

Learning a Low-coherence Dictionary to Address Spectral Variability for Hyperspectral Unmixing

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Knowledge for Tomorrow



Roadmap

- **Background & Motivation**
- **Proposed Spectral Mixing Model**
- **Experiments**
- **Conclusion & Future Work**



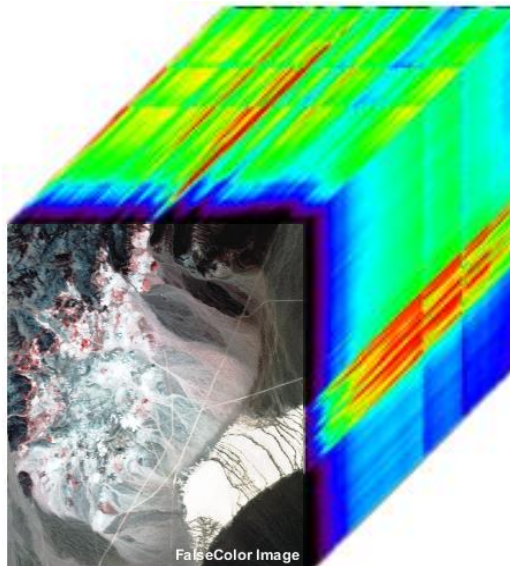
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Hyperspectral Data and Spectral Unmixing

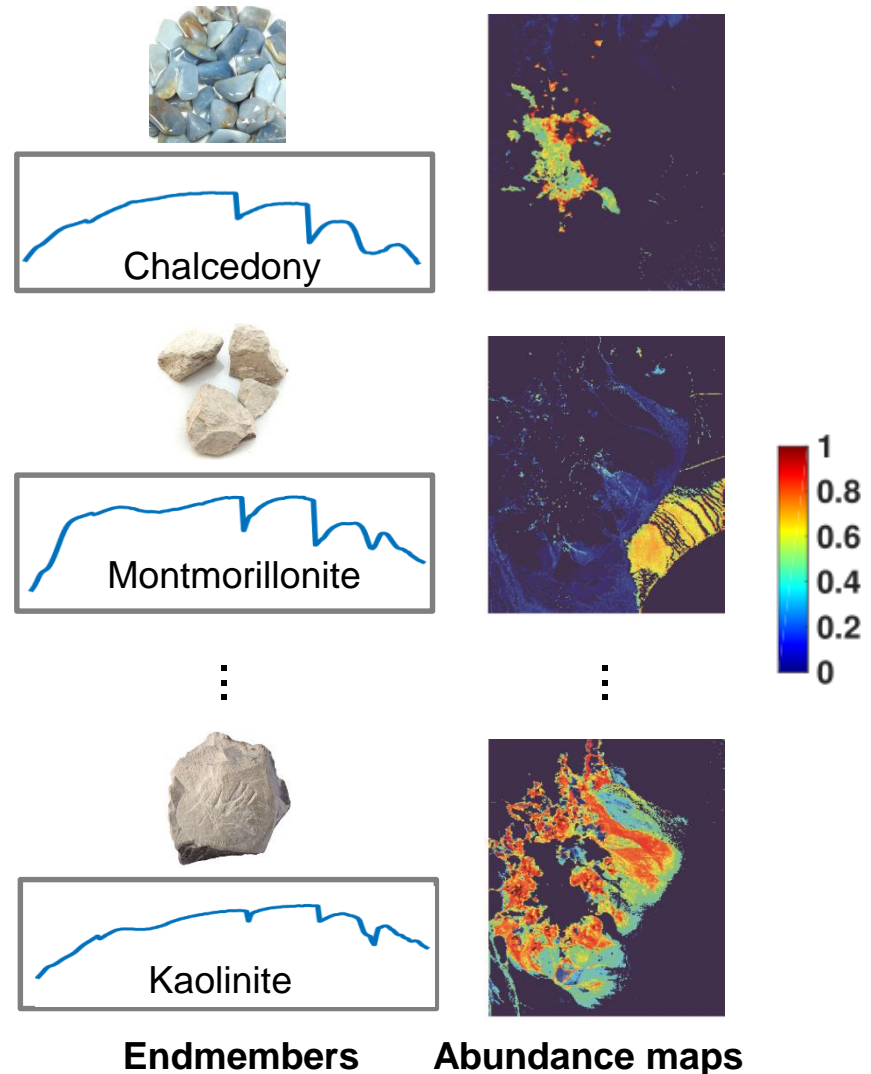
Hyperspectral data (Cuprite)



