# Appendix of Experimental Details

The experiments focus on the evolutionary algorithm (EA) selection for continuous optimization problems to validate the effectiveness of our method, which are divided into three parts. First, we conduct comparative experiments on candidate EA sets of varying sizes and different scales of training data, to compare its performance against the state-of-the-art algorithm selection methods. Subsequently, we perform ablation experiments on the model by replacing or eliminating parts of the model to validate the role of each module. Following this, there is an analysis experiment that primarily examines the impact of negative sampling on the accuracy and time efficiency of our method. Finally, we embed different large language models (LLMs) into our method to showcase and compare the capabilities of these LLMs in code comprehension and representation.

## A. PERFORMANCE COMPARISON FOR ALGORITHM SELECTION TASKS

**Comparing Algorithms**: To verify the performance of AS-LLM, we selected five comparing algorithm selection methods proposed in recent years as follows:

- Algorithm recommender for white-box problems (AR-WB) [1]. AR-WB is a algorithm selection system for continuous optimization problems. It represents problem structures using tree-based features and utilizes a deep recurrent neural network to recommend the most suitable metaheuristic algorithm.
- Implicit multi-fidelity algorithm selection (IMFAS) [2].
   IMFAS is a algorithm selection method that directly leverages multi-fidelity landmarking information from candidate algorithms, learning dataset topology and algorithm biases via LSTMs and MLP in a few-shot setting.
- Algorithm selection for multi-objective optimization problems (AS-MOP) [3]. This paper introduces an automated algorithm selection method for MOPs, which uses a SVM predictor to choose the most suitable EA based on explicit and implicit features of an MOP.
- Algorithm selection for faster Top-K retrieval (AS-TOPK) [4]. This study introduces a machine learning (decision tree / random forest) based algorithm selection method to predict the time-efficiency of different top-k query processing algorithms.
- Task-agnostic representation optimized for algorithm selection (TRIO) [5]. TRIO is a meta-learning approach that utilizes learned graphical representations of datasets to recommend the best-performing machine learning algorithms with XGBoost algorithm.

As the aforementioned models were not all explicitly designed for algorithm selection in continuous optimization problems, certain algorithms may not be suitable for feature extraction in this context. In such cases, we utilize features extracted from the objective function and employ separate learning models from the comparing methods to establish a mapping from problem features to algorithms.

**Dataset Description**: The training data utilized in this study is sourced from the literature [1]. The training data collection strategy employed in the cited work involves creating complex functions using a tree-like structure. The initial nodes of the tree consist of mean computation and x, after which a random leaf node is selected and extended with a new operator, along with the random selection of new operands based on the operator. To construct more intricate functions, various operations are introduced during the sampling process to modify the constructed tree, including the addition of noise, removal of redundant operations, and substitution of complex functions. After collecting the objective functions for the given problems, candidate algorithm sets of various sizes are deployed on each problem in the problem set, and performance for each candidate algorithm on each problem are obtained. The candidate algorithms include artificial bee colony algorithm (ABC), ant colony optimization (ACO), covariance matrix adaptation evolution strategy (CMA-ES), competitive swarm optimizer (CSO), differential evolution (DE) with rand/1/bin, fast evolutionary programming (FEP), genetic algorithm (GA) with simulated binary crossover and polynomial mutation, particle swarm optimization (PSO), simulated annealing (SA), and random search (Rand). We selected candidate algorithm sets of different sizes, denoted as Problemset  $\#i(i \in \{1, 2, \dots, 20\})$ , each containing different candidate algorithms. Tables I and II provides the number of algorithms included in each set and the proportion of each algorithms that achieved the best performance.

Experimental Settings: In addition to obtaining the aforementioned problem features, algorithm text serves as another input information for our method, from which LLM extracts features for each EA. The code used in this experiment is the original MATLAB code that generates performance data, with redundant data preprocessing stages removed from each code text, retaining only the corresponding algorithm functions. Aside from experiments comparing the code representation capabilities of different LLMs, all other experiments use Code-Bert [6] as the LLM. Since most LLMs have high-dimensional representations for code text, for example, CodeBert extracts features with dimensions ranging from hundreds of thousands to millions, a two-stage pooling layer is employed to reduce the extremely high-dimensional algorithm features in order to control their impact on the results. The final problem and algorithm representations are set to the same 10 dimensions.

