Dynamic Heterogeneous Graph Attention Neural Architecture Search

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Abstract

Dynamic heterogeneous graph neural networks (DHGNNs) have been shown to be effective in handling the ubiquitous dynamic heterogeneous graphs. However, the existing DHGNNs are hand-designed, requiring extensive human efforts and failing to adapt to diverse dynamic heterogeneous graph scenarios. In this paper, we propose to automate the design of DHGNN, which faces two major challenges: 1) how to design the search space to jointly consider the spatial-temporal dependencies and heterogeneous interactions in graphs; 2) how to design an efficient search algorithm in the potentially large and complex search space. To tackle these challenges, we propose a novel Dynamic Heterogeneous Graph Attention Search (DHGAS) method. Our proposed method can automatically discover the optimal DHGNN architecture and adapt to various dynamic heterogeneous graph scenarios without human guidance. In particular, we first propose a unified dynamic heterogeneous graph attention (DHGA) framework, which enables each node to jointly attend its heterogeneous and dynamic neighbors. Based on the framework, we design a localization space to determine where the attention should be applied and a parameterization space to determine how the attention should be parameterized. Lastly, we design a multi-stage differentiable search algorithm to efficiently explore the search space. Extensive experiments on real-world dynamic heterogeneous graph datasets demonstrate that our proposed method significantly outperforms state-of-the-art baselines for tasks including link prediction, node classification and node regression. To the best of our knowledge, DHGAS is the first dynamic heterogeneous graph neural architecture search method.

1 Introduction

Dynamic heterogeneous graphs are ubiquitous in real-world applications, including social networks, e-commerce networks, academic citation networks, etc. Compared to static homogeneous graphs, dynamic heterogeneous graphs contain richer heterogeneous information represented as node and edge types, and dynamic information like evolving graph structures over time. The modeling of heterogeneity and temporal evolutionary patterns is critical for applications of dynamic heterogeneous graphs including the prediction of future links, node labels and properties.

graph heterogeneous Dvnamic neural networks (DHGNNs) (Hu et al. 2020; Fan et al. 2022; Xue et al. 2020; Li et al. 2020) have recently achieved remarkable progress in mining graph dynamic and heterogeneous information (Huang et al. 2021; Fan et al. 2021; Luo et al. 2020). Despite their success, the existing DHGNNs are all manually designed and therefore suffer from the following problems: (1) The designs of DHGNN architectures require extensive human endeavors and expert knowledge. (2) Since the hand-designed models have a fixed architecture, they are unable to adapt to diverse dynamic heterogeneous graph scenarios. (3) The existing DHGNN architectures consider the heterogeneous and dynamic information rather separately and fail to optimally model the joint and complex interaction of heterogeneous and dynamic information.

In this paper, we propose to automate the design of DHGNN architectures on dynamic heterogeneous graphs using neural architecture search (NAS). NAS has attracted considerable attention in automated machine learning and has shown success in domains including computer vision (Wistuba, Rawat, and Pedapati 2019; Elsken, Metzen, and Hutter 2019). However, tailoring a NAS method for dynamic heterogeneous graphs is non-trivial and faces the following two challenges:

- How to design the suitable search space to jointly consider the complex spatial-temporal dependencies and heterogeneous interactions in graphs?
- How to design a tailored efficient search algorithm in the potentially large and complex search space for dynamic heterogeneous graphs?

To tackle these challenges, we propose a novel Dynamic Heterogeneous Graph Attention Search (**DHGAS**) method. Our proposed method can automatically tailor an optimal DHGNN architecture and adapt to various dynamic heterogeneous graph scenarios. In particular, we first propose a unified dynamic heterogeneous graph attention (DHGA) framework. We enable the model to simultaneously consider heterogeneous neighbors across different time stamps by extending the classic neighborhood to dynamic heterogeneous neighborhood and applying attention to the neighbor-

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hood with node and edge type-aware parameterizations. Second, we propose a localization and parameterization search space based on the DHGA framework. Localization space determines what relation types and time stamps should we apply attentions to, which can customize the connections in dynamic heterogeneous neighborhood. Parameterization space further determines the functions for calculating attentions and what types of nodes and edges and time stamps should share the same parameterization, which can customize the mapping functions in dynamic heterogeneous neighborhood. We show that our proposed search space is general and flexible, and can cover many classic DHGNN architectures as special cases. Besides, it also permits the search algorithm to make a trade-off between performance and computational resources. Lastly, we propose a multistage differentiable search algorithm to efficiently explore the search space. By relaxing the discrete choices to continuous ones in the localization and parameterization space, our proposed method can jointly optimize the architecture choices and the supernet weights in a differentiable manner, providing fast and accurate performance estimation of architecture candidates. To stabilize the training process, we further propose to train the supernet in a multi-stage manner and search the spaces sequentially based on the stage.

