

Judge and Improve: Towards a Better Reasoning of Knowledge Graphs with Large Language Models

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Abstract

Graph Neural Networks (GNNs) have shown immense potential in improving the performance of large-scale models by effectively incorporating structured relational information. However, current approaches face two key challenges: (1) achieving robust semantic alignment between graph representations and large models, and (2) ensuring interpretability in the generated outputs. To address these challenges, we propose **ExGLM** (Explainable Graph Language Model), a novel training framework designed to seamlessly integrate graph and language modalities while enhancing transparency. Our framework introduces two core components: (1) a **graph-language synergistic alignment module**, which aligns graph structures with language model to ensure semantic consistency across modalities; and (2) a **judge-and-improve paradigm**, which allows the language model to iteratively evaluate, refine, and prioritize responses with higher interpretability, thereby improving both performance and transparency. Extensive experiments conducted on three benchmark datasets—ogbn-arxiv, Cora, and PubMed—demonstrate that ExGLM not only surpasses existing methods in efficiency but also generates outputs that are significantly more interpretable, effectively addressing the primary limitations of current approaches.

1 Introduction

Large Language Models (LLMs) have demonstrated remarkable success across various natural language processing tasks, including dialogue generation (Aboussalah and Ed-dib, 2025), machine translation (Zhu et al., 2024), question answering (Zhang et al., 2024), and text summarization (Zhang et al., 2025). These models exhibit an impressive capacity for understanding and generating human-like text. However, LLMs face inherent limitations in effectively modeling structured knowledge, such as graphs, which are essential

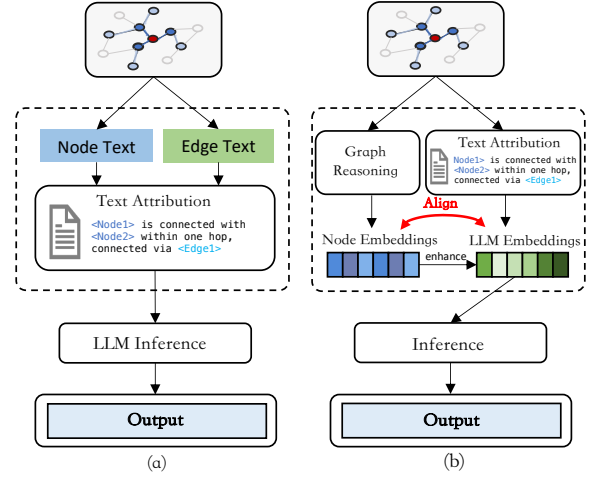


Figure 1: Overview of two mainstream methods (a) textualizing graph and inference via LLM and (b) aligning the semantic representation of LLMs and GNNs.

for capturing complex relationships and dependencies in diverse real-world domains like social networks, biological systems, and knowledge graphs. To address these limitations, recent research has explored the integration of GNNs (Kipf and Welling, 2017; Hamilton et al., 2018; Veličković et al., 2018) with LLMs (OpenAI et al., 2024; Yang et al., 2024; DeepSeek-AI et al., 2025), leveraging GNNs’ strengths in modeling structured information alongside LLMs’ powerful language capabilities, creating opportunities for enhanced performance in graph-related tasks.

Current approaches (Yang et al., 2021; Zhao et al., 2023; Xue et al., 2024) to combining GNNs and LLMs can be broadly classified into two categories. The first category involves **textualizing graph structures** and feeding them into LLMs (Figure 1(a)). For example, some methods (Zhao et al., 2023; Wang et al., 2024; Chen et al., 2024; Wu et al., 2025) describe nodes and their relationships using natural language templates to generate textual representations of subgraphs. Other approaches (Ye et al., 2024; Tang et al.,

2024) employ special tokens to represent nodes and edges, effectively converting graph structures into sequences compatible with LLM processing. However, these methods have notable pitfalls: the textualization process can result in the loss of structural information, and the sequential representations may fail to fully capture the intricate relationships within the graph. Additionally, these methods face scalability challenges due to the token length constraints of LLMs, making them unsuitable for handling large graphs with extensive neighborhood information.

