ROS: A GNN-BASED RELAX-OPTIMIZE-AND-SAMPLE FRAMEWORK FOR MAX-*k*-CUT PROBLEMS

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Abstract

The Max-k-Cut problem is a fundamental combinatorial optimization challenge that generalizes the classic \mathcal{NP} -complete Max-Cut problem. While relaxation techniques are commonly employed to tackle Max-k-Cut, they often lack guarantees of equivalence between the solutions of the original problem and its relaxation. To address this issue, we introduce the Relax-Optimize-and-Sample (ROS) framework. In particular, we begin by relaxing the discrete constraints to the continuous probability simplex form. Next, we pre-train and fine-tune a graph neural network model to efficiently optimize the relaxed problem. Subsequently, we propose a sampling-based construction algorithm to map the continuous solution back to a high-quality Max-k-Cut solution. By integrating geometric landscape analysis with statistical theory, we establish the consistency of function values between the continuous solution and its mapped counterpart. Extensive experimental results on random regular graphs and the Gset benchmark demonstrate that the proposed ROS framework effectively scales to large instances with up to 20,000 nodes in just a few seconds, outperforming state-of-the-art algorithms. Furthermore, ROS exhibits strong generalization capabilities across both in-distribution and out-ofdistribution instances, underscoring its effectiveness for large-scale optimization tasks.

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1 INTRODUCTION

The *Max-k-Cut problem* involves partitioning the vertices of a graph into k disjoint subsets in such a way that the total weight of edges between vertices in different subsets is maximized. This problem represents a significant challenge in combinatorial optimization and finds applications across various fields, including telecommunication networks (Eisenblätter, 2002; Gui et al., 2018), data clustering (Poland & Zeugmann, 2006; Ly et al., 2023), and theoretical physics (Cook et al., 2019; Coja-Oghlan et al., 2022). The Max-k-Cut problem is known to be \mathcal{NP} -complete, as it generalizes the well-known *Max-Cut problem*, which is one of the 21 classic \mathcal{NP} -complete problems identified by Karp (2010).

Significant efforts have been made to develop methods for solving Max-k-Cut problems (Nath & 040 Kuhnle, 2024). Ghaddar et al. (2011) introduced an exact branch-and-cut algorithm based on semi-041 definite programming, capable of handling graphs with up to 100 vertices. For larger instances, vari-042 ous polynomial-time approximation algorithms have been proposed. Goemans & Williamson (1995) 043 addressed the Max-Cut problem by first solving a semi-definite relaxation to obtain a fractional so-044 lution, then applying a randomization technique to convert it into a feasible solution, resulting in a 0.878-approximation algorithm. Building on this, Frieze & Jerrum (1997) extended the approach to 046 Max-k-Cut, offering feasible solutions with approximation guarantees. de Klerk et al. (2004) further 047 improved these guarantees, while Shinde et al. (2021) optimized memory usage. Despite their strong 048 theoretical performance, these approximation algorithms involve solving computationally intensive semi-definite programs, rendering them impractical for large-scale Max-k-Cut problems. A variety of heuristic methods have been developed to tackle the scalability challenge. For the Max-Cut 051 problem, Burer et al. (2002) proposed rank-two relaxation-based heuristics, and Goudet et al. (2024) introduced a meta-heuristic approach using evolutionary algorithms. For Max-k-Cut, heuristics such 052 as genetic algorithms (Li & Wang, 2016), greedy search (Gui et al., 2018), multiple operator heuristics (Ma & Hao, 2017), and local search (Garvardt et al., 2023) have been proposed. While these



Figure 1: The Relax-Optimize-and-Sample framework.

heuristics can handle much larger Max-k-Cut instances, they often struggle to balance efficiency and
 solution quality.

075 Recently, machine learning techniques have gained attention for enhancing optimization algo-076 rithms (Bengio et al., 2021; Gasse et al., 2022; Chen et al., 2024). Several studies, including Khalil 077 et al. (2017); Barrett et al. (2020); Chen et al. (2020); Barrett et al. (2022), framed the Max-Cut 078 problem as a sequential decision-making process, using reinforcement learning to train policy net-079 works for generating feasible solutions. However, RL-based methods often suffer from extensive sampling efforts and increased complexity in action space when extended to Max-k-Cut, and hence entails significantly longer training and testing time. Karalias & Loukas (2020) focuses on subset 081 selection, including Max-Cut as a special case. It trains a graph neural network (GNN) to produce a distribution over subsets of nodes of an input graph by minimizing a probabilistic penalty loss 083 function. After the network has been trained, a randomized algorithm is employed to sequentially 084 decode a valid Max-Cut solution from the learned distribution. A notable advancement by Schuetz 085 et al. (2022) reformulated Max-Cut as a quadratic unconstrained binary optimization (QUBO), removing binarity constraints to create a differentiable loss function. This loss function was used to 087 train a GNN, followed by a simple projection onto integer variables after unsupervised training. The 088 key feature of this approach is solving the Max-Cut problem during the training phase, eliminating the need for a separate testing stage. Although this method can produce high-quality solutions 090 for Max-Cut instances with millions of nodes, the computational time remains significant due to the need to optimize a parameterized GNN from scratch. The work of Tönshoff et al. (2022) first 091 formulated the Max-Cut problem as a *constraint satisfaction problem* and then proposed a novel 092 GNN-based reinforcement learning approach. This method outperforms prior neural combinatorial 093 optimization techniques and conventional search heuristics. However, to the best of our knowledge, 094 it is limited to unweighted Max-k-Cut problems. 095

In this work, we propose a GNN-based *Relax-Optimize-and-Sample* (ROS) framework for efficiently solving the Max-k-Cut problem with arbitrary edge weights. The framework is depicted in Figure 1. Initially, the Max-k-Cut problem is formulated as a discrete optimization task. To handle this, we introduce *probability simplex relaxations*, transforming the discrete problem into a continuous one. We then optimize the relaxed formulation by training parameterized GNNs in an unsupervised manner. To further improve efficiency, we apply *transfer learning*, utilizing pre-trained GNNs to warm-start the training process. Finally, we refine the continuous solution using a *random sampling algorithm*, resulting in high-quality Max-k-Cut solutions.

- 104 The key contributions of our work are summarized as follows:
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> • Novel Framework. We propose a scalable ROS framework tailored to the weighted Maxk-Cut problem with arbitrary signs, built on solving continuous relaxations using efficient learning-based techniques.

- **Theoretical Foundations.** We conduct a rigorous theoretical analysis of both the relaxation and sampling steps. By integrating geometric landscape analysis with statistical theory, we demonstrate the consistency of function values between the continuous solution and its sampled discrete counterpart.
 - **Superior Performance.** Comprehensive experiments on public benchmark datasets show that our framework produces high-quality solutions for Max-*k*-Cut instances with up to 20,000 nodes in just a few seconds. Our approach significantly outperforms state-of-the-art algorithms, while also demonstrating strong generalization across various instance types.

¹¹⁷ 2 PRELIMINARIES

119 2.1 MAX-*k*-CUT PROBLEMS

121 Let $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ represent an undirected graph with vertex set \mathcal{V} and edge set \mathcal{E} . Each edge $(i, j) \in \mathcal{E}$ 122 is assigned an arbitrary weight $W_{ij} \in \mathbb{R}$, which can have any sign. A *cut* in \mathcal{G} refers to a partition 123 of its vertex set. The Max-k-Cut problem involves finding a k-partition $(\mathcal{V}_1, \ldots, \mathcal{V}_k)$ of the vertex 124 set \mathcal{V} such that the sum of the weights of the edges between different partitions is maximized.

To represent this partitioning, we employ a k-dimensional one-hot encoding scheme. Specifically, we define a $k \times N$ matrix $X \in \mathbb{R}^{k \times N}$ where each column represents a one-hot vector. The Max-k-Cut problem can be formulated as:

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 $\max_{\boldsymbol{X} \in \mathbb{R}^{k \times N}} \quad \frac{1}{2} \sum_{i=1}^{N} \sum_{j=1}^{N} \boldsymbol{W}_{ij} \left(1 - \boldsymbol{X}_{\cdot i}^{\top} \boldsymbol{X}_{\cdot j} \right) \\
\text{s. t.} \quad \boldsymbol{X}_{\cdot j} \in \{\boldsymbol{e}_{1}, \boldsymbol{e}_{2}, \dots, \boldsymbol{e}_{k}\} \quad \forall j \in \mathcal{V},$ (1)

where $X_{.j}$ denotes the j^{th} column of X, W is a symmetric matrix with zero diagonal entries, and e_l $\in \mathbb{R}^k$ is a one-hot vector with the l^{th} entry set to 1. This formulation aims to maximize the total weight of edges between different partitions, ensuring that each node is assigned to exactly one partition, represented by the one-hot encoded vectors. We remark that weighted Max-k-Cut problems with arbitrary signs is a generalization of classic Max-Cut problems and arise in many interesting applications (De Simone et al., 1995; Poland & Zeugmann, 2006; Hojny et al., 2021).