Extensive experiments on 5 real-world dynamic heterogeneous graph datasets demonstrate that our proposed method significantly outperforms manually designed and automated state-of-the-art baselines for tasks including link prediction, node classification, and node regression. Detailed ablation studies further verify the effectiveness of our proposed search space and search strategy. The codes are publicly available¹.

In summary, the contributions of our work are as follows:

- We propose Dynamic Heterogeneous Graph Attention Search (**DHGAS**) method for dynamic heterogeneous graphs. To the best of our knowledge, **DHGAS** is the first dynamic heterogeneous graph neural architecture search method.
- We design a localization space and a parameterization space for dynamic heterogeneous graphs based on our unified dynamic heterogeneous graph attention framework. We show that our proposed space covers representative manually designed architectures as special cases.
- We propose a multi-stage differentiable search algorithm for dynamic heterogeneous graphs which can explore our proposed search space effectively and efficiently.
- Extensive experiments on real-world datasets demonstrate the superiority of our method over state-of-the-art manually designed and automated baselines.

2 Notations and Preliminaries

2.1 Dynamic Heterogeneous Graphs

Consider a graph \mathcal{G} with the node set \mathcal{V} and the edge set \mathcal{E} . Nodes are associated with a type mapping function $\phi_n : \mathcal{V} \to \mathcal{C}_n$ and edges are associated with a type mapping function $\phi_e : \mathcal{E} \to \mathcal{C}_e$, where \mathcal{C}_n and \mathcal{C}_e denote the node type set

and the edge type set, respectively. We give a formal definition for dynamic heterogeneous graphs as follows:

Definition 1 A dynamic heterogeneous graph is defined as $\mathcal{G} = (\{\mathcal{G}^t\}_{t=1}^T, \phi_n, \phi_e)$, where T is the number of time stamps, $\mathcal{G}^t = (\mathcal{V}^t, \mathcal{E}^t)$ is the graph slice at time stamp t, $\mathcal{V} = \bigcup_{t=1}^T \mathcal{V}^t$, $\mathcal{E} = \bigcup_{t=1}^T \mathcal{E}^t$, and $|\mathcal{C}_n| + |\mathcal{C}_e| \geq 2$.

Dynamic heterogeneous graphs are rather general data formats to represent relational data in real-world applications. For example, static graphs and homogeneous graphs can be considered special cases of dynamic heterogeneous graphs by setting T = 1 and $|\mathcal{C}_n| + |\mathcal{C}_e| = 2$, respectively.

2.2 Dynamic Heterogeneous Graph Neural Networks

Generally, GNNs follow a message-passing mechanism (Gilmer et al. 2017; Hamilton, Ying, and Leskovec 2017) that each node aggregates information from its neighbors. Specifically, let h_u be the representation of node u. Message-passing GNNs update the node representation by²

$$\mathbf{h}_{u} \leftarrow \text{Update}\left(\mathbf{h}_{u}, \text{Agg}(\{\text{Msg}(\mathbf{h}_{v}) : v \in \mathcal{N}(u)\}\right), \quad (1)$$

where $\mathcal{N}(u) = \{v : (u, v) \in \mathcal{E}\}$ denotes the neighborhood of the node u, $Msg(\cdot)$ extracts information from the neighbor node $v \in N(u)$, $Agg(\cdot)$ aggregates the neighborhood information, and Update(\cdot) updates the node representation. Heterogeneous GNNs further consider the heterogeneity of graphs by differentiating node and edge types and assigning different parameters for $Msg(\cdot)$, $Agg(\cdot)$ and $Update(\cdot)$ functions. The message-passing function is:

 $\mathbf{h}_{u} \leftarrow \text{Update}_{\phi(u)}(\mathbf{h}_{u}, \text{Agg}_{r}(\{\text{Msg}_{r}(\mathbf{h}_{v}) : v \in \mathcal{N}_{r}(u), r \in \mathcal{C}_{e}\}),$ (2)

where $\mathcal{N}_r(u) = \{v : (u, v) \in \mathcal{E} \land \phi_e(u, v) = r\}$ is the neighborhood of node u with the relation type r.

DHGNNs further explore the temporal information in dynamic graphs based on Eq. (2). For example, relative time encoding (Hu et al. 2020) encodes time information into edges, i.e., $\mathcal{E}' = \text{Encode}(\{\mathcal{E}^t\}_{t=1}^T)$ followed by heterogeneous message-passings. Another category of DHGNNs adopt sequence-based models to aggregate information from different time slices, i.e., $\mathbf{H} = \text{Seq}(\{\mathbf{H}^t\}_{t=1}^T)$, where \mathbf{H}^t denotes the node representations at time stamp t and **H** is the final node representation. Clearly, these existing approaches handle heterogeneous and dynamic information rather separately in a fixed form. In comparison, our proposed method can jointly aggregate spatial-temporal heterogeneous information and automatically adapt to diverse dynamic heterogeneous graph scenarios.

