↑	Testing Accuracy ↑	
BERT-base	RT	92.97%	0.875	98.48%	91.38%	9560s
	FT	96.09%	0.9219	99.56%	91.26%	5759s
	GD	96.09%	0.9219	99.58%	91.29%	5858s
	SA	92.19%	0.8906	92.52%	88.52%	5579s
	SO	92.97%	0.9063	97.69%	90.65%	832s
	SPE-SO	92.19%	0.8906	98.03%	90.72%	926s
DistilBERT	RT	90.62%	0.8203	98.52%	90.39%	6291s
	FT	98.44%	0.9453	99.61%	90.25%	3619s
	GD	96.09%	0.9297	99.65%	90.16%	3763s
	SA	95.31%	0.9141	96.80%	81.14%	3571s
	SO	92.19%	0.8594	98.23%	90.05%	415s
	SPE-SO	91.41%	0.8594	98.36%	90.12%	468s
RoBERTa-large	RT	91.41%	0.8594	99.17%	92.19%	79214s
	FT	98.44%	0.9609	99.85%	92.18%	21742s
	GD	94.53%	0.9453	99.91%	92.14%	23239s
	SA	93.75%	0.8750	98.69%	91.48%	20793s
	SO	92.97%	0.8750	98.93%	91.56%	2598s
	SPE-SO	92.87%	0.8750	98.86%	91.46%	2639s

Table 5: Overall results of unlearning performance using different unlearning methods under three fine-tuned models on SST-2 dataset.

Model	Method	Efficacy		Fidelity		Efficiency
		Unlearning Accuracy ↓	MIA ↓	Remaining Accuracy ↑	Testing Accuracy ↑	
BERT-base	RT	93.75%	0.9297	99.06%	93.00%	915s
	FT	96.88%	0.9609	99.25%	92.78%	479s
	GD	95.31%	0.8984	99.53%	92.78%	518s
	SA	95.31%	0.9062	98.82%	89.79%	450s
	SO	94.53%	0.9141	98.96%	92.89%	93s
	SPE-SO	94.53%	0.8984	98.93%	93.35%	103s
DistilBERT	RT	92.97%	0.8984	98.78%	91.40%	403s
	FT	95.31%	0.8594	97.67%	90.37%	238s
	GD	94.53%	0.8984	98.90%	90.25%	243s
	SA	94.53%	0.9141	98.28%	89.91%	235s
	SO	94.53%	0.8906	96.35%	91.63%	46s
	SPE-SO	94.53%	0.8906	96.35%	91.63%	53s
RoBERTa-large	RT	94.53%	0.9063	99.64%	96.10%	4698s
	FT	97.66%	0.9753	99.44%	96.22%	1430s
	GD	93.75%	0.9219	97.98%	95.33%	1492s
	SA	94.53%	0.9219	98.84%	95.07%	1423s
	SO	94.53%	0.9297	99.14%	94.15%	311s
	SPE-SO	94.53%	0.8984	99.45%	94.55%	320s

To implement this, we adopt Wanda (Sun et al. (2023)) as our selection mechanism. Wanda operates by analyzing the forgetting dataset, which serves as the input for the selective process. The values returned by Wanda represent the importance of each neuron to the unlearning task—higher values indicate neurons that are more critical for unlearning. After this analysis, we apply a sparsity of 90% on SPE-SO, selecting the most important parameters to retain based on their Wanda scores. These selected parameters are then targeted for unlearning. This method not only aligns with the broader goal of structural selection but also enhances the precision of unlearning by targeting specific neurons within the model. Detailed results of this selective mechanism are shown in Table 9.

However, Our experimental results indicate that incorporating the parameter selection mechanism does not improve unlearning performance in SPE-SO. We hypothesize that this outcome stems from the inherent complexity of balancing unlearning precision with model utility. While selecting individual parameters based on their Wanda scores allows for a more targeted and theoretically precise unlearning process, this fine-grained approach may inadvertently reduce the overall model’s adaptability and robustness.

756
757 Table 6: Overall results of unlearning performance using different unlearning methods under three
758 fine-tuned models on STS-B dataset.

