

APPROXIMATE CLUSTERING FOR EXTRACTING TASK RELATIONSHIPS IN MULTI-INSTRUCTION TUNING

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ABSTRACT

The development of language models involves the evaluation of a broad range of learning tasks. Recent work has shown that by using carefully designed instructions to teach a large transformer model, they can be fine-tuned on a wide range of downstream tasks. However, when the number of instructions increases, they can negatively interfere with each other if trained together. Existing works have relied on domain expertise and manual inspection to construct multi-instruction sets, which can be time-consuming and difficult to scale. To address this challenge, this paper develops a clustering algorithm to find groups of similar tasks based on a given set of task affinity scores. This is an NP-hard problem, and conventional algorithms such as spectral and Lloyd’s clustering are sensitive to variations in the scale of task losses. Our algorithm instead uses a semidefinite relaxation to maximize the average density of clusters and then rounds the solution with a threshold. We adaptively build the clusters by gradually adding tasks so that the affinities only need to be computed in the existing clusters. Then, we construct an evaluation benchmark to assess task grouping algorithms with verified group structures. The evaluation set includes 63 cases, spanning multitask instruction tuning, multi-instruction tuning, and in-context learning of multiple functions. We validate our algorithm on this evaluation set by showing that it recovers the group structure found by an exhaustive search. We also show that our approach improves performance over multi-instruction and soft-prompt tuning by up to 6% on several sentence classification and structure-to-text generative tasks.

1 INTRODUCTION

A hallmark of the recent development in language models is that they can simultaneously make predictions over a broad range of learning tasks (Roberts et al., 2019; Liang et al., 2022). The adaptation of these language models to downstream tasks is then enhanced via instruction fine-tuning (Mishra et al., 2022). Prior work has shown that fine-tuning an existing model such as T5 through multiple instructions can lead to state-of-the-art results on a diverse collection of NLP tasks (Sanh et al., 2022; Wei et al., 2022). In light of these developments, the design of instruction tuning datasets and evaluations has received much interest recently (Longpre et al., 2023). By contrast, the algorithmic problem of how to best use these instructions for fine-tuning downstream tasks remains under-explored. It is also worth noting that these sets typically involve a large number of tasks and instructions, which can lead to severe negative interference when they are trained together naively (Jang et al., 2023). Existing work on multi-instruction tuning relies on extensive domain expertise and manual selection (Chung et al., 2022). In this paper, we revisit a task grouping problem (Standley et al., 2020), which is highly relevant to a number of settings in language model fine-tuning: Given n tasks, we are interested in partitioning them into k groups so that each group of tasks can be best trained together (separately from the other groups).

A naive approach to selecting which tasks to train together in a language model is according to the category of each task. Because the datasets are collected from different sources (Wang et al., 2018; Aribandi et al., 2022), even two tasks of the same category, such as sentiment analysis, may not transfer positively to each other. Task grouping methods have been developed for jointly learning multiple datasets. For instance, Fifty et al. (2021) first compute a set of (pairwise) affinity measures and then apply optimization techniques such as branch-and-bound to find the best task combinations. The computational cost of these techniques can still be quite high for the scale of instruction fine-

tuning sets. Another natural solution is to use clustering algorithms such as spectral clustering (Ng et al., 2001) and Lloyd’s algorithm (Lloyd, 1982). We find that these methods are particularly sensitive to the scale of the varied losses across a large set of different tasks.

To address the challenge, we develop a new clustering algorithm, which involves two key steps. The first step is a semidefinite relaxation for maximizing the average density of the k groups, given an n by n task affinity matrix T . This matrix requires measuring n^2 affinity scores, which can be slow to compute when n is large. Therefore, the second step of our algorithm is an adaptive procedure, where we build the clusters gradually. This allows us to accelerate the computation of task affinities by leveraging the existing separations in the clusters. Moreover, we introduce an adaptive sampling technique to account for higher-order task relationships.

To facilitate the evaluation of task grouping methods, we curate an evaluation benchmark that contains task group structures with verified positive transfer within groups. This benchmark includes 63 evaluation cases that span three types of scenarios, including multitask (instruction) fine-tuning (over 19 NLP tasks) (Wang et al., 2018, 2019; Sanh et al., 2022), multi-instruction fine-tuning (over 100 instructions) (Bach et al., 2022; Zhou et al., 2023), and in-context learning with three function classes (Garg et al., 2022). See Tab. 1 for a summary. Based on this benchmark, we evaluate our approach by showing that the above algorithm can correctly identify the underlying groups, succeeding in all evaluation cases. Notably, the groups match the results found by exhaustive search. We also show that our approach outperforms multi-instruction and prefix tuning by **3.3%** on three sentence classification tasks from SuperGLUE (Wang et al., 2019) and two structure-to-text generative tasks from the GEM benchmark (Gehrmann et al., 2021).

In summary, in this paper, we revisit the task grouping problem for language model fine-tuning and design a new clustering algorithm that is both efficient and robust to cross-task heterogeneity. We construct an evaluation benchmark for task grouping approaches along with an easy-to-use package, spanning three scenarios of instruction fine-tuning, which can also be used for future work. Experiments show that our algorithm can correctly identify the underlying group structures and also be used to identify groups of similar instructions in multi-instruction tuning. The rest of this paper is organized as follows. Sec. 2 reviews related work. Sec. 3 provides more background information. Sec. 4 describes our algorithm. Sec. 5 presents the experiments. Additional related work and experiment details are provided in the Appendix.

2 RELATED WORK

Previous works of FLAN (Wei et al., 2022), NaturalInstructions (Mishra et al., 2022), and T0 (Sanh et al., 2022) have demonstrated that fine-tuning language models on multiple downstream tasks prompted with instructions resulting in enhanced generalization to previously unseen tasks. Moreover, there have been efforts to advance the instruction fine-tuning method, such as expanding task datasets (Chung et al., 2022; Wang et al., 2022b) and refining instruction sets (Bach et al., 2022). Furthermore, Muennighoff et al. (2023) constructed multilingual instruction fine-tuning datasets and performed instruction fine-tuning on multilingual pretrained language models to boost the generalization on unseen language and tasks. Longpre et al. (2023) study the design decisions of publicly available instruction tuning methods and find that training with mixed instructions settings yields improved performance. Wang et al. (2023b) propose multitask prompt tuning, which first learns a single transferable prompt by distilling knowledge from multiple task-specific source prompts and then learns multiplicative low-rank updates to this shared prompt to adapt it to each downstream target task efficiently. Compared to these studies, we study the algorithmic problem of how to find task structures in instruction fine-tuning sets.

