Identifying stable components of matrix/tensor factorizations via

low-rank approximation of inter-factorization similarity

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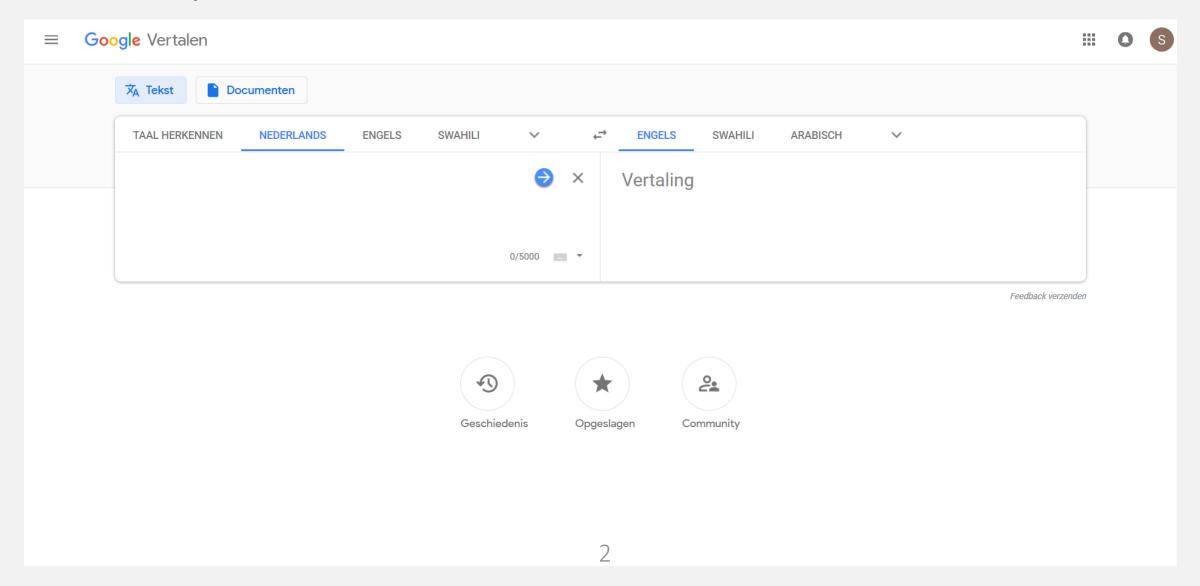
European Signal Processing Conference, A Coruña, Spain 3 September 2019