The second category of approaches (Chai et al., 2023; Tang et al., 2024; Liu et al., 2024) focuses on **aligning the representation spaces** of GNNs and LLMs in the semantic domain (Figure 1(b)). For instance, certain methods (Xia et al., 2024; Huang et al., 2023, 2024; Guo et al., 2025) project GNN-generated node embeddings into the embedding space of LLMs to achieve semantic consistency. Other techniques, such as those employing attention mechanisms (Ying et al., 2021; Kuang et al., 2022), integrate graph structure information directly into the language model’s representations. While these approaches improve the integration of graph and language modalities, challenges remain. The alignment process may not be optimal, leading to performance bottlenecks in tasks requiring a precise understanding of graph structures and language semantics. Moreover, such methods often suffer from a lack of interpretability, making it difficult to elucidate how the model leverages graph information to make decisions or derive outputs.

To address the limitations of existing approaches, we propose **ExGLM (Explainable Graph Language Model)**, a novel framework designed to effectively and interpretably integrate graph structures with LLMs. Our framework introduces a **graph-language synergistic alignment module** to achieve semantic consistency between graph structures and LLM outputs, while also maintaining interpretability. Specifically, we assign a textual attribute to each node in the graph, describing its adjacent relationships, with different nodes represented by special tokens. We then perform reasoning using the LLM and enhance its representation by incorporating the graph representation into the hidden state. To further improve interpretability, we propose a **judge-and-improve paradigm** where the LLM evaluates and selects responses with better interpretability. These optimized responses are subsequently used to refine the GNN-LLM model.

Our main contribution can be summarized as follows:

- We propose a novel graph-language synergistic alignment module that effectively bridges the gap between graph-structured data and LLM outputs, ensuring robust semantic consistency across modalities.
- We propose a judge-and-improve paradigm, enabling the model to iteratively evaluate and refine its responses for enhanced interpretability and generation quality, thereby improving both performance and transparency.
- We conduct comprehensive experiments on multiple datasets, demonstrating the superior performance and effectiveness of our approach compared to existing methods.

2 Related Work

2.1 Graph-Large Language Models

LLMs (OpenAI et al., 2024; Yang et al., 2024; DeepSeek-AI et al., 2025) achieve state-of-the-art performance on various natural language tasks, however, it lacks explicit mechanisms to effectively model structured information, such as graphs. To address this limitation, recent studies (Shu et al., 2024; Tang et al., 2024) have explored ways to integrate the benefits of GNNs into LLM-based frameworks. For instance, Zhang et al. (2020) adapts the self-attention mechanism of BERT (Devlin et al., 2019) to capture the relational structure of nodes and edges within a graph. However, its performance is highly dependent on the presence and quality of node features, which may limit its applicability when such features are sparse or noisy. InstructGLM (Ye et al., 2024) leverages the natural language modeling capabilities of LLMs to describe multi-scale geometric structures within graphs, thereby improving representation and analysis of graph data. Nonetheless, it suffers from token-length limitations, making it challenging to process large graphs with extensive neighborhood information. Jin et al. (2024) propose a framework named Graph-COT that enhances LLMs by encouraging them to perform iterative reasoning over graph structures. However, fine-tuning LLMs within this framework remains challenging, and potential misalignment between the graph structure and the text attribution can lead to inaccuracies. Another recent work, PromptGFM (Zhu et al., 2025),

explicitly prompts LLMs to mimic the workflow of GNNs within the text space, achieving naturally alignment between graph representations and textual modeling. While this approach improves graph-text integration, it struggles to differentiate between graphs with similar semantic structures. In this work, we propose a novel graph-language synergistic alignment module that aligns GNNs and LLMs at both the text attribution and semantic representation levels. This alignment enables seamless and effective incorporation of the strengths of GNNs and LLMs.

2.2 Self-Judge-and-Improve Paradigm

The self-judge-and-improve paradigm highlights the capacity of LLMs to autonomously evaluate and enhance their own performance and capabilities, thereby reducing dependence on external supervision. This approach enables models to internally refine their understanding and outputs. For instance, Self-Insturct (Wang et al., 2022) embodies this paradigm through a two-step process to improve instruction-following abilities. First, the model generates sample outputs and evaluates them using its internal mechanisms, filtering out sub-optimal results. These filtered samples are then leveraged to fine-tune the model. Similarly, Self-Refine (Madaan et al., 2023) demonstrates how LLMs can provide feedback on their own generations and use this feedback to optimize their outputs iteratively. Expanding on this concept, Yuan et al. (2025) introduced self-rewarding language models, wherein LLMs assign self-generated rewards to their outputs. Preference pairs selected based on these rewards are then utilized to optimize the models using DPO (Rafailov et al., 2023). While these approaches effectively minimize external intervention, the quality of self-judgment is inherently constrained by the performance of the LLM. To address this limitation, we propose the Judge-and-Improve paradigm, which incorporates a superior language model to evaluate the generated outputs. By introducing an external judgment mechanism, our approach enhances the reliability and accuracy of evaluations, enabling more effective refinement of the model’s outputs.