759 760 Model	761 Method	762 Efficacy		763 Fidelity		764 Efficiency Time ↓
		765 Unlearning Spearman Corr. ↓	766 MIA ↓	767 Remaining Spearman Corr. ↑	768 Testing Spearman Corr. ↑	
769 BERT-base	770 RT	771 86.60%	772 0.5156	773 97.86%	774 88.63%	775 148s
	776 FT	777 95.37%	778 0.8750	779 96.72%	780 88.49%	781 76s
	782 GD	783 91.66%	784 0.594	785 99.17%	786 88.50%	787 84s
	788 SA	789 98.70%	790 0.8750	791 99.31%	792 88.60%	793 64s
	794 SO	795 86.91%	796 0.632	797 98.00%	798 87.76%	799 9s
	800 SPE-SO	801 86.47%	802 0.5234	803 98.24%	804 87.76%	805 10s
806 DistilBERT	807 RT	808 87.31%	809 0.6563	810 93.10%	811 85.45%	812 62s
	813 FT	814 91.15%	815 0.6875	816 95.20%	817 85.16%	818 29s
	819 GD	820 89.12%	821 0.7031	822 94.81%	823 85.37%	824 30s
	825 SA	826 92.36%	827 0.7109	828 93.85%	829 85.26%	830 27s
	831 SO	832 87.61%	833 0.6875	834 91.71%	835 85.02%	836 5s
	837 SPE-SO	838 87.75%	839 0.703125	840 92.01%	841 85.26%	842 5.5s
843 RoBERTa-large	844 RT	845 90.97%	846 0.5781	847 97.95%	848 92.01%	849 671s
	850 FT	851 96.19%	852 0.7656	853 98.68%	854 91.92%	855 198s
	856 GD	857 92.18%	858 0.5703	859 96.17%	860 90.33%	861 205s
	862 SA	863 96.25%	864 0.7344	865 98.68%	866 91.57%	867 176s
	868 SO	869 91.28%	870 0.5078	871 97.46%	872 91.57%	873 31s
	874 SPE-SO	875 91.13%	876 0.484375	877 97.88%	878 91.79%	879 35s

773 Table 7: Overall results of unlearning performance using different unlearning methods under three
774 fine-tuned models on SQuAD v1.1 dataset.

776 777 Model	778 Method	779 Efficacy		780 Fidelity		781 Efficiency Time ↓
		782 Unlearning F1 ↓	783 MIA ↓	784 Remaining F1 ↑	785 Testing F1 ↑	
786 BERT-base	787 RT	788 87.62%	789 0.5938	790 95.23%	791 88.18%	792 6328s
	793 FT	794 92.36%	795 0.7578	796 96.38%	797 87.73%	798 3765s
	799 GD	800 87.27%	801 0.6797	802 96.44%	803 87.34%	804 3775s
	805 SA	806 89.75%	807 0.7031	808 91.94%	809 86.85%	810 3800s
	811 SO	812 86.26%	813 0.5625	814 94.33%	815 87.74%	816 764s
	817 SPE-SO	818 86.74%	819 0.5781	820 94.25%	821 87.60%	822 809s
823 DistilBERT	824 RT	825 84.38%	826 0.5391	827 94.34%	828 85.35%	829 3203s
	830 FT	831 92.54%	832 0.7188	833 97.49%	834 85.09%	835 1936s
	836 GD	837 87.18%	838 0.6016	839 97.54%	840 85.05%	841 1956s
	842 SA	843 89.52%	844 0.7109	845 96.42%	846 83.86%	847 1921s
	848 SO	849 85.79%	850 0.5547	851 93.51%	852 85.35%	853 763s
	854 SPE-SO	855 85.35%	856 0.5547	857 93.65%	858 85.37%	859 812s
860 RoBERTa-large	861 RT	862 90.41%	863 0.6484	864 97.92%	865 92.50%	866 18439s
	867 FT	868 94.74%	869 0.7734	870 98.97%	871 93.15%	872 11365s
	873 GD	874 91.75%	875 0.6484	876 99.15%	877 92.98%	878 11520s
	879 SA	880 91.05%	881 0.6875	882 95.16%	883 89.36%	884 11116s
	885 SO	886 90.71%	887 0.500	888 94.93%	889 90.95%	890 2008s
	891 SPE-SO	892 90.81%	893 0.5234	894 95.14%	895 91.03%	896 2141s