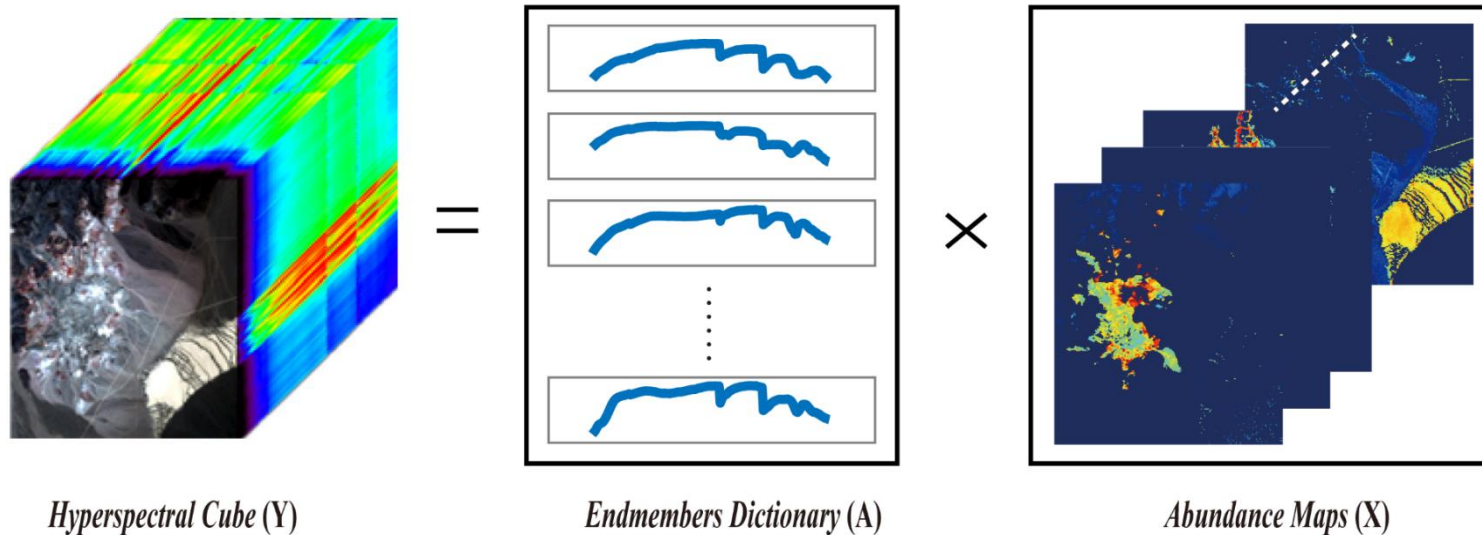
Unmixing

separate the hyperspectral data into

- Endmember signatures
- Abundance maps



Linear Mixing Model for Spectral Unmixing



Linear Mixing Model (LMM) :

$$\mathbf{Y} = \mathbf{A}\mathbf{X} + \boldsymbol{\varepsilon} \longrightarrow \text{model errors}$$

$$\mathbf{Y} \in \mathbb{R}^{D \times N} \quad \mathbf{A} \in \mathbb{R}^{D \times P} \quad \mathbf{X} \in \mathbb{R}^{P \times N} \quad \boldsymbol{\varepsilon} \in \mathbb{R}^{P \times N}$$

extracted from \mathbf{Y} by PPI [1], VCA [2]

