Decomposing User-APP Graph into Subgraphs for Effective APP and User Embedding Learning

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Abstract—APP-installation information is helpful to describe the user's characteristics. The users with similar APPs installed might share several common interests and behave similarly in some scenarios. In this work, we learn a user embedding vector based on each user's APP-installation information. Since the user APP-installation embedding is learnable without dependency on the historical intra-APP behavioral data of the user, it complements the intra-APP embedding learned within each specific APP. Thus, they considerably help improve the effectiveness of the personalized advertising in each APP, and they are particularly beneficial for the cold start of the new users in the APP. In this paper, we formulate the APP-installation user embedding learning into a bipartite graph embedding problem. The main challenge in learning an effective APP-installation user embedding is the imbalanced data distribution. In this case, graph learning tends to be dominated by the popular APPs, which billions of users have installed. In other words, some niche/specialized APPs might have a marginal influence on graph learning. To effectively exploit the valuable information from the niche APPs, we decompose the APP-installation graph into a set of subgraphs. Each subgraph contains only one APP node and the users who install the APP. For each mini-batch, we only sample the users from the same subgraph in the training process. Thus, each APP can be involved in the training process in a more balanced manner. After integrating the learned APP-installation user embedding into our online personal advertising platform, we obtained a considerable boost in CTR, CVR, and revenue.

Index Terms-advertising, search, cross-modal

I. INTRODUCTION

For different users, a personalized advertising system feeds different ads based on the estimated relevance between the ad and the user's interest. Normally, the relevance between a user and an ad is measured by the similarity between their embeddings, which are learned jointly from the users' historical behaviors on the ads. Nevertheless, for new users, there are no historical user-ad behaviors for learning effective user embedding. This issue of modeling new users is normally defined as the cold start problem. To solve the cold start problem, we usually exploit the user's demographic attributes, such as age, region, and gender. The attribute embedding has been effectively learned based on the ordinary users' rich



Fig. 1. The visualization of a user-APP undirected bipartite graph. A node with a blue dot denotes a user and a node with a green box denotes an APP. An edge exists between an APP and a user if the user installs the APP.

experience accumulated in the past and can readily generalize well to the new users. Since the attribute embedding does not rely on the historical user behaviors, they are useful for tackling the cold start problem.

This work explores a new type of attribute embedding learned from the APP-installation information. The users who install the same APP might share some common interests and tend to behave similarly. Meanwhile, a user's installed APP lists might encode much richer fine-grained information about a user than basic demographic information, such as age, gender, and location. Thus, if exploiting the users' APPinstallation information effectively, we might significantly boost the performance of our personalized advertising platform for the new users. In fact, the APP-installation information benefits not only the new users but also the regular users. This is because the learned user's APP-installation embedding complements the user's behavior embedding. Thus, when incorporating the APP-installation embedding into our model, we also observed improvements for the regular users' personalized advertising performance.

We formulate the APP-installation embedding as a bipartite graph embedding problem. The bipartite graph consists of two types of nodes, including the user nodes and the APP nodes as visualized in Figure 1. An edge exists between a user node and an APP node if the user has installed the APP in his/her mobile phone. Straightforwardly, we could utilize any existing graph learning methods such as graph convolutional neural network (GCN) to learn the user node embedding and the APP node embedding. Nevertheless, a serious issue caused by imbalanced data distribution makes the training of the graph learning model extremely challenging. Specifically, for a popular APP such as Tik Tok, it is installed by billions of users, generating billions of edges in the graph. In contrast, a niche APP installed by millions of users can only create millions of edges in the graph. In this case, graph learning is dominated by the billions of edges created by the popular APPs, and the edges from the niche APPs might be swamped. But the edges from the popular APPs might not encode useful discriminating information since everyone almost installs them. In contrast, the edges from the niche APPs might be very useful for describing a user's characteristics, but that useful information might not gain enough attention when training the graph embedding.

In this work, we propose a novel sampling approach to tackle the imbalanced data distribution issue for learning effective APP-installation user embeddings. Specifically, we decompose the user-APP graph into a set of sub-graphs. Each subgraph contains only a single APP and the users who install the APP. In the training process, we sample a subgraph for each iteration and construct training triplets based on users within the subgraph for embedding learning. In this manner, the popular APPs and the niche APPs will be involved in the training process in a fair manner. The offline and online experiments demonstrate the excellence of our method.