797 B IDENTIFY KEY STRUCTURES IN OTHER UNLEARNING OBJECTIVE

798
799 Machine unlearning typically relies on the specific unlearning objective to design optimization algo-
800 rithms. For instance, second-order unlearning is achieved by minimizing the loss on the remaining
801 dataset (i.e., Equation 2). To simplify the optimization, a Taylor expansion of the unlearning ob-
802 jective is performed on the original model. Following the optimization process, we identify the
803 influence-critical parameters using a structure-aware approach in Transformers, which then guides
804 the second-order unlearning update.

805
806 The mainstream class of existing Large Language Model (LLM) unlearning methods also follow the
807 pattern of optimization based on the objective function. Gradient Ascent (GA) (Jang et al. (2022))
808 aims to maximize the loss for the forgetting dataset. Building on this, Gradient Difference (GD) (Liu
809 et al. (2022)) further strives to maintain performance on the remaining dataset. Direct Preference
810 Optimization (DPO) (Rafailov et al. (2024)) seeks to align the model by replacing the original re-
811 sponse on forgetting dataset with the alternative answers “I don’t know”. Inspired by DPO, Negative

810
811 Table 8: Overall results of unlearning performance using different unlearning methods under three
812 fine-tuned models on SQuAD v2.0 dataset.
813

814 Model	815 Method	816 Efficacy		817 Fidelity		818 Efficiency
		819 Unlearning F1 ↓	820 MIA ↓	821 Remaining F1 ↑	822 Testing F1 ↑	
823 BERT-base	RT	73.77%	0.6484	98.72%	75.77%	9560s
	FT	88.80%	0.8047	98.84%	74.52%	5532s
	GD	81.54%	0.7344	90.28%	74.22%	5600s
	SA	79.16%	0.6797	96.03%	72.65%	5512s
	SO	78.03%	0.6797	93.66%	73.33%	1043s
	SPE-SO	77.40%	0.6563	93.90%	73.57%	1123s
824 DistilBERT	RT	71.86%	0.6641	93.75%	69.80%	4715s
	FT	89.78%	0.8047	97.55%	69.71%	2880s
	GD	79.93%	0.7188	97.28%	68.16%	2894s
	SA	80.89%	0.7497	95.76%	68.46%	2863s
	SO	77.73%	0.7109	92.10%	68.36%	415s
	SPE-SO	76.30%	0.7031	91.82%	67.95%	468s
825 RoBERTa-large	RT	87.03%	0.7734	98.42%	86.58%	27053s
	FT	89.15%	0.7891	98.01%	85.89%	16466s
	GD	88.26%	0.7734	97.93%	85.37%	16652s
	SA	84.05%	0.7343	93.21%	80.82%	13470s
	SO	87.70%	0.7188	94.68%	85.22%	3092s
	SPE-SO	87.34%	0.7188	94.76%	85.50%	3390s