2.3 Neural Architecture Search

Neural architecture search (NAS) aims at automating the design of neural architectures, which can be formulated as a bi-level optimization problem (Elsken, Metzen, and Hutter

¹https://github.com/wondergo2017/DHGAS

 $^{^{2}}$ To simplify notations, we omit the layer superscript and use arrows to show the message-passing functions in each layer. We also omit edge features, which can be easily incorporated.



Figure 1: The framework of our proposed dynamic heterogeneous graph attention search (**DHGAS**) model. For a given dynamic heterogeneous graph with multiple node and edge types and time slices, **DHGAS** can tailor an optimal architecture based on the unified Dynamic Heterogeneous Graph Attention (DHGA) framework. In particular, **DHGAS** conducts a multi-stage differentiable architecture search on the attention parameterization space and the attention localization space with several carefully designed constraints. In the localization space, we search for what types of edges and which time stamps the attention should be calculated. In the parameterization space, we search for how the attention functions should be parameterized.

2019; Wistuba, Rawat, and Pedapati 2019):

$$a^{*} = \underset{a \in \mathcal{A}}{\arg\min \mathcal{L}_{\text{val}}(a, \mathbf{w}^{*}(a))},$$

s.t. $\mathbf{w}^{*}(a) = \underset{\mathbf{w} \in \mathcal{W}(a)}{\arg\min \mathcal{L}_{\text{train}}(a, \mathbf{w})},$ (3)

where \mathcal{A} is the architecture search space, $\mathcal{W}(a)$ is the parameter space for a given architecture a, and $\mathbf{w}^*(a)$ are the optimal weights for the architecture a. In this paper, we tailor a search space, including attention localization and parameterization, and a multi-stage differentiable search algorithm for dynamic heterogeneous graphs.

3 The Proposed Method

3.1 Dynamic Heterogeneous Graph Attention

The key idea of our proposed dynamic heterogeneous graph attention (DHGA) framework is to unify the spatialtemporal aggregation and jointly integrate dynamic and heterogeneous information from neighborhoods by an attention-based message-passing mechanism. We first extend the neighborhood definition.

Definition 2 Dynamic Heterogeneous Neighborhood: for the neighborhood of each node u, we use subscripts to denote the relation type and superscripts to denote the time stamp, i.e., $\mathcal{N}_r^t(u) = \{v : (u, v) \in \mathcal{E}^t, \phi_e(u, v) = r\}$. With a slight abuse of notations, we use $\mathcal{N}(u)$ to denote all types of neighbors at all time stamps in dynamic heterogeneous graphs, i.e., $\mathcal{N}(u) = \bigcup_{r,t} \mathcal{N}_r^t(u)$. Next, we introduce our tailored message-passing framework to capture the dynamic heterogeneous neighborhood information to update node representations. Following the attention mechanism (Vaswani et al. 2017), for a node u in the time stamp t and its neighbor $v \in \mathcal{N}_r^{t'}(u)$, we calculate the Query-Key-Value vector using a set of mapping functions:

$$\mathbf{q}_{u}^{t} = \mathcal{F}_{q,\phi_{n}(u),t}^{N}(\mathbf{h}_{u}^{t}), \tag{4}$$

$$\mathbf{k}_{v}^{t'} = \mathcal{F}_{k,\phi_{n}(v),t'}^{N}(\mathbf{h}_{v}^{t'}), \tag{5}$$

$$\mathbf{v}_{v}^{t'} = \mathcal{F}_{v,\phi_{n}(v),t'}^{N}(\mathbf{h}_{v}^{t'}),\tag{6}$$

where \mathbf{h}_{u}^{t} denotes the representation of node u at the time stamp t, \mathbf{q} , \mathbf{k} , \mathbf{v} represents the query, key and value vector, respectively, and $\mathcal{F}_{q}^{N}(\cdot)$, $\mathcal{F}_{k}^{N}(\cdot)$, $\mathcal{F}_{v}^{N}(\cdot)$ denote the corresponding node mapping functions. In this paper, we adopt small fully-connected neural networks to instantiate all $\mathcal{F}^{N}(\cdot)$. Notice that the subscripts in the functions indicate that we adopt different functions, i.e., functions with different parameters, based on the node type and time stamps. Then, we calculate the attention score between u and v using a mapping function $\mathcal{F}^{R}(\cdot)$ on the query and key vector:

$$\alpha_{u,v} = \mathcal{F}^R_{\phi_e(u,v),\Delta t}(\mathbf{q}^t_u, \mathbf{k}^{t'}_v), \tag{7}$$

where $\Delta t = t - t'$, i.e., the relation mapping depends on the difference in time stamps instead of the absolute values. Inspired by HGT (Hu et al. 2020) and RGCN (Schlichtkrull et al. 2018), we adopt a relation-aware projection to instantiate $\mathcal{F}^{R}(\cdot)$, i.e.,