Fully Constrained Least Square Unmixing (FCLSU) : $\arg \min_{\mathbf{X}} \|\mathbf{Y} - \mathbf{A}\mathbf{X}\|_F^2 \quad s.t. \quad \mathbf{X} \geq 0, \mathbf{1}_{P \times 1}^T \mathbf{X} = \mathbf{1}_{1 \times N}$

[1] C. Chang et al, IEEE GRL 2006

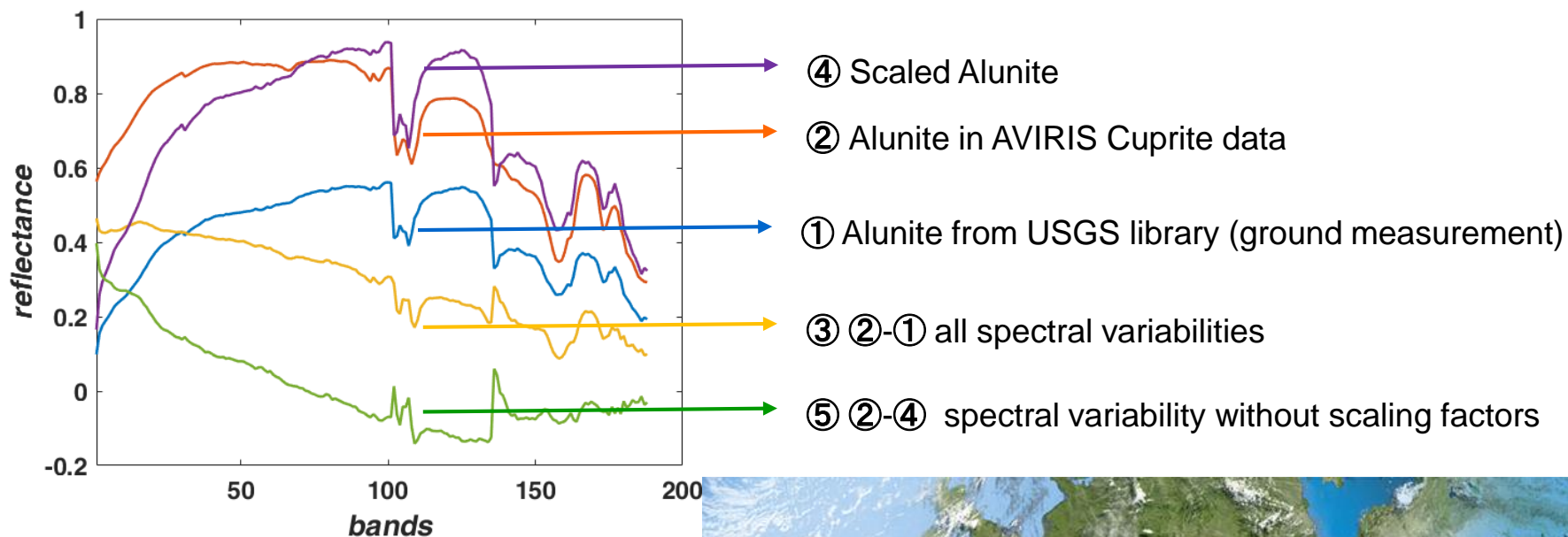
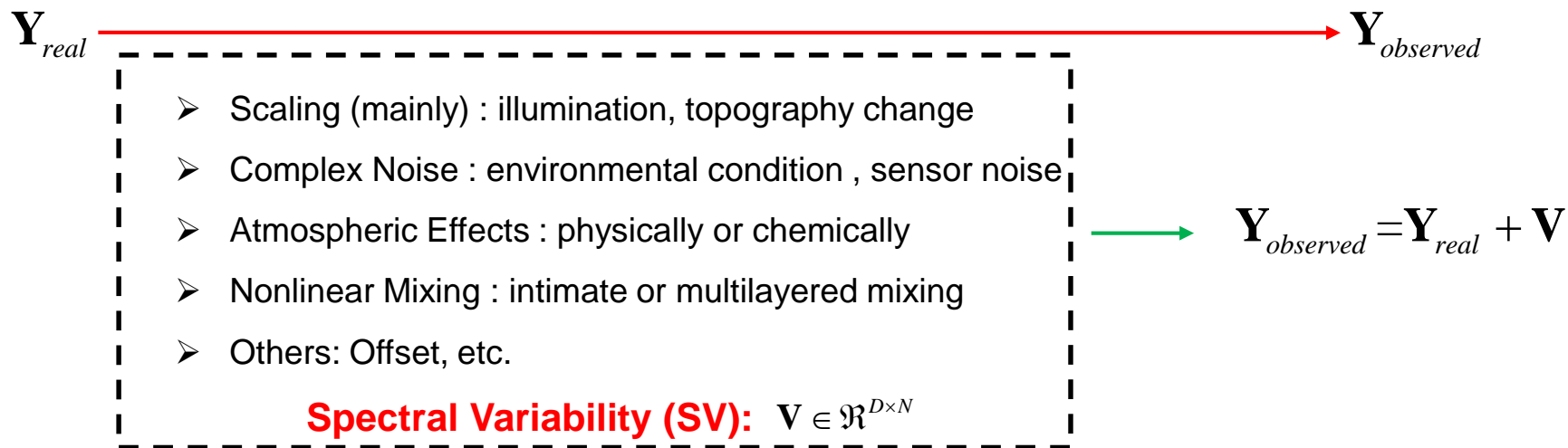
[2] J. M. P. Nascimento et al, IEEE TGRS 2005.