There is also a growing line of work on designing the best prompt to adapt pre-trained language models to downstream tasks (Shin et al., 2020; Gao et al., 2021; Zhang et al., 2022). Prefix tuning (Li and Liang, 2021) inserts continuous prompt embeddings to each layer in language models and optimizes the embeddings during fine-tuning. Prompt tuning (Lester et al., 2021) proposes to add prompt embeddings only in the inputs. PromptBoosting (Hou et al., 2023) constructs a large pool of weak learners by pairing prompts with different elements of the LM’s output distribution and then ensemble the weak learners using the AdaBoost algorithm. Instead of finding the best instruction for a downstream task, our work focuses on optimizing the average performance of a model under multiple instructions.

Table 1: We construct 63 evaluation cases with verified group structures to assess task grouping algorithms. Below is a table of summary of the category of tasks and datasets included in each evaluation set.

Multitask fine-tuning	57 evaluation cases with 2-6 (varied) groups				
	Sentiment Classification (3)	Natural Language Inference (3)	Multiple-Choice QA (4)	Open-Domain QA (3)	Coreference Resolution (3) Summarization (3)
Multi-instruction fine-tuning	5 evaluation cases with 10 groups in each				
	RTE (100)	WiC (100)	BoolQ (100)	E2E NLG (100)	Web NLG (100)
In-context learning	1 evaluation case with 3 groups				
	Linear Regression (3)	Decision Trees (3)	Neural Networks (3)		

Clustering is a fundamental aspect of machine learning. Besides SDP relaxations, linear programming relaxations are known for clustering objectives such as k -center (Guttmann-Beck and Hassin, 2000). Their approach requires pre-selecting k anchor points as the centers of each cluster. However, their approach then enumerates all k -sized subsets and thus runs in time $O(n^k)$. For the case of geometric clustering in Euclidean spaces, polynomial time approximation schemes can be achieved (De La Vega et al., 2003). Bartal et al. (2001) give a polynomial time approximation algorithm for min-sum k -clustering for arbitrary metric spaces via dynamic programming. The integrality gap of linear programming and semidefinite programming relaxations can be analyzed when there is a separation structure in the underlying clusters (Awasthi et al., 2015). These approximation guarantees typically require the underlying similarity scores to satisfy a metric condition. By contrast, the task affinity matrix, in our case, can easily violate the triangle inequality. Lastly, recent work has also looked into mixed integer programming for best subset selection (Bertsimas et al., 2016). One novel contribution of this work is to make explicit a connection between multitask/multi-instruction fine-tuning and clustering. In light of this connection, it would also be interesting to revisit hierarchical clustering (Charikar and Chatziafratis, 2017; Chami et al., 2020) and hypergraph clustering (Yin et al., 2017; Veldt, 2023) for task grouping.

3 PRELIMINARIES

Many problems in the context of language model fine-tuning are related to multitask learning (Wang et al., 2018; Aribandi et al., 2022; Sanh et al., 2022). We give three examples, which will be the focus of this paper: (1) Multitask instruction fine-tuning is an essential component of adapting language models, enabling the models with various language processing abilities, such as question answering and text summarization. (2) Multi-instruction fine-tuning involves a mix of instructions to further enhance a language model’s ability to respond to diverse instructions from users. (3) In-context learning refers to the ability of a language model to learn a function class with a few “in-context” examples; A natural question is whether these different function classes are in-context learnable simultaneously.

Task Grouping Setup. The above examples can be formulated abstractly in a multitask learning setting. Let there be n downstream tasks. The goal of task grouping (cf. Standley et al. (2020)) is to partition the n tasks into k subsets such that each subset of tasks is the best to be trained together.

For each pair of tasks u and v , let $T_{u,v}$ denote an affinity score, which quantifies the transfer effect between them. Pairwise notions of affinity scores between two tasks have been used in prior work (Fifty et al., 2021). For example, one way to quantify $T_{u,v}$ is via the performance of task u ’s validation performance evaluated on a model fine-tuned on a model trained with both u and v . Given an n by n task affinity matrix T , the extent of positive transfers within a subset of tasks S can be characterized by the density of affinity scores in the subset:

$$d_S = \sum_{u,v \in S} \frac{T_{u,v}}{|S|}. \quad (1)$$

Then, one can view task grouping as a clustering problem whose objective is to maximize the average density of all clusters. Let C_1, \dots, C_k denote a partition of the n tasks. Let v_1, \dots, v_k be a 0-1

where e is the all-ones vector. We omit the $\frac{1}{k}$ factor in the objective for simplicity.

This integer program is computationally challenging to solve, even for small values of k . To address this issue, we will relax the above integer program to a (constrained) semidefinite program (SDP), which can be solved in polynomial time. First, we note that $v_i v_i^\top$ is a rank one semidefinite variable. Let us denote the sum of them (normalized by $v_i^\top v_i$) as the following new variable

$$X = \sum_{j=1}^k \frac{v_j v_j^\top}{v_j^\top v_j}. \quad (4)$$

This matrix X has a rank equal to k because it is the sum of k rank-1 matrices, and the v_i 's are orthogonal to each other. Additionally, its trace is equal to k because $\frac{v_j v_j^\top}{v_j^\top v_j}$ has a trace of one for any j . Second, the entries of every row of X is equal to one:

$$Xe = \sum_{i=1}^k \frac{v_i (v_i^\top e)}{v_i^\top v_i} = \sum_{i=1}^k v_i = e.$$

Removing the 0-1 integer constraint, we relax Problem (3) into a rank-constrained problem:

$$\max \left\{ \langle T, X \rangle : Xe = e, \text{rank}(X) = k, \text{Tr}[X] = k, X \geq 0, X \succeq 0, X \in \mathbb{R}^{n \times n} \right\}.$$

The above program involves a rank constraint, which is still computationally challenging to solve. However, it can be further relaxed by removing the rank constraint while keeping the trace constraint:

$$\max \left\{ \langle T, X \rangle : Xe = e, \text{Tr}[X] = k, X \geq 0, X \succeq 0, X \in \mathbb{R}^{n \times n} \right\}. \quad (5)$$

The above problem can be solved efficiently using convex optimization methods. Given a solution of X , the last step is to round it into an integer solution. We set a threshold λ such that if $X_{u,v} \geq \lambda$, tasks u and v are assigned to the same cluster. In practice, we set the λ as c/n for some constant $c \geq 1$, since $X_{u,v}$ should be $\frac{1}{|C_i|}$ when they are in the same cluster C_i . In summary, we can derive an efficient clustering algorithm given a task affinity matrix; See Procedure 1 below.

