# Appendix of *GraphSAID*: *Graph Sampling via*Attention based Integer Programming Method

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# 1 Full Algorithmic Description of Attended Connected Component Generation.

**Algorithm 1** Attended Connected Component Generation

```
Require: G, \tau, E
Ensure: \mathcal{L} (the collection of connected components as node sets)
 1: List \mathcal{L} \leftarrow [\{\}]
 2: Oueue Q \leftarrow E
 3: while Q \neq null do
            Edge e \leftarrow Q.pop()
            s_1 \leftarrow (s \in \mathcal{L} \ s.t. \ e[0] \in s)
 5:
            s_2 \leftarrow (s \in \mathcal{L} \ s.t. \ e[1] \in s)
 6:
                                                                                                                     \triangleright e[0], e[1] are vertexes
 7:
            if s_1 = null and s_2 = null then
                  s^{new} \leftarrow \{e[0], e[1]\}
 8:
 9:
                  \mathcal{L}.append(s^{new})
10:
            else if s_1 \neq null or s_2 \neq null then
11:
                  s_{found} \leftarrow (s \in \{s_1, s_2\} \ s.t. \ s \neq null)
12:
                 n_{found} \leftarrow (n \in \{e[0], e[1]\} \text{ s.t. } n \in s_{found})
13:
                  n_{new} \leftarrow (n \in \{e[0], e[1]\} \text{ s.t. } n \notin s_{found})
14:
                  if |s_{found}| < \tau then
15:
                        s_{found} = s_{found} \bigcup \{n_{new}\}
                       update \mathcal{L} with s^{found}
16:
17:
                  else
                        s^{new} \leftarrow \{n_{new}\}; \ \mathcal{L}.append(s^{new})
18:
19:
                 end if
20:
            else if s_1 \neq null and s_2 \neq null then
                  s_1 \leftarrow (s \in \mathcal{L} \ s.t. \ e[0] \in s)
21:
22:
                  s_2 \leftarrow (s \in \mathcal{L} \ s.t. \ e[1] \in s)
23:
                  if s_1 \neq s_2 then
                       s^{merge} \leftarrow s_1 \cup s_2
24:
                        remove s_1, s_2 in \mathcal{L}; \mathcal{L}.append(s^{merge})
25:
26:
                  end if
27:
            end if
28: end while
```

# 2 Full Description of Datasets

**GRPG.** There are many graph generators can model a transaction network, e.g., [6, 3, 4, 1, 2, 5].

In this work, we model our synthesized network as a realization of a *Gaussian Random Partition Graph* process [2], due to the distinct nature of intra- and inter-group activities which can be used to model complex networks like cash flow. The *Gaussian Random Partition Graph* (GRPG) process can be noted as  $G(n, s, v, p_{in}, p_{out})$ , where n is the number of nodes, s is the mean of the cluster size, s/v is the variance of the cluster size,  $p_{in}$  and  $p_{out}$  are the probabilities of intra- and inter-cluster connection. Fig-







Figure 1: GRPG Example. Left: G(400, 100, 200, 0.05, 0.001); Middle: G(400, 100, 200, 0.1, 0.001); right: G(400, 100, 200, 0.1, 0.01).

ure 1 shows visualization of graph structure changes by controlling different  $p_{out}$ . In this example,

		Elliptic			Co	ora	Cite	seer	GPRG		
В	Method	$R_{attn}$	Acc	F1	$R_{attn}$	Acc	$R_{attn}$	Acc	$R_{attn}$	Acc	
60%	FF	0.156	0.940	0.474	0.248	0.747	0.628	0.618	0.434	0.832	
60%	SB	0.157	0.934	0.481	0.253	0.788	0.645	0.589	0.402	0.856	
60%	SRW	0.158	0.941	0.517	0.250	0.806	0.645	0.578	0.407	0.861	
60%	RWF	0.146	0.949	0.544	0.249	0.811	0.647	0.594	0.409	0.861	
60%	ORC-sub	oom	oom	oom	0.258	0.743	0.646	0.618	0.483	0.862	
60%	Ours	0.506	0.952	0.557	0.380	0.815	0.650	0.626	0.627	0.860	
60%	$Ours^\dagger$	0.506	0.951	0.557	0.380	0.817	0.650	0.625	0.627	0.860	
60%	$Ours^{\ddagger}$	0.506	0.952	0.559	0.380	0.815	0.650	0.626	0.627	0.860	
80%	FF	0.198	0.950	0.537	0.344	0.831	0.857	0.626	0.660	0.849	
80%	SB	0.213	0.929	0.412	0.351	0.828	0.861	0.627	0.674	0.860	
80%	SRW	0.193	0.948	0.524	0.348	0.834	0.873	0.635	0.672	0.861	
80%	RWF	0.202	0.955	0.557	0.351	0.832	0.870	0.629	0.669	0.861	
80%	ORC-sub	oom	oom	oom	0.353	0.811	0.870	0.627	0.703	0.861	
80%	Ours	0.617	0.955	0.540	0.495	0.837	0.870	0.629	0.822	0.861	
80%	$Ours^\dagger$	0.617	0.955	0.545	0.495	0.838	0.870	0.634	0.822	0.861	
80%	$Ours^{\ddagger}$	0.617	0.955	0.540	0.495	0.836	0.870	0.636	0.822	0.861	