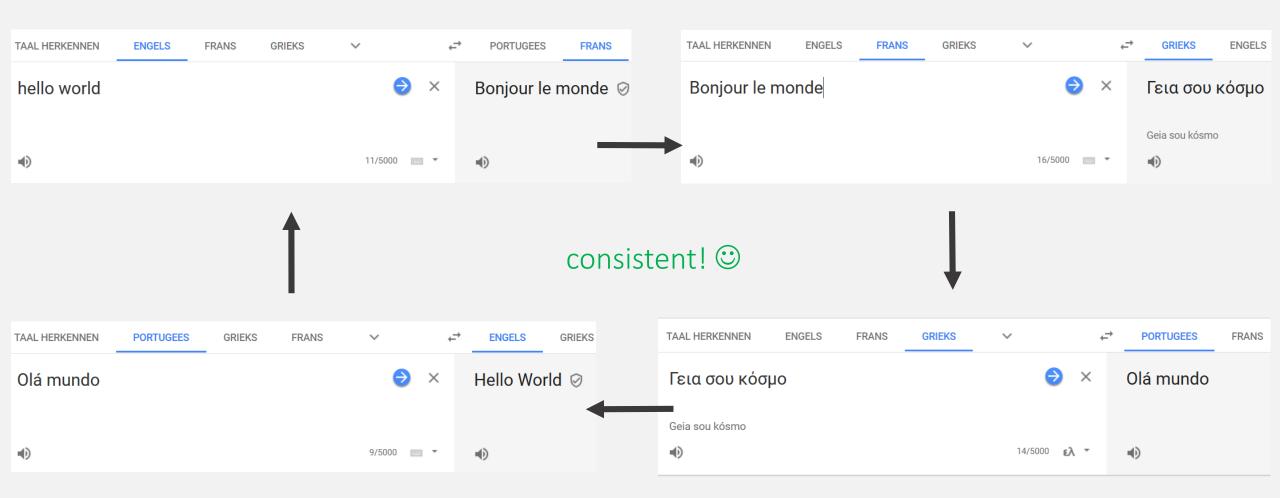




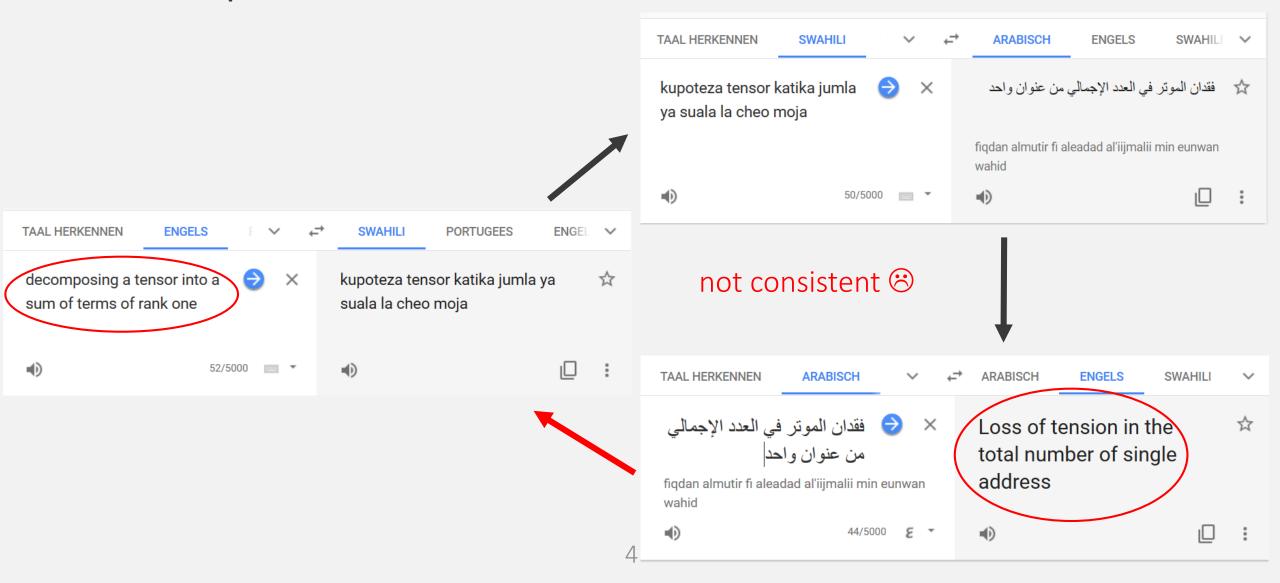
Example 1: machine translation



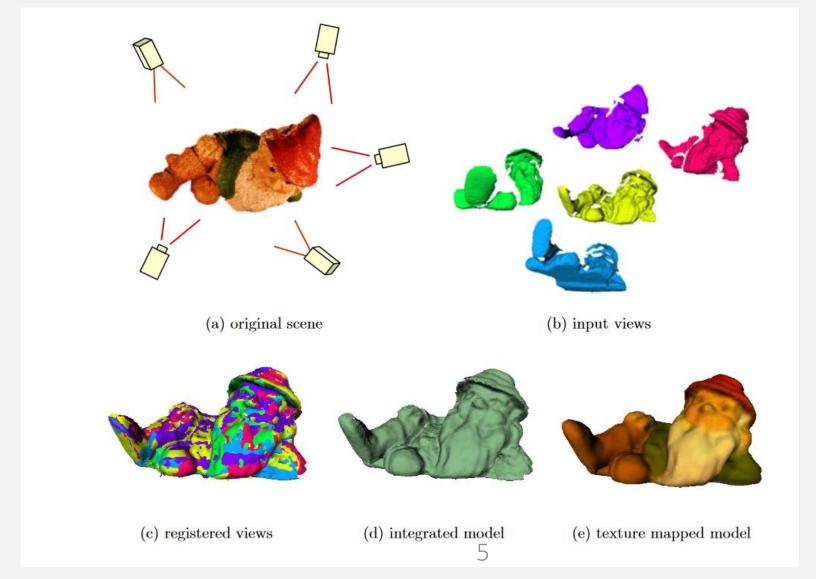
Example 1: machine translation



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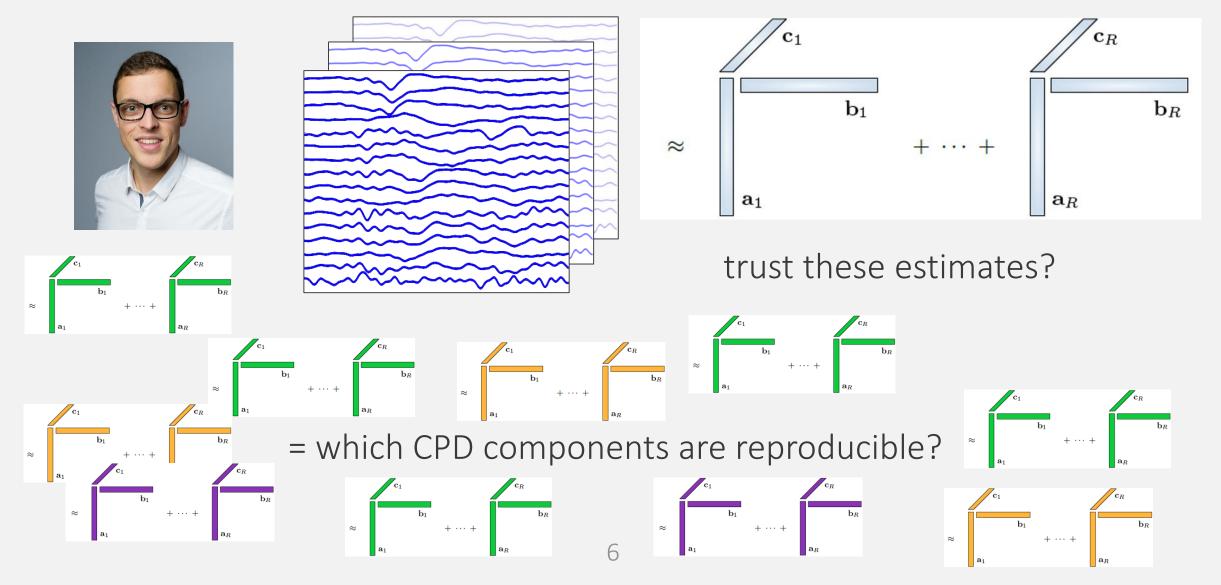