2.2 GRAPH NEURAL NETWORKS

GNNs are powerful tools for learning representations from graph-structured data. GNNs operate by
 iteratively aggregating information from a node's neighbors, enabling each node to capture increas ingly larger sub-graph structures as more layers are stacked. This process allows GNNs to learn
 complex patterns and relationships between nodes, based on their local connectivity.

At the initial layer (l = 0), each node $i \in \mathcal{V}$ is assigned a feature vector $h_i^{(0)}$, which typically originates from node features or labels. The representation of node *i* is then recursively updated at each subsequent layer through a parametric aggregation function $f_{\Phi^{(l)}}$, defined as:

$$\mathbf{p}_{i}^{(l)} = f_{\mathbf{\Phi}^{(l)}} \left(\mathbf{h}_{i}^{(l-1)}, \{ \mathbf{h}_{j}^{(l-1)} : j \in \mathcal{N}(i) \} \right),$$
(2)

where $\Phi^{(l)}$ represents the trainable parameters at layer $l, \mathcal{N}(i)$ denotes the set of neighbors of node *i*, and $h_i^{(l)}$ is the node's embedding at layer l for $l \in \{1, 2, \dots, L\}$. This iterative process enables the GNN to propagate information throughout the graph, capturing both local and global structural properties.

3 A RELAX-OPTIMIZE-AND-SAMPLE FRAMEWORK

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In this work, we leverage continuous optimization techniques to tackle Max-*k*-Cut problems, introducing a novel ROS framework. Acknowledging the inherent challenges of discrete optimization,
we begin by relaxing the problem to probability simplices and concentrate on optimizing this relaxed
version. To achieve this, we propose a machine learning-based approach. Specifically, we model the

162 relaxed problem using GNNs, pre-training the GNN on a curated graph dataset before fine-tuning 163 it on the specific target instance. After obtaining high-quality solutions to the relaxed continuous 164 problem, we employ a random sampling procedure to derive a discrete solution that preserves the 165 same objective value.

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3.1 PROBABILITY SIMPLEX RELAXATIONS

To simplify the formulation of the problem (1), we remove constant terms and negate the objective 169 170 function, yielding an equivalent formulation expressed as follows:

$$\min_{\boldsymbol{X} \in \mathcal{X}} \quad f(\boldsymbol{X}; \boldsymbol{W}) \coloneqq \operatorname{Tr}(\boldsymbol{X} \boldsymbol{W} \boldsymbol{X}^{\top}), \tag{P}$$

173 where $\mathcal{X} \coloneqq \{ \mathbf{X} \in \mathbb{R}^{k \times N} : \mathbf{X}_{.j} \in \{ \mathbf{e}_1, \mathbf{e}_2, \dots, \mathbf{e}_k \}, \forall j \in \mathcal{V} \}$. It is important to note that the matrix 174 W is indefinite due to its diagonal entries being set to zero. 175

176 Given the challenges associated with solving the discrete problem **P**, we adopt a naive relaxation approach, obtaining the convex hull of \mathcal{X} as the Cartesian product of N k-dimensional probability 177 simplices, denoted by Δ_k^N . Consequently, the discrete problem **P** is relaxed into the following 178 continuous optimization form: 179

$$\min_{\boldsymbol{X} \in \Delta_k^N} \quad f(\boldsymbol{X}; \boldsymbol{W}). \tag{\overline{P}}$$

Before optimizing problem \mathbf{P} , we will characterize its *geometric landscape*. To facilitate this, we 183 introduce the following definition. 184

Definition 1. Let \overline{X} denote a point in Δ_k^N . We define the neighborhood induced by \overline{X} as follows:

$$\mathcal{N}(\overline{X}) \coloneqq \left\{ X \in \Delta_k^N \left| \sum_{i \in \mathcal{K}(\overline{X}_{\cdot j})} X_{ij} = 1, \forall j \in \mathcal{V} \right. \right\},$$

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where $\mathcal{K}(\overline{\mathbf{X}}_{i}) \coloneqq \{i \in \{1, \ldots, k\} \mid \overline{\mathbf{X}}_{ij} > 0\}.$

The set $\mathcal{N}(\overline{X})$ represents a neighborhood around \overline{X} , where each point in $\mathcal{N}(\overline{X})$ can be derived by 192 allowing each non-zero entry of the matrix \overline{X} to vary freely, while the other entries are set to zero. 193 Utilizing this definition, we can establish the following theorem. 194

Theorem 1. Let \overline{X} denote a globally optimal solution to \overline{P} , and let $\mathcal{N}(\overline{X})$ be its induced neighbor-195 hood. Then 196

$$f(\boldsymbol{X}; \boldsymbol{W}) = f(\overline{\boldsymbol{X}}; \boldsymbol{W}), \quad \forall \boldsymbol{X} \in \mathcal{N}(\overline{\boldsymbol{X}})$$

Theorem 1 states that for a globally optimal solution \overline{X} , every point within its neighborhood $\mathcal{N}(\overline{X})$ 199 shares the same objective value as \overline{X} , thus forming a *basin* in the geometric landscape of f(X; W). 200 If $\overline{X} \in \mathcal{X}$ (i.e., an integer solution), then $\mathcal{N}(\overline{X})$ reduces to the singleton set $\{\overline{X}\}$. Conversely, if 201 $\overline{X} \notin \mathcal{X}$, there exist $\prod_{i \in \mathcal{V}} |\mathcal{K}(\overline{X}_{i})|$ unique integer solutions within $\mathcal{N}(\overline{X})$ that maintain the same 202 203 objective value as \overline{X} . This indicates that once a globally optimal solution to the relaxed problem \overline{P} 204 is identified, it becomes straightforward to construct an optimal solution for the original problem P 205 that preserves the same objective value.

206 According to Carlson & Nemhauser (1966), among all globally optimal solutions to the relaxed 207 problem $\overline{\mathbf{P}}$, there is always at least one integer solution. Theorem 1 extends this result, indicating 208 that if the globally optimal solution is fractional, we can provide a straightforward and efficient 209 method to derive its integer counterpart. We remark that it is highly non-trivial to guarantee that the 210 feasible Max-k-Cut solution obtained from the relaxation one has the same quality.

211 **Example.** Consider a Max-Cut problem (k = 2) associated with the weight matrix W. We optimize 212 its relaxation and obtain the optimal solution X^* . 213

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$$W \coloneqq \begin{pmatrix} 0 & 1 & 1 \\ 1 & 0 & 1 \\ 1 & 1 & 0 \end{pmatrix}, X^* \coloneqq \begin{pmatrix} p & 1 & 0 \\ 1 - p & 0 & 1 \end{pmatrix},$$

where $p \in [0, 1]$. From the neighborhood $\mathcal{N}(\overline{X})$, We can identify the following integer solutions that maintain the same objective value.

$$\boldsymbol{X}_{1}^{\star} = \begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 1 \end{pmatrix}, \boldsymbol{X}_{2}^{\star} = \begin{pmatrix} 1 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

Given that $\overline{\mathbf{P}}$ is a non-convex program, identifying its global minimum is challenging. Consequently, the following two critical questions arise.

- Q1. Since solving $\overline{\mathbf{P}}$ to global optimality is \mathcal{NP} -hard, how to efficiently optimize $\overline{\mathbf{P}}$ for highquality solutions?
- **Q2.** Given $\overline{\mathbf{X}} \in \Delta_k^N \setminus \mathcal{X}$ as a high-quality solution to $\overline{\mathbf{P}}$, can we construct a feasible solution $\hat{\mathbf{X}} \in \mathcal{X}$ to \mathbf{P} such that $f(\hat{\mathbf{X}}; \mathbf{W}) = f(\overline{\mathbf{X}}; \mathbf{W})$?

We provide a positive answer to **Q2** in Section 3.2, while our approach to addressing **Q1** is deferred to Section 3.3.

