

Appendix

A DATASET DETAILS

We use 5 node-level graph datasets (Cora, Citeseer (Kipf & Welling, 2017), Ogbn-arxiv (Hu et al., 2020a), Flickr (Zeng et al., 2020), and Reddit (Hamilton et al., 2017)) and 5 graph-level graph datasets (GEOM (Axelrod & Gomez-Bombarelli, 2020), BACE (Wu et al., 2018), BBBP (Martins et al., 2012), ClinTox (Gayvert et al., 2016), and SIDER (Kuhn et al., 2016)). We further provide the statics of datasets in Table A1.

Table A1: Statistics of datasets.

Level	Dataset	# Classes / # Tasks	#Nodes / #Graphs	# Edges	# Features
Node-level	Cora	7	2,708	5,429	1,433
	Citeseer	6	3,327	4,732	3,703
	Ogbn-Arxiv	40	169,343	1,166,243	128
	Flickr	7	89,250	899,756	500
	Reddit	210	232,965	57,307,946	602
Graph-level	BACE	1	1,513	-	-
	BBBP	1	2,039	-	-
	ClinTox	2	1,478	-	-
	Sider	27	1,427	-	-

B SELF-SUPERVISED TEACHER TASKS

In the paper, we utilize different self-supervised task guided models as the way we extract the “universal knowledge” from the original dataset $\mathcal{G}$. Here, we present each model we used.

For node-level classification tasks, we follow AutoSSL (Jin et al., 2022a) and adopt five classic tasks:

- **DGI** (Velickovic et al., 2019): Maximizes the different views’ representations (graph v.s. nodes).
- **CLU** (You et al., 2020b): Predicts pseudo-labels from K -means clustering on node features.
- **PAR** (You et al., 2020b): Predicts pseudo-labels from Metis graph partition (Karypis & Kumar, 1998).
- **PAIRSIM** (Jin et al., 2020): Predicts pairwise feature similarity between nodes.
- **PAIRDIS** (Peng et al., 2020): Predicts the shortest path length between nodes.

In the graph-level classification tasks, we follow WAS (Fan et al., 2024) to adopt 7 classic tasks:

- **AttrMask** (Hu et al., 2020b): Learns the regularities of node/edge attributes.
- **ContextPred** (Hu et al., 2020b): Explores graph structures by predicting the contexts.
- **EdgePred** (Hamilton et al., 2017): Predicts the connectivity of node pairs.
- **GPT-GNN** (Hu et al., 2020c): Introduces an attributed graph generation task to pre-train GNNs.
- **GraphLoG** (Xu et al., 2021): Introduces a hierarchical prototype to capture the global semantic clusters.
- **GraphCL** (You et al., 2020a): Constructs specific contrastive views of graph data.
- **InfoGraph** (Sun et al., 2019): Maximizes the mutual information between the representations of the graph and substructures.

C PROOF OF THEOREM 1

Proof. We aim to prove that:

$$I(\mathbf{Y}_s^s; \mathbf{Y}_s^h) \leq I(\mathcal{P}^T(\mathbf{X}, \mathbf{A}); \mathbf{Y}). \quad (\text{A.1})$$

Since $\mathbf{Y}_s^s$ and $\mathbf{Y}_s^h$ are obtained from $\mathcal{P}^\mathcal{T}(\mathbf{X}, \mathbf{A})$ and $\mathbf{Y}$ through kernel ridge regression—which is a deterministic mapping, they can be expressed as:

$$\mathbf{Y}_s^s = f(\mathcal{P}^\mathcal{T}(\mathbf{X}, \mathbf{A})), \quad (\text{A.2})$$

$$\mathbf{Y}_s^h = g(\mathbf{Y}), \quad (\text{A.3})$$

where f and g are the regression functions.