Example 2: Reconstructing 3D models



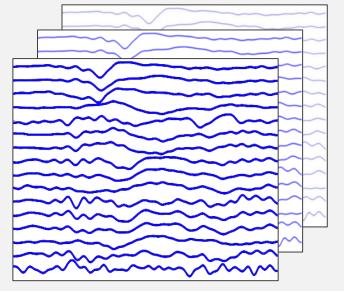
D. Huber, "Automatic three-dimensional modeling from reality",

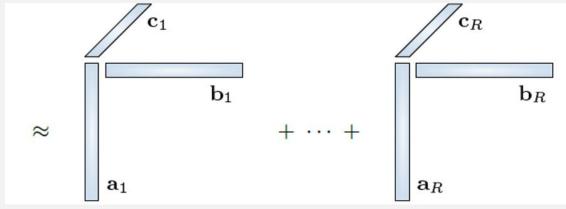
Example 3: non-convex blind source separation



Problem statement: stability analysis in BSS









Identify clusters of local solutions found by optimization.



We can use the same notion of consistency as before!

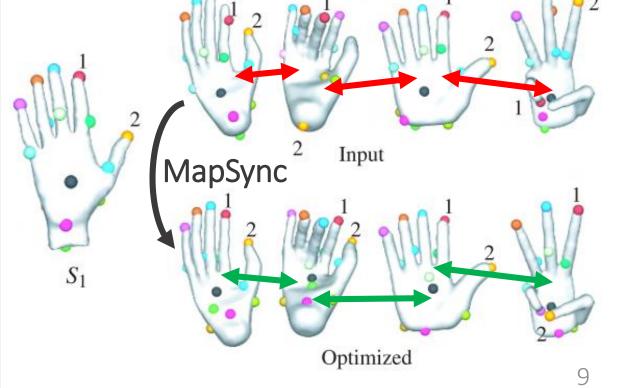
Cycle consistency

= an intuitive property of (some) graphs

Leveraging cycle consistency

from noisy, pairwise, local matching between objects... reconstruct globally consistent correspondence maps

This is called map synchronization.

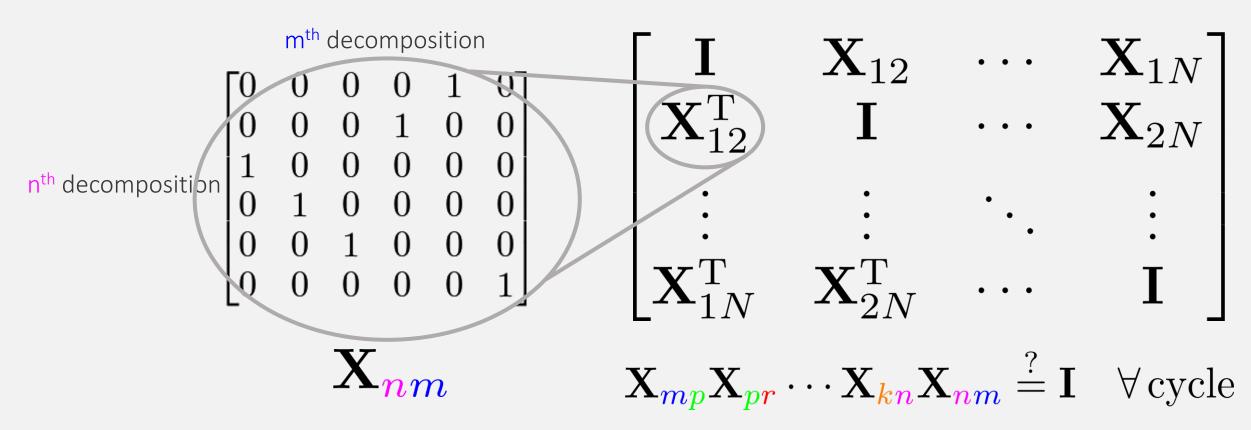


Huang, Qi-Xing, and Leonidas Guibas. "Consistent shape maps via semidefinite programming."