233 3.2 RANDOM SAMPLING

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Let $\overline{X} \in \Delta_k^N \setminus \mathcal{X}$ be a feasible solution to the relaxation $\overline{\mathbf{P}}$. Our goal is to construct a feasible 235 solution $X \in \mathcal{X}$ for the original problem **P**, ensuring that the corresponding objective values are 236 equal. Inspired by Theorem 1, we propose a random sampling procedure, outlined in Algorithm 1. 237 In this approach, we sample each column X_{i} of the matrix X from a categorical distribution char-238 acterized by the event probabilities \overline{X}_{i} (denoted as $\operatorname{Cat}(x; p = \overline{X}_{i})$ in Step 3 of Algorithm 1). 239 This randomized approach yields a feasible solution \hat{X} for **P**. However, since Algorithm 1 incorpo-240 rates randomness in generating \hat{X} from \overline{X} , the value of $f(\hat{X}; W)$ becomes random as well. This 241 raises the critical question: is this value greater or lesser than $f(\bar{X}; W)$? We address this question 242 in Theorem 2. 243

Algorithm 1 Random Sampling 245 246 1: Input: $\overline{X} \in \Delta_{h}^{N}$ \triangleright any feasible solution to \overline{P} 247 2: for i = 1 to N do ▷ each dimension is independent 248 $\hat{X}_{i} \sim \operatorname{Cat}(\boldsymbol{x}; \boldsymbol{p} = \overline{X}_{i})$ 3: ▷ sampling from a categorical distribution 249 4: end for 250 5: Output: $\hat{X} \in \mathcal{X}$ \triangleright a feasible solution to P251

Theorem 2. Let $\overline{\mathbf{X}}$ and $\hat{\mathbf{X}}$ denote the input and output of Algorithm 1, respectively. Then, we have $\mathbb{E}_{\hat{\mathbf{X}}}[f(\hat{\mathbf{X}}; \mathbf{W})] = f(\overline{\mathbf{X}}; \mathbf{W}).$

Theorem 2 states that $f(\hat{X}; W)$ is equal to $f(\overline{X}; W)$ in expectation. This implies that the random sampling procedure operates on a fractional solution, yielding Max-k-Cut feasible solutions with the same objective values in a probabilistic sense. While the Lovász-extension-based method (Bach et al., 2013) also offers a framework for continuous relaxation, achieving similar theoretical results for arbitrary k and edge weights $W_{i,j} \in \mathbb{R}$ is not always guaranteed. In practice, we execute Algorithm 1 T times and select the solution with the lowest objective value as our best result. We remark that the theoretical interpretation in Theorem 2 distinguishes our sampling algorithm from the existing ones in the literature (Toenshoff et al., 2021; Karalias & Loukas, 2020).

3.3 GNN PARAMETRIZATION-BASED OPTIMIZATION

To solve the problem $\overline{\mathbf{P}}$, we propose an efficient learning-to-optimize (L2O) method based on GNN parametrization. This approach reduces the laborious iterations typically required by classical optimization methods (e.g., mirror descent). Additionally, we introduce a "pre-train + fine-tune" strategy, where the model is endowed with prior graph knowledge during the pre-training phase, significantly decreasing the computational time required to optimize $\overline{\mathbf{P}}$. 270 GNN Parametrization. The Max-k-Cut problem can be framed as a node classification task, allow-271 ing us to leverage GNNs to aggregate node features, and obtain high-quality solutions. Initially, we 272 assign a random embedding $h_i^{(0)}$ to each node *i* in the graph \mathcal{G} , as defined in Section 2. We adopt 273 the GNN architecture proposed by Morris et al. (2019), utilizing an L-layer GNN with updates at 274 layer *l* defined as follows: 275

$$\boldsymbol{h}_{i}^{(l)} \coloneqq \sigma \left(\boldsymbol{\Phi}_{1}^{(l)} \boldsymbol{h}_{i}^{(l-1)} + \boldsymbol{\Phi}_{2}^{(l)} \sum_{j \in \mathcal{N}(i)} w_{ji} \boldsymbol{h}_{j}^{(l-1)} \right)$$

279 where $\sigma(\cdot)$ is an activation function, and $\Phi_1^{(l)}$ and $\Phi_2^{(l)}$ are the trainable parameters at layer l for 280 $l \in \{1, \dots, L\}$. This formulation facilitates efficient learning of node representations by leveraging both node features and the underlying graph structure. After processing through L layers of GNN, 282 we obtain the final output $\boldsymbol{H}_{\Phi}^{(L)} := [\boldsymbol{h}_{1}^{(L)}, \dots, \boldsymbol{h}_{N}^{(L)}] \in \mathbb{R}^{k \times N}$. A softmax activation function is then applied in the last layer to ensure $\boldsymbol{H}_{\Phi}^{(L)} \in \Delta_{k}^{N}$, making the final output feasible for $\overline{\boldsymbol{P}}$.

285 "Pre-train + Fine-tune" Optimization. We propose a "pre-train + fine-tune" framework for learn-286 ing the trainable weights of GNNs. Initially, the model is trained on a collection of pre-collected 287 datasets to produce a pre-trained model. Subsequently, we fine-tune this pre-trained model for each 288 specific problem instance. This approach equips the model with prior knowledge of graph structures during the pre-training phase, significantly reducing the overall solving time. Furthermore, it allows 289 for out-of-distribution generalization due to the fine-tuning step. 290

291 The trainable parameters $\mathbf{\Phi} := (\mathbf{\Phi}_1^{(1)}, \mathbf{\Phi}_2^{(1)}, \dots, \mathbf{\Phi}_1^{(L)}, \mathbf{\Phi}_2^{(L)})$ in the pre-training phase are optimized using the Adam optimizer with *random initialization*, targeting the objective 292 293

$$\min_{\boldsymbol{\Phi}} \quad \mathcal{L}_{\text{pre-training}}(\boldsymbol{\Phi}) \coloneqq \frac{1}{M} \sum_{m=1}^{M} f(\boldsymbol{H}_{\Phi}^{(L)}; \boldsymbol{W}_{\text{train}}^{(m)}),$$

where $\mathcal{D} \coloneqq \{ \boldsymbol{W}_{\text{train}}^{(1)}, \dots, \boldsymbol{W}_{\text{train}}^{(M)} \}$ represents the pre-training dataset. In the fine-tuning phase, for a problem instance represented by $\boldsymbol{W}_{\text{test}}$, the Adam optimizer seeks to solve

$$\min_{\boldsymbol{\Phi}} \quad \mathcal{L}_{\text{fine-tuning}}(\boldsymbol{\Phi}) \coloneqq f(\boldsymbol{H}_{\Phi}^{(L)}; \boldsymbol{W}_{\text{test}}),$$

initialized with the pre-trained parameters. 302

303 Moreover, to enable the GNN model to fully adapt to specific problem instances, the pre-training 304 phase can be omitted, enabling the model to be directly trained and tested on the same instance. 305 While this direct approach may necessitate more computational time, it often results in improved performance regarding the objective function. Consequently, users can choose to include a pre-306 307 training phase based on the specific requirements of their application scenarios.

4 **EXPERIMENTS**

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4.1 EXPERIMENTAL SETTINGS 312

313 We compare the performance of ROS against traditional methods and L2O algorithms for solving 314 the Max-k-Cut problem. Additionally, we assess the impact of the "Pre-train" stage in the GNN parametrization-based optimization. The source code is available at https://anonymous. 315 4open.science/r/ROS_anonymous-1C88/. 316

317 **Baseline Algorithms.** We denote our proposed algorithms by ROS and compare them against both 318 traditional algorithms and learning-based methods. When the pre-training step is skipped, we refer 319 to our algorithm as ROS-vanilla. The following traditional Max-k-Cut algorithms are considered 320 as baselines: (i) GW (Goemans & Williamson, 1995): an method with a 0.878-approximation guarantee based on semi-definite relaxation; (ii) BQP (Gui et al., 2018): a local search method designed for 321 binary quadratic programs; (iii) Genetic (Li & Wang, 2016): a genetic algorithm specifically for 322 Max-k-Cut problems; (iv) MD: a mirror descent algorithm that addresses the relaxed problem \mathbf{P} with 323 a convergence tolerance at 10^{-8} and adopts the same random sampling procedure; (v) LPI (Goudet



Figure 2: The computational time comparison of Max-k-Cut problems.

338 et al., 2024): an evolutionary algorithm featuring a large population organized across different is-339 lands; (vi) MOH (Ma & Hao, 2017): a heuristic algorithm based on multiple operator heuristics, 340 employing various distinct search operators within the search phase. (vii) Rank2 (Burer et al., 341 2002): a heuristic based on rank-2 relaxation. For the L2O method, we primarily examine the 342 state-of-the-art baseline: (viii) PI-GNN (Schuetz et al., 2022): A cutting-edge L2O method capable 343 of solving QUBO problems in dozens of seconds, delivering commendable performance. It is the 344 first method to eliminate the dependence on large, labeled training datasets typically required by supervised learning approaches. 345

346 Datasets. The datasets utilized in this paper comprise random regular graphs from Schuetz et al. 347 (2022) and the Gset benchmark from Ye (2003). For the random regular graphs, we employ the 348 random_regular_graph from the NetworkX library (Hagberg et al., 2008) to generate r-349 regular graphs, which are undirected graphs in which all nodes have a degree of r, with all edge weights equal to 1. The Gset benchmark is constructed using a machine-independent graph gen-350 erator, encompassing toroidal, planar, and randomly weighted graphs with vertex counts ranging 351 from 800 to 20,000 and edge densities between 2% and 6%. The edge weights in these graphs are 352 constrained to values of 1, 0, or -1. Specifically, the training dataset includes 500 3-regular graphs 353 and 500 5-regular graphs, each containing 100 nodes, tailored for the cases where k = 2 and k = 3, 354 respectively. The testing set for random regular graphs consists 20 3-regular graphs and 20 5-regular 355 graphs for both k = 2 and k = 3 tasks, with node counts of 100, 1,000, and 10,000, respectively. 356 Moreover, the testing set of Gset encompasses all instances included in the Gset benchmark. 357

Model Settings. ROS is designed as a two-layer GNN, with both the input and hidden dimensions 358 set to 100. To address the issue of gradient vanishing, we apply a graph normalization technique as 359 proposed by Cai et al. (2021). The ROS model undergoes pre-training using the Adam optimizer with 360 a learning rate of 10^{-2} for one epoch. During the fine-tuning stage, the model is further optimized 361 using the same Adam optimizer and learning rate of 10^{-2} . An early stopping strategy is employed, 362 with a tolerance of 10^{-2} and a patience of 100 iterations, terminating training if no improvement 363 is observed over this duration. Finally, in the random sampling stage, we execute Algorithm 1 for 364 T = 100 independent trials and return the best solution obtained.

