

LLM-enhanced Score Function Evolution for Causal Structure Learning

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Abstract

Causal structure learning (CSL) plays a pivotal role in causality and is often formulated as an optimization problem within score-and-search methods. Under the assumption of an infinite dataset and a predefined distribution, several well-established and consistent score functions have been shown to be both optimal and reliable for identifying ground-truth causal graphs. However, in practice, these idealized assumptions are often infeasible, which can result in CSL algorithms learning suboptimal structures. In this paper, we introduce L-SFE, a framework designed to automatically discover effective score functions by exploring the "score function space". L-SFE addresses this task from a bi-level optimization perspective. First, it leverages a Large Language Model (LLM) to interpret the characteristics of score functions and generate the corresponding code implementations. Next, L-SFE employs evolutionary algorithms along with carefully designed operators, to search for solutions with higher fitness. Additionally, we take the BIC as example and prove the consistency of the generated score functions. Experimental evaluations, conducted on discrete, continuous, and real datasets, demonstrate the high stability, generality and effectiveness of L-SFE.

1 Introduction

Causal structure learning (CSL) is a fundamental approach for understanding causality [Pearl, 2009; Spirtes *et al.*, 2001]. It uncovers causal relationships from observational or interventional data and represents them using graphical models such as Directed Acyclic Graphs (DAGs) *et al.* [Glymour *et al.*, 2019; Vowels *et al.*, 2022]. Learning an exact DAG from data is NP-hard [Chickering, 1996], and the mainstream algorithms can be broadly categorized into two types: constraint-based methods and score-based methods. Constraint-based methods reconstruct the causal graph from a statistical perspective. They first identify the skeleton by performing independence tests and then infer edge directions while adhering the acyclicity and other rules [Koller, 2009]. However, the significance of independence tests cannot be accurately

samples	10	10 ²	10 ³	10 ⁴
Plcg \perp PIP3	-18.7	-157.3	-1701.8	-16783.7
Plcg \rightarrow PIP3	-19.3	-160.8	-1662.6	-16414.7

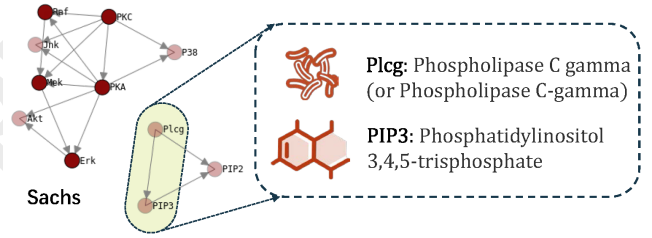


Figure 1: BIC score for subgraphs $\text{Plcg} \rightarrow \text{PIP3}$ and $\text{Plcg} \perp \text{PIP3}$, where the direct causal relationship indicating that Plcg catalyzes the conversion of PIP2 to PIP3 . However, the BIC fails to identify this causal link under the limited data (less than 10^3).

assessed, as it is influenced by variable dimensionality and the number of conditioning variables, among others. In contrast, score-based methods have gained popularity in recent research [Huang *et al.*, 2018]. They treat DAG learning as a constrained optimization problem, and define a score function to evaluate the fitness between candidate graphs and the observed data. Optimization techniques, either combinatorial or continuous, are then applied to identify the optimal DAG in the space of possible causal structures.

Although much of the current research focuses on improving the effectiveness and efficiency of search methods in more complex spaces or under weaker causal assumptions [Cheng *et al.*, 2024], these studies generally adopt consistent score functions, and when the sample size approaches infinity, CSL can reveal the true causal structure, represented by the minimal I-map that accurately captures the data distribution.

However, is the score function can be applied across various scenarios? To explore this, we still take BIC [Schwarz, 1978] as example and consider the limited sample size case, which frequently arises in medical field. As illustrated in Figure 1 [Sachs *et al.*, 2005], BIC exhibits varying preferences in model selection depending on the sample size. In fact, this gap is universal across score functions used for model selection. For instance, abnormal noise can negatively impact the BIC-Gaussian score, and the choice of priors can influence

the BDeu score. Although some improvements have been proposed under various conditions [Silander *et al.*, 2008; Huang *et al.*, 2018; Andrews *et al.*, 2018], it is impractical to manually design new scoring functions for every emerging scenario. Given these challenges, our aim is to improve the score functions in a automatic and systematic manner. In fact, this task can be also framed as an optimization problem, requiring to explore the landscape of possible scoring functions.

Previously, such optimization tasks have been hindered by the lack of a comprehensive understanding of the score function space. Fortunately, recent advancements in large language model (LLM) [Achiam *et al.*, 2023; Zhao *et al.*, 2023] offer promising tools. LLM encodes algorithms through code implementation and contextual description, enabling more effective search methods for algorithm design [Liu *et al.*, 2024b; Romera-Paredes *et al.*, 2024]. Inspired by it, we propose the L-SFE (LLM-enhanced Score Function Evolution) framework to discovery optimal score functions. Note that L-SFE still achieves the MEC-level identifiability under the causal faithfulness and causal sufficiency assumptions [Pearl, 2009]. The contribution of this paper includes:

1. We frame L-SFE as a supervised learning process and formulate it using a bi-level optimization approach. At the lower level, a greedy local search (GLS) is employed to identify the optimal DAGs based on the specific score function. At the upper level, Evolutionary algorithm (EA) is utilized to optimize the scoring function with maximal fitness, which is quantified by the structural difference between the aforementioned DAGs and the ground-truth graphs.