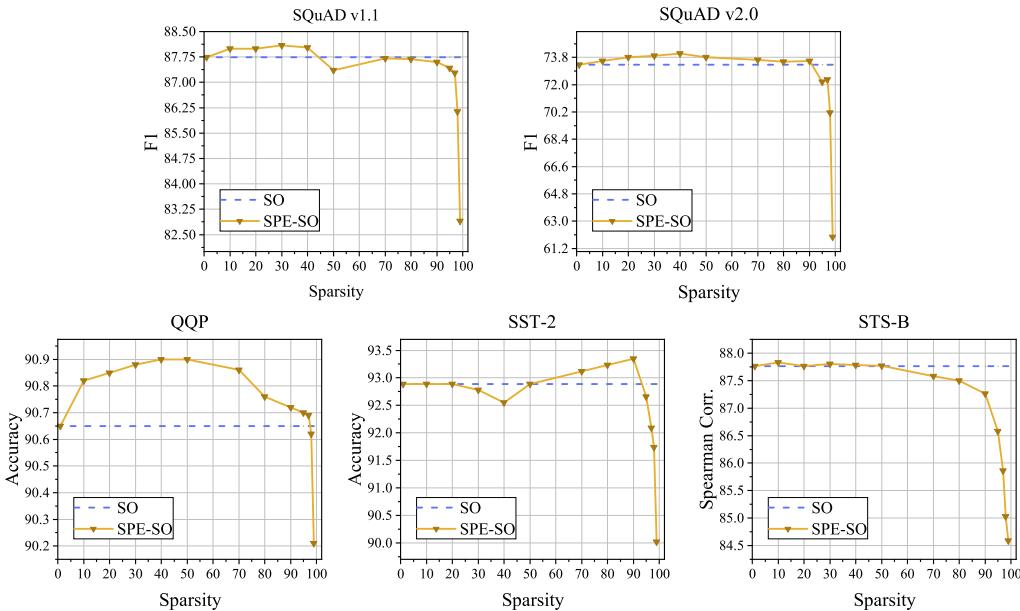


Figure 5: Accuracy of SO and SPE-SO applied to BERT-base across varying sparsity on additional datasets.

Preference Optimization (NPO) (Zhang et al. (2024a)) targets maximizing the discrepancy on the forgetting dataset between the original model and unlearned model. Although the objective functions of these methods differ, their optimization approach relies on gradient ascent, which aims to maximize the loss on the forgetting dataset. We express the unlearning objective in the following form:

$$\arg \max_{\theta} \mathcal{L}(\theta; \mathcal{D}_f), \quad (14)$$

864
 865 Table 9: Overall results of unlearning performance are presented using SPE-SO on BERT-base, both
 866 with and without the neuron selection mechanism. For clarity, SPE-SO denotes SPE-SO applied to
 867 structures at 90% sparsity, while SPE-SO(90%) indicates SPE-SO applied to both structures and
 868 parameters, each with 90% sparsity.

Datasets	Method	Efficacy		Fidelity		Efficiency
		Unlearning Accuracy ↓	MIA ↓	Remaining Accuracy ↑	Testing Accuracy ↑	
MNLI	SPE-SO	85.94%	0.7969	94.15%	84.62%	1274s
	SPE-SO(90%)	85.94%	0.7969	94.12%	84.61%	1280s
QQP	SPE-SO	92.19%	0.9062	98.03%	90.72%	926s
	SU(90%)	92.19%	0.8828	97.67%	90.46%	930s
SST-2	SPE-SO	94.53%	0.8984	98.93%	93.35%	103s
	SPE-SO(90%)	94.53%	0.9141	98.92%	93.35%	105s
STS-B	SPE-SO	86.47%	0.632	97.24%	87.26%	10s
	SPE-SO(90%)	86.87%	0.6406	96.36%	86.99%	11s
SQuAD v1.1	SPE-SO	85.74%	0.5781	94.25%	87.60%	809s
	SPE-SO(90%)	86.16%	0.5859	93.98%	87.19%	812s
SQuAD v2.0	SPE-SO	77.40%	0.6563	93.90%	73.57%	1123s
	SPE-SO(90%)	77.40%	0.6563	93.75%	73.73%	1128s

881
 882 where \mathcal{D}_f is the forgetting dataset. We observe that this objective is similar to minimizing the loss
 883 on the remaining dataset and can also identify the influence-critical parameters using a comparable
 884 approach. First, we introduce a learnable pair of masks for heads and filters:

$$m^* = \arg \max_m \mathcal{L}(m; \theta^*, \mathcal{D}_f) \quad \text{s.t. } \frac{\sum_{i=1}^{|m|} m_i}{|m|} < 1 - S, \quad (15)$$

885
 886 where $|m|$ is the number of mask variables, θ^* represents the original model, and S denotes the
 887 sparsity. We then approximate it using the second-order Taylor series around the mask variables $\mathbb{1}$:

$$\mathcal{L}(m; \theta^*, \mathcal{D}_f) \approx \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_f) - (\mathbb{1} - m) \nabla_m \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_f) + \frac{1}{2} (\mathbb{1} - m)^T \nabla_m^2 \mathcal{L}(\mathbb{1}; \theta^*, \mathcal{D}_f) (\mathbb{1} - m). \quad (16)$$