TABLE I Proportion of each best-performance EA within 10,000 problem instances.

Datasets			Pr	oportion of e	ach best-per	formance alg	orithm			Proportion of each best-performance algorithm								
Problemset #1	ABC	ACO	CMA-ES	CSO	DE	FEP	GA	PSO	SA	Rand								
	0.26%	6.62%	30.44%	7.59%	2.59%	0.08%	41.30%	9.38%	1.67%	0.07%								
Problemset #2	CMA-ES	GA																
1 TOUICHISCT #2	37.19%	62.81%																
Problemset #3	CSO	PSO																
FIODICIIISCI #3	70.07%	29.93%																
Problemset #4	ABC	ACO	CMA-ES	CSO	DE	FEP	PSO	Rand										
FIODICIIISCI #4	1.19%	10.46%	34.40%	15.55%	2.64%	13.60%	21.89%	0.27%										
Problemset #5	ABC	CMA-ES	CSO	DE	GA	PSO	SA	Rand										
FIODICIIISCI #3	2.65%	32.48%	7.63%	2.64%	43.48%	9.58%	1.47%	0.07%										
Problemset #6	ACO	CMA-ES	CSO	DE	FEP	GA	PSO	SA										
FIODICIIISCI #0	6.59%	30.86%	7.55%	2.56%	0.19%	41.34%	9.44%	1.47%										
Problemset #7	CSO	DE	PSO	SA	Rand													
1 IOυICIIIsct π/	25.46%	32.30%	17.20%	17.50%	7.54%													
Problemset #8	CMA-ES	DE	FEP	SA	Rand													
1 Ιουιαπικά πο	21.42%	11.91%	0.84%	55.48%	10.25%													
Problemset #9	ACO	CSO	FEP	GA	PSO													
1 Ιουιοπικοί π9	42.46%	9.35%	34.78%	12.30%	1.11%													
Problemset #10	ABC	CSO	FEP	PSO	SA													
1 TOURCHISEL #10	61.54%	9.08%	21.20%	7.81%	0.37%													

TABLE II PROPORTION OF EACH BEST-PERFORMANCE EA WITHIN  $30,\!000$  problem instances.

Datasets	Proportion of each best-performance algorithm									
Problemset #11	ABC	ACO	CMA-ES	CSO	DE	FEP	GA	PSO	SA	Rand
	0.20%	6.50%	30.50%	7.81%	2.41%	0.10%	41.63%	9.18%	1.57%	0.08%
Problemset #12	CMA-ES	GA								
11001cmsct #12	37.02%	62.98%								
Problemset #13	CSO	PSO								
1 Toblemset #15	70.44%	29.56%								
Problemset #14	ABC	ACO	CMA-ES	CSO	DE	FEP	PSO	Rand		
11001cmsct #14	1.20%	10.64%	34.81%	15.63%	2.92%	13.41%	21.21%	0.18%		
Problemset #15	ABC	CMA-ES	CSO	DE	GA	PSO	SA	Rand		
1100iciliset #15	2.63%	32.24%	7.85%	2.67%	43.55%	9.44%	1.56%	0.06%		
Problemset #16	ACO	CMA-ES	CSO	DE	FEP	GA	PSO	SA		
11001cmsct #10	6.89%	30.64%	7.61%	2.51%	0.19%	41.17%	9.44%	1.55%		
Problemset #17	CSO	DE	PSO	SA	Rand					
Troolemset #17	25.26%	32.57%	17.42%	17.07%	7.68%					
Problemset #18	CMA-ES	DE	FEP	SA	Rand					
11001cmsct #10	21.53%	11.47%	0.90%	55.51%	10.59%					
Problemset #19	ACO	CSO	FEP	GA	PSO					
11001cmsct #17	43.12%	9.74%	34.05%	12.07%	1.02%					
Problemset #20	ABC	CSO	FEP	PSO	SA					
riouiciliset #20	60.66%	9.33%	21.41%	8.25%	0.35%					

The loss function used is cross-entropy loss, which assesses the model's predictions of the positivity or negativity of each example. The training batch size for all deep models is set to 128. In the TRIO algorithm, the maximum tree depth for XGBoost is set to 100, and the number of weak classifiers is set to 1000. The SVM and decision tree classifiers used in the AS-MOP and AS-TOPK algorithms respectively are configured with default settings from sklearn. We conducted multiple rounds of testing on datasets of varying sizes, with 80% of the samples in each dataset allocated as training data and the remaining 20% as test data.