Non-negative constraint

Sum-to-one constraint



Spectral Variability



Variations of LMM

- LMM with spectral variability: $\mathbf{Y}_{observed} \neq \mathbf{A}\mathbf{X} + \boldsymbol{\varepsilon}$

$$= \mathbf{A}\mathbf{X} + \mathbf{V} + \boldsymbol{\varepsilon}$$

leading to **inaccurate estimation of X**, since **SV** is absorbed by **X**.

- Scaled LMM ^[1]: $\mathbf{y}_k = s_k \mathbf{A}\mathbf{x}_k + \boldsymbol{\varepsilon}_k$
- └──────────────────────────┘ a scalar shared for all endmembers

Advantage: speed up to estimate scaling factors
 obtain a relatively robust solution

Disadvantage: endmembers share the same scaling factor
 ignore other variabilities

- Extended LMM ^[2]: $\mathbf{y}_k = \mathbf{A}\mathbf{S}_k \mathbf{x}_k + \boldsymbol{\varepsilon}_k$
- └──────────────────────────┘ diagonal matrix for different endmembers

Advantage: model the scaling factor for each endmember

Disadvantage: inaccurate solution: non-convexity
 ignore other variabilities

[1] M. A. Veganzones et al, WHISPERS2014. [2] L. Drumetz, et.al, IEEE TIP, 2016.