Cycle consistency is a strong constraint

local (pairwise) map:

all pairwise maps:



⁼ inter-factorization adjacency / mapping / matching matrix

Different view of cycle consistency

- To step from m^{th} decomposition to n^{th} decomposition: premultiply with map \mathbf{X}_{nm}
- iff cycle consistent: all decompositions are (partial) instances of a latent "universe" that contains the R true components

Alternative map from m to n: 1. map from m to the universe:
$$\mathbf{X}_{um} = \mathbf{X}_{mu}^T$$
 $\mathbf{X}_{nm} = \mathbf{X}_{nu}^T \mathbf{X}_{mu}^T$ 2. map from the universe to n: \mathbf{X}_{nu}

Cycle consistent graphs are low-rank

translates to a low-rank, positive semi-definite structure!

$$\mathbf{X}_{nm} = \mathbf{X}_{nu} \mathbf{X}_{mu}^T \quad \forall n, m$$

$$\begin{bmatrix} \mathbf{I} & \mathbf{X}_{12} & \cdots & \mathbf{X}_{1N} \\ \mathbf{X}_{12}^{\mathrm{T}} & \mathbf{I} & \cdots & \mathbf{X}_{2N} \\ \vdots & \vdots & \ddots & \vdots \\ \mathbf{X}_{1N}^{\mathrm{T}} & \mathbf{X}_{2N}^{\mathrm{T}} & \cdots & \mathbf{I} \end{bmatrix} = \begin{bmatrix} \mathbf{X}_{1u} \\ \mathbf{X}_{2u} \\ \vdots \\ \mathbf{X}_{Nu} \end{bmatrix} \begin{bmatrix} \mathbf{X}_{1u}^{\mathrm{T}} & \mathbf{X}_{2u}^{\mathrm{T}} & \cdots & \mathbf{X}_{Nu}^{\mathrm{T}} \end{bmatrix}$$

 $rank(\mathbf{X})$ = size of universe = # distinct BSS components

Identifying stable BSS components via MapSync

1 Build (thresholded) inter-factorization similarity matrix

$$\mathbf{X} \in \mathbb{R}^{NR \times NR}$$

2 Determine Q, the size of the universe

= restore a denoised graph

$$Q = \arg\max_{r} \frac{\lambda_r - \lambda_{r+1}}{|\lambda_r| + |\lambda_{r+1}|}$$

Compute a symmetric rank-Q factorization of X = enforce cycle consistency

$$\mathbf{X} pprox \mathbf{X}_u \mathbf{X}_u^T$$

4 Project the obtained maps onto space of permutations

$$\tilde{\mathbf{X}} \in \{0, 1\}^{NR \times NR}$$

Identify the Q clusters of recurrent components in the approximated graph

Similarity metric

similarity between \mathbf{r}^{th} component in run \mathbf{i} ---- \mathbf{s}^{th} component in run \mathbf{j} :

$$\sigma(r_i,s_j) = \prod_{m \in \mathcal{M}} \sigma_m(r_i,s_j) = \left|\prod_{m \in \mathcal{M}} \langle \mathbf{a}_{r_i}^{(m)}, \mathbf{a}_{s_j}^{(m)}
angle \right|^{rac{1}{|\mathcal{M}_{\sigma}|}}$$

apply a threshold: e.g. 0.95

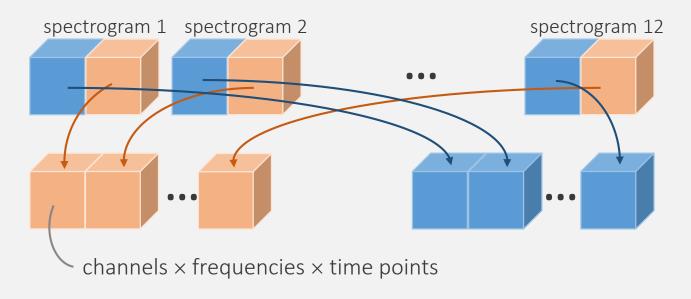
Note:

Metric can be tailored to the decomposition, e.g. KL divergence for non-negative factorizations.

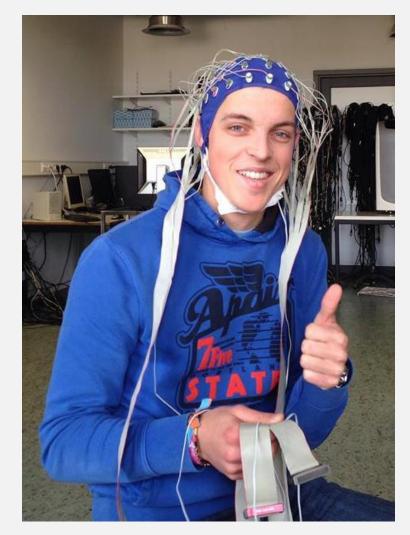
MapSync4BSS: experiment 1 Canonical Polyadic Decomposition (CPD)