According to the data processing inequality (Beaudry & Renner, 2012), applying deterministic functions to random variables does not increase mutual information. Therefore, we have:

$$I(\mathbf{Y}_s^s; \mathbf{Y}_s^h) \leq I(\mathcal{P}^\mathcal{T}(\mathbf{X}, \mathbf{A}); \mathbf{Y}). \quad (\text{A.4})$$

□

D TIME COMPLEXITY ANALYSIS

ST-GCond primarily consists of two parts: sampling sub-tasks for meta updating and involving self-supervised tasks to guide condensing. For the former, we can treat them as a composition of bi-level optimization. Following GCond (Jin et al., 2022c), we start with an L -layer GCN, where the large-scale graph has N nodes, the small yet informative condensed graph has m nodes, and the hidden dimension is d . The computation cost for a single task involves a forward pass through the GNN, which is $O(Lm^2d + Lmd)$, and through g_ϕ , which is $O(m^2d^2)$. The inner optimization of kernel ridge regression can be expressed as $O(Nmr^2 + Nm)$ (Wang et al., 2024). Therefore, the single task complexity is $O(Lm^2d + Lmd + m^2d^2 + Nmr^2 + Nm)$. Denoting the split of tasks as t , the complexity for the former part can be shown as $tO(Lm^2d + Lmd + m^2d^2 + Nmr^2 + Nm)$.

For the latter, the calculation process is similar to the former, although we introduce multiple self-supervised models during the condensing stage. Thanks to the benefits of the offline strategy, we only need the extra computation complexity of $kO(LEd + LNd^2)$, where k denotes the number of self-supervised tasks. Therefore, the overall complexity can be expressed as $(t + 1)O(Lm^2d + Lmd + m^2d^2 + Nmr^2 + Nm) + kO(LEd + LNd^2)$. Note that t and k are not set to be too large.

To intuitively demonstrate the efficiency comparison, we present the running time (in seconds) of the proposed ST-GCond and GCond over 50 epochs on a single A100 GPU. Thanks to the efficiency of kernel ridge regression, we avoid the time-consuming triple-level optimization. As a result, our method is empirically 1.14 to 2.17 times faster than the previous GCond method.

Table A2: Comparison of running time of GCond and ST-GCond(in seconds).

Ogbn-arxiv	r=0.05%	r=0.25%	r=0.5%
GCond (Jin et al., 2022c)	217.18	386.71	765.12
ST-GCond	178.27	278.44	399.15

E COMPUTATION RESOURCE

We conduct all experiments with:

- Operating System: Ubuntu 20.04 LTS.
- CPU: Intel(R) Xeon(R) Platinum 8358 CPU@2.60GHz with 1TB DDR4 of Memory.
- GPU: NVIDIA Tesla A100 SMX4 with 40GB of Memory.
- Software: CUDA 10.1, Python 3.8.12, PyTorch (Paszke et al., 2019) 1.7.0.

F PARAMETER SETTING

In our proposed ST-GCond, we mainly need to handle four hyperparameters $\mathbf{l}_r, \alpha, \beta, k$, the actual search space of them are:

lr	0.1, 0.01, 0.001
α	(0.0, 1.0)
β	(0.0, 1.0)
k	5

Note that the search space for α and β may change during training. For simplicity, we use 10 discrete points. The actual time consumption will depend on the Cartesian product of the individual runs.

G ALGORITHMS

G.1 ALGORITHM 1

Algorithm 1 ST-GCond: Self-supervised and Transferable Graph Dataset Condensation

