

# Universe Points Representation Learning for Partial Multi-Graph Matching

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## Abstract

Many challenges from natural world can be formulated as a graph matching problem. Previous deep learning-based methods mainly consider a full two-graph matching setting. In this work, we study the more general partial matching problem with multi-graph cycle consistency guarantees. Building on a recent progress in deep learning on graphs, we propose a novel data-driven method (URL) for partial multi-graph matching, which uses an object-to-universe formulation and learns latent representations of abstract universe points. The proposed approach advances the state of the art in semantic keypoint matching problem, evaluated on Pascal VOC, CUB, and Willow datasets. Moreover, the set of controlled experiments on a synthetic graph matching dataset demonstrates the scalability of our method to graphs with large number of nodes and its robustness to high partiality.

## Introduction

Many real-world problems of finding correspondences can be formalized as graph matching. The examples of tasks in computer vision and graphics, where graph matching plays a crucial role, are object tracking (Wang et al. 2015), optical flow estimation (Sun et al. 2018), shape matching (Gao et al. 2021) and image keypoint matching (Fey et al. 2020). Graph matching is also relevant in other domains for matching protein networks (Singh, Xu, and Berger 2008), comparing molecules (Kriege, Humbeck, and Koch 2019), or linking users in social networks (Zhang and Philip 2015).

The recent progress in end-to-end learning of graph matching has allowed one to find the most useful graph features instead of hand-crafting them (Zanfir and Sminchisescu 2018), and also enabled the incorporation of axiomatic graph matching solvers into learning models (Rolínek et al. 2020). In this work, we specifically focus on finding image keypoint correspondences across an image collection with the help of deep learning techniques. We consider the general setting assuming that the images in the collection depict different instances of objects from the same category, as illustrated in Fig. 1 for the object category `cat`. A strong advantage of our multi-matching approach compared to previous pairwise matching formulations (Fig. 1a) is that due

to our *object-to-universe matching representation* we obtain cycle-consistent multi-matchings. Moreover, as opposed to the previous template-based deep multi-graph matching approach (Fig. 1b), our object-to-universe matching representation does not require an explicit template shape. Instead, we propose to learn latent representations of abstract universe points through aligning them with graph node features (Fig. 1c, Fig. 3 right). The learned embeddings encode higher-order information with the help of graph neural network processing and allow to find accurate object-to-universe matchings.

In addition, the object-to-universe formulation naturally supports the challenging partial multi-matching setting, as demonstrated in Fig. 1d. This is in stark contrast to a large portion of recent deep graph matching approaches that only tackle the full (bijective) matching problem. However, in a real-world collection of images, not all keypoints may be observed in each image, e.g. due to ambiguities of 2D image projections stemming from occlusions, camera-view change, object deformation, shape variability, etc. Hence, these methods require a keypoint pre-filtering before matching. As such, the partial matching setting is the most natural one and has thus the highest relevance in practical scenarios, since it can be applied to *unfiltered keypoints*.

Since image keypoints represent 2D projections of 3D points, we would like to expand the dimensionality of the matched keypoints. Therefore, we propose to learn a third *virtual* coordinate lifting (Fig. 4), and to incorporate it into the graph neural network architecture (3D SplineCNN (Fey et al. 2018)). The third dimension is learned with respect to the task of graph matching, thus it is a virtual coordinate. As we demonstrate in our ablation study in Sec. , the use of 3D coordinates leads to an improved matching performance.

Overall, our method advances the state of the art (SOTA) on various keypoint matching benchmarks. Namely, it achieves a 71.7% F1-score on the Pascal VOC dataset and a 95.1% F1-score on the CUB-200-2011 dataset on partial multi-graph matching without keypoint filtering. Additionally, it gives 81.8% accuracy on Pascal VOC dataset on full two-graph matching with intersection-filtered keypoints. Moreover, we show that the learned node features taken from a model pre-trained on (reduced) Pascal VOC can be transferred without fine-tuning to the Willow ObjectClass dataset, yielding 98.9% accuracy. On a set of controlled ex-