**Performance Evaluation**: Tables III and IV demonstrate the average accuracy of decision-making by each comparing method. In both Table III and Table IV, we can observe that our method consistently outperforms the other algorithm selection methods in terms of accuracy across most of the problem sets. This trend indicates the robustness and effectiveness of our method in handling both smaller (10,000 problem instances) and larger (30,000 problem instances) tasks. We can also observe variations in the performance of other methods. TRIO, AS-TOPK, and IMFAS show competitive results in specific datasets, especially on Problemset #4 with 10,000 problems and Problemset #9 with 30,000problems. By comparing the results from the two tables, we have discovered that our method gains a significant performance advantage by leveraging additional information sources, i.e., algorithm features, particularly when the training data is limited. For example, when 10,000 samples are insufficient to model

TABLE III
PERFORMANCE OF DIFFERENT ALGORITHM SELECTION METHODS ON 10000 PROBLEM INSTANCES.

Algorithm	TRIO	AS-TOPK	AS-MOP	AR-WB	IMFAS	Ours
Problemset #1	0.9161	0.8814	0.8346	0.9275	0.9271	0.9304
Problemset #2	0.8021	0.7585	0.6255	0.8219	0.8348	0.9357
Problemset #3	0.8145	0.7568	0.6953	0.864	0.8651	0.9554
Problemset #4	0.8704	0.8391	0.7625	0.8751	0.8720	0.8639
Problemset #5	0.9054	0.8573	0.8031	0.9168	0.9205	0.9319
Problemset #6	0.8964	0.8541	0.7923	0.9157	0.9120	0.9244
Problemset #7	0.8122	0.7680	0.6635	0.8234	0.8362	0.8617
Problemset #8	0.8112	0.8135	0.7782	0.9004	0.8941	0.9126
Problemset #9	0.8383	0.7912	0.7163	0.8650	0.8563	0.8788
Problemset #10	0.8778	0.8235	0.7954	0.8921	0.8916	0.9123

TABLE IV
PERFORMANCE OF DIFFERENT ALGORITHM SELECTION METHODS ON 30000 PROBLEM INSTANCES.

Algorithm	TRIO	AS-TOPK	AS-MOP	AR-WB	IMFAS	Ours
Problemset #11	0.9448	0.9499	0.8423	0.9587	0.9523	0.9662
Problemset #12	0.8963	0.8897	0.6308	0.8931	0.8993	0.9658
Problemset #13	0.9135	0.8987	0.7042	0.9135	0.9167	0.9850
Problemset #14	0.9218	0.9301	0.7670	0.9090	0.8650	0.9280
Problemset #15	0.9396	0.9412	0.7988	0.9287	0.9269	0.9681
Problemset #16	0.9361	0.9392	0.7954	0.9369	0.9296	0.9513
Problemset #17	0.9026	0.9056	0.6602	0.8768	0.8699	0.9113
Problemset #18	0.9325	0.9297	0.7777	0.9441	0.9106	0.9588
Problemset #19	0.9096	0.9161	0.7113	0.9114	0.9220	0.8969
Problemset #20	0.9286	0.9249	0.8031	0.9283	0.9271	0.9578

the mapping from problem features to algorithm features, our method, with the assistance of algorithm features, can converge to a superior model more effectively. The results suggest that the ability to leverage algorithm-specific features and model bidirectional relationships in algorithm selection is particularly advantageous. It exhibits robustness across various datasets, making it a reliable choice for algorithm selection tasks, especially especially it is difficult to directly model the mapping from problem features to algorithms under low-quality training data.

#### **B.** ABLATION EXPERIMENTS

In this subsection, we conduct ablation experiments to assess the impact of the four modules: (1) the significance of cosine similarity calculations, (2) the effectiveness of algorithm track for algorithm representation, (3) the efficacy of average pooling layer in algorithm track, and (4) the function of the MLP layer used for output generation. Specifically, 'Ours' represents the original model, while its four variant models are defined as follows: (1) Ours-Si utilizes both algorithm and problem features but does not calculate the degree of matching degree between algorithm and problem representations; instead, it passes the representations of algorithms and problems through an MLP layer. (2) Ours-AF removes algorithm track from the original model, retaining only problem features, and directly learns the mapping from problem representations. (3) Ours-Pool replaces the average pooling in the original model with a maximum pooling layer. (4) Ours-MLP employs both algorithm and problem features and computes the matching degree between algorithm and problem representations, but the final MLP layer does not consider algorithm features.