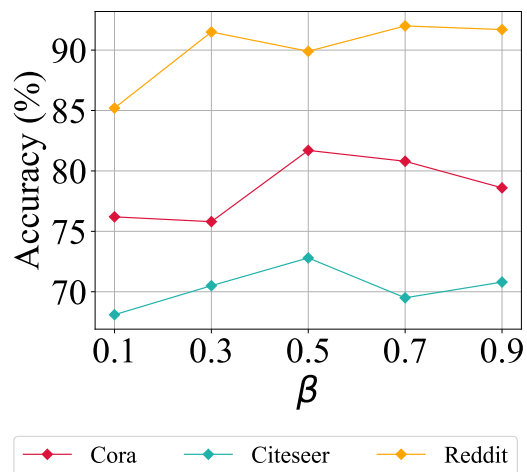
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1: Input: Graph Dataset  $\mathcal{G} = (\mathbf{X}, \mathbf{A}, \mathbf{Y})$ , pre-trained teachers  $\{f_1^T(\cdot), \dots, f_K^T(\cdot)\}$ , steps  $T$ , condensation
   ratio  $r$ .
2: Output: Condensed graph dataset  $\mathcal{G}_s = (\mathbf{X}_s, \mathbf{A}_s, \mathbf{Y}_s^h, \mathbf{Y}_s^s)$ .
3: Initialize weights  $\{\lambda_i = \frac{1}{K}\}_{i=1}^K$ ,  $\mathbf{X}_s$  by selecting  $r\%$  features/class,  $\mathbf{Y}_s^h$  with labels.
4: Initialize  $\mathbf{A}_s = g_\phi(\mathbf{X}_s)$ ,  $\mathbf{Y}_s^s = \frac{1}{K} \sum \lambda_i f_i^T(\mathbf{X}_s, \mathbf{A}_s)$ .
5: for  $t = 0, \dots, T - 1$  do
6:   Initialize  $\theta \sim P_\theta$ .
7:   while not converge do
8:      $D' = 0$ .
9:     Sample tasks  $\mathcal{T}_c \sim p(\mathcal{T}_Y)$ .
10:    for  $\mathcal{T}_i$  do
11:      Sample  $\mathcal{G}^{\mathcal{T}_i} \sim \mathcal{G}, \mathcal{G}_s^{\mathcal{T}_i} \sim \mathcal{G}_s$ .
12:      // Meta-training
13:      Adapt parameters with  $\mathcal{L}_{self}$  on  $\mathcal{G}_s^{\mathcal{T}_i}$ :  $\theta'_i \leftarrow \theta - \lambda_1 \nabla_\theta \mathcal{L}_{self}^{\mathcal{T}_i}(GNN_\theta, \mathcal{G}_s^{\mathcal{T}_i})$ .
14:      // Meta-updating
15:      Combine representation:  $\hat{\mathbf{Y}}^{\mathcal{T}_i} = \sum_{i=1}^K \lambda_i f_i^T(\mathcal{G}^{\mathcal{T}_i})$ .
16:      Compute  $\mathcal{L}_{cls}$ ,  $\mathcal{L}_{self}$ , and  $\mathcal{L}_{MI}$  on  $\mathcal{G}^{\mathcal{T}_i}$ :
17:       $D' \leftarrow D' + \left( \nabla_{\theta'_i} \mathcal{L}_{cls}^{\mathcal{T}_i}(GNN_{\theta'_i}, \mathcal{G}^{\mathcal{T}_i}) + \alpha \nabla_{\theta'_i} \mathcal{L}_{self}^{\mathcal{T}_i}(\hat{\mathbf{Y}}^{\mathcal{T}_i}, \mathcal{G}^{\mathcal{T}_i}) + \beta \nabla_{\theta'_i} \mathcal{L}_{MI}^{\mathcal{T}_i}(\mathbf{Y}_s^s, \mathbf{Y}_s^h) \right)$ .
18:    end for
19:    Update  $\{\lambda_i\}_{i=1}^K$ ,  $\mathbf{X}_s$ ,  $\mathbf{Y}_s^s$ ,  $\phi$ , and  $\theta$ .
20:  end while
21: end for
22: Generate the condensed graph:  $\mathbf{A}_s = \text{ReLU}(g_\phi(\mathbf{X}_s) - \delta)$ ,  $\mathcal{G}_s = (\mathbf{X}_s, \mathbf{A}_s, \mathbf{Y}_s^h, \mathbf{Y}_s^s)$ 

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H MORE EXPERIMENTS

H.1 PARAMETER SENSITIVITY

Figure A1: Sensitive of β