II. RELATED WORK

Factorization based methods. Factorization-based methods rely on an affinity matrix encoding the connections between nodes in the graph. They factorize the affinity matrix to obtain the embedding vectors for nodes. A pioneering work, Laplacian Eigenmaps [3] aims to keep the embedding of two nodes close when the weight of the edge connecting these two nodes is high. It seeks to minimize the weighted summation of squares of distance between nodes while the weight of each item in the summation corresponds to the weight of the edge. It is formulated into an eigenproblem. Nevertheless, it is extremely slow when solving the eigen problem in the scenario when the number of nodes is huge. Ahmed et al. [1] propose a framework for large-scale graph decomposition. They partition a graph based on minimizing the number of neighboring vertices. GraRep [5] integrates global structural information of the graph into the graph learning process. HOPE [17] learns the graph embedding for nodes with the asymmetric transitivity, which is a critical property of the directed graph.

Random walks based methods. Random walks are very useful when we have only access to a part of the graph or the graph is too large to be modeled globally. DeepWalk [18] creates multiple random walks, and maximizes the sum of log-likelihoods for each random walk. It preserves higher-order proximity between nodes in the graph. *node2vec* [10] also encodes higher-order proximity between nodes by maximizing the probability of occurrence of subsequent nodes. It conducts a trade-off between breadth-first searches (BFS) and depth-first searches (DFS) on the graph to generate a more effective graph embedding than DeepWalk. Walklets [19] additionally incorporates explicit modeling in random walks. Hierarchical Representation Learning for Networks (HARP) [7] proposes a better initialization strategy to avoid the local optima in optimization.

Neural network based methods. SDNE [21] stacks multiple layers of non-linear functions to preserve highly non-linear network structure. It adopts an auto-encoder structure which uses the embedding to reconstruct its neighbors. DNGR [6] feeds the positive point-wise mutual information matrix into a stacked denoising autoencoder to capture higher-order proximity in the learned graph embedding. Nevertheless, SDNE and DNGR consider the whole graph and take as input the global neighborhood of each node, which are not efficient for large-scale graphs. Recently, graph convolution neural network (GCN) provides an effective and efficient solution by adopting a configuration with local constraints. These methods can be categorizes into spatial-based methods [2], [8], [11], [16], [20], [22] and spectral-based methods [4], [9], [12]-[15]. Spatialbased methods directly conduct convolution on the original graph. In contrast, spectral-based methods conduct convolution on the spectrum of the adjacent matrix of the graph.

III. METHOD

In this section, we introduce graph-based embedding learning for modeling the APP-installation information of users.

A. Graph Decomposition



Fig. 2. The visualization of decomposing a user-APP graph into subgraphs. The user-APP graph \mathcal{G} consists of the node set $\mathcal{V} = \{p_i\}_{i=1}^3 \cup \{u_i\}_{i=1}^8$ where p_i denotes an APP node and u_i is a user node. \mathcal{G} is decomposed into three subgraphs $\{\mathcal{G}_i\}_{i=1}^3$. \mathcal{G}_i consists of the node set \mathcal{V}_i . In this example, $\mathcal{V}_1 = \{p_1, u_4, u_5, u_6\}, \mathcal{V}_2 = \{p_2, u_1, u_2, u_3\}, \text{ and } \mathcal{V}_3 = \{p_3, u_7, u_8, u_1\}.$ Note that, the user u_1 is connected with two APPs p_2 and p_3 . Thus, u_1 is included in two subgraphs, \mathcal{G}_1 and \mathcal{G}_2 .

Definition. We denote the set of APPs used for building the graph by $\{s_i\}_{i=1}^N$, and denote the set of users by $\{t_j\}_{j=1}^M$. They constitute the node set $\mathcal{V} = \{p_1, \dots, p_N, u_1, \dots, u_M\}$. Meanwhile, the edge set \mathcal{E} contains all edges connecting two nodes $\{e_{s_i,t_i}\}_{i=1}^L$, where s_i denotes the index of the user and t_i denotes the index of the user in the *i*-th edge, e_{s_i,t_i} . That is, the existence of edge e_{s_i,t_i} means that the user u_{s_i} has installed the APP a_{t_i} . The user-APP graph \mathcal{G} is constructed based on the node set and the edge set $\{\mathcal{E}, \mathcal{V}\}$. We further define the subgraph \mathcal{G}_i with the node set \mathcal{V}_i and the edge set \mathcal{E}_i . \mathcal{V}_i contains only one APP node p_i and the user nodes connected to p_i . \mathcal{E}_i contains all edges which connect the APP node p_i . We visualize the process of decomposing a graph into a set of subgraphs in Figure 3.

B. Graph Learning

Initialization. We denote the embedding of the user u_i by \mathbf{u}_i and the embedding of an APP p_j by \mathbf{p}_j . We denote the indices of users installing the APP u_i by \mathcal{I}_i . The user embeddings are randomly initialized. In parallel, an APP embedding \mathbf{p}_j is initialized by averaging the embeddings of users installing the APP:

$$\mathbf{p}_j = \frac{\sum_{k \in \mathcal{P}_i} \mathbf{u}_k}{|\mathcal{P}_i|},\tag{1}$$

where $|\mathcal{P}_i|$ denotes the cardinality of the set \mathcal{P}_i , *i.e.*, the number of users installing the APP p_i .

