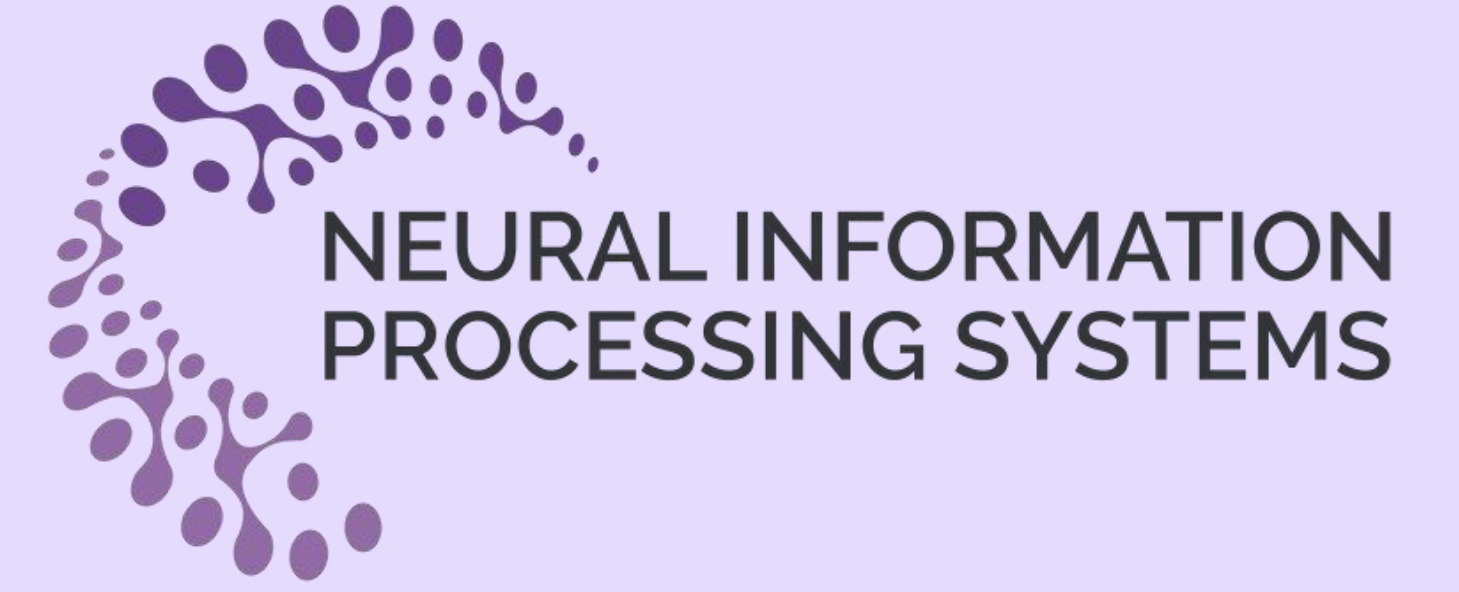


# Boosting Spectral Clustering on Incomplete Data via Kernel Correction and Affinity Learning

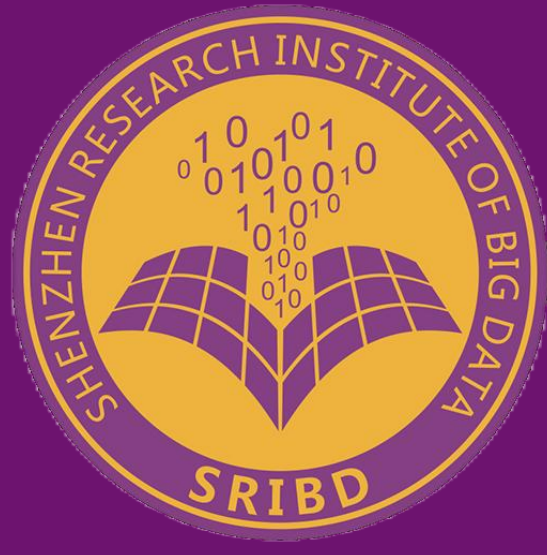
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## 1. INTRODUCTION

### Spectral Clustering on Complete Data:

#### Step 1. Construct A Similarity Matrix: $S = K$ or $C$

- Calculate a (Gaussian) kernel  $K$  from  $X$
- Learn a self-expressive affinity matrix  $C$  from  $X$

#### Step 2. Perform Normalized Cut Algorithm [1]

- Calculate the Laplacian matrix  $L$
- Select the first  $k$  eigenvectors of  $L$  as  $W$
- Conduct k-means clustering on  $W$

### Q: Spectral Clustering on Incomplete Data?

## 2. AIM

When dealing with incomplete data, how to

- 1) Estimate a high-quality kernel;
- 2) Learn a high-quality self-expressive affinity, benefiting the spectral clustering task.