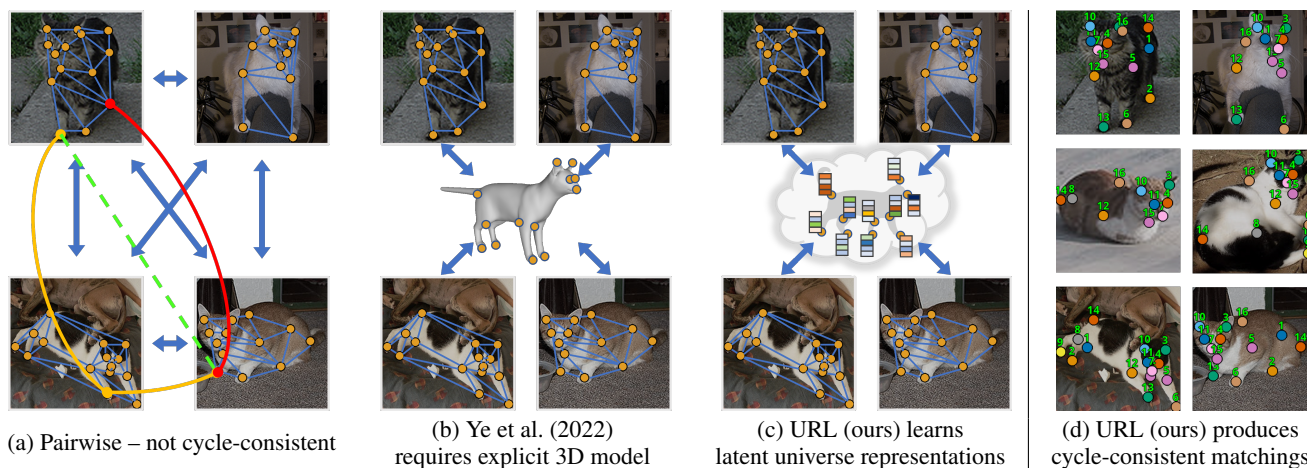


Figure 1: Overview of pairwise and multi-graph matching methods. (a) Pairwise matchings from prior works do not guarantee cycle consistency. The *red* arc shows an incorrect matching that violates the cycle consistency constraint, indicated by the *green dashed* line. (b) Previous cycle-consistent approach by Ye et al. (2022) requires an explicit template of sparse 3D geometry of the universe object. (c) Our approach learns latent representations of abstract universe points containing geometric and visual information, which allows to produce accurate cycle-consistent matchings under non-rigid deformations and occlusions (d).

periments, we demonstrate the robustness of the proposed approach to the partiality of the input data, its scalability to large graphs, and computational efficiency compared to the strong baseline BBGM-Multi (Rolínek et al. 2020).

We summarize our main contributions as follows:

- We propose the deep learning approach for *partial* multi-graph matching based on universe points representation learning;
- Building on the recent progress in graph-based learning, we introduce a virtual coordinate lifting, thus improving 2D image keypoint matching performance;
- The proposed method achieves state-of-the-art results in partial multi-graph matching with unfiltered keypoints on Pascal VOC and CUB-200-2011 datasets.

## Related Work

In this section, we introduce prior works on deep graph matching and discuss related methods.

**Deep graph matching** Graph matching consists of two main parts: feature extraction from raw input data and a matching algorithm. The seminal paper (Zanfir and Sminchisescu 2018) on two-graph matching has demonstrated how learnable features can be combined with a differentiable graph matching solver. Recent progress in this area allows differentiating through general combinatorial solvers (Rolínek et al. 2020). Another direction considers solving the simple linear assignment problem (LAP) (Munkres 1957) (via differentiable Sinkhorn iterations (Adams and Zemel 2011)) instead of the quadratic assignment problem QAP (Lawler 1963). The higher-order information is incorporated into the node features via graph neural network processing (Yu et al. 2019). This also allows an iterative improvement of soft matchings (Fey et al. 2020). Another way to solve a graph matching problem is

by learning a solver on a product graph (Wang et al. 2020). While these methods allow end-to-end learning, they focus on matching pairs of graphs, and for most of the solvers, the unrealistic setting of full (bijective) matchings is assumed. In contrast, we match multiple graphs in a cycle-consistent way and naturally allow for the realistic setting of partial matching.

**Multi-Graph Matching** In previous deep learning works that addressed multi-graph matching, cycle consistency was obtained via permutation synchronization of predicted pairwise matchings (Wang, Yan, and Yang 2020, 2021). In contrast, when using an object-to-universe matching formulation, cycle consistency can be directly obtained by construction (Schmidt et al. 2007; Huang and Guibas 2013). The deep learning-based approach (Ye et al. 2022) operates in this setting, but explicitly reconstructs a 3D model of a universe graph. While this gives the desired cycle consistency, the method requires 3D reconstruction, and it does not perform on par with the state-of-the-art pairwise matching methods (Rolínek et al. 2020; Wang, Yan, and Yang 2021). In our work, we propose to find an object-to-universe matching without explicitly reconstructing a 3D model of the universe. Instead, we learn latent representations of universe points by matching them with node features.

**Graph Neural Networks** One of the main sources of the recent progress in deep graph matching is the rise of graph neural network (GNN) architectures (Scarselli et al. 2008). Previous works have implemented various GNNs with intra- and cross-graph communication to optimally learn interconnections between nodes (Wang, Yan, and Yang 2019). Particularly, for the problem of keypoint matching, a SplineCNN architecture (Fey et al. 2018, 2020) has shown promising performance (Wang, Yan, and Yang 2021). The reason is that it allows to efficiently incorporate geomet-