3 Method

The training framework of our method, illustrated in Figure 2, is composed of two key modules: (1) Graph-language synergistic alignment mod-

ule and (2) Judge-and-improve paradigm. The graph-language synergistic alignment module ensures effective integration between the GNNs and the LLMs by aligning textual attributes and semantic representations, thereby maintaining consistency across modalities. The judge-and-improve paradigm operates in two stages: first, it generates and selects accurate and explainable results through prompting, creating a supervised fine-tuning (SFT) and preference dataset; second, it uses these two datasets to optimize the model, progressively enhancing both performance and interpretability.

3.1 Problem Setup

Graph structure. Generally, a graph can be formally defined as $G = (V, E, X)$, where $V = \{v_1, v_2, \dots, v_n\}$ represents the set of nodes, $E \subseteq V \times V$ represents the set of edges, encoding pairwise relationships between nodes, and $X \in \mathbb{R}^{n \times d}$ is the node feature matrix. Each $x_i \in \mathbb{R}^d$ corresponds to the feature vector of node v_i , where d represents the dimensionality of the node features.

Node classification with LLM. Consider a node classification problem over a graph $G = (V, E, X)$, where the goal is to assign one of k discrete class labels to each node. Let $Y = \{1, 2, \dots, k\}$ denote the set of class labels. The training data consists of labeled examples (x_i, y_i) , where $x_i \in \mathbb{R}^d$ represents the graph feature vector of node $v_i \in V$, and $y_i \in Y$ is the corresponding class label. The objective is to learn a classifier $f : X \xrightarrow{G} Y$, such that $f(x_i)$ accurately predicts the class label y_i for each node. In this work, we first derive textual attributions T_v of each node v , capturing its structural and feature information in a textual format. We then leverage both the graph structure and an LLM to perform reasoning. Consequently, the classification objective is refined to learning a classifier $f : T \xrightarrow{G-LLM} Y$ where T represents the textual descriptions derived from the graph’s structural and feature information. This approach integrates the representational strengths of both GNNs and LLMs, enabling a more interpretable and semantically rich node classification framework.

Classification with interpretability. In real-world applications where interpretability is paramount, it is essential for classification models to not only make accurate decisions but also provide clear explanations for those decisions. Therefore, our ultimate goal is to train a classifier that not only performs classification tasks but also generates an-

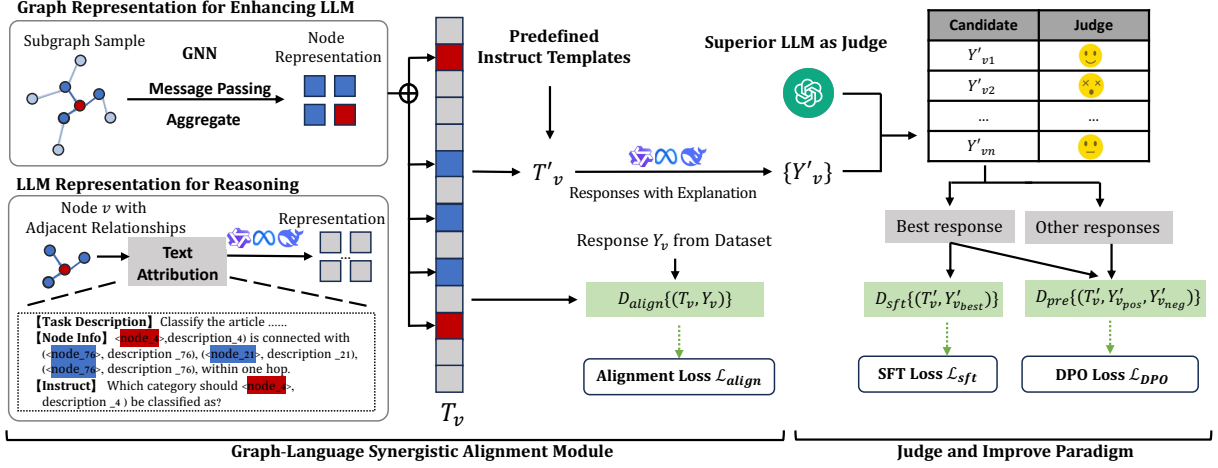


Figure 2: The training framework of ExGLM.