365 Evaluation Configuration. All our experiments were conducted on an NVIDIA RTX 3090 GPU, 366 using Python 3.8.19 and PyTorch 2.2.0. 367

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> 4.2 PERFORMANCE COMPARISON AGAINST BASELINES

371 4.2.1COMPUTATIONAL TIME 372

373 We evaluated the performance of ROS against baseline algorithms GW, BQP, Genetic, MD, and 374 PI-GNN on random regular graphs, focusing on computational time for both the Max-Cut and Max-375 3-Cut tasks. The experiments were conducted across three problem sizes: N = 100, N = 1,000,and N = 10,000, as illustrated in Figure 2a. Additionally, Figure 2b compares the scalable methods 376 MD, Rank2, and PI-GNN on problem instances from the Gset benchmark with $N \ge 10,000$. "N/A" 377 denotes a failure to return a solution within 30 minutes. A comprehensive summary of the results for

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Table 1: Objective value comparison of Max-k-Cut problems on random regular graphs.

Methods	N=1	.00	N=1,	000	N=10	,000
in como dis	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3
GW	$130.20_{\pm 2.79}$	_	N/A	-	N/A	-
BQP	$131.55_{\pm 2.42}$	$239.70_{\pm 1.82}$	$1324.45_{\pm 6.34}$	2419.15 ± 6.78	N/A	N/A
Genetic	127.55 ± 2.82	$235.50_{\pm 3.15}$	1136.65 ± 10.37	$2130.30_{\pm 8.49}$	N/A	N/A
MD	127.20 ± 2.16	235.50 ± 3.29	1250.35 ± 11.21	2344.85 ± 9.86	12428.85 ± 26.13	$23341.20 {\scriptstyle \pm 32.87}$
PI-GNN	122.75 ± 4.36	-	1263.95 ± 21.59	-	$12655.05_{\pm 94.25}$	-
ROS	128.20 ± 2.82	240.30 ± 2.59	1283.75 ± 6.89	2405.75 ± 5.72	12856.85 ± 26.50	$24085.95 {\scriptstyle \pm 21.88}$

Table 2: Objective value comparison of Max-k-Cut problems on Gset instances.

Methods	G70 (N=	10,000)	G72 (N=	10,000)	G77 (N=	14,000)	G81 (N=	20,000)
methods	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3
MD	8551	9728	5638	6612	7934	9294	11226	13098
Rank2	9529	-	6820	-	9670	-	13662	-
PI-GNN	8956	-	4544	-	6406	-	8970	-
ROS	8916	9971	6102	7297	8740	10329	12332	14464

all Gset instances on Max-Cut and Max-3-Cut, including comparisons with state-of-the-art methods LPI and MOH, is presented in Table 3 and Table 4 in the Appendix.

The results depicted in Figure 2a indicate that ROS efficiently solves all problem instances within 402 seconds, even for large problem sizes of N = 10,000. In terms of baseline performance, the 403 approximation algorithm GW performs efficiently on instances with N = 100, but it struggles with 404 larger sizes such as N = 1,000 and N = 10,000 due to the substantial computational burden 405 associated with solving the underlying semi-definite programming problem. Heuristic methods such 406 as BQP and Genetic can manage cases up to N = 1,000 in a few hundred seconds, yet they 407 fail to solve larger instances with N = 10,000 because of the high computational cost of each 408 iteration. Notably, MD is the only method capable of solving large instances within a reasonable 409 time frame; however, when N reaches 10,000, the computational time for MD approaches 15 times 410 that of ROS. Regarding learning-based methods, PI-GNN necessitates retraining and prediction for 411 each test instance, with test times exceeding dozens of seconds even for N = 100. In contrast, ROS 412 solves these large instances in merely a few seconds. Throughout the experiments, ROS consistently completes its tasks in under 10 seconds, requiring only 10% of the computational time utilized by 413 PI-GNN. Figure 2b illustrates the results for the Gset benchmark, where ROS efficiently solves the 414 largest instances in just a few seconds, while other methods, such as Rank2, take tens to hundreds 415 of seconds for equivalent tasks. Remarkably, ROS utilizes only about 1% of the computational time 416 required by PI-GNN. 417

418 419 4.2.2 OBJECTIVE VALUE

We also evaluate the performance of ROS on random regular graphs and the Gset benchmark concerning the objective values of Problem (1). The results for the random regular graphs and Gset are presented in Tables 1 and 2, respectively. Note that "–" indicates that the method is unable to handle Max-k-Cut problems.

424 The results for random regular graphs, presented in Table 1, indicate that ROS effectively addresses 425 both k = 2 and k = 3 cases, producing high-quality solutions even for large-scale problem in-426 stances. In contrast, traditional methods such as GW and the L2O method PI-GNN are restricted to 427 k = 2 and fail to generalize to the general k, i.e., k = 3. While GW achieves high-quality solu-428 tions for the Max-Cut problem with an instance size of N = 100, it cannot generalize to arbitrary k without integrating additional randomized algorithms to yield discrete solutions. Similarly, the 429 L2O method PI-GNN cannot manage k = 3 because the Max-k-Cut problem cannot be modeled 430 as a QUBO problem. Furthermore, its heuristic rounding lacks theoretical guarantees, which results 431 in sub-optimal performance regarding objective function values. Traditional methods such as BQP



Figure 3: The ratio of computational time and objective value comparison of Max-*k*-Cut problems between ROS-vanilla and ROS.

and Genetic can accommodate k = 3, but they often become trapped in sub-optimal solutions. Among all the baselines, only MD can handle general k while producing solutions of comparable quality to ROS. However, MD consistently exhibits inferior performance compared to ROS across all experiments. The results for the Gset benchmark, shown in Table 2, offer similar insights: ROS demonstrates better generalizability compared to the traditional Rank2 method and the L2O method PI-GNN. Moreover, ROS yields higher-quality solutions than MD in terms of objective function values.

4.3 EFFECT OF THE "PRE-TRAIN" STAGE IN ROS

To evaluate the impact of the pre-training stage in ROS, we compared it with ROS-vanilla, a variant that omits the pre-training phase (see Section 3.3). We assessed both methods based on objective function values and computational time. Figure 3 illustrates the ratios of these metrics between ROS-vanilla and ROS. In this figure, the horizontal axis represents the problem instances, while the left vertical axis (green bars) displays the ratio of objective function values, and the right vertical axis (red curve) indicates the ratio of computational times.

462 As shown in Figure 3a, during experiments on regular graphs, ROS-vanilla achieves higher 463 objective function values in most settings; however, its computational time is approximately 1.5 464 times greater than that of ROS. Thus, ROS demonstrates a faster solving speed compared to 465 ROS-vanilla. Similarly, in experiments conducted on the Gset benchmark (Figure 3b), ROS 466 reduces computational time by around 40% while maintaining performance comparable to that of 467 ROS-vanilla. Notably, in the Max-3-Cut problem for the largest instance, G81, ROS effectively halves the solving time, showcasing the significant acceleration effect of pre-training. It is worth 468 mentioning that the ROS model was pre-trained on random regular graphs with N = 100 and gen-469 eralized well to regular graphs with N = 1,000 and N = 10,000, as well as to Gset problem 470 instances of varying sizes and types. This illustrates ROS's capability to generalize and accelerate 471 the solving of large-scale problems across diverse graph types and sizes, emphasizing the strong 472 out-of-distribution generalization afforded by pre-training. 473

In summary, while ROS-vanilla achieves slightly higher objective function values on individual instances, it requires longer solving times and struggles to generalize to other problem instances. This observation highlights the trade-off between a model's ability to generalize and its capacity to fit specific instances. Specifically, a model that fits individual instances exceptionally well may fail to generalize to new data, resulting in longer solving times. Conversely, a model that generalizes effectively may exhibit slightly weaker performance on specific instances, leading to a marginal decrease in objective function values. Therefore, the choice between these two training modes should be guided by the specific requirements of the application.