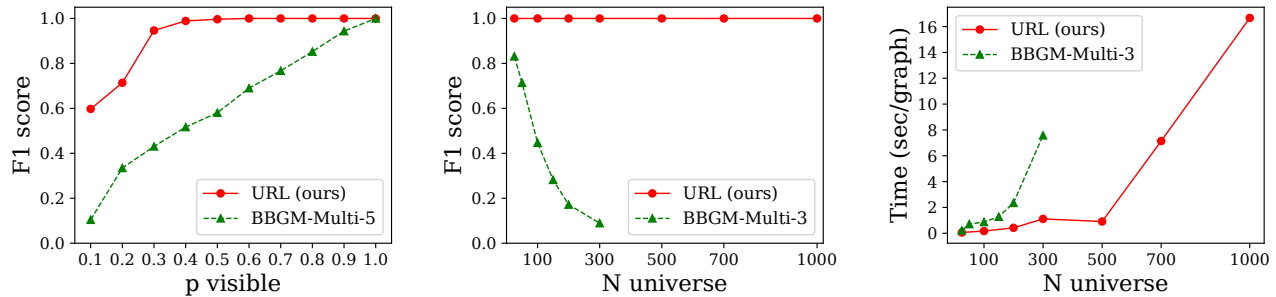


Figure 2: Partial multi-graph matching performance evaluation on synthetic dataset. Our method (URL) handles well high partiality in input points (left plot, low visibility). Its accuracy does not degrade on larger problems (middle plot, large number of points). And it scales more efficiently (right plot, processing time) than previous method BBGM-Multi.

ric information directly into the node features. Similarly to these methods, we utilize the SplineCNN architecture for our image keypoint matching problem. However, we propose to also predict a third virtual coordinate, which helps to resolve 2D keypoint location conflicts. As we later demonstrate, this leads to an improved matching performance.

### Background

In the following, we provide the general background for the multi-graph matching problem. First, we introduce the notation concerning graphs and matchings. Then we revisit the theory behind the object-to-universe multi-graph matching.

#### Notation

We assume that  $k \in \mathbb{N}$  different graphs  $\mathcal{G}_1, \dots, \mathcal{G}_k$  are given, where for each  $i \in [k] := \{1, \dots, k\}$  the graph  $\mathcal{G}_i = (\mathcal{V}_i, \mathcal{E}_i, F_i, E_i)$  comprises of a node set  $\mathcal{V}_i$ , an edge set  $\mathcal{E}_i \subset \mathcal{V}_i \times \mathcal{V}_i$ , node features  $F_i : \mathcal{V}_i \rightarrow \mathcal{F}_1$ , and edge features  $E_i : \mathcal{E}_i \rightarrow \mathcal{F}_2$ . The feature spaces  $\mathcal{F}_1$  and  $\mathcal{F}_2$  are some finite-dimensional vector spaces and we write  $m_i = |\mathcal{V}_i|$  for the amount of nodes in  $\mathcal{G}_i$ .

In this work, we consider the task of *partial* multi-graph matching, in which we want to find correspondences between nodes of the graphs  $\mathcal{G}_1, \dots, \mathcal{G}_k$ . Note that the number of nodes  $m_i$  can differ among the graphs  $\mathcal{G}_i$ , i.e. not all of the matched nodes are present in each graph, which backs the term ‘partial’. Before we introduce our formalism for multi-graph matching, we introduce the correspondence  $X_{ij} \in \mathbb{P}_{m_i m_j}$  between a pair of graphs  $(\mathcal{G}_i, \mathcal{G}_j)$ , where the set of *partial permutation matrices* is given as

$$\mathbb{P}_{m_i m_j} = \{X \in \{0, 1\}^{m_i \times m_j} : X \mathbf{1}_{m_j} \leq \mathbf{1}_{m_i}, \mathbf{1}_{m_i}^\top X \leq \mathbf{1}_{m_j}^\top\}, \quad (1)$$

for  $\mathbf{1}_m$  being the  $m$ -dimensional vector of ones.

#### Multi-graph matching

When considering multiple ( $k > 2$ ) graphs, it is necessary to ensure that the pairwise matchings  $\{X_{ij}\}_{i,j=1}^k$  do not contradict each other. Namely, the matchings should be *cycle-consistent* (Bernard et al. 2019a):

**Definition 1** (Cycle consistency). *The set  $\mathcal{X} = \{X_{ij}\}_{i,j=1}^k$  of pairwise matchings is said to be cycle-consistent if for all  $i, j, \ell \in [k]$  it holds that*

$$X_{ii} = I_{m_i}, \quad (\text{Identity})$$

$$X_{ij} = X_{ji}^\top, \quad (\text{Symmetry})$$

$$X_{ij} X_{j\ell} \leq X_{i\ell}. \quad (\text{Partial Transitivity})$$

The property of (Partial Transitivity) introduces a cubic number of non-convex quadratic constraints. A more efficient way to characterize cycle-consistent matchings is by using the notion of *universe matchings*.