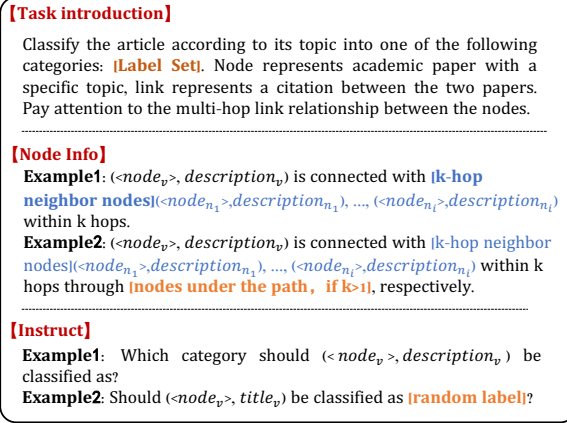


Figure 3: Example of text attribution.

analytical content to elucidate its decision-making process. This can be represented as a function $f: X \xrightarrow{G} Y, Analysis$, where *Analysis* provides the explanatory content.

3.2 Graph-Language Synergistic Alignment Module

To effectively leverage both structural information from graphs and textual attributes from LLMs, we integrate GNNs and LLMs to obtain node representations. To bridge the gap between these two modalities, we propose a graph-language synergistic alignment module. This module consists of two core components: (1) Textual attribution of adjacency relationships, which captures the textual representation of graph structures. (2) Incorporating graph semantic information into textual attribution, which enriches textual descriptions with graph-based semantics. We detail these components below.

Textual attribution of adjacency relationships.

We derive the textual attribution of each node through a two-step process: (1) Subgraph sampling for node information. In the context of large graphs, subgraph sampling is crucial to mitigate computational complexity and enable scalable processing. In this work, we adopt a k -hop sampling strategy to extract localized subgraphs centered around each node. Specifically, for a central node v , we sample its neighbors within k hops, and represent it as $\mathcal{N}_v^{(n)}$. v and $\mathcal{N}_v^{(n)}$ are then further utilized to derive the textual attribution of adjacency relationships.

(2) Node description via text. For each central node v , we construct multiple text descriptions. Each description is represented as a tuple T_v : (task introduction, <node info>, instruction). Here, task introduction provides a brief overview of the task, <node info> contains textual descriptions of the central node's neighbors sampled from i -hop neighborhoods (where $0 < i \leq k$, selected randomly), and instruction specifies a task-related question tailored to the node and its neighborhood information. Specifically, in the <node info> part, each individual graph node is represented by a special token $node_i$ with its brief text descriptions $description_i$. Detailed examples are provided in Figure 3. This approach enables the attribution of each graph node to be naturally expressed in textual form, bridging the structural information of graphs with the representational capabilities of LLMs.

Incorporating graph semantic information into textual attribution. Since the aforementioned special tokens for each node cannot be effectively represented by LLMs alone, we integrate them with representations derived from GNNs. To obtain

the representations of the GNNs, we adopt GraphSAGE, which primarily consists of three steps: neighbor sampling, message aggregation, and node updating. For each node v , we sample its m -hop neighbors and denote them as $\mathcal{N}^{(m)}(v)$. In this paper, we set $m = 2$ in all scenarios. For the aggregation representation is calculated via mean-pooling of neighborhood features:

$$h_{\text{agg}}^{(l)} = \frac{1}{|\mathcal{N}^{(m)}(v)| + 1} \left(h_v^{(l-1)} + \sum_{u \in \mathcal{N}^{(m)}(v)} h_u^{(l-1)} \right) \quad (1)$$

in which $h_v^{(l-1)}$ denotes the representation of node v at layer $(l-1)$. Then each node is updated via Non-linear projection with learnable parameters which is denoted as:

$$h_v^{(l)} = \sigma \left(W^{(l)} \cdot h_{\text{agg}}^{(l)} \right), \quad (2)$$

$W^{(l)}$ is the layer-specific weight matrix, $\sigma(\cdot)$ denotes the ReLU activation function, $d^{(l)}$ is the dimensionality at layer l .

After obtaining $h_v^{(l)}$, we directly add it to the LLM’s hidden states corresponding to the special token v which is shown in Figure 2 left.