CPD of resting-state EEG data

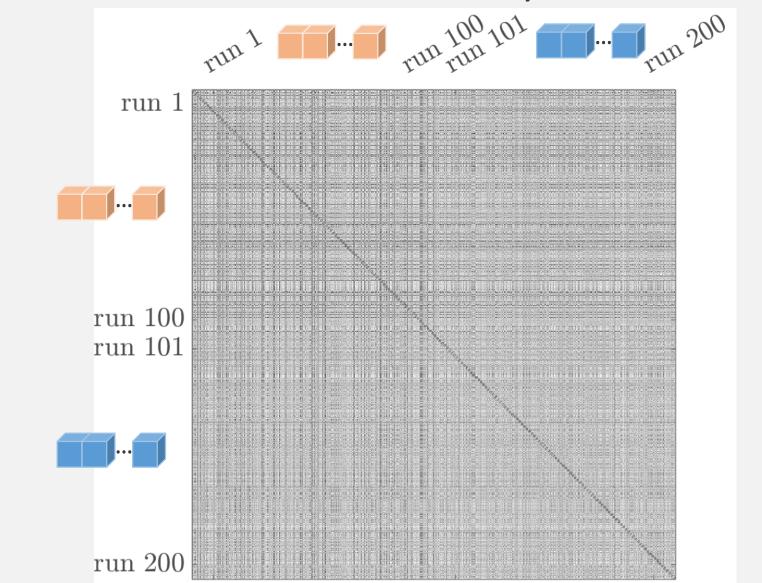
We concatenated EEG spectrograms from 12 patients, and divided the data in 2 halves



We Which spatial spectral patterns are commons to both halves, and which ones are specific?