Assuming that we are given an upper bound of the *universe size*  $d \in \mathbb{N}$ , for graph  $\mathcal{G}$  with  $m$  nodes ( $m \leq d$ ) the space of *object-to-universe* matchings is denoted as

$$\mathbb{U}_{m,d} = \{X \in \{0, 1\}^{m \times d} : X \mathbf{1}_d = \mathbf{1}_m, \mathbf{1}_m^\top X \leq \mathbf{1}_d^\top\}. \quad (2)$$

Then the cycle-consistency constraints can be reformulated in the following manner (Huang and Guibas 2013):

**Lemma 1** (Universe matchings). *The set  $\mathcal{X} = \{X_{ij}\}_{i,j=1}^k$  of pairwise matchings is cycle-consistent, iff there exists a universe of size  $d$  and a collection of object-to-universe matchings  $\{X_i \in \mathbb{U}_{m_i,d}\}_{i=1}^k$ , such that for each  $X_{ij} \in \mathcal{X}$  it holds that  $X_{ij} = X_i X_j^\top$ .*

In other words, by matching each node of a graph with exactly one abstract *universe point*, we can find object-to-universe matchings  $\{X_i \in \mathbb{U}_{m_i,d}\}_{i=1}^k$ , which by construction ensures the cycle consistency of pairwise matchings.

### Proposed Method

Below, we introduce our method for partial multi-graph matching by **universe points representation learning (URL)**. We first formulate the general problem setup and clarify the assumptions of our method. Then we formalize the proposed approach as a deep learning problem and introduce our proposed neural network architecture with a novel virtual coordinate lifting block.

#### Problem Setup and Assumptions

Our method accepts as input a set of graphs  $\{\mathcal{G}_i\}_{i=1}^k$  that are to be matched. It outputs globally cycle-consistent pairwise matchings from object-to-universe matchings, i.e.  $\{\hat{X}_{ij} := \hat{X}_i \hat{X}_j^\top \in \mathbb{P}_{m_i m_j}\}_{i=1}^k$ .

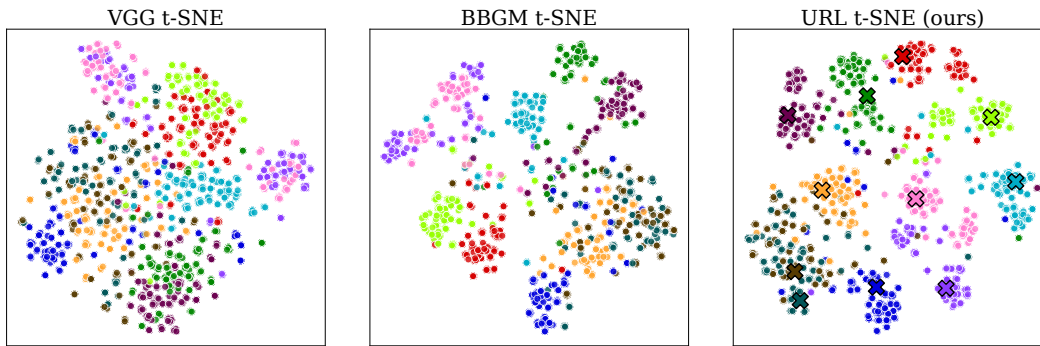


Figure 3: Two dimensional t-SNE embeddings of latent keypoint representations from different neural networks (on `bicycle` object class from test set of Pascal VOC dataset). The colours represent the ground-truth matchings of keypoints. Both BBGM (middle) and URL (right, ours) models start with the node features from the ImageNet pre-trained VGG network (left), and further improve them. We empirically demonstrate that matching keypoints with learnable latent representations of *universe points* (centroids  $\times$ ) yields more accurate results than matching keypoints with each other (see Table 1, BBGM vs URL).

**Object classes** Since the set of matched graphs  $\{\mathcal{G}_i\}_{i=1}^k$  share common nodes, we expect that the graphs correspond to the same abstract class, called object class. In other words, each graph represents an object from the same object class. We *assume* that the input graphs are from a known object class during inference. Therefore, we consider matchings for different object classes independently.

**Supervision** Further, we *require* the ground truth matchings for the training set as a supervisory signal. We note that any set of correct pairwise matchings can be transformed into object-to-universe matching form via matrix factorization (Bernard et al. 2019a). Therefore, we can always extract the ground truth universe matchings  $\{X_i^{\text{gt}}\}_{i=1}^k$  from pairwise matchings  $\{X_{i,j}^{\text{gt}}\}_{i,j=1}^k$  during training. From this, we *estimate* an upper bound of the number of universe points  $d$ .

The above-stated assumptions hold for multi-image keypoint matching problems with predefined sets of detected keypoints sharing similar semantic meanings. For example, commonly used datasets, such as Pascal VOC, Willow ObjectClass, CUB-200-2011, provide keypoint annotations as node labels, i.e. in object-to-universe matching form, and give the ground truth number of universe points  $d$ .

## Deep Object-to-Universe Matching

Given the problem setup, we propose to utilize the object-to-universe matching formulation from Lemma 1. This allows to match a set of graphs  $\{\mathcal{G}_i\}_{i=1}^k$  with abstract universe points individually. Thus, it reduces the memory and time consumption compared to multi-graph matching approaches based on synchronization of all pairwise matchings.