To achieve better alignment between the LLM and GNN in the semantic space, we perform joint training. First, we construct a dataset $\mathcal{D}_{\text{align}} = \{(T_v, Y_v)\}, v \in V$, where $Y_v = [c_1, c_2, \dots, c_n]$ denotes the label sequence associated with node v . The alignment is achieved using the Negative Log-Likelihood (NLL) loss function:

$$\mathcal{L}_{\text{align}} = - \sum_{t=1}^n \log P(c_t | c_{<t}, T_v; \theta_{\text{LLM}}, \theta_{\text{GNN}}), \quad (3)$$

where θ_{LLM} denotes the parameters of the LLM, and θ_{GNN} denotes the parameters of the GNN encoder.

3.3 Judge and Improve

Building upon the dual-projection constrained mechanism, we achieve a deep collaboration between Graph Neural Networks (GNNs) and Large Language Models (LLMs). Beyond mere decision-making, providing reasonable and trustworthy analyses significantly enhances the interpretability of these decisions, which is crucial for various real-world applications. To ensure the interpretability of model decisions, we require the LLMs to not only generate accurate answers but also provide comprehensive explanations for their decisions.

However, we have observed that the explanations generated by the LLMs are often suboptimal, indicating a need for further training. Considering the challenge of obtaining training data with annotated explanations, we adopt a judge-and-improve paradigm (Yuan et al., 2025) to enhance the interpretability of the LLMs. Specifically, our approach involves the following steps:

(1) Generating multiple responses: For a given input, the LLM generates multiple responses, each accompanied by an explanation.

(2) Judging quality: Superior LLM acts as a judge to evaluate these responses, selecting the one that is not only accurate but also provides a reasonable explanation.

(3) Optimizing through annotated data: The generated responses and explanations are then used to optimize the LLM, thereby improving the quality of its explanations.

Responses generation. As illustrated in Figure 3, we construct the text attributes of node v using a tuple T_v (task introduction, <node info>, instruction). To assemble a high-quality and diverse dataset, we first replace the instruction with several predefined instruction templates that convey the same intent, denoted as T'_v . Subsequently, we generate a response Y'_v or each T'_v a set $\{(T'_v, Y'_v)\}$, for all $v \in V$.

Judging quality. We require the superior LLM such as GPT-4 to evaluate the generated responses based on three criteria: correctness of the response, adherence to instructions, and reasonableness of the explanation. If none of the samples meet all three criteria, we repeat the response generation procedure. Ultimately, for each T'_v , we obtain a set of candidates $(Y'_{v_{\text{best}}}, Y'_{v_1}, \dots, Y'_{v_{n-1}})$, where n denotes the number of generated responses.

Optimizing through annotated data. Building upon the generated responses, we construct a supervised fine-tuning (SFT) dataset: $\mathcal{D}_{\text{sft}} = \{(T'_v, Y'_{v_{\text{best}}})\}, v \in V$, which aims to teach the LLM to learn the pattern of the best response. The loss is computed as follows:

$$\mathcal{L}_{\text{sft}} = - \sum_{t=1}^n \log P(c'_t | c'_{<t}, T'_v; \theta_{\text{LLM}}, \theta_{\text{GNN}}), \quad (4)$$

where $Y'_{v_{\text{best}}} = [c'_{\text{best}_1}, \dots, c'_{\text{best}_n}]$ and best_i denotes the i -th token of the best response.

This procedure is trained alongside the alignment process, and the overall loss becomes:

$$\mathcal{L}_{\text{total}} = \lambda_1 \cdot \mathcal{L}_{\text{align}} + \lambda_2 \cdot \mathcal{L}_{\text{sft}} \quad (5)$$

where λ_1 and λ_2 denotes the hyperparameters.

Moreover, to enable the LLM to distinguish between good and bad responses, we construct a preference dataset: $\mathcal{D}_{\text{pre}} = \{(T'_v, Y'_{v_{\text{pos}}}, Y'_{v_{\text{neg}}})\}, v \in V$ where $Y'_{v_{\text{pos}}}$ denotes the best response (positive example), and $Y'_{v_{\text{neg}}}$ denotes any other response (negative example) corresponding to the same input T'_v . This dataset pairs each best response with its corresponding non-optimal responses for every node $v \in V$.