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5 CONCLUSIONS

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In this paper, we propose ROS, an efficient method for addressing the Max-*k*-Cut problem with any arbitrary edge weights. Our approach begins by relaxing the constraints of the original dis-

486 crete problem to probabilistic simplices. To effectively solve this relaxed problem, we propose an 487 optimization algorithm based on GNN parametrization and incorporate transfer learning by lever-488 aging pre-trained GNNs to warm-start the training process. After resolving the relaxed problem, 489 we present a novel random sampling algorithm that maps the continuous solution back to a discrete 490 form. By integrating geometric landscape analysis with statistical theory, we establish the consistency of function values between the continuous and discrete solutions. Experiments conducted on 491 random regular graphs and the Gset benchmark demonstrate that our method is highly efficient for 492 solving large-scale Max-k-Cut problems, requiring only a few seconds, even for instances with tens 493 of thousands of variables. Furthermore, it exhibits robust generalization capabilities across both 494 in-distribution and out-of-distribution instances, highlighting its effectiveness for large-scale opti-495 mization tasks. Exploring other sampling algorithms to further boost ROS performance is a future 496 research direction. Moreover, the ROS framework with theoretical insights could be potentially ex-497 tended to other graph-related combinatorial problems, and this direction is also worth investigating 498 as future work. 499

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A PROOF OF THEOREM 1

Proof. Before proceeding with the proof of Theorem 1, we first define the neighborhood of a vector $\bar{x} \in \Delta_k$, and establish results of Lemma 1 and Lemma 2.

Definition 2. Let $\bar{x} = (\bar{x}_1, \dots, \bar{x}_k)$ denote a point in Δ_k . We define the neighborhood induced by \bar{x} as follows:

$$\widetilde{\mathcal{N}}(ar{m{x}}) \coloneqq \left\{ (m{x}_1, \cdots, m{x}_k) \in \Delta_k \left| \sum_{j \in \mathcal{K}(ar{m{x}})} m{x}_j = 1 \right.
ight\},$$

where $\mathcal{K}(\bar{\boldsymbol{x}}) = \{j \in \{1, \cdots, k\} \mid \bar{\boldsymbol{x}}_j > 0\}.$

Lemma 1. Given $X_{\cdot i} \in \widetilde{\mathcal{N}}(\overline{X}_{\cdot i})$, it follows that

$$\mathcal{C}(\boldsymbol{X}_{\cdot i}) \subseteq \mathcal{K}(\overline{\boldsymbol{X}}_{\cdot i}).$$

Proof. Suppose there exists $j \in \mathcal{K}(\mathbf{X}_{\cdot i})$ such that $j \notin \mathcal{K}(\overline{\mathbf{X}}_{\cdot i})$, implying $\mathbf{X}_{ji} > 0$ and $\overline{\mathbf{X}}_{ji} = 0$. We then have

$$\sum_{l \in \mathcal{K}(\overline{\mathbf{X}}_{\cdot,i})} \mathbf{X}_{li} + \mathbf{X}_{ji} \leq \sum_{l=1}^{k} \mathbf{X}_{li} = 1,$$

which leads to

$$\sum_{\in \mathcal{K}(\overline{\boldsymbol{X}}_{\cdot i})} \boldsymbol{X}_{li} \leq 1 - \boldsymbol{X}_{ji} < 1$$

contradicting with the fact that $X_{i} \in \widetilde{\mathcal{N}}(\overline{X}_{i})$.

Lemma 2. Let \overline{X} be a globally optimal solution to \overline{P} , then

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 $f(\boldsymbol{X}; \boldsymbol{W}) = f(\overline{\boldsymbol{X}}; \boldsymbol{W}),$

where X has only the i^{th} column $X_{i} \in \widetilde{\mathcal{N}}(\overline{X}_{i})$, and other columns are identical to those of \overline{X} . Moreover, X is also a globally optimal solution to \overline{P} .

Proof. The fact that X is a globally optimal solution to \overline{P} follows directly from the equality $f(X; W) = f(\overline{X}; W)$. Thus, it suffices to prove this equality. Consider that \overline{X} and X differ only in the i^{th} column, and $X_{i} \in \widetilde{\mathcal{N}}(\overline{X}_{i})$. We can rewrite the objective value function as

$$f(\boldsymbol{X}; \boldsymbol{W}) = g(\boldsymbol{X}_{\cdot i}; \boldsymbol{X}_{\cdot - i}) + h(\boldsymbol{X}_{\cdot - i}),$$

where $X_{.-i}$ represents all column vectors of X except the i^{th} column. The functions g and h are defined as follows:

$$g(\boldsymbol{X}_{\cdot i}; \boldsymbol{X}_{\cdot - i}) = \sum_{j=1}^{N} \boldsymbol{W}_{ij} \boldsymbol{X}_{\cdot i}^{\top} \boldsymbol{X}_{\cdot j} + \sum_{j=1}^{N} \boldsymbol{W}_{ji} \boldsymbol{X}_{\cdot j}^{\top} \boldsymbol{X}_{\cdot i} - \boldsymbol{W}_{ii} \boldsymbol{X}_{\cdot i}^{\top} \boldsymbol{X}_{\cdot i},$$
$$h(\boldsymbol{X}_{\cdot - i}) = \sum_{l=1, l \neq i}^{N} \sum_{j=1, j \neq i}^{N} \boldsymbol{W}_{lj} \boldsymbol{X}_{\cdot l}^{\top} \boldsymbol{X}_{\cdot j}$$

To establish that $f(X; W) = f(\overline{X}; W)$, it suffices to show that

$$g(\boldsymbol{X}_{\cdot i}; \boldsymbol{X}_{\cdot -i}) = g(\overline{\boldsymbol{X}}_{\cdot i}; \boldsymbol{X}_{\cdot -i})$$

as $X_{\cdot-i} = \overline{X}_{\cdot-i}$.

Rewriting $q(\mathbf{X}_{\cdot i}; \mathbf{X}_{\cdot - i})$, we obtain

$$g(\boldsymbol{X}_{\cdot i}; \boldsymbol{X}_{\cdot - i}) = \sum_{j=1}^{N} \boldsymbol{W}_{ij} \boldsymbol{X}_{\cdot i}^{\top} \boldsymbol{X}_{\cdot j} + \sum_{j=1}^{N} \boldsymbol{W}_{ji} \boldsymbol{X}_{\cdot j}^{\top} \boldsymbol{X}_{\cdot i}$$
$$= 2 \sum_{j=1}^{N} \boldsymbol{W}_{ij} \boldsymbol{X}_{\cdot i}^{\top} \boldsymbol{X}_{\cdot j}$$
$$= 2 \boldsymbol{X}_{\cdot i}^{\top} \sum_{j=1, j \neq i}^{N} \boldsymbol{W}_{ij} \boldsymbol{X}_{\cdot j}$$

where $\boldsymbol{Y}_{i} := \sum_{j=1, j \neq i}^{N} \boldsymbol{W}_{ij} \boldsymbol{X}_{j}$.

If $|\mathcal{K}(\overline{X}_{i})| = 1$, then there is only one non-zero element in \overline{X}_{i} equal to one. Therefore, $g(\overline{X}_{\cdot i}; X_{\cdot - i}) = g(X_{\cdot i}; X_{\cdot - i})$ since $X_{\cdot i} = \overline{X}_{\cdot i}$.

 $= 2 \boldsymbol{X}_{i}^{\mathsf{T}} \boldsymbol{Y}_{i}$

For the case where $|\mathcal{K}(\overline{X}_{\cdot i})| > 1$, we consider any indices $j, l \in \mathcal{K}(\overline{X}_{\cdot i})$ such that $\overline{X}_{ji}, \overline{X}_{li} > 0$. Then, there exists $\epsilon > 0$ such that we can construct a point $\tilde{x} \in \Delta_k$ where the j^{th} element is set to $\overline{X}_{ji} - \epsilon$, the l^{th} element is set to $\overline{X}_{li} + \epsilon$, and all other elements remain the same as in \overline{X}_{i} . Since \overline{X} is a globally optimum of the function f(X; W), it follows that \overline{X}_{i} is also a global optimum for the function $q(\overline{X}_{i}; X_{i-i})$. Thus, we have

$$egin{aligned} g(\overline{oldsymbol{X}}_{\cdot i};oldsymbol{X}_{\cdot -i}) &\leq g(\widetilde{oldsymbol{x}};oldsymbol{X}_{\cdot -i}) \ \overline{oldsymbol{X}}_{\cdot i}^{ op}oldsymbol{Y}_{\cdot i} &\leq \widetilde{oldsymbol{x}}^{ op}oldsymbol{Y}_{\cdot i} \end{aligned}$$

$$= \overline{oldsymbol{X}}_{\cdot i}^{ op} oldsymbol{Y}_{\cdot i} - \epsilon oldsymbol{Y}_{ji} + \epsilon oldsymbol{Y}_{lij}$$

which leads to the inequality

 $Y_{ii} \leq Y_{li}$. (3)

Next, we can similarly construct another point $\hat{x} \in \Delta_k$ with its j^{th} element equal to $\overline{X}_{ji} + \epsilon$, the k^{th} element equal to $\overline{X}_{ki} - \epsilon$, and all other elements remain the same as in \overline{X}_{i} . Subsequently, we can also derive that

$$egin{aligned} g(oldsymbol{X}_{\cdot i};oldsymbol{X}_{\cdot - i}) & \leq g(\hat{oldsymbol{x}};oldsymbol{X}_{\cdot - i}) \ & = \overline{oldsymbol{X}}_{\cdot i}^{ op}oldsymbol{Y}_{\cdot i} + \epsilonoldsymbol{Y}_{li} - \epsilonoldsymbol{Y}_{li} \end{aligned}$$

which leads to another inequality

 $Y_{li} \leq Y_{ii}$. (4)

Consequently, combined inequalities (3) and (4), we have

 $Y_{ii} = Y_{li}$,

for $j, l \in \mathcal{K}(\overline{X}_{\cdot i})$.