To facilitate the data-driven approach, we propose to simultaneously learn the feature encoder of graph nodes and the latent representations of universe points. Formally, given a graph  $\mathcal{G}_i$  with its initial node and edge features  $F_i$  and  $E_i$ , we update them to produce new node and edge features  $\hat{F}_i$  and  $\hat{E}_i$ . This is achieved with a parameterized graph-processing function  $\Psi_\theta$ , where  $\theta$  denotes the learnable pa-

rameters:

$$\Psi_\theta(\mathcal{G}_i) =: \hat{\mathcal{G}}_i = (\mathcal{V}_i, \mathcal{E}_i, \hat{F}_i, \hat{E}_i) \quad (3)$$

In practice,  $\Psi_\theta$  is implemented as a graph neural network. As a consequence, the node features contain the neighborhood (including edge) information. Thus, we will only use the improved node features  $\hat{F}_i \in \mathbb{R}^{m_i \times h}$  for matching. The features are from  $h$ -dimensional vector space.

Concurrently, we learn the latent representations of  $d$  universe points  $\hat{U} \in \mathbb{R}^{d \times h}$  in the same  $h$ -dimensional vector space. The *soft* (non-binary, continuous) object-to-universe matching is then found by the normalization of feature similarities as

$$\hat{S}_i := \text{Norm}(\hat{F}_i \hat{U}^\top) \in [0, 1]^{m_i \times d}. \quad (4)$$

The normalization can be performed row-wise, i.e. for each node individually, since every node of the graph has exactly one match among the universe points. So for each row (node)  $v$ , we apply

$$\text{Norm}(\hat{F}_i \hat{U}^\top)_v := \text{SoftMax}\left(\left(\hat{F}_i \hat{U}^\top\right)_v\right), \quad (5)$$

where  $\text{SoftMax}(z)_j = \frac{e^{z_j}}{\sum_k e^{z_k}}$ . Note, that in the general case of partial pairwise matching the simple normalization is not applicable, because not every node can be matched.

We predict the discretized object-to-universe matching  $\hat{X}_i$  from soft matching  $\hat{S}_i$  by solving a linear assignment problem (LAP) with an efficient auction algorithm (Bertsekas 1998)

$$\hat{X}_i := \underset{X \in \mathbb{U}_{m_i d}}{\text{argmax}} \langle X, \hat{S}_i \rangle = \text{LAP}(\hat{S}_i). \quad (6)$$

Finally, to obtain cycle-consistent pairwise matchings, we combine corresponding individual object-to-universe matchings, i.e.  $\{\hat{X}_{ij} = \hat{X}_i \hat{X}_j^\top\}_{i,j=1}^k$ .

For training graph feature encoder  $\Psi_\theta$  and universe representations  $\hat{U}$ , we minimize the discrepancy between soft matchings  $\hat{S}_i$  and ground truth universe matchings  $X_i^{\text{gt}}$ . We

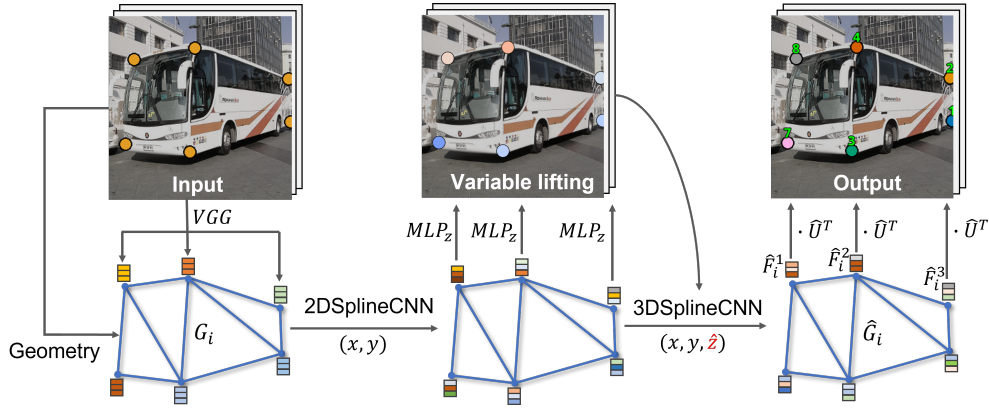


Figure 4: Proposed pipeline for image keypoint matching

propose to use a CrossEntropy loss ( $L_{CE}$ ) for each node individually

$$L_i := \frac{1}{m_i} \sum_{v=1}^{m_i} L_{CE} \left( (\hat{S}_i)_v, (X_i^{gt})_v \right), \quad (7)$$

where the  $L_{CE}(\vec{z}, \vec{y}) = -\sum_j y_j \log z_j$ . The total loss is a sum of graph-wise losses, i.e.  $L := \sum_{i=1}^k L_i$ .