We then utilize the preference dataset to optimize the LLM using DPO loss:

$$\mathcal{L}_{\text{DPO}} = -\mathbb{E}_{(T'_v, Y'_{v_{\text{pos}}}, Y'_{v_{\text{neg}}}) \sim \mathcal{D}_{\text{pre}}} \left[\log \sigma \left(\beta \log \frac{\pi_{\theta}(Y'_{v_{\text{pos}}}|T'_v)}{\pi_{\text{ref}}(Y'_{v_{\text{pos}}}|T'_v)} - \beta \log \frac{\pi_{\theta}(Y'_{v_{\text{neg}}}|T'_v)}{\pi_{\text{ref}}(Y'_{v_{\text{neg}}}|T'_v)} \right) \right], \quad (6)$$

where π_{ref} denotes the reference model, which we adopt as the model before DPO training, and β is a hyperparameter.

4 Experiments

4.1 Experimental Setup

Datasets. We utilize three graph datasets of varying scales: Cora, comprising 2,708 nodes and 5,429 edges (Yang et al., 2016); Pubmed, containing 19,717 nodes and 44,338 edges (Namata et al., 2012); and ogbn-arxiv, consisting of 169,343 nodes and 1,166,243 edges (Hu et al., 2020). For our experiments, we adopt the same dataset partitioning strategy as proposed in (Ye et al., 2024).

Metrics. Following (Namata et al., 2012), we use accuracy as the primary metric to evaluate node classification performance. To assess the interpretability of the generated outputs, we utilize GPT-4 (Brown et al., 2020) as an automated evaluator. Additionally, to ensure a more robust and reliable assessment of interpretability, we complement this with a questionnaire-based survey, which provides valuable human-centered insights (Sperrle et al., 2021).

Baselines. We compare the proposed method against three categories of existing approaches: (1) GNN-based models, including GCN (Kipf and Welling, 2017), GraphSAGE (Hamilton et al., 2018), GAT (Veličković et al., 2018), TransGAT, (Louis et al., 2020) etc.; (2) Transformer-based models, such as Graphormer (Ying et al.,

2021), GT (Dwivedi and Bresson, 2021) and CoarFormer (Kuang et al., 2022); and (3) LLM-based models, such as InstructGLM (Ye et al., 2024).

Implementations. In our implementation, we utilize Llama-7B (Touvron et al., 2023) and Llama3.1-8B-Instruct (Touvron et al., 2023) as the LLM backbones. For Llama3.1-8B-Instruct, when it serves as the backbone for InstructGLM, we make minimal adjustments to the inputs to align with its requirements, such as embedding dialogue templates and mapping node IDs to token IDs. During the data generation and annotation phase, we use the Qwen2.5-7B-Instruct model (Bai et al., 2023) to generate decision-analysis content for node classification tasks. To ensure high-quality annotations, we further employ the more powerful Qwen2.5-72B-Instruct model (Bai et al., 2023) as a "super annotator," automatically refining and validating the generated analyses. The model training is performed on 8 A100 GPUs, and all experiments are conducted over 1-3 epochs.

4.2 Performance Comparison

Tables 1 compares the performance of various models on the Cora and PubMed datasets, showcasing the effectiveness of different approaches.

Accuracy on Cora dataset. Among the GNN-based methods, ACM-GCN+ achieves the best accuracy on the Cora dataset (89.75%). Transformers-based methods, on the other hand, generally exhibit relatively lower performance. Notably, the hybrid InstructGLM approach, which combines GNN and LLM techniques, is the most comparable to our method, achieving competitive performance with an accuracy of 87.08% on Cora. In contrast, our proposed method achieves 88.8% accuracy, surpassing all existing Transformers-based and GNN-LLM-based approaches.

Accuracy on PubMed dataset. On the PubMed dataset, InstructGLM sets a strong baseline with the best performance among prior methods. Our method outperforms all baselines, achieving a new state-of-the-art accuracy of 94.6%. These results highlight the superiority of our training framework.

Accuracy on Ogbn-Arxiv dataset. Table 2 summarizes the performance of various models on the Ogbn-Arxiv dataset. Among traditional GNN-based approaches, DRGAT achieves the highest accuracy at 76.11%, outperforming simpler architectures such as GraphSAGE (74.35%) and GAT (74.15%), which exhibit moderate performance. Notably, methods that integrate large language

Table 1: Accuracy on Cora and PubMed datasets.