CPD of EEG data: similarity



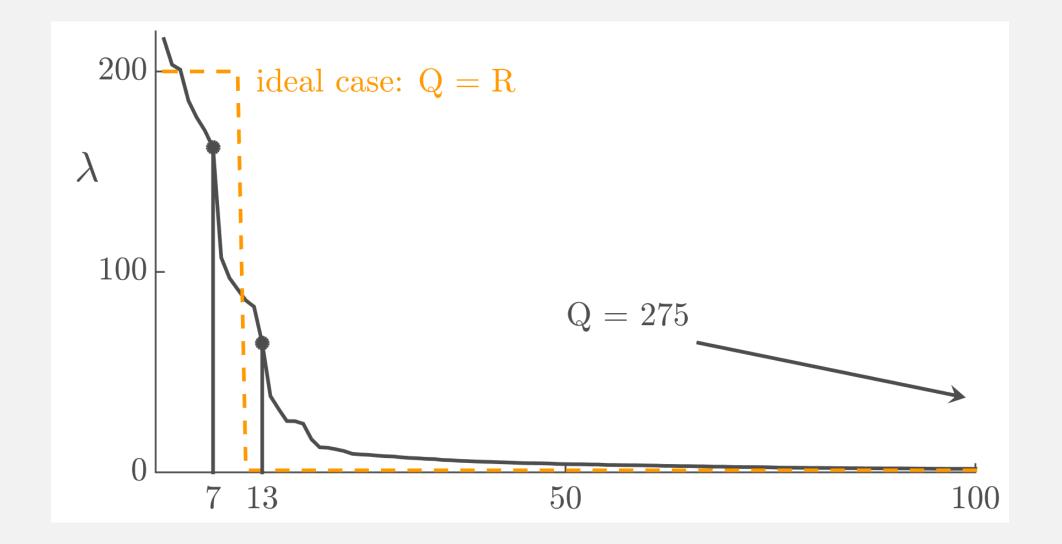
CPD of EEG data: similarity

run 1 run 100 run 101 run 200

1

CPD of EEG data: variability

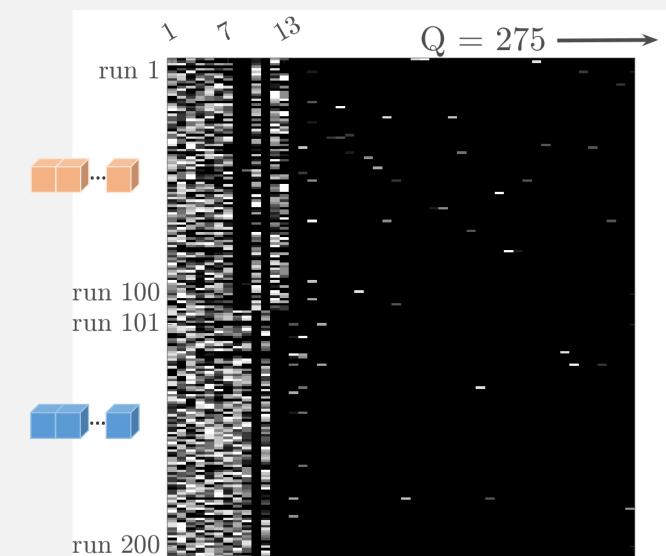
2



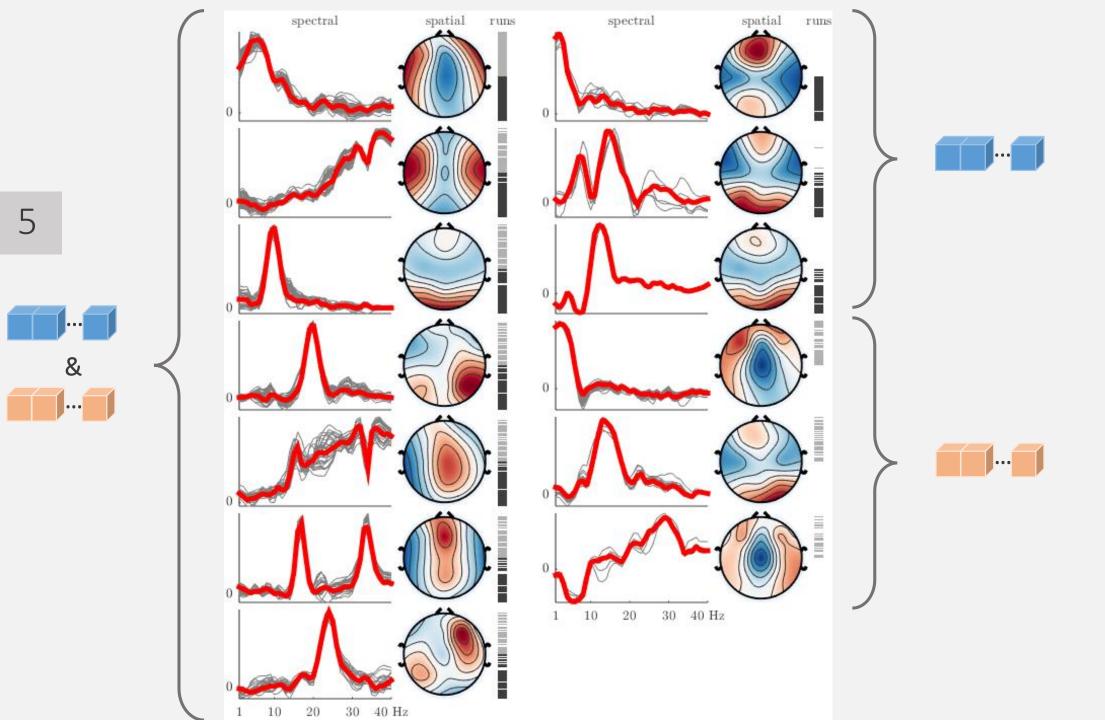
CPD of EEG data: enforcing consistency

```
3 >> [ V , D ] = eig(X)
>> % sort eigenvalues and vectors
>> Xu = V(:,1:Q) * sqrt(D(1:Q,1:Q))
```

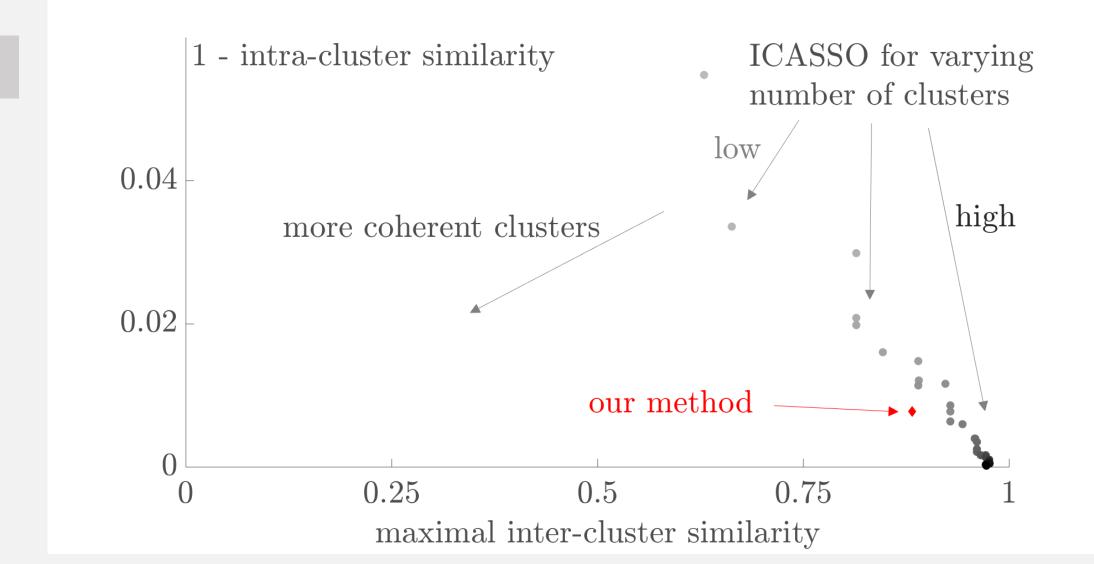
CPD of EEG data: denoise / cluster



assignment matrix of components to clusters



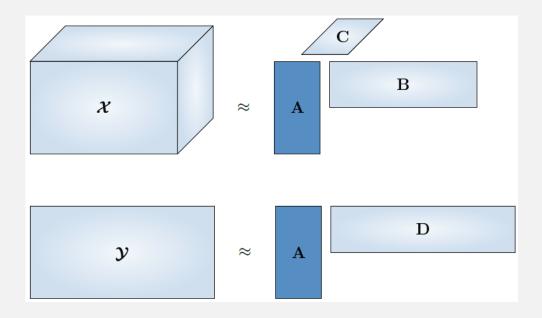
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MapSync4BSS: experiment 2 Coupled matrix-tensor factorization (CMTF)

CMTF of synthetic data

We coupled a non-negative tensor with a matrix, both of rank 7.