**Universe Points Representations** For the fixed node features  $\{\hat{F}_i\}_{i=1}^k$  and given ground truth object-to-universe matchings  $\{X_i^{gt}\}_{i=1}^k$ , it is possible to analytically derive the universe points representations, optimal in  $L_2$ -distance manner, as the average vectors, i.e. *centroids*,

$$U^* := \operatorname{argmin}_U \sum_{i=1}^k \left\| X_i^{gt \top} \hat{F}_i - U \right\|^2 = \frac{1}{k} \sum_{i=1}^k X_i^{gt \top} \hat{F}_i. \quad (8)$$

The visualization of t-SNE embeddings (in cosine distance) of the learned keypoint features and the universe points representations in Fig. 3 (right plot) shows that the universe points are located at the regions with a high density of the matched keypoints, confirming the above intuition. Note that in the visualization the keypoints are taken from the test set, while universe points are learned from the training set.

## Architecture Design

In this section, we introduce a neural network architecture for the image keypoint matching problem.

Following previous works (Rolínek et al. 2020), we initialize the node features by interpolating ImageNet pre-trained VGG features (Simonyan and Zisserman 2015; Deng et al. 2009) at the keypoint locations (see left plot in Fig. 3), and the edge features – by relative keypoint locations. Further, we assume that the graph structure is fixed by the geometrical properties of keypoints, e.g. by Delaunay triangulation (Delaunay 1934). The graph processing network  $\Psi_\theta$  is implemented using SplineCNN (Fey et al. 2018) blocks.

In addition to previous graph neural network architecture from BBGM (Rolínek et al. 2020), we propose to utilize a predicted *third virtual coordinate*  $\hat{z}$  in the 3DSplineCNN

block. The intuition behind this virtual variable is the nature of keypoints in images, which represent 2D projections of 3D points. The value of the third coordinate is driven by the matching objective.

Overall, our graph processing module  $\Psi_\theta$  takes VGG features as input and combines 2DSplineCNN and 3DSplineCNN blocks. The third virtual coordinate  $\hat{z}$  is predicted by  $MLP_z$  using the output node features of 2DSplineCNN. The edge features are set as relative keypoint locations. The proposed architecture is illustrated in Fig. 4.

## Experiments

We evaluate the proposed method on various keypoint matching benchmarks. First, we introduce compared methods and metrics. Then we assess different aspects of our method, such as robustness to partiality and scalability, on a controlled synthetic dataset. We evaluate the real-world performance on Pascal VOC and CUB-200-2011 datasets. We also estimate the transferability of the learned node feature encoder to other datasets (Pascal  $\rightarrow$  Willow). Finally, we analyze the contribution of each component of the proposed approach by ablation studies.

### Metrics and Peer Methods

In *partial* multi-graph matching (i.e. all keypoints are included,  $\cup$ ), the pairwise matchings  $\{X_{ij} \in \mathbb{P}_{m_i m_j}\}$  contain both inlier matches and non-matches (zero rows/columns in  $X_{ij}$ ). This leads to both types of false positive and false negative errors. To measure the balanced performance of partial matching, we use **F1** score, the harmonic mean of precision and recall. In this setting, we compare our method with recent deep learning-based multi-graph matching methods BBGM-Multi (Rolínek et al. 2020) and JointMGM3D (Ye et al. 2022). We also evaluate learning-free multi-graph matching solvers HiPPI (Bernard et al. 2019b) and MGM-Floyd (Jiang, Wang, and Yan 2020) on Willow dataset.

If graphs contain only (the same number of) inliers (i.e. the intersection of all keypoints,  $\cap$ ), the matching problem becomes *full*, and is a special case of partial multi-graph matching. Since in this setting there are no non-matches, we

measure the performance of models by computing the accuracy (**Acc**) of predicted full pairwise matchings.

The majority of previous deep graph matching methods consider the full pairwise matching problem (i.e. the intersection of keypoints from two graphs,  $\cdot \cap \cdot$ ): GMN (Zanfir and Sminchisescu 2018), SuperGlue (Sarlin et al. 2020), PCA-GM (Wang, Yan, and Yang 2019), LCS (Wang et al. 2020), CIE-H (Yu et al. 2019), GANN-MGM (Wang, Yan, and Yang 2020), DGMC (Fey et al. 2020), BBGM (Rolínek et al. 2020), NHGM-v2 (Wang, Yan, and Yang 2021), GLAM (Liu et al. 2021). Our method can also be evaluated on this restricted setting, as it is a special case of partial multi-matching.

We categorize the methods according to their multi-graph *cycle-consistency* property into 3 categories. First, the methods that consider only two graphs for matching (**no**). Next, the methods that take as input  $N$  graphs, and synchronize the matchings between them – the cycle consistency is provided **locally** for the considered graphs. Lastly, the methods that provide cycle consistency without depending on the number of considered input graphs, e.g. by matching each individual graph with a common universe, are categorized as fully cycle-consistent (**yes**).