Method	Type	Cora (%)	PubMed (%)
MixHop	GNN	75.65	90.04
GAT	GNN	76.70	83.28
Geom-GCN	GNN	85.27	90.05
SGC-v2	GNN	85.48	85.36
GraphSAGE	GNN	86.58	86.85
GCN	GNN	87.78	88.90
BernNet	GNN	88.52	88.48
FAGCN	GNN	88.85	89.98
GCNII	GNN	88.93	89.80
RevGAT	GNN	89.11	88.50
Snowball-V3	GNN	89.59	91.44
ACM-GCN+	GNN	89.75	90.96
Graphormer	Transformers	80.41	88.24
GT	Transformers	86.42	88.75
CoarFormer	Transformers	88.69	89.75
InstructGLM	GNN-LLM	87.08	93.84
ExGLM	GNN-LLM	88.8	94.6

models (LLMs) with GNN frameworks surpass all conventional GNN models, demonstrating the potential of combining structured graph data with the rich semantic understanding of LLMs. For instance, InstructGLM achieves an accuracy of 76.42%, further highlighting the effectiveness of this hybrid approach. Our proposed method achieves the highest overall accuracy at 77.4%, setting a new state-of-the-art performance on this task. This result underscores the advantages of our framework in effectively leveraging both graph structures and textual information to improve predictive performance.

Table 2: Accuracy on Ogbn-Arxiv dataset.

Method	Type	Accuracy (%)
GAT	GNN	74.15
GraphSAGE	GNN	74.35
GCN	GNN	73.29
AGDN	GNN	76.02
RvGAT	GNN	75.90
DRGAT	GNN	76.11
InstructGLM	GNN-LLM	76.42
ExGLM	GNN-LLM	77.4

Accuracy with Other LLMs. Table 3 compares the performance of our proposed method against InstructGLM on the Cora and PubMed datasets, utilizing two different LLM backbones: LLaMA and LLaMA3. Two key observations can be drawn from the results: (1) Our method consistently outperforms the baseline InstructGLM across both datasets, regardless of the underlying LLM backbone. This demonstrates the robustness and ef-

Table 3: Performance comparison with different LLMs.

Method	Cora (%)	PubMed (%)
InstructGLM (LLaMA)	87.08	93.84
Ours (LLaMA)	88.8	94.6
InstructGLM (LLaMA3)	88.01	94.17
ExGLM (LLaMA3)	89.30	94.42

fectiveness of our approach. (2) The use of a more advanced backbone does not always guarantee a significant performance improvement. While both methods perform slightly better with LLaMA3 compared to LLaMA, the relative gain is marginal. Notably, when applying LLaMA3, the performance on the PubMed dataset drops slightly from 94.6% to 94.42%. This indicates that the integration mechanism and model design play a more critical role than simply using a stronger LLM.

GPT-4 evaluation. To evaluate the different methods more comprehensively, we use GPT-4 as a proxy for human judgment. Specifically, we task GPT-4 with performing pairwise evaluations to select the better response based on three key criteria: correctness of the response, adherence to instructions, and reasonableness of the explanation. The evaluation results, presented in Table 4, demonstrate that our method outperforms InstructGLM on both datasets. The low performance of InstructGLM may be attributed to its overfitting on the dataset, which can lead to less fluent or less adaptable language generation. Additionally, the integration of DPO enhances overall performance on both datasets.

Table 4: GPT-4 evaluation results with LLaMA3 as base model.

ExGLM vs.	Dataset	Win (%)	Lose (%)
InstructGLM	Cora	81.61	0
	PubMed	92.45	0
ExGLM (w/o DPO)	Cora	8.46	6.80
	PubMed	5.81	4.42

4.3 Interpretability

DPO influence for accuracy. In the judge-and-improve paradigm, DPO is utilized to prioritize generations that exhibit better interpretability. However, it remains essential to evaluate how this prioritization affects reasoning accuracy. The results presented in Table 5 demonstrate that enhancing interpretability does not compromise accuracy