Procedure 1 Approximate Task Clustering through SDP Relaxations

Input: Task affinity matrix $T \in \mathbb{R}^{n \times n}$

Require: Number of clusters k ; A threshold λ for rounding

Output: A list of clusters \mathcal{C}

- 1: Obtain X by solving problem (5)
 - 2: Generate a list of clusters \mathcal{C} by assigning u and v into a cluster if $X_{u,v} > \lambda$
-

Illustrative Example. A naive way to maximize the clustering objective is using algorithms such as spectral clustering or Lloyd’s algorithm on the task affinity matrix T . Curiously, we observe that these algorithms are not robust in multitask learning as the scale of different tasks’ losses varies dramatically. In Fig. 2, we illustrate the clustering results with these methods. We use a planted model by generating a random matrix including one low-density cluster and two high-density clusters.

- In spectral clustering, the eigenvector values remain constant on the high-density clusters with the presence of the low-density cluster.
- Lloyd’s algorithm iteratively selects the cluster centroids and updates assignments to each cluster. With higher values in high-density clusters, the centroids are assigned to them, and the algorithm does not separate the low-density cluster.

4.2 ADAPTIVELY ESTIMATING TASK AFFINITIES AND BUILDING TASK CLUSTERS

Next, we design an algorithm to speed up the clustering process. The above clustering algorithm requires access to the pairwise task affinity matrix. For n tasks, computing the pairwise affinity scores between every pair of tasks is time-consuming, as it requires training n^2 models. Furthermore, it ignores higher-order task relationships beyond the combinations of two tasks, which adds more

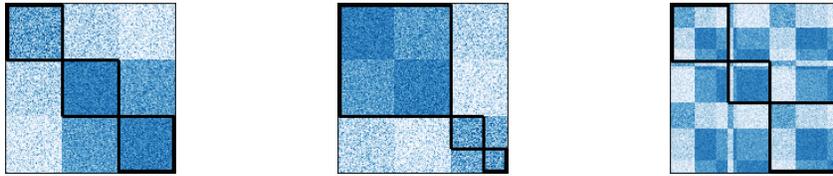


Figure 2: We illustrate the SDP relaxation compared to spectral clustering and Lloyd’s algorithm for recovering three hidden clusters, as shown by the densities in the figure. Spectral clustering groups the two higher-density clusters together, while Lloyd’s algorithm mixes the three clusters. By contrast, the SDP relaxation manages to identify all three hidden clusters.

complexity to multitask relationships. We first introduce a task affinity score that captures higher-order task relationships over task subsets. Then, we design an efficient sampling procedure by iteratively computing part of the affinity matrix and growing the clusters adaptively.

Higher-Order Task Affinity. We consider a higher-order task affinity score estimated from subsets of more than two tasks. First, sample m subsets of tasks $\{1, 2, \dots, n\}$ uniformly over subsets of size α , denoted as S_1, S_2, \dots, S_m . Then, compute the value of $f_i(S_j)$ by fine-tuning a model on tasks in every subset S_j for $j = 1, \dots, m$. Lastly, compute $T_{i,j}$ as the average multitask performance over all subsets that include task i and j :

$$T_{i,j} = \frac{1}{n_{i,j}} \sum_{1 \leq k \leq n: \{i,j\} \subseteq S_k} f_i(S_k), \text{ for all } 1 \leq i, j \leq n, \quad (6)$$

where $n_{i,j}$ be the number of subsets that include both i, j . This sampling is analogous to the sampling of features in random forests (due to space limit, a detailed justification is stated in App. B.3).

Adaptive Sampling. The next step is an adaptive sampling procedure to accelerate the above estimation. The idea is to divide tasks into small batches and iteratively estimate affinity scores for a new batch of tasks. In each iteration, we have existing cluster structures and a new batch of unclustered tasks. We pick one cluster to estimate task affinity scores between the chosen cluster and the new batch of tasks. This uses the existing separations, as described in Procedure 2.

Procedure 2 Adaptive Estimation of Task Affinity Scores

Input: n tasks, training and validation sets of each task, cluster structure \mathcal{C}_0 for the first n_0 tasks

Require: Number of subsets m ; Size of each subset α ; Multitask learning algorithm f

Output: An n by n task affinity matrix T

- 1: **for** $i = 1, 2, \dots, m$ **do**
 - 2: Randomly choose a group C from cluster \mathcal{C}_0
 - 3: Sample a random subset S_i from $\{n_0 + 1, n_0 + 2, \dots, n\} \cup C$ with size α
 - 4: Evaluate multitask performance $f(S_i)$ for every task in S_i
 - 5: **end for**
 - 6: Calculate the affinity score matrix via Eq. (6)
-

After estimating the affinity scores for the new batch of tasks, we update the clusters by solving the relaxed SDP in Eq. (5). We initialize the task assignment variable X by assigning $X_{u,v}$ as $\frac{1}{|C|}$ if u and v are in a cluster C with size $|C|$. Then, we solve the SDP again to re-generate the clusters. At iteration t , we search the number of clusters within a range of $|C^{(t)}|$ to k and choose the one that maximizes the objective $\langle T, X \rangle$. The complete procedure is described in Algorithm 3.

Runtime. We examine the runtime of our algorithm. There are s iterations. During each iteration:

- We estimate task affinity scores for b tasks. We train m models on sampled subsets to compute the scores. In practice, we notice that collecting $m = 5b = \frac{5n}{s}$ subsets suffices for estimating the affinity scores until convergence. For $n = 100$ tasks, we take $s = 10$ steps. Each step trains 50 models on sampled subsets and takes 23 hours using a single GPU.
- We solve a convex program on an affinity matrix of size n by n . In practice, this step typically runs quickly in our experiments, taking less than 1.5 seconds for n up to 100.