891
 892 We then use the diagonal FIM to approximate the Hessian matrix and omit constant terms, resulting
 893 in a simplified optimization objective:

$$m^* \approx \arg \max_m (\mathbb{1} - m) \sum_{x \in \mathcal{D}_r} \nabla_m \ell(\mathbb{1}; \theta^*, x) + \frac{1}{2} (\mathbb{1} - m)^2 \widehat{\mathcal{I}}(\mathbb{1}; \theta^*, \mathcal{D}_f). \quad (17)$$

894 Since the mask can only take values of 0 or 1, we can derive the importance evaluation function:

$$m^* \approx \arg \max_m \sum_i \left[(1 - m_i) \left[\sum_{x \in \mathcal{D}_r} \nabla_m \ell(\mathbb{1}; \theta^*, x) \right]_i + \frac{1}{2} (1 - m_i)^2 [\widehat{\mathcal{I}}(\mathbb{1}; \theta^*, \mathcal{D}_f)]_i \right]. \quad (18)$$

903 After obtaining the initial mask, we further optimize the objective using the block diagonal FIM to
 904 rearrange mask:

$$m_l^* \approx \arg \max_{m_l} (\mathbb{1} - m_l) \left[\sum_{x \in \mathcal{D}_r} \nabla_m \ell(\mathbb{1}; \theta^*, x) \right]_l + \frac{1}{2} (\mathbb{1} - m_l)^2 [\widehat{\mathcal{I}}(\mathbb{1}; \theta^*, \mathcal{D}_f)]_l. \quad (19)$$

905
 906 where l represents the layer being optimized. Equipped with the identified key structures, we facilitate
 907 four LLM unlearning methods.

911 B.1 EXPERIMENTS

912
 913 We evaluate unlearning methods on the Task of Fictitious Unlearning (TOFU) dataset (Maini et al.
 914 (2024)) using LLama2-7b-chat model (Touvron et al. (2023)). The unlearning scenarios of TOFU
 915 can be divided into three types: Forget01, Forget05, and Forget10, which represent forgetting dataset
 916 proportions of 1%, 5%, and 10% of the total dataset, respectively. The baseline includes seven
 917 methods: Retraining (RT), Fine-tuning (FT), Sparsity-Aware Unlearning (SA), GA, GD, DPO and
 NPO. We apply three structure-aware parameter-efficient unlearning methods into GA, GD, DPO

918 and NPO for comparison. These methods include: 1) maximize the loss on the forgotten dataset
 919 (MLF) as the unlearning objective, 2) minimize the loss of the remaining dataset (MLR) as the
 920 forgetting objective (i.e. the original method in Section 3.1), and 3) use the norm of the gradients
 921 associated with the structure to evaluate its importance (NORM).
 922

923 **Experimental details.** We use AdamW with a weight decay of 0.01 and a learning rate of 10^{-5} in
 924 RT, FT, SA, GA, GD, DPO and NPO. Besides, we set the learning rate for structure-aware parameter-
 925 efficient methods to $2 \cdot 10^{-4}$ or $3 \cdot 10^{-4}$. In addition, the sparsity of structure-aware parameter-
 926 efficient methods is 90%. All the experiments run for 5 epochs. We also use three main aspects
 927 (i.e., efficacy, fidelity and efficiency) to evaluate the unlearning performance. We use Rouge scores,
 928 normalized probabilities, and the True Ratios on the forgotten dataset to measure efficacy, and those
 929 metrics on the real authors, world facts, and remaining dataset to measure fidelity. We still use
 930 unlearning time to evaluate efficiency. Note that smaller values do not necessarily indicate better
 931 forgetting performance. The goal for unlearning is to closely match that achieved through retraining.
 932

933 **Results.** We find that performing FT and SA only on the remaining dataset does not meet the
 934 unlearning requirements. Although the original GA, GD, and DPO methods can achieve unlearning,
 935 they all exhibit severe catastrophic forgetting on the Forget10 dataset. In contrast, NPO is the most
 936 efficient among these methods. Furthermore, our experiments indicate that sparse updates are better
 937 suited for unlearning than full updates, as they offer a stronger guarantee of unlearning while more
 938 effectively preserving performance, even on the Forget10 dataset. Additionally, the NORM-based
 939 method significantly reduces computation time, but it is less effective than the MLF-based and MLR-
 940 based methods. We observe that the MLR-based method offers a robust balanced trade-off among
 941 unlearning efficacy, model fidelity, and computational efficiency.
 942