### Synthetic Partial Multi-Matching

We test various aspects of our method on a set of controlled experiments. For this, we create synthetic graphs following the setup from previous works (Wang, Yan, and Yang 2019; Cho, Lee, and Lee 2010). See details in appendix. We enforce *partiality* in generated graphs  $\{G_i\}$  by randomly sampling keypoints among total  $N_{\text{univ}}$  universe points for each graph from Binomial distribution with visibility rate  $p_{\text{vis}}$ , i.e.  $m_i \sim \text{Bin}(p_{\text{vis}}, N_{\text{univ}})$ .

We compare our method with the strongest peer method BBGM-Multi- $K$ . To assess the robustness to partiality, we vary the visibility rate  $p_{\text{vis}}$  (with fixed  $N_{\text{univ}} = 25$ ). Further we evaluate scalability and computational efficiency by altering number of universe points  $N_{\text{univ}}$  with fixed partiality  $p_{\text{vis}} = 0.8$ . As demonstrated in Fig. 2, our method scales well to large problems (graphs with up to 1000 nodes) and handles high partiality rates better than the method based on a multi-matching solver. Training of BBGM-Multi-3 on problems with more than 300 universe points does not fit into GPU memory (48 GB).

### Real-World Partial Multi-Matching

To evaluate our method on a real-world problem, we consider multi-image keypoint matching on challenging datasets with natural partiality: the Pascal VOC (Everingham et al. 2010; Bourdev and Malik 2009), and the CUB datasets (Wah et al. 2011). These datasets contain different object classes, and up to 23 universe points depending on object category (Fig. 5). The average visibility rate of Pascal VOC is  $p_{\text{vis}}^{\text{VOC}} = 0.72$ , and of CUB is  $p_{\text{vis}}^{\text{CUB}} = 0.8$ . We follow the data pre-processing procedures of previous methods (Wang, Yan, and Yang 2019).

The comparison of matching performance is presented in Tables 1-2 for Pascal VOC and CUB datasets respectively. So in the real-world setting without any keypoint

Method	F1, $\mathbf{U}$	Acc, $\cdot \cap \cdot$	Cycle
GMN	n/a	55.3	no
SuperGlue	n/a	58.1	no
PCA-GM	n/a	63.8	no
LCS	n/a	68.5	no
CIE-H	n/a	68.9	no
DGMC	n/a	73.0	no
NGM-v2	n/a	80.1	no
NHGM-v2	n/a	80.4	no
GLAM	n/a	<b>86.2</b>	no
BBGM	61.4	80.1	no
BBGM-Multi-5	62.8	-	locally
JointMGM3D	42.9 (59.4 <sup>†</sup> )	59.0 (67.1 <sup>†</sup> )	<b>yes</b>
URL ( <b>ours</b> )	<b>71.7 (82.9<sup>†</sup>)</b>	<b>81.8 (88.7<sup>†</sup>)</b>	<b>yes</b>

Table 1: F1 score (%) of partial multi-graph matching and accuracy of full two-graph matching on the Pascal VOC dataset (averaged over object classes). Since the majority of previous works consider full graph matching, their results are not available (*n/a*) in a partial multi-matching setup.

<sup>†</sup>Scores in parentheses are directly computed on object-to-universe matchings  $\{X_i\}$ .

Method	F1, $\mathbf{U}$	Cycle
GMN	79	no
PCA-GM	83.2	no
CIE-H	83	no
GANN-MGM	82.6	locally
URL (ours)	<b>95.1</b>	<b>yes</b>

Table 2: F1 scores (%) of partial multi-graph matching on CUB-200-2011 dataset. Our method produces the most accurate matchings with built-in cycle consistency.

pre-filtering ( $\mathbf{U}$ ), our method achieves state-of-the-art results on both datasets: 71.7% F1 on Pascal VOC, and 95.1% F1 on CUB-200-2011. On the object-to-universe matchings, our method achieves an F1-score of 82.9%, and thus outperforms the previous object-to-universe matching method JointMGM3D (Ye et al. 2022) by a large margin. Moreover, our method shows strong performance with 81.8% accuracy on a restricted setting of full two-graph matching ( $\cdot \cap \cdot$ ). The recent method GLAM (Liu et al. 2021) performs best on this setting. However, in contrast to our method, it learns the graph structures of given points. Thus, the improvement from GLAM is orthogonal to our main contribution.

### Node Feature Transfer

We study the generalization ability of our method by transferring a node feature encoder model  $\Psi_\theta$ , pre-trained on Pascal VOC dataset (without overlapping images), to Willow dataset (Cho, Alahari, and Ponce 2013). The Willow dataset contains images with consistent object orientations and the same number of keypoints, i.e. all matchings are bijective (see Fig. 5). We evaluate our model by computing cosine similarities of node features with a subsequent discretization by a combinatorial solver. This variation of our model does