$$\mathcal{F}^{R}_{\phi_{e}(u,v),\Delta t}(\mathbf{q},\mathbf{k}) = \frac{\mathbf{q}\mathbf{W}_{\phi_{e}(u,v),\Delta t}\mathbf{k}^{\top}}{\sqrt{d}},\qquad(8)$$

where $\mathbf{W}_{\phi_e(u,v),\Delta t} \in \mathbb{R}^{d \times d}$ denote the learnable parameters for a specific edge type and time stamps and d is the dimensionality. Finally, we normalize the attention scores using the softmax function and aggregate all neighborhoods, i.e.,

$$\mathbf{h}_{u}^{t} \leftarrow \text{Update}(\mathbf{h}_{u}^{t}, \sum_{v \in \mathcal{N}(u)} \hat{\alpha}_{u,v} \mathbf{v}_{v}^{t'})),$$
$$\hat{\alpha}_{u,v} = \frac{\exp(\alpha_{u,v})}{\sum_{v' \in \mathcal{N}(u)} \exp(\alpha_{u,v'})}.$$
(9)

Note that we can easily extend our method into multi-head attention (Vaswani et al. 2017) to stabilize the training process and improve the model's expressiveness. We omit the detailed formulations for brevity.

In summary, we can jointly aggregate information across different types of neighborhoods in all time stamps using one layer of DHGA. Compared to previous works which aggregate spatial and temporal information separately, our proposed method can capture more flexible heterogeneous spatial-temporal graph structures.

Besides, we explicitly consider different types of relations, i.e., node types in the node mapping functions $\mathcal{F}^{N}(\cdot)$ and edges types in the relation mapping functions $\mathcal{F}^{R}(\cdot)$, and different time stamps by setting different parameters. Thus, our proposed DHGA can learn to adaptively assign different attention scores to handle different dynamic heterogeneous graph applications.

Though DHGA is flexible and expressive in modeling dynamic heterogeneous graphs, naively searching architectures based on DHGA can incur high complexity when solving the bi-level optimization in Eq. (3). In the following, we introduce our tailored search space and search algorithm to reduce the complexity while maintaining the expressiveness of the model.

3.2 Attention Localization and Parameterization Search Space

The full version of our proposed DHGA introduced in Section 3.1 calculates attention across all types of neighbors and all time stamps. While being the most expressive architecture, the computation cost is also extensive. To sparsify the attention and enable more lightweight and efficient architectures, we propose the localization space and the parameterization space based on the full DHGA.

Localization Space: Locate where to apply attention. First, we introduce the localization space which determines what types of edges and time stamps should the attention be calculated. In specific, we denote the localization space as $\mathcal{A}^{Lo} = \{0,1\}^{T \times T \times |\mathcal{C}_e|}$. For $\mathbf{A}^{Lo} \in \mathcal{A}^{Lo}$, $\mathbf{A}^{Lo}_{t,t',r}$ denotes whether calculating the representation of node u at the time stamp t should attend to its neighbors $\mathcal{N}_r^{t'}(u)$ in the message-passing. Therefore, \mathbf{A}^{Lo} completely determines where the attention functions are applied. Notice that our proposed localization space is general and flexible since it can cover many existing architectures as special cases. For example, the full DHGA equivalents to every value of \mathbf{A}^{Lo} equals to one. It can also cover other architectures including GAT (Veličković et al. 2018), Temporal self-attention (Fan et al. 2022), Masked temporal selfattention (Xue et al. 2020; Sankar et al. 2020), identity mapping to support skip connections, etc. (please refer to Appendix to the detailed correspondence).

Besides being general and flexible, we also greatly reduce the complexity using the localization space. Specifically, it is easy to see that the full DHGA has a time complexity

$$O(\sum_{t=1}^{T} \sum_{t'=1}^{T} \sum_{r \in \mathcal{C}_e} |\mathcal{E}_r^{t'}|) = O(T^2 |\mathcal{C}_e| \max_{1 \le t \le T, r \in \mathcal{C}_e} |\mathcal{E}_r^{t}|)$$
(10)

In comparison, the time complexity of using \mathbf{A}^{Lo} is:

$$O(\sum_{t=1}^{T}\sum_{t'=1}^{T}\sum_{r\in\mathcal{C}_{e}}\mathbf{A}_{t,t',r}^{Lo}|\mathcal{E}_{r}^{t'}|) = O(|\mathbf{A}^{Lo}|\max_{1\leq t\leq T,r\in\mathcal{C}_{e}}|\mathcal{E}_{r}^{t}|)$$
(11)

where $|\mathbf{A}^{Lo}|$ denotes the number of non-zero values in \mathbf{A}^{Lo} . By constraining the size of $|\mathbf{A}^{Lo}|$, we can reduce DHGA time complexity to be independent of the number of time stamps T and the number of relation types $|\mathcal{C}_e|$.