943 Table 10: Overall results of unlearning performance using different unlearning methods under
 944 LLama2-7b-chat on TOFU Forget01. ‘Prob.’ indicates the normalized probabilities, ‘TR’ repre-
 945 sents the True Ratios. Forget quality (FQ) and Model Utility (MU) are also used to evaluate the
 946 efficacy and fidelity respectively.
 947

Method	Efficacy			Fidelity												Efficiency
	Forgetting Dataset			Real Authors			World Facts			Remaining Dataset			MU↑	Time↓		
	Rouge	Prob.	TR	Rouge↑	Prob.↑	TR↑	Rouge↑	Prob.↑	TR↑	Rouge↑	Prob.↑	TR↑				
RT	0.39	0.18	0.69	1.0	0.93	0.45	0.58	0.88	0.41	0.54	0.99	0.99	0.47	0.62	-	
FT	0.96	0.99	0.53	5.04e-4	0.94	0.45	0.58	0.87	0.42	0.55	0.97	0.99	0.48	0.62	95.19s	
SA	0.95	0.99	0.53	1.88e-4	0.93	0.45	0.58	0.87	0.42	0.56	0.98	0.99	0.48	0.62	94.88s	
GA	0.49	0.23	0.54	1.27e-3	0.92	0.42	0.55	0.89	0.41	0.54	0.92	0.95	0.49	0.60	96.13s	
MLF-GA	0.64	0.83	0.54	1.27e-3	0.93	0.45	0.58	0.88	0.43	0.57	0.97	0.98	0.48	0.63	149.46s	
MLR-GA	0.43	0.55	0.56	1.27e-3	0.93	0.45	0.58	0.89	0.44	0.57	0.93	0.96	0.48	0.63	147.53s	
NORM-GA	0.57	0.79	0.54	1.27e-3	0.93	0.45	0.58	0.88	0.43	0.57	0.96	0.98	0.48	0.63	77.07s	
GD	0.55	0.53	0.53	1.27e-3	0.94	0.44	0.57	0.86	0.42	0.55	0.96	0.98	0.48	0.61	220.57s	
MLF-GD	0.64	0.83	0.53	1.27e-3	0.94	0.45	0.59	0.88	0.43	0.56	0.96	0.98	0.48	0.63	174.66s	
MLR-GD	0.48	0.61	0.54	1.27e-3	0.94	0.45	0.57	0.88	0.43	0.56	0.94	0.98	0.48	0.62	172.73s	
NORM-GD	0.64	0.83	0.53	1.27e-3	0.94	0.45	0.59	0.88	0.43	0.56	0.96	0.98	0.48	0.63	102.12s	
DPO	0.69	0.92	0.58	5.04e-4	0.93	0.48	0.62	0.88	0.45	0.56	0.94	0.98	0.46	0.64	380.96s	
MLF-DPO	0.69	0.83	0.54	5.04e-4	0.94	0.45	0.58	0.88	0.43	0.56	0.96	0.98	0.48	0.63	169.90s	
MLR-DPO	0.65	0.81	0.54	1.88e-4	0.94	0.45	0.58	0.88	0.43	0.56	0.96	0.98	0.48	0.63	237.73s	
NORM-DPO	0.69	0.83	0.54	5.04e-4	0.94	0.45	0.58	0.87	0.43	0.56	0.96	0.98	0.48	0.63	167.12s	
NPO	0.52	0.27	0.55	3.02e-3	0.92	0.42	0.55	0.87	0.41	0.54	0.94	0.95	0.49	0.61	253.88s	
MLF-NPO	0.59	0.68	0.54	1.27e-3	0.94	0.45	0.58	0.89	0.44	0.57	0.96	0.98	0.48	0.63	174.66s	
MLR-NPO	0.55	0.75	0.54	1.27e-3	0.93	0.45	0.58	0.89	0.44	0.56	0.95	0.98	0.48	0.63	196.31s	
NORM-NPO	0.59	0.78	0.54	1.27e-3	0.93	0.45	0.58	0.89	0.44	0.56	0.95	0.98	0.48	0.63	125.70s	

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976 Table 11: Overall results of unlearning performance using different unlearning methods under
977 LLama2-7b-chat on TOFU Forget05. ‘Prob.’ indicates the normalized probabilities, ‘TR’ repre-
978 sents the True Ratios. Forget quality (FQ) and Model Utility (MU) are also used to evaluate the
979 efficacy and fidelity respectively.