From this, we can deduce that

 $\boldsymbol{Y}_{j_1i} = \boldsymbol{Y}_{j_2i} = \cdots = \boldsymbol{Y}_{j_{|\mathcal{K}(\overline{\boldsymbol{X}}_{-i})|}i} = t,$

where $j_1, \cdots, j_{|\mathcal{K}(\overline{X}_i)|} \in \mathcal{K}(\overline{X}_i)$.

 $g(\overline{\boldsymbol{X}}_{\cdot i}; \boldsymbol{X}_{\cdot - i}) = 2\overline{\boldsymbol{X}}_{\cdot i}^{\top} \boldsymbol{Y}_{\cdot i}$

 $=2\sum_{j=1}^{k}\overline{X}_{ji}Y_{ji}$

 $=2\sum_{j=1,j\in\mathcal{K}(\overline{\boldsymbol{X}}_{\cdot i})}^{N}\overline{\boldsymbol{X}}_{ji}\boldsymbol{Y}_{ji}$

 $=2t\sum_{j=1,j\in\mathcal{K}(\overline{\boldsymbol{X}}_{\cdot i})}^{N}\overline{\boldsymbol{X}}_{ji}$

= 2t.

 $=2\sum_{j=1}^k \boldsymbol{X}_{ji}\boldsymbol{Y}_{ji}$

 $= 2 \sum_{j=1,j\in\mathcal{K}(\boldsymbol{X}_{\cdot i})} \boldsymbol{X}_{ji} \boldsymbol{Y}_{ji}$

 $\stackrel{\text{Lemma I}}{=} 2t \sum_{j=1, j \in \mathcal{K}(\boldsymbol{X}_{\cdot i})} \boldsymbol{X}_{ji}$

 $g(\boldsymbol{X}_{\cdot i}; \boldsymbol{X}_{\cdot -i}) = 2\boldsymbol{X}_{\cdot i}^{\top} \boldsymbol{Y}_{\cdot i}$

Next, we find that

Accordingly, we conclude that

Similarly, we have

which leads us to the result

 $f(\boldsymbol{X}; \boldsymbol{W}) = f(\overline{\boldsymbol{X}}; \boldsymbol{W}),$

 $q(\boldsymbol{X}_{\cdot i}; \boldsymbol{X}_{\cdot - i}) = q(\overline{\boldsymbol{X}}_{\cdot i}; \boldsymbol{X}_{\cdot - i}),$

= 2t

 $= q(\overline{X}_{\cdot i})$

where $X_{i} \in \widetilde{\mathcal{N}}(\overline{X}_{i}), X_{i-i} = \overline{X}_{i-i}$.

> Accordingly, for any $X \in \mathcal{N}(\overline{X})$, we iteratively apply Lemma 2 to each column of \overline{X} while holding the other columns fixed, thereby proving Theorem 1.

PROOF OF THEOREM 2 В

Proof. Based on \overline{X} , we can construct the random variable \widetilde{X} , where $\widetilde{X}_{i} \sim \operatorname{Cat}(x; p = \overline{X}_{i})$. The probability mass function is given by

> $\mathbf{P}(\widetilde{\mathbf{X}}_{i}=\mathbf{e}_{\ell})=\overline{\mathbf{X}}_{\ell i},$ (5)

where $\ell = 1, \cdots, k$.

Next, we have $\mathbb{E}_{\widetilde{\mathbf{X}}}[f(\widetilde{\mathbf{X}}; \mathbf{W})] = \mathbb{E}_{\widetilde{\mathbf{X}}}[\widetilde{\mathbf{X}} \mathbf{W} \widetilde{\mathbf{X}}^{\top}] = \mathbb{E}_{\widetilde{\mathbf{X}}}[\sum_{i=1}^{N} \sum_{j=1}^{N} \mathbf{W}_{ij} \widetilde{\mathbf{X}}_{\cdot i}^{\top} \widetilde{\mathbf{X}}_{\cdot j}]$ $= \sum_{i=1}^{N} \sum_{j=1}^{N} \mathbf{W}_{ij} \mathbb{E}_{\widetilde{\mathbf{X}}_{\cdot i}, \widetilde{\mathbf{X}}_{\cdot j}}[\widetilde{\mathbf{X}}_{\cdot i}^{\top} \widetilde{\mathbf{X}}_{\cdot j}]$ $= \sum_{i=1}^{N} \sum_{j=1}^{N} \mathbf{W}_{ij} \mathbb{E}_{\widetilde{\mathbf{X}}_{\cdot i}, \widetilde{\mathbf{X}}_{\cdot j}}[\mathbb{1}(\widetilde{\mathbf{X}}_{\cdot i} = \widetilde{\mathbf{X}}_{\cdot j})]$ $= \sum_{i=1}^{N} \sum_{j=1}^{N} \mathbf{W}_{ij} \mathbb{P}(\widetilde{\mathbf{X}}_{\cdot i} = \widetilde{\mathbf{X}}_{\cdot j})$ $= \sum_{i=1}^{N} \sum_{j=1, j \neq i}^{N} \mathbf{W}_{ij} \mathbb{P}(\widetilde{\mathbf{X}}_{\cdot i} = \widetilde{\mathbf{X}}_{\cdot j}).$ Since $\widetilde{\mathbf{X}}_{\cdot i}$ and $\widetilde{\mathbf{X}}_{\cdot j}$ are independent for $i \neq j$, we have $\mathbb{P}(\widetilde{\mathbf{X}}_{\cdot i} = \widetilde{\mathbf{X}}_{\cdot j}) = \sum_{\ell=1}^{k} \mathbb{P}(\widetilde{\mathbf{X}}_{\cdot i} = \widetilde{\mathbf{X}}_{\cdot j} = \mathbf{e}_{\ell})$ $= \sum_{\ell=1}^{k} \mathbb{P}(\widetilde{\mathbf{X}}_{\cdot i} = \mathbf{e}_{\ell}, \widetilde{\mathbf{X}}_{\cdot j} = \mathbf{e}_{\ell})$ $= \sum_{\ell=1}^{k} \mathbb{P}(\widetilde{\mathbf{X}}_{\cdot i} = \mathbf{e}_{\ell}) \mathbb{P}(\widetilde{\mathbf{X}}_{\cdot j} = \mathbf{e}_{\ell})$

 $=\sum_{\ell=1}^{k}\overline{X}_{\ell i}\overline{X}_{\ell j}$

 $=\overline{X}_{i}^{\top}\overline{X}_{i}$

Substitute (7) into (6), we obtain

$$\mathbb{E}_{\widetilde{\boldsymbol{X}}}[f(\widetilde{\boldsymbol{X}};\boldsymbol{W})] = \sum_{i=1}^{N} \sum_{j=1}^{N} \boldsymbol{W}_{ij} \overline{\boldsymbol{X}}_{\cdot i}^{\top} \overline{\boldsymbol{X}}_{\cdot j} = f(\overline{\boldsymbol{X}};\boldsymbol{W}).$$
(8)

(7)

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C THE COMPLETE RESULTS ON GSET INSTANCES