Parameterization Space: How to parameterize attention. To reduce the number of parameters, we propose an parameterization space to search for how the attention functions should be calculated. Specifically, we denote the parameterization space as $\mathcal{A}^{Pa} = \mathcal{A}^N \times \mathcal{A}^R$, where $\mathcal{A}^N = \{1, ..., K_N\}^{T \times |\mathcal{C}_n|}$ is the parameterization matrix for the node mapping functions $\mathcal{F}^N(\cdot)$ and $\mathcal{A}^R =$ $\{1, ..., K_R\}^{2T \times |\mathcal{C}_e|}$ is the parameterization matrix for the relation mapping functions $\mathcal{F}^R(\cdot)$, and K_N and K_R are two hyper-parameters. In a nutshell, we store K_N mapping functions for $\mathcal{F}_N(\cdot)$ and K_R mapping functions for $\mathcal{F}_R(\cdot)$ as prototypes, and each attention function can choose from the corresponding prototypes. Concretely, let $\mathbf{A}^N \in \mathcal{A}^N$. $\mathbf{A}_{t,c}^N = k$ indicates that the node mapping function node u with $\phi_n(u) = c$ and time stamp t, i.e., $\mathcal{F}_{q,c,t}^N(\cdot)$, $\mathcal{F}_{k,c,t}^N(\cdot)$, $\mathcal{F}^N(\cdot)$. Similarly, for $\mathbf{A}^R \in \mathcal{A}^R$, $\mathbf{A}_{\Delta t,c}^R = k$ denotes that the relation mapping function $\mathcal{F}_{c,\Delta t}^R$ for relation $\phi_e(u, v) = c$ should choose the k^{th} prototype for $\mathcal{F}^R(\cdot)$.

Using the parameterization space, our proposed method can flexibly determine which mapping functions, including both node mapping functions and relation mapping functions, should share parameters. Intuitively, some node types, relation types, or time stamps share similar patterns and therefore can enjoy parameter sharing without affecting the performance. Since these patterns may depend on specific dynamic heterogeneous graph datasets and tasks, we propose to search and learn these patterns adaptively instead of manually setting the parameter sharing rules.

Similar to the localization space, the parameterization space is general and covers diverse existing architectures. For example, when $K_N = T \times |\mathcal{C}_n|$ and $K_R = 2T \times |\mathcal{C}_e|$, we can search for a unique prototype vector for each function and recover the full DHGA. When $K_N = 1$ and $K_R = 1$, we recover the existing homogeneous attention-based GNNs.

Using the parameterization space, we can reduce the number of learnable parameters. It is easy to see that the number of learnable parameters for the full DHGA is of $O(T(|\mathcal{C}_n|+|\mathcal{C}_e|))$. Using the parameterization space, we can reduce it to $O(K_N + K_R)$. When constraining K_N and K_R as constants, the number of learnable parameters is also a constant, i.e., unrelated to the number of edges $|\mathcal{E}|$, the number of time stamps T, or the number of node and edge types $|\mathcal{C}_n|$ and $|\mathcal{C}_e|$.

In short, the localization space and parameterization space balance the model complexity and model expressiveness by determining the edge types and time stamps in calculating the attentions and the parameterization of attentions.

3.3 Multi-Stage Differentiable Search

With our proposed localization space and parameterization space, we introduce our proposed search strategy. Denote the whole search space as $\mathcal{A} = \mathcal{A}^{Lo} \times \mathcal{A}^{Pa}$. It is easy to see that the space can contain up to $2^{T^2|\mathcal{C}_e|}K_N^{T|\mathcal{C}_n|}K_R^{2T|\mathcal{C}_e|}$ possible choices, which is considerably large and it is infeasible to enumerate all possible choices in practice. To reduce the complexity of searching, we first propose heuristic constraints on the search space to remove invalid or ineffective architectures, and then adopt the one-shot neural architecture search algorithm to speed up the search process.

Space Constraint. Inspired by Masked temporal selfattention (Xue et al. 2020; Sankar et al. 2020), we constrain the searched localization to respect the chronological order of graph slices, i.e., the representation of node u at time stamp t can only receive messages from neighborhood $\mathcal{N}_r^{t'}(u)$ with $t' \leq t$. This constraint has clear explanations, since in practice it is infeasible to predict the current situation using future information. Besides, we add another constraint as $|\mathbf{A}_t^{Lo}| \leq K_{Lo}, 1 \leq t \leq T$, where K_{Lo} is a hyperparameter. In this way, we constrain the sparsity of the attention connections in each time slice and reduce the complexity, as shown in paragraph 3.2. Assuming the continuity of mapping functions in the temporal domain, we further break time slices into consecutive patches, where functions within one patch share the same parameters. Lastly, we constrain the last layer of the architecture to only contain connections to the last time slice T so that we can utilize these representations for downstream tasks.