Method	Efficacy			Fidelity												Efficiency	
	Forgetting Dataset			FQ↑	Real Authors			World Facts			Remaining Dataset			MU↑	Time↓		
	Rouge	Prob.	TR		Rouge↑	Prob.↑	TR↑	Rouge↑	Prob.↑	TR↑	Rouge↑	Prob.↑	TR↑				
RT	0.39	0.15	0.67	1.0	0.96	0.42	0.55	0.90	0.40	0.53	0.98	0.99	0.46	0.62	-		
FT	0.92	0.97	0.51	8.33e-16	0.94	0.47	0.61	0.89	0.44	0.57	0.93	0.96	0.48	0.63	404.03s		
SA	0.97	0.99	0.51	3.43e-16	0.94	0.45	0.58	0.87	0.42	0.55	0.98	0.99	0.48	0.62	404.03s		
GA	0.10	3.62e-3	0.65	4.31e-4	0.63	0.35	0.49	0.85	0.40	0.53	0.17	0.02	0.46	0.11	404.27s		
MLF-GA	0.20	3.93e-3	0.62	1.18e-4	0.86	0.41	0.56	0.88	0.41	0.57	0.40	0.31	0.46	0.48	346.03s		
MLR-GA	0.17	2.31e-3	0.61	4.75e-5	0.83	0.42	0.58	0.86	0.42	0.58	0.33	0.17	0.45	0.42	342.01s		
NORM-GA	0.18	1.08e-3	0.56	1.21e-10	0.85	0.40	0.55	0.86	0.40	0.57	0.36	0.22	0.44	0.44	252.45s		
GD	0.30	1.79e-2	0.54	2.83e-4	0.79	0.35	0.49	0.87	0.38	0.53	0.46	0.42	0.50	0.49	1009.22s		
MLF-GD	0.37	0.15	0.61	2.83e-4	0.94	0.44	0.59	0.86	0.43	0.57	0.81	0.91	0.48	0.61	463.18s		
MLR-GD	0.33	2.31e-2	0.64	0.63	0.90	0.46	0.60	0.87	0.43	0.57	0.79	0.88	0.48	0.61	460.68s		
NORM-GD	0.37	0.15	0.61	2.83e-4	0.94	0.44	0.59	0.86	0.43	0.57	0.81	0.91	0.48	0.61	371.12s		
DPO	4.57e-2	0.64	0.62	6.57e-12	0.57	0.46	0.60	0.83	0.46	0.57	0.23	0.73	0.40	0.47	1800.04s		
MLF-DPO	0.30	0.19	0.60	8.06e-7	0.92	0.46	0.59	0.86	0.45	0.58	0.76	0.88	0.48	0.61	782.51s		
MLR-DPO	0.28	9.35e-2	0.62	1.84e-4	0.90	0.46	0.60	0.87	0.45	0.58	0.69	0.88	0.48	0.61	780.65s		
NORM-DPO	0.19	0.14	0.61	4.75e-5	0.88	0.46	0.60	0.82	0.45	0.58	0.66	0.78	0.48	0.60	691.08s		
NPO	0.34	0.11	0.66	1.18e-4	0.94	0.33	0.42	0.89	0.38	0.49	0.42	0.36	0.46	0.46	1183.55s		
MLF-NPO	0.33	0.12	0.61	1.11e-5	0.91	0.44	0.58	0.88	0.42	0.56	0.74	0.83	0.48	0.60	568.75s		
MLR-NPO	0.34	0.15	0.59	8.11e-8	0.90	0.44	0.58	0.87	0.42	0.57	0.76	0.85	0.48	0.60	570.66s		
NORM-NPO	0.32	0.12	0.61	1.11e-5	0.91	0.43	0.58	0.88	0.42	0.56	0.74	0.82	0.48	0.59	481.09s		

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1000 Table 12: Overall results of unlearning performance using different unlearning methods under
1001 LLama2-7b-chat on TOFU Forget10. ‘Prob.’ indicates the normalized probabilities, ‘TR’ repre-
1002 sents the True Ratios. Forget quality (FQ) and Model Utility (MU) are also used to evaluate the
1003 efficacy and fidelity respectively.