Table 1: Learning performance comparison using different sampled subgraphs. The best performance is in **bold**. "oom" means out-of-memory issue during the computation.

four partitions are labeled in colors, while from left to right the difficulty to detect the partitions by graph structure increases.

**Citeseer.** The CiteSeer dataset consists of 3312 scientific publications classified into one of six classes. The citation network consists of 4732 links. Each publication in the dataset is described by a 0/1-valued word vector indicating the absence/presence of the corresponding word from the dictionary. The dictionary consists of 3703 unique words.

**Cora.** The Cora dataset consists of 2708 scientific publications classified into one of seven classes. The citation network consists of 5429 links. Each publication in the dataset is described by a 0/1-valued word vector indicating the absence/presence of the corresponding word from the dictionary. The dictionary consists of 1433 unique words.

**Elliptic.** The Elliptic dataset maps Bitcoin transactions to real entities belonging to licit categories (exchanges, wallet providers, miners, licit services, etc.) versus illicit ones (scams, malware, terrorist organizations, ransomware, Ponzi schemes, etc.). It consists of 203,769 nodes and 234,355 transactions as edges. In addition, 157,205 nodes are unlabelled.

## 2.1 Full Description of GRPG dataset synthesis

We synthesize our GRPG dataset by setting: (1) G(1000, 200, 100, 0.05, 0.005); and (2) let the 5 expected clusters have a  $\mathcal{N}_{k=3}(\mu_i, \Sigma_i)$ , where  $i \in \{1, 2, 3, 4, 5\}$ ,  $\mu_i = (i-3) \cdot \mathbf{1}_k^T$  is the mean vector for the *ith* partition and  $\Sigma_i = diag(\mathbf{1}_k^T)$  for all partitions. We denote node feature for node j as  $v_j$ , and  $p_j \in \{1, 2, 3, 4, 5\}$  is the the partition for node j. To make our toy dataset more realistic, we set 25% of nodes as "unknown" (i.e., "-1"). Then we define the node labels of the other 75% nodes as:

$$label_j = \begin{cases} 0, & \text{if } \mathbf{1}^T v_j \le p_j \\ 1, & \text{otherwise} \end{cases}$$
 (1)

Table 2 summarizes the label distribution in our GRPG dataset.

block	1		2		3			4			5				
label	-1	0	1	-1	0	1	-1	0	1	-1	0	1	-1	0	1
count	43	156	2	63	119	20	46	74	80	58	15	125	40	2	156

Table 2: Statistics of toy datast *GRPG*.

### References

- [1] Albert-László Barabási and Réka Albert. Emergence of scaling in random networks. *science*, 1999.
- [2] Ulrik Brandes, Marco Gaertler, and Dorothea Wagner. Experiments on graph clustering algorithms. In *European symposium on algorithms*, 2003.
- [3] Sergey N Dorogovtsev, Alexander V Goltsev, and José Ferreira F Mendes. Pseudofractal scale-free web. *Physical review E*, 2002.
- [4] Paul Erdos, Alfréd Rényi, et al. On the evolution of random graphs. *Publ. Math. Inst. Hung. Acad. Sci*, 1960.
- [5] Paul W Holland, Kathryn Blackmond Laskey, and Samuel Leinhardt. Stochastic blockmodels: First steps. *Social networks*, 1983.
- [6] Jaewon Yang and Jure Leskovec. Community-affiliation graph model for overlapping network community detection. In *Data Mining*, 2012.