Supernet Construction. Following the recent advancements of NAS (Liu, Simonyan, and Yang 2019; Xie et al. 2018; Guo et al. 2020), we transform the bi-level optimization in Eq. (3) into an one-shot NAS problem using a supernet: Since every possible architecture $a \in \mathcal{A}$ is contained in the supernet, its performance can be quickly evaluated using the corresponding weights in the supernet. Specifically, in the supernet, the categorical choice of a particular operation is relaxed into a softmax overall all possible operations: $\bar{\mathcal{F}}(\mathbf{x}) = \sum_{i=1}^{|\mathcal{A}|} \frac{\exp(\beta_i)}{\sum_{j=1}^{|\mathcal{A}|} \exp(\beta_j)} \mathcal{F}_i(\mathbf{x})$, where **x** is the input, $\bar{\mathcal{F}}(\mathbf{x})$ is the output, $|\mathcal{A}|$ denotes the number of possible operation.

erations, and β_i denotes the mixing weights for the *i*th possible function $\mathcal{F}_i(\cdot)$. For the localization space, operations indicate whether the attention function is applied. For the pa-

rameterization space, operations represent different node/relation prototype mapping functions. Using the supernet, we can jointly optimize the mixing weights β and all parameters in the mapping functions in a differentiable manner:

$$\mathbf{w} \leftarrow \mathbf{w} - \eta_{\mathbf{w}} \frac{\partial \mathcal{L}_{\text{train}}}{\partial \mathbf{w}}, \beta \leftarrow \beta - \eta_{\beta} \frac{\partial \mathcal{L}_{\text{val}}}{\partial \beta},$$
 (12)

where η_{β} and η_{w} are the learning rate for model weights and architecture weights, respectively.

Multi-stage Supernet Training. To stabilize the training of the supernet, we divide the training process into three stages: node parameterization, relation parameterization, and localization. In the node parameterization stage, we force the attention location as fully-connected, and force the relation mapping functions to share the same parameterizations. In the relation parameterization stage, we select and fix the choice in the node parameterization space and focus on searching for the relations. Similarly, in the localization stage, we focus on searching in the localization space while fixing the choices in the other two spaces. When the training is finished, we obtain an optimal architecture by discretizing the operation choices.

4 Experiments

In this section, we evaluate the proposed method through tasks including link prediction, node classification, and node regression. We compare static homogeneous GNNs **GCN** (Kipf and Welling 2017),**GAT** (Veličković et al. 2018) ; static heterogenous GNNs **RGCN** (Schlichtkrull et al. 2018),**HGT** (Hu et al. 2020); dynamic heterogeneous GNNs **DyHATR**(Xue et al. 2020),**HGT+** (Hu et al. 2020),**HTGNN** (Fan et al. 2022) as hand-designed baselines. We also compare to a state-of-the-art static homogeneous graph NAS method **GraphNAS** (Gao et al. 2020) and a heterogeneous graph NAS method **DiffMG** (Ding et al. 2021). More details about the experimental setup can be found in Appendix, including datasets, baselines, training protocol, hyperparameters, task setup, loss functions, etc.

4.1 Main results

Link Prediction. First, we conduct experiments for the link prediction task on two datasets: an academic citation dataset Aminer (Ji et al. 2021) and a recommendation dataset **Ecomm** (Xue et al. 2020). The results are shown in Table 1. We have the following findings. (1) DHGAS achieves the best result on both datasets with a large margin, i.e., improving the AUC by approximately 2.5% and 4% over the most competent baseline, respectively. The results demonstrate that DHGAS can effectively handle the link prediction task on dynamic heterogeneous graph datasets by tailoring the most suitable architecture. (2) DiffMG reports reasonably good results and outperforms most manually designed heterogeneous methods, demonstrating the importance and potentials of automatically designing neural architectures. However, there still exists a large performance gap between DiffMG and our proposed DHGAS, especially in the Ecomm dataset. We attribute this difference to that our Table 1: The overall results for different methods for tasks including link prediction, node classification, and node regression. The evaluation metrics are in parentheses, and $\uparrow(\downarrow)$ means that higher (lower) value indicate better results. The best results are in bold and the second-best results are underlined."-" indicates the method is not applicable.