Method	Efficacy			FQ↑	Fidelity												Efficiency	
	Forgetting Dataset				Real Authors			World Facts			Remaining Dataset			MU↑	Time↓			
	Rouge	Prob.	TR		Rouge↑	Prob.↑	TR↑	Rouge↑	Prob.↑	TR↑	Rouge↑	Prob.↑	TR↑					
RT	0.41	0.15	0.67	1.0	0.92	0.43	0.57	0.90	0.41	0.54	0.98	0.99	0.47	0.61	-			
FT	0.89	0.96	0.51	2.43e-19	0.94	0.48	0.62	0.89	0.45	0.58	0.89	0.96	0.47	0.64	827.26s			
SA	0.98	0.99	0.50	1.69e-15	0.92	0.44	0.58	0.86	0.41	0.55	0.98	0.99	0.49	0.62	832.32s			
GA	1.19e-3	6.26e-33	0.79	5.40e-18	0.0	0.25	0.21	0	0.25	0.20	0.01	1.57e-32	0.12	0.0	822.49s			
MLF-GA	0.14	3.25e-4	0.56	2.06e-13	0.72	0.47	0.67	0.73	0.46	0.61	0.21	2.51e-2	0.40	0.16	590.08s			
MLR-GA	0.22	2.79e-2	0.62	0.34	0.84	0.49	0.66	0.87	0.46	0.59	0.36	0.33	0.46	0.51	588.77s			
NORM-GA	0.15	3.02e-4	0.54	1.45e-14	0.68	0.47	0.67	0.73	0.46	0.60	0.20	3.31e-2	0.41	0.19	464.81s			
GD	1.31e-2	3.01e-18	0.70	1.07e-13	0.49	0.46	0.63	0.82	0.44	0.58	0.25	0.24	0.48	0.42	2042.64s			
MLF-GD	0.31	1.85e-2	0.59	7.31e-3	0.89	0.48	0.61	0.87	0.46	0.59	0.61	0.75	0.47	0.60	847.88s			
MLR-GD	0.31	1.85e-2	0.59	7.31e-3	0.89	0.48	0.61	0.87	0.46	0.59	0.61	0.75	0.49	0.61	843.89s			
NORM-GD	0.30	1.53e-2	0.59	3.11e-3	0.89	0.51	0.67	0.85	0.46	0.61	0.53	0.69	0.48	0.60	719.93s			
DPO	1.05e-2	0.51	0.66	1.49e-9	5.33e-3	0.43	0.57	0.17	0.43	0.53	1.17e-2	0.57	0.37	3.08e-2	3420.63s			
MLF-DPO	0.37	0.27	0.59	2.55e-9	0.89	0.46	0.60	0.85	0.44	0.57	0.79	0.93	0.49	0.62	1511.82s			
MLR-DPO	0.25	0.21	0.61	3.63e-7	0.81	0.46	0.60	0.79	0.44	0.58	0.69	0.92	0.49	0.60	1509.28s			
NORM-DPO	0.17	0.19	0.61	1.40e-6	0.77	0.46	0.60	0.66	0.45	0.57	0.58	0.89	0.48	0.58	1385.32s			
NPO	0.27	0.11	0.72	3.36e-2	0.72	0.46	0.62	0.86	0.45	0.59	0.35	0.29	0.36	0.47	2407.90s			
MLF-NPO	0.33	7.25e-2	0.62	6.54e-4	0.94	0.46	0.62	0.89	0.45	0.59	0.63	0.73	0.47	0.60	1063.28s			
MLR-NPO	0.33	0.12	0.61	3.63e-7	0.91	0.47	0.62	0.86	0.45	0.59	0.61	0.72	0.47	0.60	1060.32s			
NORM-NPO	0.32	4.10e-2	0.63	9.96e-3	0.90	0.47	0.63	0.86	0.45	0.59	0.60	0.67	0.47	0.59	932.37s			

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