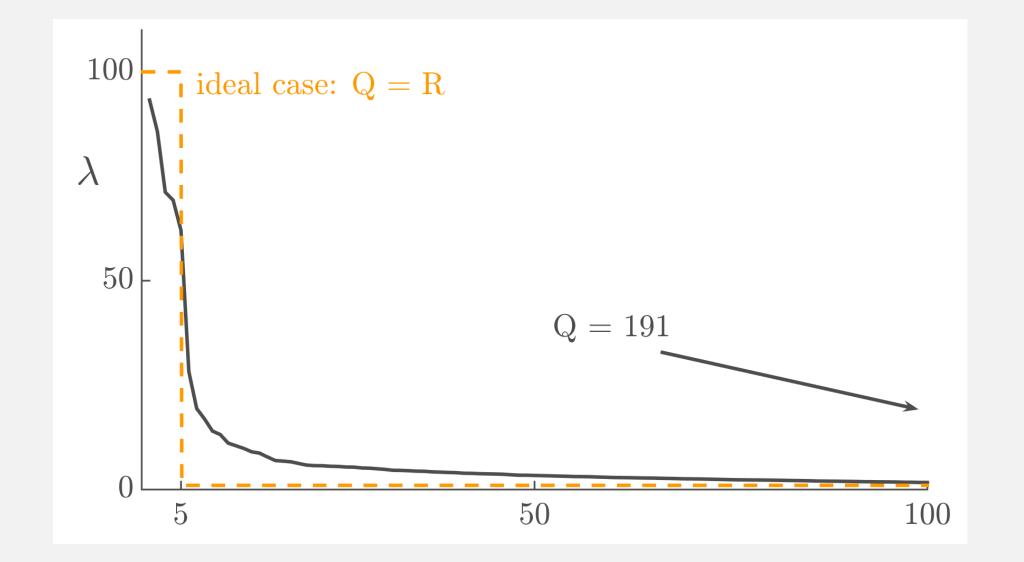


We compute 100× a CMTF with 5 components (underfitting the data),

Q: How stable is the coupled matrix-tensor factorization?

CMTF of synthetic data: variability

2

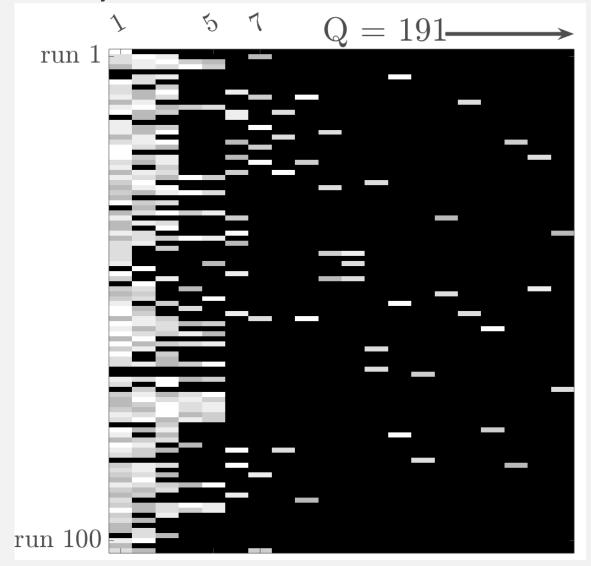


CMTF of synthetic data: denoise / cluster

4

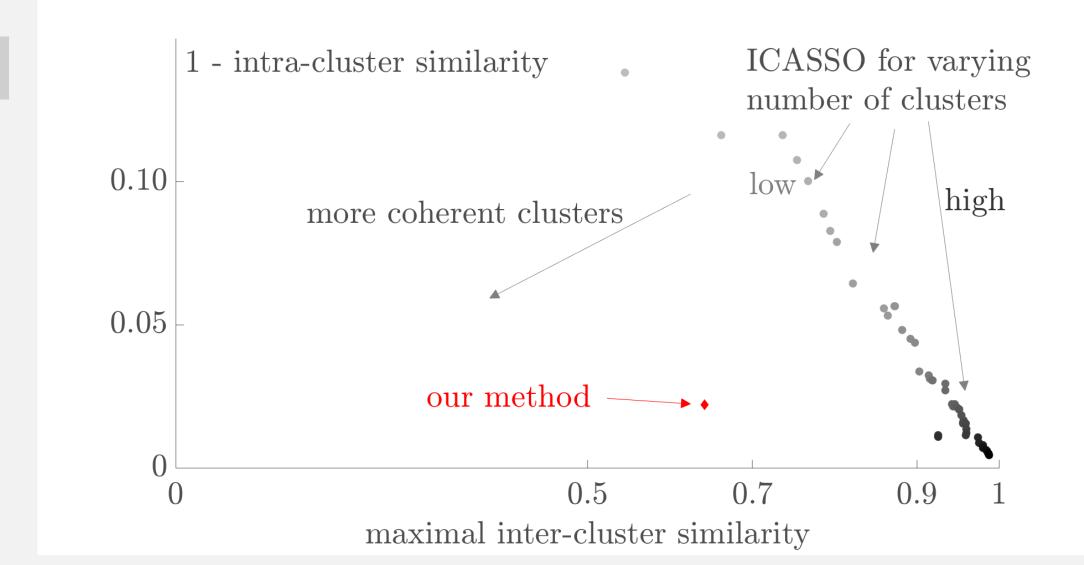
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CMTF of synthetic data: cluster

5



MapSync4BSS is a diagnostic tool for practitioners

accurate

principled

intuitive

generic

easier than ICASSO

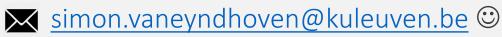
PS: Interested in the MATLAB implementation?