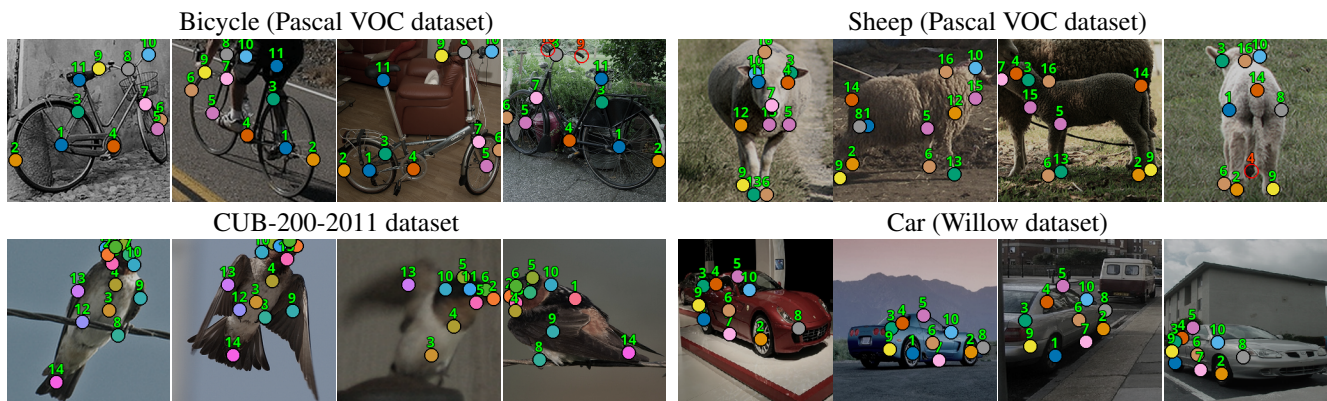


Figure 5: Qualitative results of our method. The *numbers* represent the ground-truth ordering of universe points. Correct predictions are marked with *green* and incorrect ones with *red* numbers. The datasets cover various poses, object deformations, and shape variability of different object classes. The keypoints are matched without any pre-filtering.

Data	Method	Acc, $\cap$	Cycle
—	HiPPI	88.2	locally
	MGM-Floyd	88.3	locally
Pascal $\uparrow$ Willow	GMN	76.2	no
	PCA-GM	85.2	no
	CIE-H	83.5	no
	DGMC	84.1	no
	BBGM	94.5	no
	URL-LAP	95.0	no
	URL-HiPPI	<b>98.9</b>	locally

Table 3: Accuracy of full matchings on Willow dataset. Models, pre-trained on (reduced) Pascal VOC dataset, are evaluated on the Willow dataset *without fine-tuning*. Our learned features in URL-HiPPI are better than the hand-crafted ones in a learning-free HiPPI solver, which allows us to also outperform previous learning-based methods.

not use universe points and depends on an external solver (denoted as URL-Solver).

From Table 3, we observe that our model URL-HiPPI with a multi-matching solver HiPPI outperforms all other methods, achieving 98.9% matching accuracy. Notably, our approach with the simple linear assignment solver, URL-LAP, produces results comparable with the strong baseline BBGM, which points to the successful transferability of node features in this experiment. The extended results with different training strategies on Pascal VOC and Willow datasets are presented in the supplementary materials.

### Ablation Study

We evaluate the effectiveness of each component of our proposed method via an ablation study, presented in Table 4. In this experiment, we solve partial multi-matching on the Pascal VOC dataset. We observe that using the 3DSplineCNN block with predicted virtual coordinate gives 4.1% F1-score improvement (rows 1, 4). The graph neural network processing  $\Psi_\theta$  (2D and 3D SplineCNN blocks, row 2) corresponds

VGG	2D	3D	URL	F1	Cycle
✓	✓	✗	✓	67.6±1.0	yes
✓	✗	✗	✓	55.1±0.6	yes
✓	✓	✓	✗	55.4±0.6	locally
✓	✓	✓	✓	<b>71.7±0.5</b>	yes

Table 4: Ablation studies of different components of the proposed method on Pascal VOC dataset. Each building block of the neural network (3D and 2D SplineCNN) significantly contributes to the final performance. The universe representation learning (URL) component has the highest impact both on the F1 score and on the cycle consistency.

to 16.6% of performance increase, identifying the value of graph-level feature encoding. More importantly, our learned universe points representations (URL, row 3) show a critical influence in providing cycle consistency and in achieving the optimal level of efficacy, accounting for 16.3% gain in F1-score. The reported improvements are statistically significant according to the non-parametric Mann-Whitney U test. The studied methods are run 5 times with random seeds.

### Conclusion

We have presented a deep learning approach for partial multi-graph matching. It builds upon an object-to-universe formulation via learning latent universe point representations. Our important conceptual contribution is that we demonstrate the effectiveness of matching against abstract universe representations, in contrast to pairwise matching, achieving the state-of-the-art performance in practically relevant settings. On a technical level, this is made possible by combining the recent progress in graph-based learning and the architectural improvement with a virtual coordinate lifting. The insights about universe points representations open up new future directions based on exploring unsupervised learning methods for deep multi-graph matching problems. We believe that the observations from our work will be beneficial in other domains with correspondence problems.

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