Task Metric	Link Prediction (AUC%)↑		Node Classification (F1%) \uparrow		Node Regression (MAE) \downarrow
Dataset	Aminer	Ecomm	Yelp	Drugs	COVID-19
GCN	73.84 ± 0.06	77.94 ± 0.22	37.02 ± 0.00	56.43 ± 0.21	846 ± 101
GAT	80.84 ± 0.96	78.49 ± 0.31	35.54 ± 0.00	57.06 ± 0.00	821 ± 91
RGCN	82.75 ± 0.12	82.27 ± 0.51	37.75 ± 0.00	57.97 ± 0.14	833 ± 95
HGT	78.43 ± 1.81	$\overline{81.09\pm0.52}$	34.62 ± 0.00	57.65 ± 0.01	805 ± 88
DyHATR	74.24 ± 2.09	71.69 ± 0.90	34.49 ± 0.16	55.51 ± 0.09	643 ± 36
HGT+	85.60 ± 0.12	76.68 ± 0.85	38.33 ± 0.00	59.09 ± 0.00	-
HTGNN	$\overline{78.08\pm0.80}$	76.78 ± 6.37	36.33 ± 0.07	$\overline{56.24\pm0.34}$	555 ± 34
GraphNAS	81.61 ± 0.98	79.37 ± 0.21	37.73 ± 0.00	57.13 ± 0.52	$\overline{820\pm43}$
DiffMG	85.04 ± 0.30	81.69 ± 0.06	$\underline{38.65\pm0.00}$	58.45 ± 0.15	629 ± 63
DHGAS	$\textbf{88.13} \pm \textbf{0.18}$	$\textbf{86.56} \pm \textbf{0.58}$	$\overline{\textbf{41.99} \pm \textbf{0.18}}$	$\textbf{62.35} \pm \textbf{0.03}$	536 ± 43



Figure 2: Visualization of the searched architecture on Aminer with $K_{Lo} = 8$.

proposed method can effectively jointly capture the temporal and heterogeneous information, as opposed to DiffMG which only models the heterogeneous information. (3) In general, modeling heterogeneous and temporal information are both critical to boosting the performance of manually designed baselines. For example, HGT+, which adopts the relative time encoding technique, reports the second-best result on Aminer. However, HGT+ fails to handle Ecomm and even performs worse than HGT. The results re-validate that different datasets may require different GNN architectures and manually designed methods may fail to adaptively handle such diverse application scenarios.

Node Classification. Next, we compare different methods for the node classification task adopting two datasets: a business review dataset Yelp (Ji et al. 2021) and an e-commerce risk management dataset **Drugs**³. From the results also shown in Table 1, we have the following observations: (1) Our proposed method DHGAS again reports the best results on both datasets by improving the Macro-F1 score by more than 3%. The results demonstrate that we can effectively handle the node classification task for dynamic heterogeneous graphs by automatically designing architectures using DHGAS. (2) The automated baseline DiffMG and manually designed dynamic heterogeneous method HGT+ reports the second-best result on Yelp and Drugs, showing the effectiveness of NAS and importance of capturing dynamic heterogeneous information. Nevertheless, failing to merging the best of two worlds, their performance gap compared to DHGAS is still considerable.

Node Regression. For the node regression task, we adopt an epidemic disease dataset **COVID-19** (Fan et al. 2022). We report the results in Table 1 and observe the following findings: (1) Similar to the other two tasks, **DHGAS** again achieves the best performance. The results demonstrate that **DHGAS** can adaptively handle diverse applications of heterogeneous dynamic graphs. (2) For this task, manually designed dynamic baselines (i.e., DyHATR and HTGNN) greatly outperforms static methods, showing that modeling temporal information is critical to predicting the COVID-19 cases, which is consistent with the literature. (3) Though not considering dynamic information, DiffMG again shows competitive performance, illustrating the great potential of NAS methods. **DHGAS** can fully utilize such potentials by our tailored search space and search algorithm for dynamic heterogeneous graphs.

4.2 Ablation Studies

Search Space. To test the effectiveness of our proposed localization space and parameterization space, we compare the full version with two ablated versions: "**DHGAS** w/o temporal" and "**DHGAS** w/o temporal & heterogeneous". The former removes any attention localization in with different time slices and the latter further forces the method to use the same parameterization for all types of nodes and edges. For simplicity, we only report the results on Aminer when the space constraint hyper-parameters K_{Lo} are set as 20 and 40, while other datasets and settings show similar patterns.

Figure 3 shows that removing temporal connections in localization space and heterogeneous parameterization will reduce the performance of the searched model. The results verify the effectiveness of **DHGAS** in exploiting temporal and heterogeneous information in our tailored search space.

Search algorithm. We verify the design of our proposed multi-stage supernet training. We compare **DHGAS** with random search and DARTS (Liu, Simonyan, and Yang 2019) based on our proposed search space. We report the results on the Aminer dataset when the localization constraint hyperparameters K_{Lo} is chosen from $\{4, 8, 10, 20, 40\}$ while other results indicate similar conclusions.

As shown in Table 4, our proposed search algorithm outperforms DARTS and random search for all localization constraints. In particular, as K_{Lo} grows larger, the performance of all methods increases, showing a clear tradeoff between efficiency and effectiveness. When K_{Lo} is small, i.e., tight localization constraints, **DHGAS** can automatically search important attention locations and maintain impressive performance. In contrast, random search and DARTS fail in these cases.