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Instance G1 G2 G3 G4					Table 4: (Complete res	sults on G	set instances	s for Max-	-3-Cut.				
Instance G1 G2 G4 G4														
G1 G2 G4	$\overline{\Sigma}$	3		MD	Ger	letic	[BQP		HOM	ROS-	vanilla	[SOS
G1 G2 G4	-	-	Obj. ↑	Time (s) \downarrow	Obj.↑	Time (s) \downarrow	Obj. ↑	Time (s) \downarrow	Obj.↑	Time (s) \downarrow	Obj.↑	Time (s) \downarrow	Obj.↑	Time (s) \downarrow
G2 G3 G4	800	19176	14735	9.6	14075	595.3	14880	16.5	15165	557.3	14949	2.8	14961	1.9
G3 G4	800	19176	14787	8.4	14035	595.3	14845	17.0	15172	333.3	15033	2.8	14932	2.3
G4	800	19176	14663	6.5	14105	588.6	14872	17.0	15173	269.6	15016	2.9	14914	1.9
	800	19176	14716	6.9	14055	588.7	14886	17.1	15184	300.6	14984	3.3	14961	1.9
G5	800	19176	14681	8.1	14104	591.9	14847	17.3	15193	98.2	15006	3.2	14962	2.9
G6	800	19176	2161	7.8	1504	604.4	2302	25.0	2632	307.3	2436	2.8	2361	1.8
G7	800	19176	2017	8.9	1260	589.9	2081	16.6	2409	381.0	2188	2.1	2188	2.4
G8	800	19176	1938	T.T	1252	589.7	2096	19.3	2428	456.5	2237	2.8	2171	2.1
G9	800	19176	2031	8.2	1326	604.4	2099	16.5	2478	282.0	2246	2.8	2185	2.2
G10	800	19176	1961	7.5	1266	593.3	2055	18.2	2407	569.3	2201	2.9	2181	2.3
G11	800	1600	553	4.0	414	554.5	624	16.4	699	143.8	616	2	591	1.4
G12	800	1600	530	4.4	388	543.6	608	17.4	660	100.7	604	7	582	1.5
G13	800	1600	558	4.0	425	550.8	638	18.9	686	459.4	617	7	629	1.4
G14	800	4694	3844	5.0	3679	571.1	3900	16.9	4012	88.2	3914	2.8	3892	2.1
G15	800	4661	3815	4.8	3625	567.6	3885	17.3	3984	80.3	3817	1.9	3838	2
G16	800	4672	3825	5.3	3642	561.5	3896	18.2	3991	1.3	3843	2.3	3845	1.6
G17	800	4667	3815	5.3	3640	558.7	3886	20.2	3983	7.8	3841	2.4	3852	1.6
G18	800	4694	992	4.5	704	584.0	1083	18.7	1207	0.3	1094	2.2	1067	1.7
G19	800	4661	869	4.4	595	584.2	962	17.0	1081	0.2	972	2.1	967	1.7
G20	800	4672	928	4.5	589	576.8	776	17.0	1122	13.3	1006	2.2	993	1.8
G21	800	4667	936	4.9	612	576.3	984	17.5	1109	55.8	1011	2.2	975	1.5
G22	2000	19990	16402	15.2	N/A	N/A	16599	135.5	17167	28.5	16790	3.3	16601	2.2
G23	2000	19990	16422	15.0	N/A	N/A	16626	135.6	17168	45.1	16819	3.9	16702	2.1
G24	2000	19990	16452	16.1	N/A	N/A	16591	137.7	17162	16.3	16801	3.6	16754	б
G25	2000	19990	16407	16.2	N/A	N/A	16661	141.8	17163	64.8	16795	2.1	16673	1.8
G26	2000	19990	16422	15.3	N/A	N/A	16608	136.3	17154	44.8	16758	3.1	16665	2
G27	2000	19990	3250	16.4	N/A	N/A	3475	134.3	4020	53.2	3517	1.7	3532	0
G28	2000	19990	3198	16.1	N/A	N/A	3433	136.4	3973	38.9	3507	3	3414	2.1
G29	2000	19990	3324	16.0	N/A	N/A	3582	136.2	4106	68.2	3634	3.4	3596	2
G30	2000	19990	3320	16.2	N/A	N/A	3578	133.6	4119	150.4	3656	3.1	3654	3.4
G31	2000	19990	3243	17.0	N/A	N/A	3439	131.0	4003	124.7	3596	3	3525	2.5
G32	2000	4000	1342	11.1	N/A	N/A	1545	129.3	1653	160.1	1488	2.5	1482	1.7
G33	2000	4000	1284	10.7	N/A	N/A	1517	126.2	1625	62.6	1449	2.5	1454	2
G34	2000	4000	1292	10.9	N/A	N/A	1499	126.0	1607	88.9	1418	2.4	1435	1.7
G35	2000	11778	9644	14.2	N/A	N/A	9816	138.1	10046	66.2	9225	2	9536	1.7

1077 1078 1079	1074 1075 1076	1071 1072 1073	1068 1069 1070	1063 1064 1065 1066 1067	1060 1061 1062	1056 1057 1058 1059	1053 1054 1055	1048 1049 1050 1051 1052	1045 1046 1047	1041 1042 1043 1044	1038 1039 1040	1034 1035 1036 1037	1030 1031 1032 1033	1026 1027 1028 1029
						Та	lble 4: Co	ntinued.						
Instance	2	3		MD	Gei	netic		BQP		НОМ	ROS-	vanilla	[ROS
	-	-	Obj. ↑	Time (s) \downarrow	Obj. ↑	Time (s) \downarrow	Obj. ↑	Time (s) \downarrow	Obj. ↑	Time (s) \downarrow	Obj. ↑	Time (s) \downarrow	Obj. ↑	Time (s) \downarrow
G36	2000	11766	9600	13.6	N/A	N/A	9786	138.6	10039	74.3	9372	2.1	9581	2.3
G37	2000	11785	9632	14.9	N/A	N/A	9821	139.2	10052	3.4	8893	1.4	9422	1.5
G38	2000	11779	9629	14.0	N/A	N/A	9775	142.3	10040	116.6	9489	2.5	9370	1.5
G39	2000	11778	2368	13.4	N/A	NA	2600	132.8	2903	9.0	2621	2.5	2557	2.2
G40	2000	11766	2315	13.3	N/A	NA	2568	131.2	2870	82.8	2474	2	2524	2.4
G41	2000	11785	2386	12.7	N/A	N/A	2606	129.9	2887	87.7	2521	3.2	2584	2.5
G42	2000	11779	2490	13.1	N/A	NA	2682	129.2	2980	2.5	2638	2.7	2613	2.2
G43	1000	0666	8214	8.1	7624	926.7	8329	29.9	8573	380.3	8414	2.6	8349	2.3
G44	1000	0666	8187	7.0	7617	919.0	8326	27.7	8571	616.8	8369	2.6	8311	1.7
G45	1000	0666	8226	<i>T.T</i>	7602	926.7	8296	34.2	8566	186.2	8397	2.9	8342	1.8
G46	1000	0666	8229	7.5	7635	918.7	8312	27.8	8568	215.3	8409	2.6	8339	1.7
G47	1000	0666	8211	7.2	7619	928.0	8322	27.3	8572	239.4	8386	2.6	8357	2.2
G48	3000	6000	5806	14.7	N/A	NA	5998	394.8	6000	0.4	5954	2.8	5912	2
G49	3000	6000	5794	14.4	N/A	NA	5998	404.0	6000	0.9	5938	2.8	5914	1.8
G50	3000	0009	5823	14.5	N/A	NA	6000	427.1	0009	119.2	5938	2.9	5918	1.8
G51	1000	5909	4805	6.6	4582	889.5	4922	28.6	5037	47.9	4814	2.4	4820	1.7
G52	1000	5916	4849	6.4	4571	908.1	4910	27.8	5040	0.7	4796	1.9	4866	1.9
G53	1000	5914	4845	6.8	4568	898.6	4920	27.6	5039	223.9	4846	2.6	4808	1.6
G54	1000	5916	4836	6.4	4562	911.7	4921	30.1	5036	134.0	4833	2.2	4785	1.4
G55	5000	12498	11612	37.9	N/A	NA	12042	1506.0	12429	383.1	12010	2.1	11965	2.6
G56	5000	12498	3716	38.5	N/A	NA	4205	1341.5	4752	569.2	4085	3.3	4037	2.1
G57	5000	10000	3246	33.0	N/A	N/A	3817	1317.2	4083	535.6	3597	3.3	3595	2.8
G58	5000	29570	24099	47.1	N/A	N/A	24603	1468.3	25195	576.0	22748	2.1	23274	1.9
G59	5000	29570	6057	46.3	N/A	NA	6631	1377.1	7262	27.5	6133	1.7	6448	3.5
G60	7000	17148	15993	58.5	N/A	NA	NA	N/A	17076	683.0	16467	2.6	16398	2.3
G61	7000	17148	5374	57.7	N/A	N/A	N/A	N/A	6853	503.1	5881	2.5	5861	3.6
G62	7000	14000	4497	49.7	N/A	NA	NA	N/A	5685	242.4	4983	3.4	5086	2.7
G63	7000	41459	33861	73.4	N/A	NA	NA	N/A	35322	658.5	32868	4	31926	1.9
G64	7000	41459	8773	73.4	N/A	NA	NA	N/A	10443	186.9	8911	2.8	9171	2.5
G65	8000	16000	5212	59.6	N/A	NA	N/A	N/A	6490	324.7	5735	3.5	5775	2.6
G66	0006	18000	5948	69.0	N/A	NA	NA	N/A	7416	542.5	6501	5.4	6610	3.9
G67	10000	20000	6545	79.0	N/A	NA	NA	N/A	8086	756.7	7001	3.5	7259	4.1
G70	10000	6666	9718	74.8	N/A	NA	NA	N/A	6666	7.8	9982	4.2	9971	2.5
G72	10000	20000	6612	79.2	N/A	NA	NA	N/A	8192	271.2	7210	5.1	7297	3.5
G77	14000	28000	9294	142.3	N/A	NA	NA	N/A	11578	154.9	10191	8.6	10329	8.5
G81	20000	40000	13098	241.1	N/A	N/A	N/A	N/A	16321	331.2	14418	20.2	14464	9.7

1080 D EVALUATION ON GRAPH COLORING DATASET

To further verify the performance of ROS, we conduct numerical experiments on the publicly available COLOR dataset (three benchmark instances: anna, david, and huck). The COLOR dataset provides dense problem instances with relatively large known chromatic numbers ($\chi \sim 10$), which is suitable for testing the performance on Max-k-Cut tasks. As reported in Tables 5 and 6, ROS achieves superior performances across nearly all settings with the least computational time (in seconds).