We choose 8 problem sets with different numbers of problems and candidate algorithms, to conduct experiments on four variants of the original model, and comprehensively and fairly assess the impact of each module. The results of the ablation experiments are presented in Tables V and VI. (1) The variant with the most significant difference from the original model is Ours-AF. It exhibited the poorest performance in all experiments when algorithm features were omitted. This underscores the crucial influence of algorithm features and algorithm track on algorithm selection problems. (2) Ours-Si displayed a substantial performance difference compared to the original model, indicating that the matching degree between problem features and algorithm features in our model is effective and has a significant impact on the final results. (3) Ours-MLP showed performance differences compared to the original model, but the disparities were not as pronounced as in the first two variants. This module is not a core component of the model, as it is placed after similarity computation and mainly facilitates gradient backpropagation, thus positively influencing the model's accuracy but exerting a limited impact on it. (4) The variant with the smallest difference from the original model is Ours-Pool. Despite selecting the performance-enhancing average pooling, different pooling methods had less noticeable effects on performance in the algorithm selection scenario. The main reason is the preprocessing of high-dimensional algorithm features, which diminishes the differences between various pooling methods.

In summary, the core contributions highlighted in this study,

 $\label{table v} \mbox{Table V} \\ \mbox{Ablation Experiment on } 10,000 \mbox{ problem instances}.$ 

	Ours-Si	Ours-AF	Ours-MLP	Ours-Pool	Ours
Problemset #1	0.918	0.797	0.929	0.929	0.931
Problemset #2	0.944	0.341	0.945	0.945	0.945
Problemset #4	0.836	0.749	0.866	0.869	0.872
Problemset #9	0.857	0.656	0.888	0.887	0.892

	Ours-Si	Ours-AF	Ours-MLP	Ours-Pool	Ours
Problemset #11	0.948	0.802	0.949	0.952	0.953
Problemset #12	0.968	0.402	0.974	0.979	0.982
Problemset #14	0.904	0.754	0.919	0.932	0.932
Problemset #19	0.922	0.657	0.935	0.939	0.943

including algorithm representation and modeling the matching degree between problems and algorithms, had a significant impact on the model's results. The original model accurately collects training information through problem track and algorithm track, effectively models the matching mechanism between problems and algorithms through similarity calculation, and thus demonstrates the state-of-the-art performance.

#### C. IMPACT OF NEGATIVE SAMPLING

To explore the relationship between the probability of negative sampling and model performance as well as training time, a series of experiments were conducted on Problemset #1 and Problemset #11, which contain the most candidate algorithms. The probability of negative sampling was incrementally increased from 10% to 90%, and sampling was performed under a uniform distribution according to Section 3.4 in the main text. The experiments recorded the best performance achieved within 300 epochs and the average runtime for each epoch. The experimental results are presented in Figure 1.

From the graph, it is evident that as the probability of negative sampling increases, the algorithm's training time also increases sequentially. This is attributed to the higher probability resulting in the generation of more samples, thereby increasing training time. However, the changes in performance were not consistently increasing with the rise in sampling rates. In both datasets, there were performance degradation phenomenon under an increasing sampling rate. This suggests that the impact of sampling rates on accuracy should be considered in conjunction with various factors, including the specific scenario, the scale of training data, the quality of the samples, and so on. From this experiment, two conclusions can be drawn: (1) The choice of negative sampling probability should strike a balance between accuracy and time efficiency. (2) Careful adjustment of the negative sampling probability is essential to avoid performance degradation.