³Collected from Alibaba.com



Figure 3: The results of ablation study on the localization and parameterization space on the Aminer dataset (AUC%).

4.3 Additional Analyses

The efficiency of the searched architectures. Figure 5 shows that as we gradually increase the computational budget K_{Lo} , **DHGAS** can obtain architectures with better performance. The results show that **DHGAS** can search architectures tailored to the datasets as well as balance the computational budgets and model performance.

Visualization of the searched architecture. Figure 2 visualizes the search architecture on Aminer with $K_{Lo} = 8$, where the letter and number denote the node type and time, respectively, and the colors denote the choices of node and relation mapping functions. It verfies that **DHGAS** can flexibly tailor localizations and mapping functions, demonstrating that our method can automate DHGNN designs and save human endeavors tackling graph heterogenity and dynamics.

More results and visualizations are shown in Appendix.

5 Related Works

Dynamic Heterogeneous Graph Neural Networks. Graph-structured data are ubiquitous in the real-world (Wu et al. 2020; Zhou et al. 2020; Zhang, Cui, and Zhu 2020; Li et al. 2022a,b,c,d, 2021a,b; Zhang et al. 2022c). To generalize the success of GNNs in homogeneous graphs, considerable research attention have been devoted to heterogeneous GNNs (Yang et al. 2020; Wang et al. 2022; Schlichtkrull et al. 2018; Zhang et al. 2019; Wang et al. 2019; Fu et al. 2020; Hu et al. 2020). Some works attempt to consider dynamic information (Skarding, Gabrys, and Musial 2021; Zhu et al. 2022; Yang et al. 2021; Zhang et al. 2022b; Sankar et al. 2020; Wang et al. 2021; Xu et al. 2020; Rossi et al. 2020), and study dynamic heterogeneous graphs (Kazemi et al. 2020; Xue et al. 2022, 2020; Barros et al. 2021; Yuan et al. 2020; Hu et al. 2020; Fan et al. 2022). Despite the success of these existing approaches, they are all manually designed with a fixed architecture. Besides, the spatial and temporal information are processed relatively independently. In comparison, our proposed method can jointly attend dynamic and heterogeneous neighborhoods and automatically adapt to diverse dynamic heterogeneous graph tasks and datasets.

Graph Neural Architecture Search. To automate the design of GNNs, graph NAS has drawn increasing popularity in the last two years (Zhang, Wang, and Zhu 2021), including reinforcement learning based methods (Gao et al. 2020; Zhou et al. 2019; Qin et al. 2021a, 2022b; Zhou et al.



Figure 4: Comparison of the search algorithms on the Aminer dataset using **DHGAS** search space (AUC%).



Figure 5: Comparison of searched architectures under different computational budgets K_{Lo} in terms of inference time and performance on Aminer. **DHGAS**(k) means **DHGAS** with budget $K_{Lo} = k$.

2022; Guan et al. 2021), evolutionary learning based methods (Nunes and Pappa 2020; Li and King 2020; Shi et al. 2022; Guan, Wang, and Zhu 2021; Guan et al. 2022; Zhang et al. 2022a), bayesian optimization based methods (Hou et al. 2021) and differentiable methods (Zhao et al. 2020; Huan, Quanming, and Weiwei 2021; Li et al. 2021c; Cai et al. 2021; Qin et al. 2021b, 2022a) have also been studied. However, all aforementioned works focus on static homogeneous graphs. More relevant to our work, DiffMG (Ding et al. 2021) and HGNAS (Gao et al. 2021) propose to search heterogeneous GNN architectures using meta-paths (Sun et al. 2011) to differentiate node and edge types. However, they cannot capture the temporal information in dynamic graphs. Additionally, AutoSTG (Pan et al. 2021) proposes to search GNN architectures for homogeneous spatialtemporal graphs, neglecting the heterogeneous interactions.

In summary, the existing graph NAS methods cannot fully capture the complex spatial-temporal information in real dynamic heterogeneous graphs. Our proposed **DHGAS** is the first tailored dynamic heterogeneous graph neural architecture search method, to the best of our knowledge.

6 Conclusion

In this paper, we propose a novel Dynamic Heterogeneous Graph Attention Search (**DHGAS**) method to automate the design of DHGNN. We propose a unified dynamic heterogeneous graph attention framework to jointly consider heterogeneous and dynamic neighbors of nodes. Based on the framework, we design a localization space to determine where the attention should be applied and a parameterization space to determine how the attention should be parameterized. We further design a multi-stage differentiable search algorithm to efficiently explore the search space. Extensive experiments on real-world dynamic heterogeneous graph datasets demonstrate the superiority of our method.

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