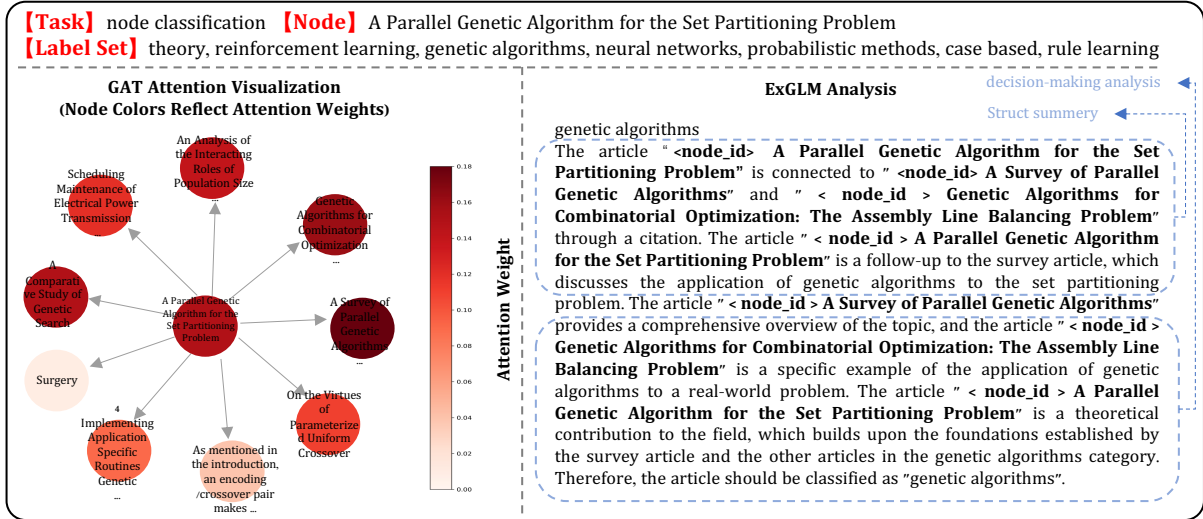


Figure 4: A show case of explanation provided by (left) GNN-based method (2) ExGLM.

and may even lead to improvements in reasoning performance.

Table 5: Ablation study of DPO with LLaMA as base model.

Method	Cora (%)	PubMed (%)
ExGLM w/o dpo	88.92	94.85
ExGLM	90.03	94.75

A showcase. We present a showcase in Figure 4 to illustrate the interpretability of our method in comparison with GNN-based approaches. While GNN-based methods provide explanations for their reasoning through attention weights, these weights may not accurately capture the underlying inference process and can be challenging for humans to interpret. In contrast, our method generates natural language explanations directly, thereby enhancing comprehensibility and interpretability.

Human evaluation. We aim to evaluate whether the use of DPO in the judge-and-improve paradigm enhances interpretability. However, assessing interpretability is challenging due to the lack of a standardized metric. To address this, we conducted a human evaluation. Specifically, we designed a questionnaire involving 20 human participants, each answering 20 questions. Participants were asked to select the response they deemed more interpretable based on three key criteria: coherency, logical consistency, and factuality. The results of this evaluation, presented in Table 6, demonstrate the effectiveness of our approach. The baseline InstructGLM suffers from overfitting on the training dataset, which harms its language generation capa-

bilities and limits its ability to provide meaningful explanations.

Table 6: Human evaluation results with LLaMA3 as base model.

ExGLM vs.	Dataset	Win (%)	Lose (%)
InstructGLM	Cora	100.00	0
	PubMed	100.00	0
ExGLM (w/o DPO)	Cora	23.25	13.75
	PubMed	9.75	9.50

5 Conclusion

This work investigates how to better leverage Large Language Models (LLMs) for reasoning with structured data. Concretely, we aim to address two main limitations identified in recent studies: cross-modality alignment and interpretability. We propose a novel training framework named **ExGLM**, within which a graph-language synergistic alignment module is introduced to ensure semantic consistency across modalities. Additionally, we introduce a judge-and-improve paradigm that adopts a superior language model to evaluate and select generated responses with better interpretability. The selected data is subsequently utilized to optimize the reasoning model. Experiments across various scenarios demonstrate the effectiveness of our approach, showcasing its potential to advance reasoning with structured data.

6 Limitations

While our work achieves promising results, there are several limitations that warrant attention. First, the effectiveness of the judge-and-improve module depends heavily on the performance of the superior language model used for evaluation. If the evaluating model introduces biases or provides inaccurate assessments, the refinement process may be suboptimal, potentially constraining the overall improvement of the target model’s outputs. Second, the current framework does not implement the judgment-and-improvement process iteratively. Iterative refinement, which involves multiple rounds of evaluation and optimization, could further enhance the quality and robustness of the model’s outputs. However, this remains an unexplored avenue and is left for future work.

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