## D.EVALUATING THE CODE COMPREHENSION CAPABILITIES OF LLMS

LLM assists our model in extracting algorithm representations from code text. Conversely, the algorithm selection

task based on our framework can serve as a benchmark task for evaluating LLM's code comprehension abilities. To validate this assertion, this subsection conducts experiments where different pre-trained LLMs are deployed within the code feature extraction module to compare their performance differences. Two categories of large models participate in this experiment: one comprises general-purpose LLMs trained on massive textual data (including code text and other text), while the other consists of LLMs trained specifically for code-related tasks. Their respective descriptions are as follows:

### (1) General-purpose LLMs:

- XLNet [7]: XLNet is a self-regressive pre-trained language model that employs the permutation language modeling approach, which allows for better handling of long texts and multi-document tasks. Its primary functionalities include text generation, text classification, and question-answering systems, among others.
- BART [8]: BART is a sequence-to-sequence generative model that utilizes both autoregressive and autoencoding methods, making it suitable for various tasks such as text summarization, translation, and dialogue generation.
- BGE [9]: BGE is a recently released large-scale pretrained model based on natural language processing.
   It has demonstrated superior performance in various tasks, including text generation, text classification, and question-answering systems, in both Chinese and English.
- (2) Code-specific LLMs trained for code-related tasks:
- CodeBERT [6]: CodeBERT is a pre-trained model specifically designed for source code. It can be applied to tasks like code completion, code recommendation, and code searching.
- GraphCodeBERT [10]: GraphCodeBERT is a model that combines graph neural networks with self-regressive pretraining. It can be used for tasks like code completion, code recommendation, and code searching.
- UniXCoder [11]: UniXcoder is a unified cross-modal pre-trained model that leverages multimodal data (i.e. code comment and abstract syntax tree) to pre-train code representation.

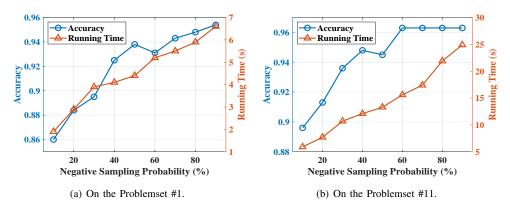


Fig. 1. The relationship between the probability of negative sampling and model performance as well as time efficiency.

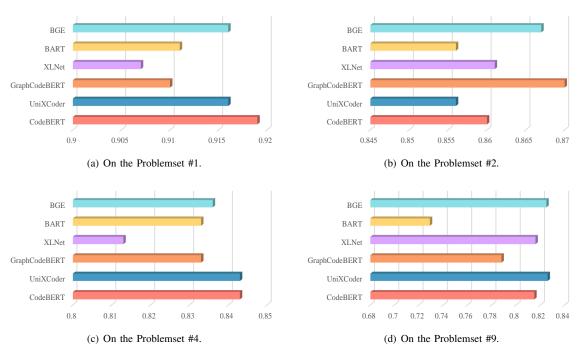


Fig. 2. The performance of different LLMs on 10,000 problem instances.

It is reasonable to evaluate LLMs using the algorithm selection framework because the algorithm selection process can take into account code text from various programming language and across multiple algorithms. We tested the six LLMs on datasets with different candidate algorithm scales, as depicted in Figures 2 and 3. Overall, LLMs pre-trained on code text, such as UniXCoder and CodeBERT, exhibit greater stability, with UniXCoder achieving the best or highly competitive performance on almost all datasets. Although large models pre-trained on common text also possess code representation capabilities and occasionally perform exceptionally well, their stability falls short of LLMs pre-trained on code text. For instance, BART even outperformed all code-pretrained LLMs on Problemset #12. However, as the number of algorithms increases, LLMs pre-trained on code text exhibit greater advantages. The recently introduced BGE model also demonstrates relatively stable performance in code representation. In summary, this experiment indicates the potential of the algorithm selection framework as a benchmark for evaluating LLMs' code representation capabilities.

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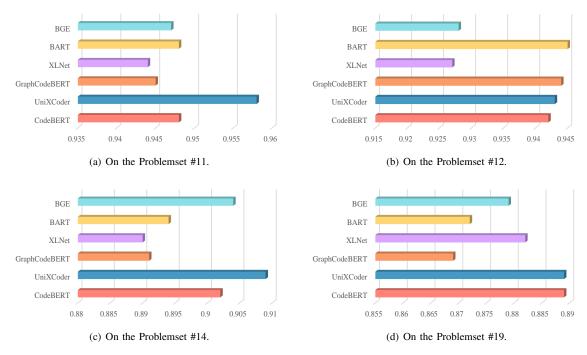


Fig. 3. The performance of different LLMs on 30,000 problem instances.

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