Algorithm 3 Adaptive Task Grouping (AdaGroup)**Input:** n tasks, training and validation datasets of each task**Require:** # final clusters k ; # adaptive steps s ; # sampled subsets in each step m ; Size of subset α **Output:** k groups of tasks

- 1: Initialize the clusters as $\mathcal{C}^{(0)} = \{\}$. Let $b = \frac{n}{s}$ be the number of additional tasks in each step
- 2: **for** $t = 0, 1, \dots, s - 1$ **do**
- 3: Choose b tasks from the remaining tasks
- 4: Estimate the task affinity matrix $T^{(t+1)}$ by Procedure (2) with current cluster structure $\mathcal{C}^{(t)}$
- 5: Generate clusters $\mathcal{C}^{(t+1)}$ following Procedure (1)
- 6: **end for**
- 7: **return** $\mathcal{C}^{(s)}$

5 EXPERIMENTS

We describe experiments to apply our algorithm to three problems relevant to language model fine-tuning, including multitask fine-tuning, multi-instruction tuning, and in-context learning. We will discuss the evaluation datasets used in the experiments. Then, we describe the setup along with the comparative results. Lastly, we give ablation studies to justify our algorithm design and end this section with a discussion for future work.

5.1 EVALUATION OF TASK GROUPING

The evaluation of task grouping algorithms requires a clear specification of task grouping structures. A naive way to conduct evaluations is using existing multitask learning benchmarks such as GLUE (Wang et al., 2018) and SuperGLUE (Wang et al., 2019). These benchmarks come with pre-defined groups. Curiously, we noticed that nearly 40% of pairwise transfers are negative even within these groups, as shown in Fig. 1. With this context in mind, the first aim of our experiments is to collect and then construct an evaluation benchmark that is more suitable for assessing task grouping algorithms. Ideally, such an evaluation set should have clearly defined group structures.

Multitask Instruction Fine-Tuning. We collect a list of NLP datasets under different (human-labeled) categories, such as sentiment analysis, question answering, summarization, etc. Then, we measure the pairwise transfers between each pair of tasks from the same category. We use T5-Base as the base model (Raffel et al., 2023).

After getting all the pairwise effects for each category, we select the subsets whose ratio of positive effects is higher than 90%. This leads to an evaluation set of six groups of six task categories. These include sentiment analysis, natural language inference, multiple-choice QA, open-domain QA, coreference solution, and summarization tasks. Each category contains three or four tasks, leading to 19 tasks in total. We display a complete list in Table 3 (App. B).

Multi-Instruction Tuning. We consider three datasets from SuperGLUE, including RTE, WiC, and BoolQ, and two structure-to-text generation datasets from the GEM benchmark (Gehrmann et al., 2021), including E2E NLG challenge and Web NLG. Each dataset contains 100 instructions, including ten instructions from Bach et al. (2022) and 90 instructions that we generate with an automatic instruction generation method from Zhou et al. (2023).

In-Context Learning. We are interested in three types of functions, including linear regression (LR), decision trees (DT), and two-layer ReLU neural networks (NN). For each type, we define three function classes with different distributions. For example, for linear regression, we specify a Gaussian distribution over their weight parameters for one function class. We regard each function class as one task, leading to a total of nine tasks. For each task, we generate training prompts containing d in-context examples, denoted as $(x_1, \phi(x_1), x_2, \phi(x_2), \dots, x_d, \phi(x_d))$ where ϕ is a random function sampled from the function distribution.

5.2 IMPLEMENTATION AND BASELINES

For multitask instruction fine-tuning, we create evaluation cases and verify the group structure inside each case. Altogether, we have 15 cases with two groups, 20 cases with three groups, 15 cases with

Table 2: Test performance averaged over all instructions on three sentence classification tasks from SuperGLUE and two structure-to-text generative tasks from GEM. We compare our approach with multi-instruction tuning, prefix tuning, and prompt tuning. we report the average results over three random seeds.

Dataset Task Type (Metric)	RTE Classification tasks (Accuracy)	WiC Classification tasks (Accuracy)	BoolQ Classification tasks (Accuracy)	E2E NLG Generative tasks (ROGUE-1)	Web NLG Generative tasks (ROGUE-1)
Multi-Instruction Tuning	75.09±0.68	66.44±0.98	78.16±0.77	71.46±0.27	80.80±0.19
Prefix Tuning	72.74±2.40	62.29±2.93	76.19±0.98	70.23±0.40	78.69±0.26
Prompt Tuning	73.12±1.26	62.88± 2.19	75.51±0.85	70.72±0.81	77.42±0.31
Our Approach	80.96±0.85	69.89±0.87	81.76±0.62	73.03±0.67	82.95±0.75

four groups, 6 cases with five groups, and 1 case with six groups. To verify that the group structure is correct, we use an exhaustive search to enumerate all task combinations that optimize the clustering objective (cf. Eq. (3)) and make sure that the group structure indeed achieves the optimum for the clustering objective.

For multi-instruction fine-tuning, we use T5-Base as the base model. For classification tasks, we report the accuracy as the performance. For generative tasks, we report the Rouge1 score as the performance. For each dataset, we evaluate the average performance over all 100 instructions. In our approach, we view one instruction as one task. We apply our approach to find groups of instructions and then fine-tune one model for each group of instructions. Our approach requires three hyper-parameters: the number of adaptive steps, the number of subsets in each step, and the size of subsets. We select the size between 3, 5, and 10. We select adaptive steps between 10, 5, and 3. We then set the number of subsets as five times the number of new tasks in each step. We select the number of clusters from a range between 10, 15, and 20.

We compare our approach with multi-instruction and report the results of two soft-instruction tuning baselines in terms of relative improvement, including Prefix Tuning (Li and Liang, 2021) and Prompt Tuning (Lester et al., 2021). We use LoRA fine-tuning (Hu et al., 2022) for our approach and multi-instruction to match the same amount of training parameters as soft-instruction tuning. The information for the training details is included in Appendix B.