2. For LLM, we take the BIC as standard and design the prompt integrating both conceptual and code-level information to initialize and evolve the scoring function. Additionally, we refine several operators, including mutation, crossover, and injection, to effectively balance convergence and diversity of EA.

3. We evaluate the generated scoring functions from both theoretical and experimental analysis. The former focuses on equivalence and consistency, while latter is performed across synthetic (including discrete and continuous) and real-world data. The results demonstrate the effectiveness, stability, and generality of L-SFE.

2 Related Works

Causal Structure Learning. The score-based CSL algorithms can be categorized into combination-based and continuous-based types. The former execute approximate or exact search in: (1) DAG space \mathcal{G} (e.g., HC [Heckerman *et al.*, 1995] and MAHC [Constantinou *et al.*, 2022]), which explore plausible DAGs by heuristically manipulating single or multiple directed edges; (2) EC (Equivalent space) \mathcal{E} (e.g., GES [Chickering, 2002] and fGES [Ramsey *et al.*, 2017]), which greedy search the completely partially directed acyclic graph (CPDAG) via forward equivalent search and backward equivalent search. (3) Permutation space \mathcal{O} (e.g., OBS [Teyssier and Koller, 2012], GRASP [Lam *et al.*, 2022], and BOSS [Andrews *et al.*, 2023]), which identify the best causal structure by discovering the topological ordering

that maximizes ancestral information. Continuous-based algorithms make assumptions about the data distribution and learn causal graphs from a structural equation model (SEM), such as LiNGAM [Shimizu *et al.*, 2011], NOTEARS [Zheng *et al.*, 2018], DAG-RL [Zhu *et al.*, 2019] and DAGMA [Bello *et al.*, 2022] *et al.*

LLM for CSL. Given the exceptional text comprehension capabilities, numerous studies have leveraged LLM to research the causality [Kıcıman *et al.*, 2023; Takayama *et al.*, 2024]. These approaches regard LLM as sole determiners of pairwise causal relationships [Zhiheng *et al.*, 2022; Wan *et al.*, 2024] or domain experts. The latter harness meta-information embedded in the LLM’s training data to enhance the causal discovery, including initialization [Ban *et al.*, 2023; Li *et al.*, 2024], post-door adjustments [Khatibi *et al.*, 2024], or structural constraints fusion [Ban *et al.*, 2023; Zhou *et al.*, 2024; Jiralerspong *et al.*, 2024; Zhang *et al.*, 2024].

LLM for Algorithm Design. Finally, we provide a brief overview of using LLM for auto algorithm design. Most existing frameworks employ evolutionary approaches and leverage operators defined within the algorithmic space for algorithm generation, such as FunSearch [Romera-Paredes *et al.*, 2024], EoH [Liu *et al.*, 2024b], and ReEov [Ye *et al.*, 2024]. Additionally, these frameworks have been applied to a variety of scenarios, including the capacitated vehicle routing problem [Liu *et al.*, 2024a], critical node discovery [Mao *et al.*, 2024], and tensor network search [Zeng *et al.*, 2024].

3 Background

DAG can be described as a tuple $\mathcal{G} = (\mathbf{V}, \mathbf{E})$, where $\mathbf{V} = \{X_1, X_2, \dots, X_n\}$ represents the collection of variables, and $\mathbf{E} = \{X_i \rightarrow X_j | X_i, X_j \in \mathbf{V}\}$ denotes the directed edges between the variables. CSL can be typically formulated as the optimization problem

$$\begin{aligned} \mathcal{G}^* &= \arg \min_{\mathcal{G} \in \mathcal{G}} \Psi(\mathcal{G} | \mathcal{D}) \\ \text{s.t. } \mathcal{G} &\text{ is acyclic.} \end{aligned} \quad (1)$$

$\Psi(\cdot)$ is a score function that evaluates the fitness of DAG \mathcal{G} on i.i.d dataset $\mathcal{D} = \{\mathcal{D}_1, \dots, \mathcal{D}_m\}$, and it is usually decomposable and equivalence [Koller, 2009].

Definition 1. (Decomposable) If the score function $\Psi(\mathcal{G} | \mathcal{D})$ can be written as $\Psi(\mathcal{G} | \mathcal{D}) = \sum_{X_i \in \mathbf{V}} \Psi(X_i, Pa_i^{\mathcal{G}} | \mathcal{D})$, then Ψ is decomposable.