## 3. METHOD

### I. Kernel Correction (KC) Algorithm

- A valid kernel should be positive semi-definite [2].
- **Step 1.** Estimate a naive kernel  $K^0$  on incomplete data;
- **Step 2.** Find a nearest PSD kernel as the new estimate.

$$\hat{K} = \underset{K \in \mathbb{R}^{n \times n}}{\operatorname{argmin}} \|K - K^0\|_F^2 \text{ s.t. } K \succeq 0, k_{ij} = k_{ji} \in [0, 1]$$

- $\|K^* - \hat{K}\|_F \leq \|K^* - K^0\|_F$  ( $K^*$  is the ground-truth) [3]

### II. Affinity Learning Algorithms

- Kernel Self-expressive Affinity Learning with  $\ell_p$  norm:

- **Proximal P-norm:**  $\|C\|_p^p := \sum_{i,j=1}^n |c_{ij}|^p$

$$\min_{C \in \mathbb{R}^{n \times n}} \|\phi(X) - \phi(X)C\|_F^2 + \lambda \|C\|_p^p \text{ s.t. } C_{ij} \in [0, 1]$$

- **Schatten P-norm:**  $\|C\|_{S_p} := (\sum_{i=1}^n \sigma_i^p(C))^{1/p}$

$$\min_{C \in \mathbb{R}^{n \times n}} \|\phi(X) - \phi(X)C\|_F^2 + \lambda \|C\|_{S_p} \text{ s.t. } C_{ij} \in [0, 1]$$

- Adaptive Kernel Least-Squares Representation:

$$\min_{K \succeq 0, C} \|K - K^0\|_F^2 + \operatorname{Tr}(K - 2KC + C^T KC) + \lambda \|C\|_F^2$$

## 4. RESULTS

### I. Performance on Gaussian Kernel Estimation

Relative error:  $\text{RE} = \|\hat{K} - K^*\|_F / \|K^*\|_F$ ; Recall: measures the accuracy of top-10 neighbors.

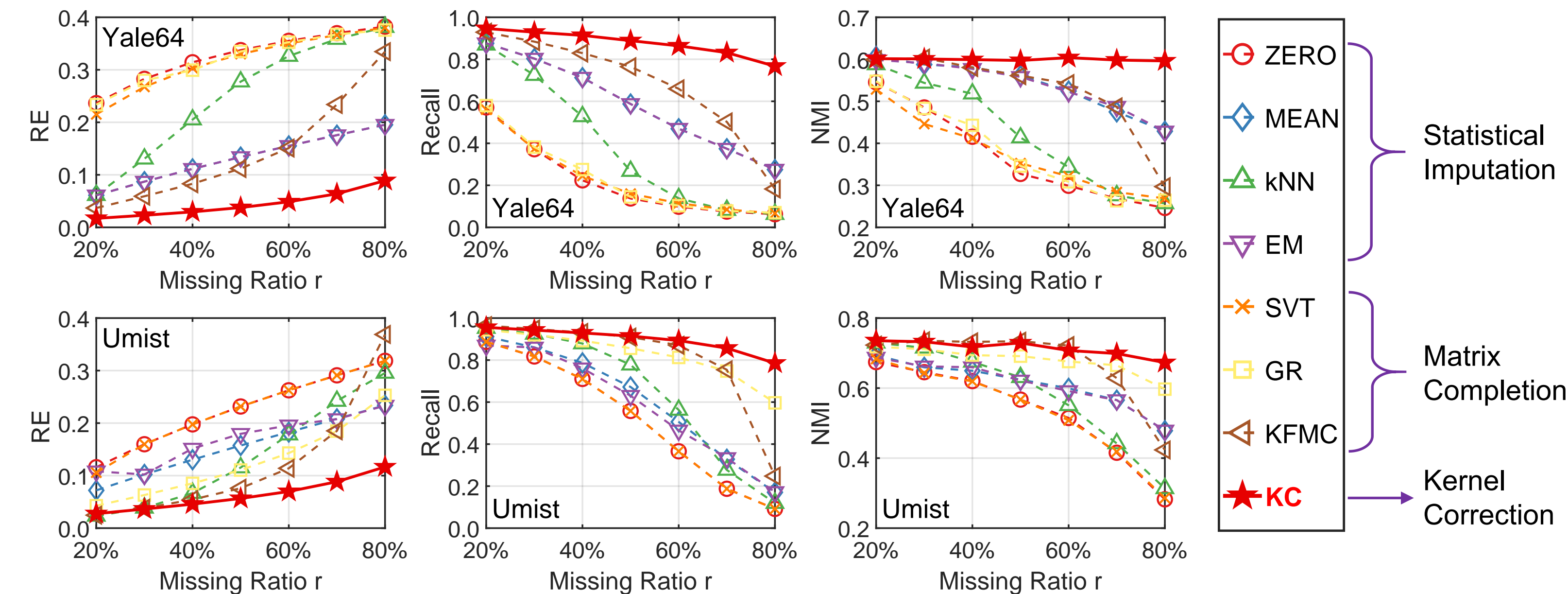
Table 1. Gaussian kernel estimation on the Yale64 dataset with 80% random missing.