For grouping function classes during in-context learning, a transformer is trained to predict $\phi(x_i)$ for a given x_i based on the preceding in-context examples. We evaluate the prediction loss as the squared error of predictions averaged over $d = 100$ in-context learning steps and use this loss as the MTL performance $f(S)$. We estimate task affinity scores by sampling subsets of three tasks and compute the MTL performance of a transformer trained on prompts from three tasks jointly.

5.3 EXPERIMENTAL RESULTS

Multitask Instruction Fine-Tuning Results. We evaluate our approach on the 57 evaluation cases ranging from two to six groups of tasks. Our approach correctly identifies the underlying groups under all cases, obtaining the same results as the exhaustive search. In contrast, using spectral and Lloyd’s clustering correctly identifies the group structures in 16/4 out of the 57 cases.

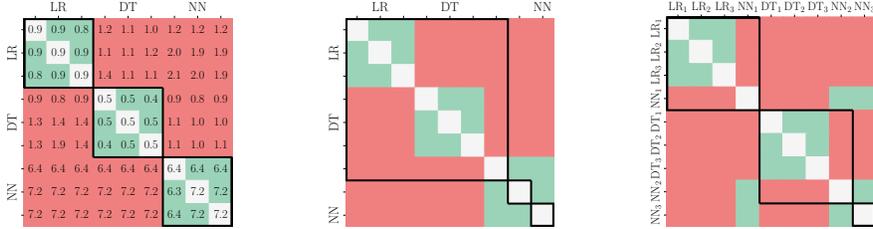
Multi-Instruction Tuning Results. Table 2 shows the results of the average performance on the development set evaluated with 100 instructions. We observe that our approach improves over the baseline methods by 3.3% on average, suggesting the benefit of separating instructions to reduce their negative interference.

In-Context Learning Results. We observe that transformers trained on different function classes perform worse than being trained on a single function class, except for neural networks. We illustrate the task affinity scores between the function classes in Figure 3 (Left). Our approach recovers the cluster structure for three types of function classes. In contrast, spectral clustering and Lloyd’s clustering yield clusters mixed between different function classes, shown in Figure 3 (Right).

5.4 ABLATION STUDIES

We provide two ablation studies of our algorithm, including the clustering step and task affinity. Then, we illustrate an intriguing transfer between function classes during in-context learning.

Figure 3: Clusters of function classes generated by our approach (Left), spectral clustering (Middle), and Lloyd’s clustering (Right). Each entry corresponds to an affinity score using the mean squared loss as the MTL performance (green means a positive transfer, while red means a negative transfer).



Instruction Selection. We compare our clustering algorithm with alternative selection methods, including (greedy) forward selection, spectral clustering, and Lloyd’s clustering. Forward selection starts from an empty group, enumerates all tasks, and adds one task to an existing group that results in the highest increase of densities within groups. In multi-instruction tuning, we find that our algorithm outperforms these alternative selections by 1.2%, averaged over the datasets.

Task Affinity. We compare alternative methods to estimate task affinity scores and validate the benefit of using higher-order task affinity. We compare the higher-order task affinity with two pairwise task affinity scores, including loss-based pairwise affinity (Standley et al., 2020), and gradient-based affinity score (as the ratio of task i ’s loss before and after applying the gradient of task j on the model parameters) (Fifty et al., 2021). We find that using higher-order task affinity improves the performance of grouping instructions by 1.7% over the two pairwise affinity scores on average.

In-Context Transferability. We examine the test MSE of transformers for predicting $f(x_{n+1})$, given n in-context examples $x_1, \phi(x_1), x_2, \phi(x_2), \dots, x_{n+1}$. We first train a transformer only on neural network functions (STL). Then, we combine the training examples with another function class, including linear regression or decision trees, and train a transformer on examples of both function classes (MTL). We compare the error between MTL and STL in Fig. 4. Curiously, we find that a transformer trained to predict neural networks with linear regression or decision trees compares comparably to a transformer trained only on neural networks. On the other hand, if the target class is decision tree or linear regression, then learning it with the other two classes will significantly degrade MSE (see Fig. 6 of App. B.2).

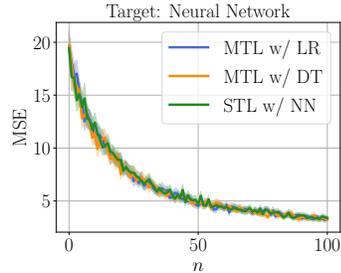


Figure 4: Test MSE of in-context learning neural networks.

5.5 DISCUSSIONS

Our findings provide some evidence that modeling task relationships can also enhance language modeling (particularly instruction tuning). It might be interesting to investigate if this perspective applies in other contexts, such as modeling the relationship between generating different programming languages or algorithmic reasoning. It may also be worth investigating hierarchical relationships: The above in-context learning example shows learning NNs implies the learning of linear regression and decision trees. It is plausible to revisit curriculum learning for tuning instructions with increasing complexities. To facilitate the discussion, we provide an easy-to-use package to make our evaluation sets accessible to researchers. Future work could use these evaluation cases to design and evaluate clustering algorithms for instruction fine-tuning. An anonymized link to our experiment code is at <https://anonymous.4open.science/r/AdaGroup4InstructionTuning>.

6 CONCLUSION

This paper developed an approximate clustering algorithm to extract task group structures so that the most related tasks are trained together. We construct a new evaluation benchmark for this clustering problem, spanning three use cases of language model fine-tuning, with a total of 63 evaluation cases. A package is developed for reusing this evaluation set to facilitate future discussions.

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Table 3: Dataset description and statistics of six groups of text datasets.