Definition 2. (Equivalence) $\forall \mathcal{G}_1, \mathcal{G}_2 \in \mathcal{G}$, if $\mathcal{G}_1, \mathcal{G}_2$ is I-equivalent, and $\Psi(\mathcal{G}_1 | \mathcal{D}) = \Psi(\mathcal{G}_2 | \mathcal{D})$ holds, then Ψ is equivalence.

where $Pa_i^{\mathcal{G}}$ represent the parents of X_i in \mathcal{G} . For discrete datasets, the most commonly used information theoretic score functions include BIC and AIC [Kitson *et al.*, 2023], which evaluate model based on the multi-information content and entropy of the variables. Other Bayesian scores, such as BDeu, BDs, K2, and BDe [Kitson *et al.*, 2023], place priors on the parameters of the tabular conditional probability distribution. For continuous datasets, BIC remains applicable when using conditional covariance matrix. Thus, we take BIC as

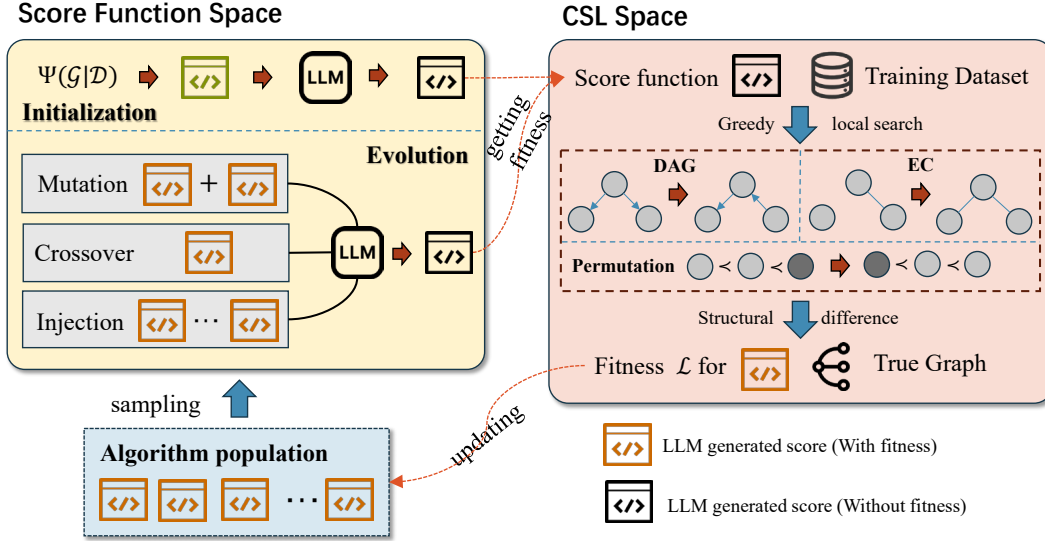


Figure 2: The workflow of the L-SFE. In score function space, EA is employed to guide the search, while well-designed prompts simulate mutation, crossover, and injection operators to push LLM generate new score functions. In CSL space, every generate score function should be combined with various GLS strategies to find the optimal DAGs, which used to compute the fitness.

benchmark for evolution [Andrews *et al.*, 2023]

$$\Psi(\mathcal{G}|\mathcal{D}) = \sum_{X_i \in \mathcal{V}} \ell_{Pa_i^{\mathcal{G}} \rightarrow X_i}(\hat{\theta}_{\text{mle}}|\mathcal{D}) - \frac{\lambda}{2} |\hat{\theta}_{\text{mle}}| \log(m), \quad (2)$$

where $\ell(\cdot)$ is a log-likelihood function, and $\hat{\theta}_{\text{mle}}$ represents the maximum likelihood estimation (MLE) for the parameters of subgraph $\mathcal{G}_i : Pa_i^{\mathcal{G}} \rightarrow X_i$. Furthermore, BIC is also consistent if the underlying distribution \mathcal{P} behind \mathcal{D} belongs to a curved exponential family [Haughton, 1988].

Definition 3. (Consistent) Suppose \mathcal{G}^* is a P -map under the distribution \mathcal{P}^* . A score function Ψ is said to be consistent if as the sample size $m \rightarrow \infty$, the following conditions hold:

- The structure \mathcal{G}^* maximum the $\Psi(\mathcal{G}^*|\mathcal{D})$.
- $\forall \mathcal{G} \in \mathcal{G}$, if \mathcal{G} is not I -equivalent to \mathcal{G}^* , then $\Psi(\mathcal{G}|\mathcal{D}) < \Psi(\mathcal{G}^*|\mathcal{D})$.

4 Framework

The search in score function space can be formulated as a bi-level optimization problem

$$\begin{aligned} \Psi^* &= \arg \min_{\Psi \in \Psi} \mathbb{E}_{\mathcal{G}^*, \mathcal{D}^{\mathcal{G}^*}} (\text{Dis}(\mathcal{G}_{\Psi, \mathcal{D}^{\mathcal{G}^*}}^{\dagger}, \mathcal{G}^*)) \\ \mathcal{G}_{\Psi, \mathcal{D}^{\mathcal{G}^*}}^{\dagger} &= \arg \max_{\mathcal{G} \in \mathcal{G}} \Psi(\mathcal{G}|\mathcal{D}^{\mathcal{G}^*}) \\ \text{s.t. } \mathcal{G} &\text{ is acyclic,} \end{aligned} \quad (3)$$

where \mathcal{G}^* denotes the arbitrary DAG within \mathcal{G} , and $\mathcal{D}^{\mathcal{G}^*}$ represents the dataset sampled from \mathcal{G}^* under the gaussian or multinomial distribution. $\text{Dis}(\mathcal{G}_a, \mathcal{G}_b)$ is a structural distance measure for $\mathcal{G}_a, \mathcal{G}_b$. Note that in equation 3, the upper level optimization identify the optimal Ψ by minimizing the expectation of structural loss over all possible \mathcal{G}^* assigned with