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Thank you!



PS: Interested in the MATLAB implementation?

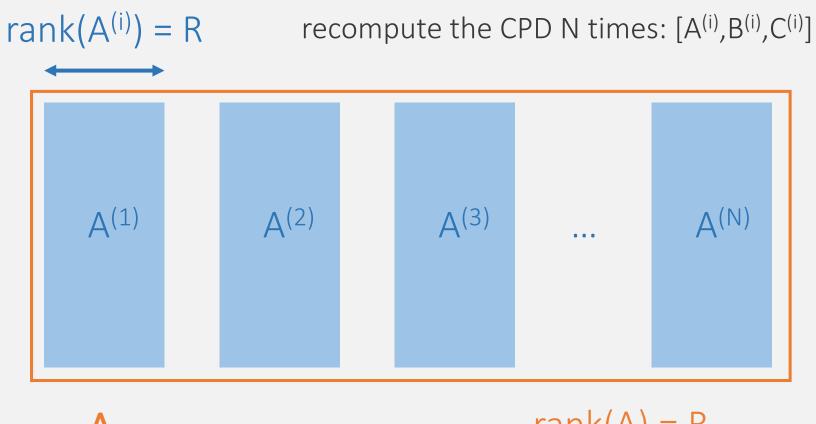








Different view of cycle consistency



rank(A) = Riff factor $A^{(i)}$ is reproducible (analogous for mode 2 and 3)

Existing work

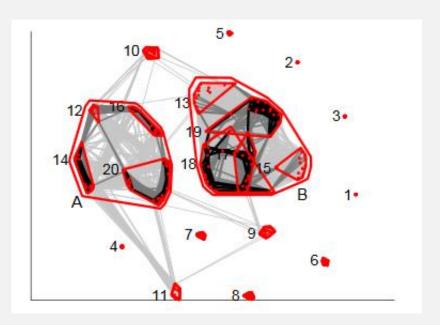
ICASSO: SOFTWARE FOR INVESTIGATING THE RELIABILITY OF ICA ESTIMATES BY CLUSTERING AND VISUALIZATION

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similar algorithm found in ICA lore

extension to tensors not explicit

different parametrization



N samples of k-dimensional vectors. The estimates of demixing matrices $\hat{\mathbf{W}}_i$ from each run $i=1,2,\ldots,M$ are collected into a single matrix $\hat{\mathbf{W}}=[\hat{\mathbf{W}}_1^T \ \hat{\mathbf{W}}_2^T \cdots \hat{\mathbf{W}}_M^T]^T$. If n_i independent components are estimated on each round, we get $K = \sum_i n_i$ estimates, and the size of $\hat{\mathbf{W}}$ will be $K \times k$.

We can resample independent component estimates by a) Randomizing the initial condition: FastICA is run M times for the same data X, so that for each run the algorithm starts from a new random initial condition; b) Bootstrapping: FastICA is run M times. The initial condition is kept the same in every run, but the data is bootstrapped every time; and c) Bootstrapping with randomized initial condition as a combination of a) and b).

A natural measure of similarity between the estimated independent components is the absolute value of their mutual correlation coefficients r_{ij} , i, j = 1, ..., K. Straightforward calculations show that they can be obtained as elements of $\mathbf{R} = \hat{\mathbf{W}} \boldsymbol{\Sigma} \hat{\mathbf{W}}^T$ where $\boldsymbol{\Sigma}$ is the covariance matrix for \mathbf{X} . The final similarity matrix has then elements

$$\sigma_{ij} = |r_{ij}|. \tag{1}$$

Selected references

- 1. Bajaj, Chandrajit, et al. "*SMAC: simultaneous mapping and clustering using spectral decompositions*." *International Conference on Machine Learning*. 2018.
- 2. Huang, Qi-Xing, and Leonidas Guibas. "*Consistent shape maps via semidefinite programming*." *Proceedings of the Eleventh Eurographics/ACMSIGGRAPH Symposium on Geometry Processing*. Eurographics Association, 2013.
- 3. Shen, Yanyao, et al. "*Normalized spectral map synchronization*." Advances in Neural Information Processing Systems. 2016.
- 4. Himberg, Johan, Aapo Hyvärinen, and Fabrizio Esposito. "Validating the independent components of neuroimaging time series via clustering and visualization." Neuroimage 22.3 (2004): 1214-1222.