Subgraph sampling. As we mentioned, for a subgraph G_i , it contains an APP node (p_i) and the users installing the APP p_i . Let us denote the probability of sampling the subgraph G_i as $P(G_i)$. A native sampling approach is sampling the sub-graph with a probability proportional to the number of user nodes in the subgraph. That is,

$$P(\mathcal{G}_i) = \frac{|\mathcal{P}_i|}{\sum_{j=1}^N |\mathcal{P}_j|}, \forall i \in [1, N],$$
(2)

where N is total number of APPs and $|\mathcal{P}_i|$ denotes the number of users in the subgraph \mathcal{G}_i . In this case, each edge connecting a user and an APP will be involved in the training process with an equal probability. Nevertheless, this strategy will make the embedding learning dominated by the popular APPs with a huge number of users and the contributions from some niche APPs with a small number of users will be underestimated. To make the contributions from different APPs balanced, we can devise that the sampling probability of each sub-graph to be equal. That is,

$$P(\mathcal{G}_i) = \frac{1}{N}, \forall i \in [1, N].$$
(3)

In this case, the edges based on niche APPs with a small number of users will be over-sampled, and the edges based on the popular APPs with a huge number of users will be undersampled. Nevertheless, it might lead to repeatedly sampling for edges from niche APPs, and some edges from the popular APPs might have little chance to be involved in the training process. It tends to make the learned embedding prone to overfitting due to a lack of diversity in the training samples. To achieve a balanced sampling and meanwhile suppress overfitting, we adopt a trade-off sampling approach. It devises the probability as

$$P(\mathcal{G}_i) = \frac{|\mathcal{P}_i|^{\tau}}{\sum_{j=1}^N |\mathcal{P}_j|^{\tau}}, \forall i \in [1, N],$$
(4)

where τ is a pre-defined positive constant. Normally, we set $0 < \tau < 1$. It assigns a higher sampling probability to the subgraph containing more nodes for suppressing over-fitting and meanwhile achieving a good balance among different APPs. When $\tau = 1$, it degenerates to the naive sampling approach defined in Eq. (2). On the other hand, when $\tau = 0$, it degenerates to the balanced sampling approach defined in Eq. (3). By default, we set $\tau = 0.5$ in our experiments.

Embedding learning within a subgraph. Let denote the app embedding with a subgraph by \mathbf{p} , the embedding of a user installing the APP by \mathbf{u}_i^+ and that of a user who does not install the APP by \mathbf{u}_j^- . The user and APP embedding learning seeks to keep a large similarity between \mathbf{p} and \mathbf{u}_i^+ . At the same time, it seeks to maintain a small similarity between \mathbf{p} and \mathbf{u}_j^- . Straightforwardly, we can learn the user and the APP embedding through a pairwise loss:



Fig. 3. The visualization of user and APP embedding learning within a subgraph. In this example, the APP embedding is \mathbf{p} (green box). There are four users installing the APP, $\{\mathbf{u}_i^+\}_{i=1}^4$ (blue dots) with the centroid \mathbf{u}^c (yellow dot) and four users not installing the APP, $\{\mathbf{u}_i^-\}_{i=1}^4$ (red dots).

$$\mathcal{L}_{\text{pair}} = \frac{1}{n^+} \sum_{i=1}^{n^+} \log(1 + e^{-\beta s(\mathbf{p}, \mathbf{u}_i^+)}) - \frac{1}{n^-} \sum_{i=1}^{n^-} \log(1 + e^{-\beta s(\mathbf{p}, \mathbf{u}_i^-)})$$
(5)

where n_+ denotes the number of users installing the APP and n_- denotes the number of users who do not install, β is a predefined positive constant controlling the softness, and $s(\cdot, \cdot)$ measures the cosine similarity between two vectors. In parallel to the pairwise loss defined above, we devise an additional centroid loss to further enhance the effectiveness of the learned embedding. To be specific, we first compute the centroid of the embeddings of users installing the APP:

$$\mathbf{u}_{c} = \frac{1}{n_{+}} \sum_{i=1}^{n^{+}} \mathbf{u}_{i}^{+}.$$
 (6)

Then the centroid loss is computed by

$$\mathcal{L}_{\text{centroid}} = \log(1 + e^{-\beta s(\mathbf{p}, \mathbf{u}_c^+)}).$$
(7)

To stabilize the training, we update the user embedding and the APP embedding in an alternating manner:

- Fix user embedding {u_i⁺}^{N₊}_{i=1} and {u_i⁻}^{N₋}_{i=1}, and update the APP embedding p using the centroid loss L_{centroid}.
- 2) Fix the APP embedding **p**, and update the positive user embedding $\{\mathbf{u}_i^+\}_{i=1}^{N_+}$ using the pairwise loss $\mathcal{L}_{\text{pair}}$.