$\mathcal{D}^{\mathcal{G}^*}$. This objective can be achieved through lower level optimization, which finds the best $\mathcal{G}_{\Psi, \mathcal{D}^{\mathcal{G}^*}}^{\dagger}$ for a given score function Ψ . Furthermore, since the score function space is abstract, L-SFE employs EA to solve equation 3, as it does not require a clear characterization for the mathematical properties of the solution. As shown in figure 2, the workflow of the L-SFE can be divided into two stages. Firstly, EA searches the score function space to identify an improved Ψ ; then, Ψ is combined with the GLS to solve the lower level optimization problem, ultimately leading to a best DAG for evaluating the fitness of Ψ .

4.1 Score Function Generation

In this subsection, we provide a detailed explanation of how the score function is discovered using LLM.

Prompts Design. Inspired by EoH [Liu *et al.*, 2024b], all prompts pt are designed using a three-level hierarchical structure as shown in figure 3. *Task Description* informs the LLM of its intended role and the objective of CSL. *Code Snippets* presents few-shot examples to heuristic the LLM designing analogous score functions, and it contains two key components: **Idea**, which provides a textual explanation of the core thought behind the score function; **Code**, which presents the detailed Python implementation. In addition, there is an optional item that demonstrates the fitness of the score function. *Task Assignment* applies the chain-of-thought (CoT) reasoning [Wei *et al.*, 2022] to derive better scoring functions from the given examples, and it includes four steps:

1. First, the LLM carefully reads and interprets the core idea provided in the code comments;
2. Second, if the input score function is labeled with a fitness, the LLM analyzes the reasons behind its effectiveness or shortcomings;

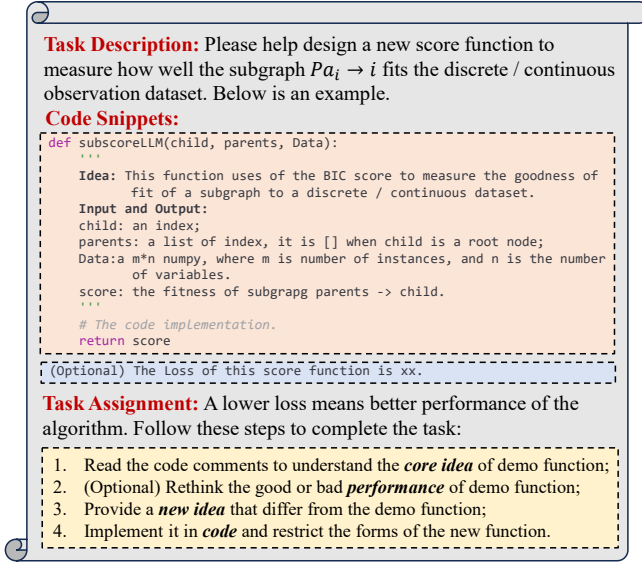


Figure 3: The prompt paradigm in L-SFE. Including a basic task description, few-shot examples, and detailed reasoning instructions.

3. Third, based on the selected operators, different prompts are generated to inspire new ideas that diverge from the given examples;
4. Finally, the LLM implements the new idea in code, ensuring that the function name, input, and output match those in the example code.

Fitness Evaluation in CSL Space. Based on equation 3, L-SFE employs the Monte Carlo method [Metropolis and Ulam, 1949] to randomly sample n_p ground-truth graphs \mathcal{G}^* and corresponding training datasets $\mathcal{D}^{\mathcal{G}^*}$ for training. For convenience, the weight matrix $W^k \in \{0, 1\}^{n \times n}$ is used to represent the k -th \mathcal{G}^* , where $W_{ij}^k = 1$ implies the directed edge $X_i \rightarrow X_j$, and the sampled dataset is denoted as \mathcal{D}^k . Then, the Normalized Hamming Distance [Kıcıman *et al.*, 2023] is used to quantify the $\text{Dis}(\mathcal{G}_a, \mathcal{G}_b)$. Thus, the fitness function can be formulated as

$$\mathcal{L}(\Psi) = \frac{1}{n_p} \sum_{k=1}^{n_p} \frac{\|W_{\Psi, \mathcal{D}^k}^\dagger - W^k\|_1}{d_{W^k}^2}, \quad (4)$$

where d_{W^k} denotes the number of variables in W^k . $W_{\Psi, \mathcal{D}^k}^\dagger$ represents the best DAG learned by the pre-selected GLS under Ψ and \mathcal{D}^k . Obviously, a lower $\mathcal{L}(\Psi)$ indicates a higher fitness. Note that GLS can source from different CSL spaces, such as HC in DAG space, GES in EC space and BOSS in permutation space.

EA in Score Function Space. The initial algorithm population $\Psi = \{\Psi_1, \dots, \Psi_{n_a}\}$ is generated with standard BIC, and mutation, crossover, and injection are employed for evolution. More detailed prompt design can be referred in Supplementary material 1.1¹.