Dataset	Benchmark	Train. Set	Dev. Set	Task Category
Sentiment Classification				
SST-2	GLUE	67k	1.8k	Sentiment classification
IMDB Reviews	-	25k	25k	Sentiment classification
Yelp Reviews	-	650k	50k	Sentiment classification
Natural Language Inference				
MNLI	GLUE	393k	20k	NLI
RTE	GLUE	2.5k	3k	NLI
WNLI	GLUE	634	146	NLI
Multiple-Choice QA				
BoolQ	SuperGLUE	9.4k	3.3k	Question answering
COPA	SuperGLUE	400	100	Question answering
MultiRC	SuperGLUE	5.1k	953	Question answering
ReCoRD	SuperGLUE	101k	10k	Question answering
Open-Domain QA				
SocialiQA	-	33.4k	1.95k	Social commonsense QA
WikiQA	-	20.4k	2.2k	Wikipedia-based question answering
HotpotQA	-	90.4k	7.4k	Wikipedia-based question answering
Coreference Resolution				
WSC	SuperGLUE	554	104	Pronoun Coreference
Winogrande	-	9.3k	1.8k	Word Coreference
Quoref	-	19.4k	2.42k	Coreferential reasoning
Summarization				
XSum	-	204k	11.3k	Document summarization
SamSum	-	14k	818	Conversation summarization
MultiNews	-	45k	5k	News summarization

A DETAILED DISCUSSION OF RELATED WORKS

Instruction fine-tuning. Many recent papers have begun looking into algorithmic aspects of instruction fine-tuning. Wang et al. (2023a) propose CocktailSGD for distributed training of large language models, which is a novel communication-efficient training framework that combines three distinct compression techniques (random sparsification, top-K sparsification, and quantization) to achieve much greater compression than each individual technique alone. Tanwisuth et al. (2023) propose an unsupervised fine-tuning framework to fine-tune the model directly or prompt on the unlabeled target data.

Besides, Lee et al. (2023) found that selectively fine-tuning a subset of layers outperforms full fine-tuning when transferring to tasks with various distribution shifts. Jang et al. (2023) find that an expert LM trained on a single task can outperform a multitask LM trained with different tasks on unseen datasets.

There have also been efforts to design more efficient fine-tuning methods. Hu et al. (2022) proposes Low-Rank Adaptation (LoRA) that freezes the pre-trained model weights and injects trainable rank decomposition matrices into each layer of the Transformer architecture. Built on this approach, Zhang et al. (2023a) adaptively allocates the parameter budget among weight matrices according to their importance score. Furthermore, Chen et al. (2023) introduce parameter-efficient fine-tuning design spaces that parameterize tuning structures characterized by four components: layer grouping, trainable parameter allocation, tunable groups, and strategy assignment.

At the same time, prompt generation has received growing interest in recent literature. Wang et al. (2022a) propose Self-Instruct to improve the instruction-following capabilities of pretrained language models by bootstrapping off their generations. Zhou et al. (2023) propose an automatic instruction generation method by searching over a pool of instruction candidates proposed by an LLM

to maximize a chosen score function. Ye et al. (2023) trains the LM to generate the task instruction given the input instance and label. Zhang et al. (2023b) design an action space for editing the prompts and use reinforcement learning to optimize prompts during test time. Si et al. (2023) design simple and effective prompts that improve LLM’s reliability.

Understanding Language Model Fine-tuning. Malladi et al. (2023) analyze fine-tuning large language models through neural tangent kernels and provide empirical evidence that the prompt-based LM fine-tuning exhibits kernel behavior characterized by linearization and fixed features. Wei et al. (2023) investigate the neural tangent kernel (NTK) to reveal the gradient descent dynamics of the multilayer perceptron modules in an LM and propose to coin a lightweight LM through NTK-approximating MLP fusion. Panigrahi et al. (2023) study the learned skills that reside inside the fine-tuned model by identifying a very small subset of parameters responsible for the model’s performance.

B EXPERIMENT DETAILS

B.1 IMPLEMENTATION DETAILS

Planted Model for Generating Clusters with Varied Densities. We use a planted model to compare against different clustering algorithms. In the model, we consider a matrix with k clusters with high weights within clusters and relatively low weights between clusters. We consider sampling random entries in $T \in \mathbb{R}^{n \times n}$ with the following distributions.

- For $1 \leq i \leq k$, the entries of each diagonal block T_{C_i, C_i} , i.e., within a cluster, are independently sampled from a Gaussian distribution $\Omega_i = \mathcal{N}(\alpha_i, \sigma_i^2)$.
- For $1 \leq i, j \leq k$, the remaining entries between two clusters i and j are independently sampled from a Gaussian distribution $\Omega_{i,j} = \mathcal{N}(\beta_{i,j}, \sigma_{i,j}^2)$.
- We define that the mean α_i is larger than the means $\beta_{i,j}$ to ensure weights within clusters are higher than weights between clusters. Then, We create clusters with different densities by defining various values of α_i for different i .

Specifically, we consider three clusters $k = 3$ and 50 points within each cluster, i.e., $n = 150$. We consider one low-density cluster with $\alpha_1 = 12$ and $\sigma_1 = 4$, and two high-density cluster with $\alpha_2 = \alpha_3 = 14$ and $\sigma_2 = \sigma_3 = 2$. The values between clusters are defined as follows: $\beta_{1,2} = 8, \beta_{1,3} = 7, \beta_{2,3} = 13$ and $\sigma_{1,2} = 4, \sigma_{1,3} = 4, \sigma_{2,3} = 2$.

Multitask Learning on the GLUE Benchmark. We study the pairwise relationship between all the tasks in the GLUE benchmark. More specifically, we consider all pair combinations of the tasks in the GLUE benchmark, including CoLA, SST-2, MRPC, STS-B, QQP, MNLI, QNLI, RTE, and WNLI, which are nine tasks and 36 combinations in total. For all the combinations, we combined all datasets inside each combination to fine-tune a pre-trained model respectively.

In this fine-tuning experiment, we use RoBERT-Base from hugging face as the pre-trained model, set the batch size of 16, and use the learning rate $5e-5$. Therefore, we will have 36 different fine-tuned models for different combinations. Also, we run the single task fine-tuning with the same setup as the baseline. The next step is to obtain the evaluation accuracy of each task we used to fine-tune the model. For example, if we have a combination (CoLA, SST-2) and a model fine-tuned by these two tasks, we will get the evaluation accuracy of CoLA and SST-2, respectively. After getting all the results above, we calculate the affinity score between all pairs and compare them with the single-task performance.

Evaluation on NLP Task Groups. After collecting the groups of NLP tasks that exhibit positive transfer within each group, we evaluate whether our approach can identify the underlying group structures with the following process. For each evaluation case, such as two groups in Table 3, we randomly sample 3 tasks inside these two groups. Then, we fine-tune a pre-trained model on the sampled subsets. We use T5-Base as the base model for fine-tuning with an AdamW optimizer using a constant learning rate. We set the learning rate as $5e-5$ and the batch size as 16.