Metric	Naive	ZERO	MEAN	kNN	EM	SVT	GR	KFMC	DC	TRF	EE	KC
RE↓	0.113	0.382	0.195	0.381	0.195	0.380	0.376	0.335	0.180	0.112	0.097	<b>0.089</b>
Recall↑	0.721	0.063	0.275	0.063	0.275	0.066	0.070	0.183	0.571	0.722	0.751	<b>0.767</b>

### II. Performance on Standard Spectral Clustering

Table 2. NMI of standard spectral clustering on Yale64 and Umist with 80% missing.

Dataset	Naive	ZERO	MEAN	kNN	EM	SVT	GR	KFMC	DC	TRF	EE	KC
Yale64	0.588	0.246	0.429	0.257	0.428	0.269	0.264	0.297	0.551	0.587	0.593	<b>0.596</b>
Umist	0.669	0.282	0.478	0.314	0.479	0.286	0.597	0.423	0.488	0.667	0.669	<b>0.673</b>



### III. Performance on Self-expressive Affinity Learning

KSL-Sp and AKLSR employ corrected kernels to yield dependable affinity matrices.

Table 3. NMI of spectral clustering using learned affinity on Yale64 with 80% missing.

Method	Naive	ZERO	MEAN	kNN	EM	SVT	GR	KFMC	DC	TRF	EE	KC
KSSC [4]:	0.219	0.215	0.167	0.173	0.177	0.218	0.208	0.259	0.588	0.210	0.209	<b>0.616</b>
KLSR [5]:	0.606	0.311	0.604	0.320	0.609	0.321	0.327	0.318	0.597	0.603	0.604	<b>0.616</b>
KSL-Sp (Ours):	0.370	0.315	0.581	0.303	0.579	0.305	0.304	0.295	0.555	0.364	0.599	<b>0.619</b>
AKLSR (Ours):	0.452	0.327	0.606	0.338	0.605	0.308	0.338	0.312	0.570	0.464	0.575	<b>0.614</b>

Three ways to estimate  $\hat{K}$ :

- **Data Imputation:**  $X^0 \rightarrow \hat{X} \rightarrow \hat{K}$
  - **Distance Calibration:**  $X^0 \rightarrow D^0 \rightarrow \hat{D} \rightarrow \hat{K}$
  - **Kernel Correction (KC):**  $X^0 \rightarrow K^0 \rightarrow \hat{K}$
- Smallest RE & Highest Recall**

- Missing completely at random
- an estimated Gaussian kernel  $\rightarrow$  kNN graph  $\rightarrow$  normalized cut algorithm
- Normalized Mutual Information

## 5. CONCLUSIONS

Q: How to boost spectral clustering on incomplete data via effective kernel and affinity learning?

A: We propose an imputation-free framework:

- Start with a naive kernel instead of imputing;
- Learn a high-quality **kernel matrix** by the kernel correction method with a theoretical guarantee;
- Learn a high-quality **affinity matrix** by the kernel self-expressive affinity learning algorithms with an adaptive extension using joint optimization.

Experimental results show the superiority:

- **Experiments I & II:** show our kernel correction method significantly outperforms data imputation and distance calibration approaches, especially for a large missing ratio (i.e., 80%).
- **Experiment III:** shows the effectiveness of KSL-Sp and AKLSR with comparable performance.

## 6. REFERENCES

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## 7. MORE INFORMATION



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