¹<https://github.com/wzd2502/L-SFE>

The crossover operator facilitates the random combination of two score functions, generating a new one that exhibits different structural and conceptual features. Specifically, Ψ_1 and Ψ_2 are randomly selected from the Ψ . Based on their code implementations and the associated fitness $\mathcal{L}(\Psi_1)$ and $\mathcal{L}(\Psi_2)$, LLM analyses and assimilates the ideas of parents to creates a new score function Ψ' .

The mutation operator modifies an existing score function to generate a variant with potentially altered characteristics. Specifically, Ψ_1 is randomly selected from the Ψ with its corresponding fitness $\mathcal{L}(\Psi_1)$. Based on the extent of the modifications, mutation can be classified into two types: (1) structural mutation, which involves significant alterations to the score function’s underlying structure, and (2) parameter mutation, which only modifies the parameters of the score function without altering its fundamental structure.

The injection operator is designed to maintain diversity within the algorithm population, preventing EA convergence to a local optima. Specifically, a new score function Ψ' is generated, which is entirely distinct from all previous $\Psi \in \Psi$. $\mathcal{L}(\Psi')$ is not evaluated at this stage, meaning Ψ' is directly incorporated into the mutation and crossover operations. Ψ' will only be discarded at the end of the iteration if it does not contribute positively to the evolutionary.

In each iteration, all of the aforementioned operators are applied repeatedly, and the generated score functions are tested for legality, ensuring that they do not introduce cyclic graphs and can produce results within a specified time limit. After completing an iteration, the population Ψ would be updated. Only the top n_a algorithms are retained for the next iteration. The evolution continues until a predefined stopping condition is met. To provide a clear understanding of L-SFE, we outline the pseudo-code in Alg. 1 in Supplementary material 1.2.

4.2 Score Function Analysis

L-SFE is repeated n_l times on different training datasets. To avoid ambiguity, we introduce the notation $\text{L-SFE}_{b_k}^a$, $a \in \{D, C\}$, $b \in \{1, \dots, n_l\}$, $k \in \{1, \dots, n_a\}$. This represents the k -th score function learned by L-SFE over the last iteration in b -th repetitions under the a -type dataset (D for Discrete and C for Continuous). Here we take the L-SFE_1^D as example and analyze its evolution in figure 4, and more detailed results can be found in Supplementary material 1.3.

At the first iteration, L-SFE enhances the CSL by incorporating extra penalty terms that account for the length of the parent sets

$$\begin{aligned} \Psi(\mathcal{G}|\mathcal{D}) &= \sum_{i=1}^n \ell_{Pa_i^{\mathcal{G}} \rightarrow X_i}(\hat{\theta}_{mle}|\mathcal{D}) \\ &\quad - \frac{\lambda}{2} (|\hat{\theta}_{mle}| + \sum_{i=1}^n |Pa_i^{\mathcal{G}}|) \log(m). \end{aligned} \quad (5)$$

Equation 5 modifies only the structure of the penalty, so it remains a consistent and equivalent score. At the end of the evolution, L-SFE identifies an optimal Ψ by incorporating priors into the BIC score, which can effectively mitigate over-fitting under limited data. Furthermore, a quadratic penalty is

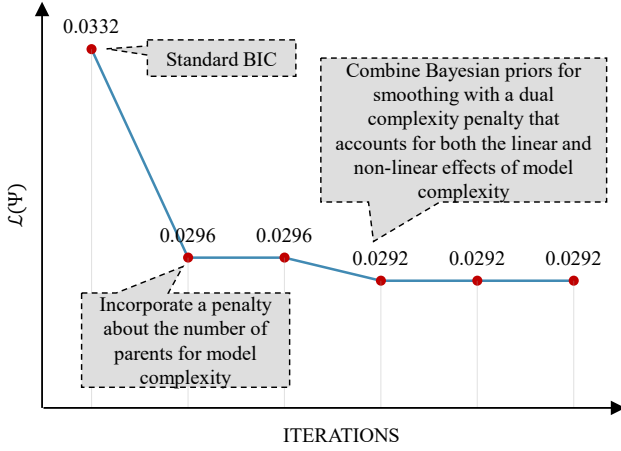


Figure 4: The variation of fitness with respect to the iterations, along with the corresponding evolution of the algorithmic ideas.

introduced to enhance the model’s preference for simplicity, and the penalty term related to the number of data is modified from $\log(m)$ to $1/m$. This formulation is expressed as

$$\Psi(\mathcal{G}|\mathcal{D}) = \sum_{i=1}^n \ell_{Pa_i^{\mathcal{G}} \rightarrow X_i}(\hat{\theta}_{MAP}|\mathcal{D}) - \frac{1}{2m}(\beta|\hat{\theta}_{MAP}| + \gamma|\hat{\theta}_{MAP}|^2), \quad (6)$$

where $\hat{\theta}_{MAP}$ represents the maximum a posterior (MAP) estimate, which is based on Dirichlet prior. Notably, the score function in Equation 6 is decomposable and possesses several advantageous properties

Theorem 1. *Score function Equation 6 is still consistent but not equivalence.*

The proof can be found in supplementary material 1.3. Actually, all LLM learned scores provide similar insights: they leverage the prior to enhance robustness under limited data and use $1/m$ rather than $\log(m)$ to penalize the graph’s complexity in a less restrictive manner.