Figure 5: The cluster structures in terms of affinity scores between NLP tasks of different categories.

We compute the affinity scores of every task and apply a clustering algorithm to the affinity matrix to evaluate whether the tasks in different groups can be separated. Take the Natural Language Inference (MNLI RTE WNLI) and Multiple-Choice QA (BoolQ COPA MultiRC ReCoRD). we obtain an affinity matrix with size 7×7 after conducting the process above. Then, we apply a clustering algorithm to the affinity matrix: the tasks of two groups could be separated where the tasks inside each group are related. More examples are illustrated in Figure 5.

Distribution of Function Classes. For linear regression, we consider the class of linear functions denoted as $\{f|f(x) = w^\top x, w \in \mathbb{R}^d\}$. Each class of functions is associated with a Gaussian distribution which the function weight w is sampled from. For decision trees, we consider the class of functions represented as depth-4 binary decision trees where each non-leaf node is associated with a coordinate in x and each leaf node is associated with a target value, which is sampled from the normal distribution. For neural networks, we consider two-layer ReLU neural networks. The weight parameters are sampled from Gaussian distributions.

B.2 ADDITIONAL IN-CONTEXT LEARNING RESULTS

Training a Transformer on Pairs of Function Classes. To begin with, we train a transformer on the sampled prompt sequences from two types of function classes. We observe that those three different function classes are affected by other functions for different extents. The results are illustrated in Figure 6, where we plot the test mean squared error of predicting the function values using a different number of in-context examples.

Based on our observation of the linear regression experiments, we noticed that the in-context learning performance became worse after training the model on both linear functions and other function classes. Combining with neural networks leads to even poorer performance than combining with decision trees. For decision trees, we noticed that the in-context learning performance is dramatically decreased when training a transformer on decision trees with other function classes. In the case of neural networks, the in-context learning performance remains comparable between training a transformer only on neural network examples and training with linear functions or decision trees.

Spectral Clustering Results. We find that the loss of NN function classes is significantly larger than the other two. As spectral clustering uses eigenvectors of the Laplacian matrix of T , we found that the tasks of linear regression and decision trees are not separable due to the higher loss value of NNs, as shown in Figure 7.

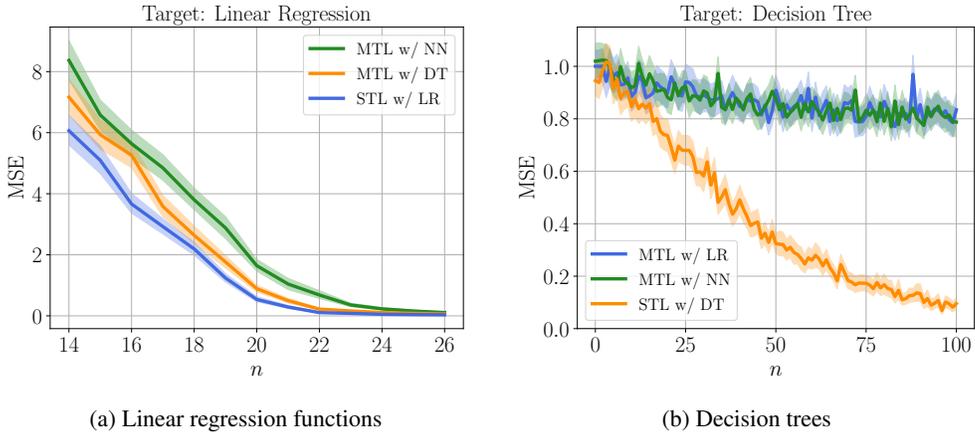


Figure 6: We plot the test mean squared error of the transformer for predicting $f(x_{k+1})$ given $(x_1, f(x_1), x_2, f(x_2), \dots, x_{k+1})$, w.r.t. the number of in-context examples k . For each figure, we fix one function class as the target function class. Then, we pick another type of function class and train a transformer on prompts from both function classes (MTL). We compare the prediction error of MTL with training a transformer only on the target function class (STL). (Left) A transformer trained on linear functions with decision trees or neural networks performs worse than a transformer only trained on linear functions; (Right) A transformer trained on decision trees with linear functions or neural networks becomes much worse than a transformer only trained on decision trees.

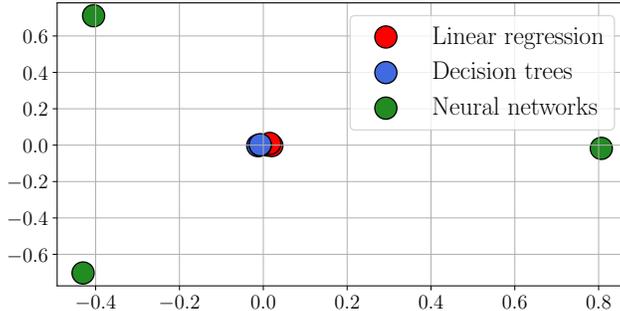


Figure 7: We show that the linear regression and decision tree function classes are not separable in the eigenvector space of the affinity score matrix due to the large loss value of the neural network function class. We illustrate the eigenvectors corresponding to the second and third largest eigenvalues of the normalized Laplacian $D^{-1}T$ of the affinity score matrix T .

B.3 HIGHER-ORDER TASK AFFINITY FOR INSTRUCTION FINE-TUNING

As the widespread negative transfer is observed in fine-tuning multiple tasks, we study a related question of deciding negative or positive transfer effects from a combination of tasks to a target task. The most thorough solution to answer this question is to enumerate all combinations of tasks. However, this solution is expensive, as there are $O(2^n)$ possible combinations given n tasks. We use task affinity measures to capture the negative and positive transfer effects among tasks.

The first method for the above question is to compute *pairwise task affinity scores* for every pair of tasks and approximate the transfer effects of a subset by averaging the pairwise affinity scores in the subset. We consider two ways to estimate the pairwise task affinities: (1) the loss-based pairwise affinity score as the task i 's loss evaluated on the model fine-tuned on both task i and another task j (as in Standley et al. (2020)). (2) the gradient-based pairwise affinity score as the ratio of task i 's loss before and after applying the gradient of task j on the model parameters (as in Fifty et al. (2021)).