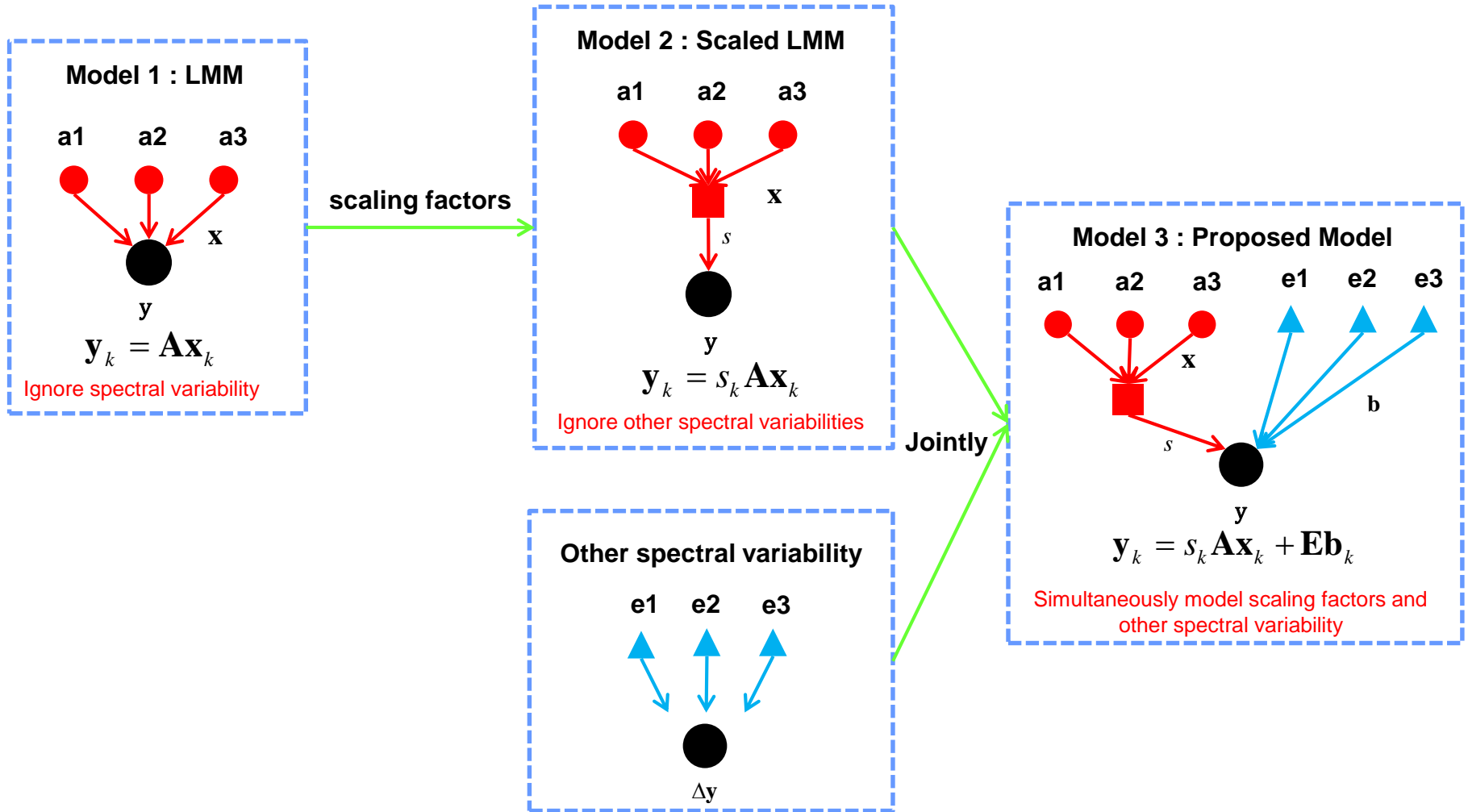


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Proposed Spectral Mixture Model



Model Formulation and Optimization

The object function of spectral variability dictionary learning can be formulated as

$$\arg \min_{\mathbf{X}, \mathbf{B}, \mathbf{S}, \mathbf{E}} \frac{1}{2} \|\mathbf{Y} - \mathbf{A}\mathbf{X}\mathbf{S} - \mathbf{E}\mathbf{B}\|_F^2 + \alpha\Phi(\mathbf{X}) + \beta\Psi(\mathbf{B}) + \gamma\Upsilon(\mathbf{E})$$

$$s.t. \quad \mathbf{X} \geq \mathbf{0}, \quad \mathbf{1}_{P \times 1}^T \mathbf{X} = \mathbf{1}_{1 \times N}, \quad \mathbf{S} \geq \mathbf{0}$$

➤ Sub-problem (1) : Pixel-wise spectral unmixing

$$\arg \min_{\mathbf{X}, \mathbf{B}, \mathbf{S}} \sum_{k=1}^N \left(\frac{1}{2} \|\mathbf{y}_k - s_k \mathbf{A}\mathbf{x}_k - \mathbf{E}\mathbf{b}_k\|_2^2 \right) + \alpha\Phi(\mathbf{X}) + \beta\Psi(\mathbf{B})$$

$$s.t. \quad \mathbf{X} \geq \mathbf{0}, \quad \mathbf{1}_{P \times 1}^T \mathbf{X} = \mathbf{1}_{1 \times N}, \quad \mathbf{S} \geq \mathbf{0}$$