5 Experiment

5.1 Settings

Datasets. GPT-4o mini is utilized for score function discovery in L-SFE, and synthetic datasets generated from *pytetrat* are employed for training and testing². For the discrete dataset, L-SFE is trained on ten RandomGraphs with $n = 30$, where the variables follow a multinomial distribution, and tested on ErdosRenyi (ER) and ScaleFree (SF) graphs. For the continuous datasets, L-SFE is trained on ten RandomGraphs under the linear Gaussian SEM assumption with $n = 30$, and evaluation is performed on the linear exponential SEM and the linear gumbel SEM. The GLS used for training is HC with a tabu search. Each test is repeated 10 times with $m = 5000$. Additional details of experimental setup are provided in Supplementary material 2.1.

²Code is available on <https://github.com/wzd2502/L-SFE>

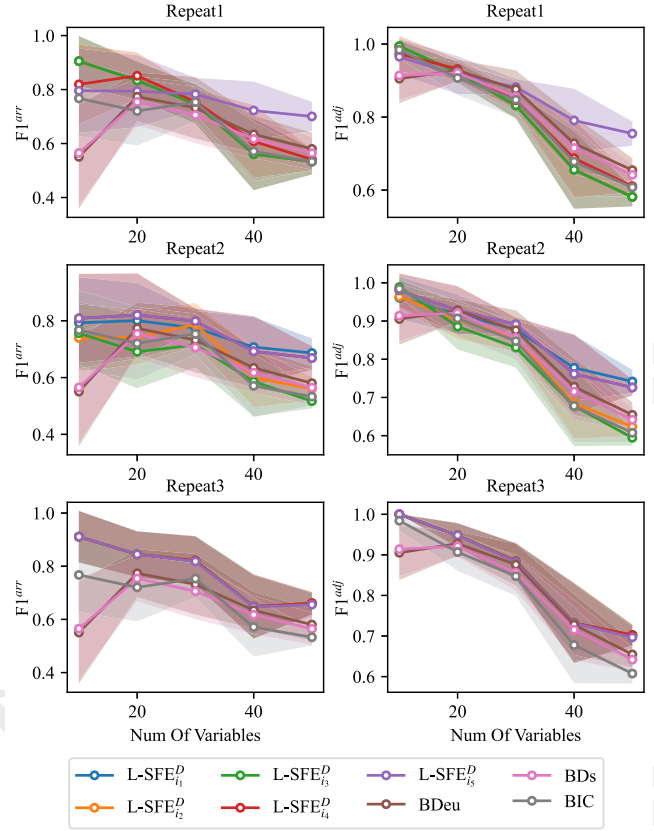


Figure 5: $F1^{arr}(\uparrow)$ and $F1^{adj}(\uparrow)$ comparison of L-SFE against human-designed scores on ER graphs. $L-SFE_{ij}^D$ represent the j -th score function in last iteration of i -th repeat.

Metrics. We evaluate the structural accuracy of L-SFE learned graph against the ground-truth graphs using the Structural Hamming Distance (SHD), the F1 score for adjacency accuracy ($F1^{adj}$), and the F1 score for arrow accuracy ($F1^{arr}$).

Baselines. The score function generated by L-SFE is compared with commonly used scores, including BIC, BDeu, and BDs for discrete data [Kitson *et al.*, 2023], and BIC-Gaussian for continuous data [Kitson *et al.*, 2023]. Additionally, these methods are integrated with GLS across various search spaces and compared to baseline algorithms, including:

- **Discrete.** HC [Heckerman *et al.*, 1995], PC [Colombo *et al.*, 2014], BOSS [Andrews *et al.*, 2023], fGES [Ramsey *et al.*, 2017], GRaSP [Lam *et al.*, 2022].
- **Continuous.** LinGAM [Shimizu *et al.*, 2011], DAGMA [Bello *et al.*, 2022], BOSS, fGES, HC.

5.2 Overall Analysis

In this subsection, we want to answer the following questions: **Stability: Are the all score functions identified by L-SFE consistently superior to human-designed scores?** In this topic, we conduct $n_l = 3$ repetitions of L-SFE and record the all n_a algorithms identified in the last iteration of each repetition. HC is employed as the search method. Figure 5

	ER-10	ER-20	ER-30	ER-40	ER-50	SF-10	SF-20	SF-30	SF-40	SF-50
PC	1.1	14.4	/	112.3	200.8	8.6	34.2	50.0	136.4	163.6
HC	1.3	13.0	39.3	99.4	158.0	11.6	33.6	42.0	126.6	152.1
BOSS	1.2	6.9	31.9	87.6	149.3	10.3	35.9	44.1	130.8	156.8
GRaSP	0.8	11.1	33.2	86.7	151.8	7.8	32.8	42.0	128.3	155.4
fGES	0.5	8.6	32.2	86.7	159.2	10.5	34.7	43.2	130.4	155.2
L-SFE _* ^D (\mathcal{G})	4.7	10.8	21.4	69.1	117.5	8.0	25.9	33.8	123.3	144.8
L-SFE _{**} ^D (\mathcal{E})	2.8	7.3	14.9	45.8	85.3	5.8	25.9	41.0	115.5	142.0
L-SFE _* ^D (\mathcal{O})	3.4	14.1	42.9	102.7	/	5.2	21.0	48.4	130.9	156.1

Table 1: The SHD (\downarrow) comparison with baseline methods on discrete datasets. L-SFE_{*}^D refers to the **best-of- n_l** score function among the L-SFE _{i *}^D ($i = 1, 2, 3$). \mathcal{G} , \mathcal{E} and \mathcal{O} correspond to the GLS used. / indicates that the algorithm cannot find acyclic graphs within 2 hours.