To improve the training efficiency, we achieve this iterative training manner in a parallel way by utilizing the stop-gradient trick. That is, we devise the final loss $\mathcal{L} = \mathcal{L}_{centroid} + \mathcal{L}_{pair}$. In the meanwhile, we stop the gradient derived by $\mathcal{L}_{centroid}$ backpropagating to $\{\mathbf{u}_i^+\}_{i=1}^{N_+}$ and $\{\mathbf{u}_i^-\}_{i=1}^{n_-}$ and meanwhile stop the gradient from \mathcal{L}_{pair} back-propagating to \mathbf{p} and $\{\mathbf{u}_i^-\}_{i=1}^{n_-}$.

IV. EXPERIMENTS

Dataset. To train the model, we collect the information of 80 million users and 50 thousand APPs. On average, each user installs around 30 APPs.

A. Offline experiments

Memory. For each APP, we randomly sample 96 users who have already installed the APP and 96 users not installing the APP. Note that these APP installation has been involved in the training process. For each APP-user pair, we compute the cosine similarity between their embeddings. Then we threshold the cosine similarity to 0 or 1 to predict whether the user has installed the APP or not. In Table I, we show the experimental result. As shown in the table, in the training data, the learned embedding can achieve a 0.953 precision and 0.981 AUC, which demonstrates the powerful fitting capability of the learned embeddings.

 TABLE I

 The memory performance of the learned user and APP

 embeddings. We report the prediction precision and

 area-under-curve (AUC) for the APP installation.

Precision	0.953
AUC	0.981

Inference. To evaluate the inference performance of the learned user and APP embedding, we report the classification AUC on the user side and that on the APP side. The user-side AUC is averaged over users. For each user, we test the prediction accuracy using several APPs the user has installed and several APPs the user does not install. The APP-side AUC is measured in a similar manner but is averaged over APPs.

TABLE II

The inference performance of the learned user and APP embeddings. AUC⁺ denotes the AUC excluding APPs with more than 8% users and AUC^{*} denotes the AUC excluding APPs with more than 2.5% users.

	AUC	AUC ⁺	AUC*
APP-side	0.797	0.840	0.854
User-side	0.786	0.829	0.844

Note that the testing cases for inference are not involved in the training process. To be specific, our whole data is collected during N days. We use the data in the first N - 5 days for training and that from the last 5 days for testing. Meanwhile, we also report the AUC without excluding APPs with a huge number of users. To be specific, we report AUC^{*}, which excludes APPs with more than 2.5% users. We also report AUC⁺, which excludes that with more than 8% users. As shown in Table II, the AUC achieved in the inference is lower than that in Table I. In the meanwhile, by excluding some APPs with a huge number of users, AUC⁺ and AUC^{*} are larger than AUC.

Ablation study Here, we investigate the influence of removing $\mathcal{L}_{centroid}$ or the stop-gradient strategy through ablation study. As shown in Table IV, when removing $\mathcal{L}_{centroid}$, the AUC drops from 0.981 to 0.977 and the precision decreases from 0.953 to 0.948. Meanwhile, without the stop-gradient strategy, both the AUC and the precision decrease considerably.

TABLE III The influence of removing $\mathcal{L}_{centroid}$ or the stop-gradient strategy in inference.

	Ours	w/o $\mathcal{L}_{ ext{centroid}}$	w/o stop-gradient
Precision	0.953	0.948	0.942
AUC	0.981	0.977	0.973

B. Online experiments

We have integrated the user embedding learned from the APP-installation information as a feature which complements the existing user embedding learned from historical behaviors. After launching in our online personalized advertising platform, it has achieved a +1.1% CTR improvement, a +1.7% CVR boost, a +2.6% increase in the revenue.

TABLE IV The online experiments in our online advertising platform during one week.

CTR	CVR	Revenue
+1.1%	+1.7%	+2.6%

V. CONCLUSION

In this paper, we exploit the APP-installation information to assist in modeling the user's characteristics for personalized advertising. To this end, we build a user-APP bipartite graph and adopt a graph convolution network to learn the user embedding. We use the learned user embedding from our user-APP graph as the complementary information to the existing user representation learned from the user profile and the user's historical behaviors. After deploying it in our advertising platform, both CTR and CVR improve considerably.

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