Table 5: Objective values returned by each method on the COLOR dataset.

Methods	an	na	da	vid	hu	ck
	k = 2	k = 3	k = 2	k=3	k = 2	k = 3
MD	339	421	259	329	184	242
PI-GNN	322	-	218	-	170	-
ecord	351	-	267	-	191	-
ANYCSP	351	-	267	-	191	-
ROS	351	421	266	338	191	244

Table 6: Computational time for each method on the COLOR dataset.

Methods	an	na	dav	vid	hu	ck
112011005	k = 2	k = 3	k=2	k = 3	k=2	k = 3
MD	2.75	2.08	2.78	2.79	2.62	2.82
PI-GNN	93.40	-	86.84	-	102.57	-
ecord	4.87	-	4.74	-	4.88	-
ANYCSP	159.35	-	138.14	-	127.36	-
ROS	1.21	1.23	1.18	1.15	1.11	1.10

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1111 E ABLATION STUDY

1113 E.1 MODEL ABLATION

We conducted additional ablation studies to clarify the contributions of different modules.

Effect of Neural Networks: We consider two cases: (i) replace GNNs by multi-layer perceptrons
(denoted by ROS-MLP) in our ROS framework and (ii) solve the relaxation via mirror descent (denoted by MD). Experiments on the Gset dataset show that ROS consistently outperforms ROS-MLP
and MD, highlighting the benefits of using GNNs for the relaxation step.

 Effect of Random Sampling: We compared ROS with PI-GNN, which employs heuristic rounding instead of our random sampling algorithm. Results indicate that ROS generally outperforms
 PI-GNN, demonstrating the importance of the sampling procedure.

These comparisons, detailed in Tables 7 and 8, confirm that both the GNN-based optimization and the random sampling algorithm contribute significantly to the overall performance.

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 SAMPLE EFFECT ABLATION

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 Sample Effect Ablation

We investigated the effect of the number of sampling iterations and report the results in Tables 9, 10,11, and 12.

Objective Value (Table 9, Table 11): The objective values stabilize after approximately 5 sampling iterations, demonstrating strong performance without requiring extensive sampling.

Sampling Time (Table 10, Table 12): The time spent on sampling remains negligible compared to the total computational time, even with an increased number of samples.

Table 7: Objective values returned by each method on Gset.

Methods	G	70	G	72	G	77	G8	31
Wiethous	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3	k=2	k = 3
ROS-MLP	8867	9943	6052	6854	8287	9302	12238	12298
PI-GNN	8956	_	4544	_	6406	_	8970	_
MD	8551	9728	5638	6612	7934	9294	11226	13098
ROS	8916	9971	6102	7297	8740	10329	12332	14464

Table 8: Computational time for each method on Gset.

Methods	G	70	G7	2	G	77	G	31
	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3
ROS-MLP	3.49	3.71	3.93	4.06	8.39	9.29	11.98	16.97
PI-GNN	34.50	_	253.00	_	349.40	_	557.70	_
MD	54.30	74.80	44.20	79.20	66.00	142.30	130.80	241.10
ROS	3.40	2.50	3.90	3.50	8.10	8.50	9.30	9.70

These results highlight the efficiency of our sampling method, achieving stable and robust performance with little computational cost.

F COMPARISON AGAINST ADDITIONAL BASELINES ON GSET

We have conducted additional experiments comparing ROS against ANYCSP and ECORD on the Gset benchmark for Max-Cut, focusing on both solution quality and computational efficiency. The results are presented below.

1163Results on Gset (unweighted) with Edge Weights of ± 1 : Tables 13 and 14 present the comparison1164of objective values and inference times for each method on unweighted Gset instances. Although1165ANYCSP achieves marginally better objective values, its computational time is considerably longer.1166ECORD, on the other hand, fails to generate competitive solutions. In contrast, our ROS framework1167strikes an optimal balance, delivering high-quality solutions in a fraction of the time required by
ANYCSP.

Weighted Max-Cut has numerous applications, including but not limited to physics (De Simone et al., 1995), power networks (Hojny et al., 2021), and data clustering (Poland & Zeugmann, 2006).
To demonstrate the capability of our ROS framework in solving general Max-k-Cut problems, we evaluate the performance of ROS, ANYCSP, and ECORD in this context.

Results on Gset with Arbitrary Edge Weights. We modified the four largest Gset instances (G70, G72, G77, and G81) to incorporate arbitrary edge weights. Specifically, we perturb the edge weights uniformly within the range [-10%, 10%] for each Gset benchmark and generate 10 instances. The averaged results, along with their standard deviations, summarized in Tables 15 and 16, reveal the limitations of both ANYCSP and ECORD in this setting. ANYCSP fails to produce meaningful so-lutions due to its CSP-based formulation, which overlooks edge weights, and ECORD again demon-strates poor performance. In contrast, ROS consistently generates high-quality solutions while main-taining computational efficiency, showcasing its robustness and versatility across different scenarios.

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T	G	70	G	72	G	77	G8	31
-	k = 2	k = 3	k=2	k = 3	k=2	k = 3	k=2	k = 3
1	8911	9968	6100	7305	8736	10321	12328	14460
5	8915	9969	6102	7304	8740	10326	12332	14462
10	8915	9971	6102	7305	8740	10324	12332	14459
25	8915	9971	6102	7307	8740	10326	12332	14460
50	8915	9971	6102	7307	8740	10327	12332	14461
100	8916	9971	6102	7308	8740	10327	12332	14462

Table 9: Objective value results corresponding to the times of sample T on Gset.

Table 10: Sampling time results corresponding to the times of sample T on Gset.

Т	G	70	G	72	G	77	G8	31
-	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3
1	0.0011	0.0006	0.0011	0.0006	0.0020	0.0010	0.0039	0.0020
5	0.0030	0.0029	0.0029	0.0030	0.0053	0.0053	0.0099	0.0098
10	0.0058	0.0059	0.0058	0.0058	0.0104	0.0104	0.0196	0.0196
25	0.0144	0.0145	0.0145	0.0145	0.0259	0.0260	0.0489	0.0489
50	0.0289	0.0289	0.0288	0.0289	0.0517	0.0518	0.0975	0.0977
100	0.0577	0.0577	0.0576	0.0578	0.1033	0.1037	0.1949	0.1953

Table 11: Objective value results corresponding to the times of sample T on random regular graphs.

Т	n = 100		n =	1000	n = 10000		
-	k = 2	k = 3	k=2	k = 3	k = 2	k = 3	
1	127	245	1293	2408	12856	24103	
5	127	245	1293	2410	12863	24103	
10	127	245	1293	2410	12862	24103	
25	127	245	1293	2410	12864	24103	
50	127	245	1293	2410	12864	24103	
100	127	245	1293	2410	12864	24103	

Table 12: Sampling time results corresponding to the times of sample T on random regular graphs.

Т	n =	100	n =	1000	n = 10000		
-	k = 2	k = 3	k = 2	k = 3	k = 2	k = 3	
1	0.0001	0.0001	0.0001	0.0001	0.0006	0.0006	
5	0.0006	0.0006	0.0007	0.0007	0.0030	0.0030	
10	0.0011	0.0011	0.0014	0.0013	0.0059	0.0059	
25	0.0026	0.0026	0.0033	0.0031	0.0145	0.0145	
50	0.0052	0.0052	0.0065	0.0060	0.0289	0.0289	
100	0.0103	0.0103	0.0128	0.0122	0.0577	0.0578	

Table 13: Objective values returned by each method on Gset for k = 2.

1237	Methods	G70	G72	G77	G81
1239	Ecord	5137	206	382	358
1240	ANYCSP	9417	6826	9694	13684
1241	ROS	8916	6102	8740	12332

		Ecord ANYCSP ROS	1.4 180.0 3.4	1.2 180.0 3.9	1.7 180.0 8.1	2.4 180.0 9.3			
		ANYCSP ROS	1.4 180.0 3.4	1.2 180.0 3.9	180.0 8.1	2.4 180.0 9.3			
		ROS	3.4	3.9	8.1	9.3			
	Table 15: O	bjective valu	ue on Gset	with art	oitrary ed	ge weights	for $k = 2$.		
Methods	G70		G72			G77		G81	
Ecord	$5154.28 \pm$	28.26	28.26 254.46 ± 37.2		344.79 ± 29.73		280.0	280.09 ± 33	
ANYCSP	$5198.87 \pm$	69.76	-15.57 ± 5	57.88	81.7	6 ± 69.97	33.4)±	
ROS	$8941.80 \pm$	17.79 6	$165.62\pm$	50.81	8737.5	59 ± 114.2	12325.	85	
Table	16: Computat	ional time co	omparison G73	on Gset	with arb	itrary edge	weights for G81	k =	
Table	16: Computat Methods Ecord 3.	ional time concerning $\overline{G70}$ $\overline{39 \pm 0.11}$	omparison G7: 3.43 ±	on Gset 2 0.12	with arb G7 3.89 ±	itrary edge	weights for $r_{\rm G81}$	k = 	