	Exp-10	Exp-20	Exp-30	Exp-40	Exp-50	Gum-10	Gum-20	Gum-30	Gum-40	Gum-50
PC	14.5	23.1	35.8	39.9	37.5	12.6	18.3	27.9	37.3	50.4
fGES	12.9	28.5	38.7	33.2	17.1	13.3	17.6	21.7	27.9	28.2
BOSS	7.1	14.6	16.1	15.8	7.8	6.2	4.5	9.3	12.1	9.2
DAGMA	17.4	40.8	56.2	61.7	76.3	22.6	39.1	54.3	65.4	79.3
LiNGAM	3.2	6.9	9.2	13.8	17.1	4.7	11.0	21.3	24.0	32.9
L-SFE _* ^C (\mathcal{G})	14.0	22.7	59.5	31.0	23.2	14.5	19.0	32.6	46.3	38.4
L-SFE _* ^C (\mathcal{E})	12.3	29.9	37.9	34.6	20.7	12.3	15.2	19.6	21.3	27.1
L-SFE _* ^C (\mathcal{O})	9.1	3.7	7.0	8.0	8.7	0.8	2.7	5.8	8.3	/

Table 2: The SHD (\downarrow) comparison with baseline methods on continuous datasets.

presents the comparison with three human-designed scores on ER graphs under the discrete dataset, and detailed test results on other settings can be found in Supplementary material 2.2. From figure 5, most of LLM learned score functions are either comparable to or outperform the human-designed scores in each repetition. For instance, L-SFE₁₅^D performs best in Repetition 1, while L-SFE₂₁^D and L-SFE₂₅^D show superior performance in Repetition 2, and L-SFE₃₁^D and L-SFE₃₂^D lead in Repetition 3. A similar trend is also observed in the SF graphs. In the case of continuous datasets, all score functions learned by the LLM outperform the human-designed BIC-gaussian score, highlighting the stability of the L-SFE.

Therefore, we select the best score function from i -th repetition, denoted as L-SFE _{i *}^D and evaluate their generality.

Generality: Can the score functions discovered by L-SFE be applied across different search methods? In this topic, we explore whether the score function, trained using HC in the DAG (\mathcal{G}) space, remains effective in the EC (\mathcal{E}) and permutation (\mathcal{O}) spaces. To investigate this, we employ the GES and BOSS, along with L-SFE _{i *}^D for testing. The results on ER graphs are displayed in figure 6, with additional results for other settings provided in Supplementary material 2.2. LLM learned score functions still demonstrate strong performance in the EC space as well compared with BIC and BDeu. However, in the permutation space, this superiority narrows significantly. This is likely due to the smaller scope of the permutation space ($O(n2^{n-1})$) compared to the graphical space ($O(n!2^{C_n^2})$), where GLS are more likely to find the optimal solution, with less sensitive to the choice of

score function. Notably, the computation cost of the Grow-Shrink Tree (GST) for BDeu is prohibitively high, preventing BOSS from identifying a DAG in test datasets when $n \geq 40$. For the continuous dataset, L-SFE maintains excellent generality in the EC space. However, in the permutation space, due to higher capability of BOSS, L-SFE_{1*}^D and L-SFE_{2*}^D slightly under-perform compared to BIC-gaussian, with only L-SFE_{3*}^D achieving a marginal advantage.

Effectiveness: Does L-SFE improve the performance of CSL compared to state-of-the-art algorithms? Finally, we investigate whether the LLM generated score function, when combined with GLS in different search spaces, can outperform SOTA causal learning methods in tables 1 and 2. Overall, L-SFE achieves the lowest SHD in 8/10 settings on discrete datasets and in 7/10 settings on continuous datasets. Counterexamples are observed in some small-scale networks, such as ER-10, ER-20, and Exp-10. These occurrences are mainly because 5000 instances provide a sufficiently sample size for the BIC to yield optimal results. However, as the problem scale increases, the advantages of L-SFE become more pronounced. For discrete datasets, L-SFE in \mathcal{E} space identifies the most optimal DAGs among the eight algorithms, and achieves similar results on continuous datasets in \mathcal{O} space. This observation aligns with our understanding that graphical-based GLS are more suited for discrete datasets, whereas permutation-based GLS can explore causal structures in larger steps, making them more effective for SEM causal models. For further results, including $F1^{arr}$ and $F1^{adj}$ scores, please refer to Supplementary material 2.2.