Given a target task i and a subset of tasks S containing i , we aim to predict whether the subset S has a negative transfer effect to task i , i.e., the MTL prediction loss $f_i(S)$ is worse than the STL loss $f_i(\{i\})$. For each target task i , we view the prediction of negative transfers as binary classification. We evaluate the average F_1 -score over the n tasks. We build affinity measures between tasks to approach this question. We define an affinity score between task i and j as $T_{i,j}$, which measures how well task j transfers to task i when combined in fine-tuning.

Another method is learning task affinity scores to approximate the MTL performance on task subsets directly. Concretely, we adopt the *surrogate modeling* setting (Ilyas et al., 2022; Li et al., 2023) where a surrogate model $g_i(S)$ is used to approximate the MTL performance $f_i(S)$ of task i on any subset S . We specify the surrogate model as a linear model and parameterize it by the affinity scores $\theta_i = [\theta_{i,1}, \dots, \theta_{i,n}]$, i.e., $g(S; \theta_i) = \sum_{j \in S} \theta_{i,j}$. This choice of $g(S; \theta_i)$ allows us to estimate affinity scores more efficiently than more complex ones. The procedure to estimate the affinity scores is as follows. First, sample m subsets of tasks $\{1, 2, \dots, n\}$ uniformly over subsets of size α , denoted as S_1, S_2, \dots, S_m . Then, compute the value of $f_i(S_j)$ by fine-tuning a model on tasks in every subset S_j for $j = 1, \dots, m$. Lastly, minimize the mean squared error between the surrogate model $g_i(S)$ and $f_i(S)$ over the m subsets:

$$\min_{\theta_{i,1}, \dots, \theta_{i,n}} \frac{1}{m} \sum_{j=1}^m (g(S; \theta_i) - f_i(S_j))$$

The affinity scores $\{\hat{\theta}_{i,1}, \dots, \hat{\theta}_{i,n}\}$ are the minimizer of the above objective. After estimating the affinity scores for an unseen subset of tasks S , we use $g_i(S)$ to predict whether S has a negative transfer effect on task i .

Results. We evaluate the accuracy of using the affinity scores to predict the negative transfer effects from a task subset to a target task. We use the ten instructions to conduct instruction fine-tuning on the RTE dataset as an example. We vary the subset size $|S|$ ranging from 2 to 6 and evaluate the prediction results on a holdout set of subsets.

As shown in Figure 8, we found that linear surrogate modeling yields more accurate predictions than the other two pairwise task affinity scores. While the F_1 -score of the predictions using the two pairwise affinity scores gradually gets worse, using linear surrogate modeling remains accurate as increasing the subset size.

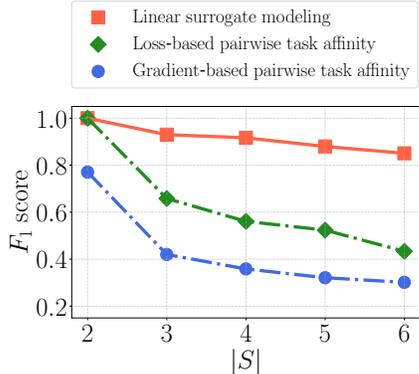


Figure 8: Linear surrogate modeling can consistently predict positive or negative transfers from task subsets to a target task with various task sizes.

Analysis. Next, we analyze the task affinity scores estimated in the linear surrogate model. We find that the task affinity score between two tasks is proportional to the average MTL performance of all subsets that include the two tasks. This suggests an explicit way to estimate a higher-order task affinity between tasks as the average MTL performance over subsets including two tasks.

For task i and task j , suppose that there are $n_{i,j}$ subsets that include both task i and j . Let $T_{i,j}$ be the average MTL performance of all subsets that include them:

$$T_{i,j} = \frac{1}{n_{i,j}} \sum_{1 \leq k \leq n: \{i,j\} \subseteq S} f_i(S_k), \text{ for all } 1 \leq i, j \leq n. \quad (7)$$

We show the connection between $\hat{\theta}_{i,j}$ and $T_{i,j}$ as follows. Let m be the number of sampled subsets and α be the size of subsets less than $\frac{k}{2}$. With probability $1 - \delta$, for any $\delta > 0$, the following holds:

$$\left| \frac{1}{m} (\hat{\theta}_{i,j} - \hat{\theta}_{i,k}) - (T_{i,j} - T_{i,k}) \right| \lesssim \frac{\log(\delta^{-1}n)}{\sqrt{m}}, \text{ for any } 1 \leq j, k \leq n. \quad (8)$$

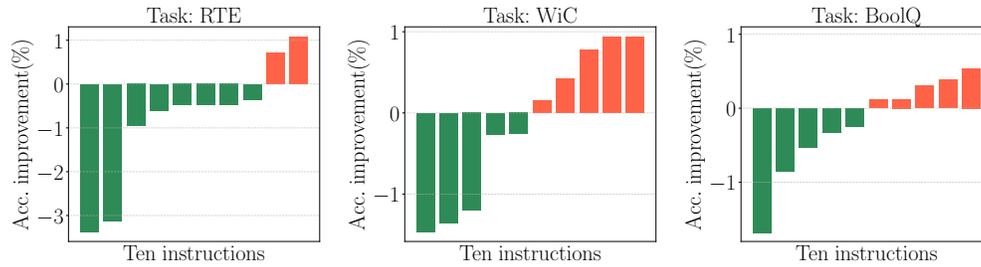


Figure 9: This figure illustrates the negative interference between instructions, i.e., tuning a model with multiple instructions can be worse than tuning with a single instruction for four datasets. Each plot shows the results from one dataset. For each plot, we first fine-tune a model on all instructions. Then, we also fine-tune a model with each single instruction separately. We evaluate the test performance on each instruction i and report the accuracy difference between fine-tuning with multiple instructions and fine-tuning with instruction i . Bars below zero indicate using all instructions results in poorer performance compared to using a single instruction..

Given the above results, we treat $T_{i,j}$ as a higher-order notion of task affinity score that measures how well task j transfers to task i , accounting for the presence of other tasks combined in fine-tuning. A higher value of $T_{i,j}$ indicates task j transfers better to task i . This justifies this notion of higher-order task affinity in Sec. 4.2.