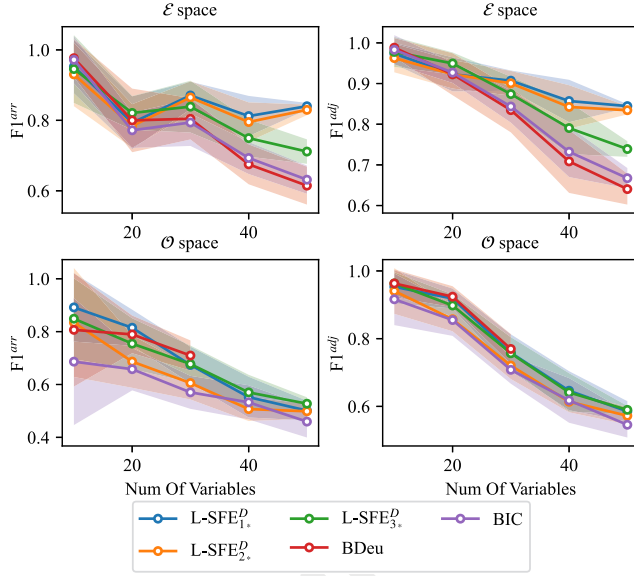


Figure 6: HC trained L-SFE for \mathcal{E} and \mathcal{O} spaces on ER graphs. BDeu + BOSS cannot output a DAG within 1 hour when $n \geq 40$.

5.3 Other Experiments

Case Study. We present a case study using data from the real world COVID-19 pandemic in the UK³. The data set comprises 866 samples that encompass eight categories, including viral tests, infections, hospitalizations, and related factors. The expert-designed benchmark network includes 17 variables, 37 directed edges, and a maximum node degree of 10. The dataset is discretized using k-means and quartiles discretization, as introduced by [Constantinou *et al.*, 2023]. A graphical accuracy analysis is provided in Table 3. Due to the absence of a predefined distribution, the expert-designed benchmark network is not a P-map for dataset. As a result, most existing algorithms fail to produce high F1 score graphs. Even then, the L-SFE-guided GLS approach can consistently identify more correct edges and accurately orient more arrows across different settings.

Ablation Study. We investigate how different types of standard score function hints impact the performance of L-SFE under three configurations: 1) *Seed + Prompt*: The BIC is included in both the population seeds and the code snippets within the prompts; 2) *Seed*: The initialization code snippets provided only the basic inputs and outputs information, with the BIC placed in the seed; 3) *Random Walk*: BIC is absent from both the seed and the prompts. Table 3 evaluates their test performance on ER-20 and SF-20. As expected, L-SFE(Prompt + Seed) yields the best $F1^{arr}$ and $F1^{adj}$, indicating that evolution helps refine the BIC to generate even better score functions, and adding the seed within the population also achieves this goal. However, allowing the LLM to randomly search the score function space leads to suboptimal results. Interestingly, when we extended the number of iterations for L-SFE(Random Walk) to 10, L-SFE still learn the

³<https://bayesian-ai.eecs.qmul.ac.uk/bayesys/>

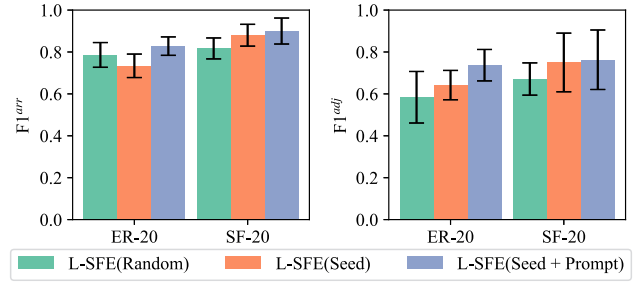


Figure 7: The accuracy of L-SFE trained under three modes.

	k-means		quartiles	
	$F1^{arr}$	$F1^{adj}$	$F1^{arr}$	$F1^{adj}$
PC	0.235	0.441	/	/
HC	0.419	0.581	0.194	0.451
BOSS	0.172	0.483	0.201	0.448
GRaSP	0.142	0.500	0.250	0.463
fGES	0.264	0.415	*0.302	0.453
L-SFE ^D (\mathcal{G})	0.253	0.507	0.347	0.533
L-SFE ^D (\mathcal{E})	0.257	*0.571	0.225	*0.507
L-SFE ^D (\mathcal{O})	*0.338	0.479	0.250	0.417

Table 3: The $F1^{arr}$ and $F1^{adj}$ comparison with baseline methods on COVID-19 dataset. Bold for best and * for second-best performance.

similar thought of BIC. This observation further supports that BIC represents an optimal solution within the score function space. Furthermore, we also analyze the performance of L-SFE across varying sample sizes in Supplementary material 2.2.

6 Conclusion

In this paper, we introduce the L-SFE, which leverages LLM and EA to explore the score function space through a bi-level optimization framework. Three key insights from this work are concluded as follows:

- LLM can effectively learn valuable components, such as Dirichlet priors and relaxed penalties. These components are theoretically well-founded and align with human-like reasoning.
- EA are crucial for finding optimal solutions. Even in the absence of explicit guidance, LLM can discover similar principles through iterative refinement.
- The score function space is highly abstract and potentially multimodal, making best-of-n search a valuable approach for yielding improved results.

Although L-SFE focuses primarily on exploring the structure of score functions, additional work is required to optimize their parameters using LLM. Furthermore, future research could focus on guiding LLM in exploring more efficient causal search strategies or expanding the search space beyond conventional frameworks.

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