➤ Sub-problem (2) : Spectral variability dictionary updating

$$\arg \min_{\mathbf{E}} \frac{1}{2} \|\mathbf{Y} - \mathbf{A}\mathbf{X}\mathbf{S} - \mathbf{E}\mathbf{B}\|_F^2 + \gamma\Upsilon(\mathbf{E})$$

➤ Regularization terms :

$$\Phi(\mathbf{X}) = \frac{1}{2} \|\mathbf{X}\|_{1,1} \quad \Upsilon(\mathbf{E}) = \frac{1}{2} \left(\|\mathbf{A}^T \mathbf{E}\|_F^2 + \|\mathbf{E}^T \mathbf{E} - \mathbf{I}\|_F^2 \right)$$

$$\Psi(\mathbf{B}) = \frac{1}{2} \|\mathbf{B}\|_F^2$$


Pseudocode

Learning Spectral Variability Dictionary[↵]

Input: Endmember dictionary (**A**), Mixture spectral signature (**Y**), and parameters (α, β, γ).[↵]

Initialize: Variability dictionary (**E**), Abundance map (**X**), Coefficient for Variability dictionary (**B=0**)[↵]
Scaling factor (**S=1**)[↵]

Main steps:[↵]

while *not converged* or $t > \text{maxIter}$ **do**[↵]

Dictionary Updating[↵]

 fix **X**, **B**, **S** and update **E** by solving the problem (2) using ADMM[↵]

if conditions are satisfied [↵]

then[↵]

Spectral Unmixing[↵]

 fix **E** and update **X**, **B**, **S** by solving the problem (1) using ADMM [↵]

if conditions are satisfied [↵]

then[↵]

 Stop iteration;[↵]

else[↵]

$t=t+1$ [↵]

end[↵]

end[↵]

Output : **X**, **B**, **E**, **S**[↵]

Note: **E**: Random Orthogonal Matrix ; Initial **X** : obtained by Scaled LMM[↵]



Regularization Terms

$$\Phi(\mathbf{X}) = \frac{1}{2} \|\mathbf{X}\|_{1,1}$$

A limited number of materials $\xrightarrow{\text{promote}}$ Sparsity

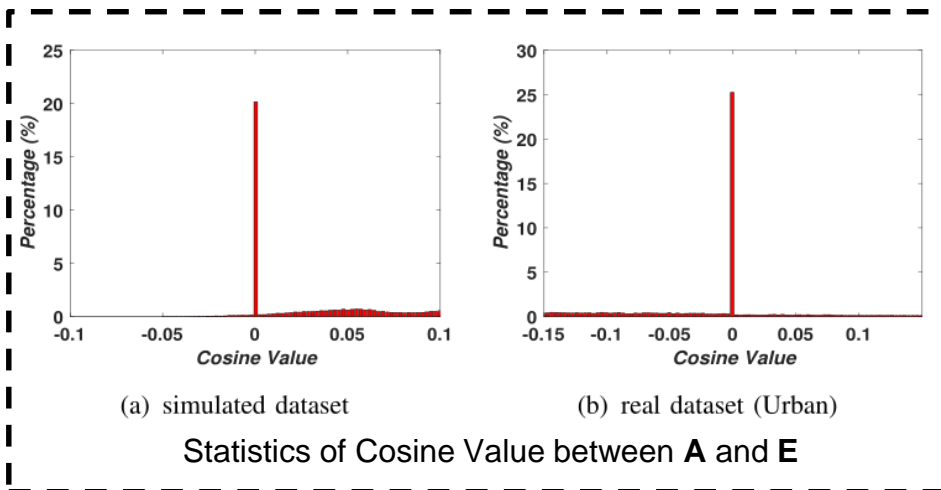
$$\Psi(\mathbf{B}) = \frac{1}{2} \|\mathbf{B}\|_F^2$$

Many atoms in \mathbf{E} should be activated $\xrightarrow{\text{promote}}$ Density

$$\Upsilon(\mathbf{E}) = \frac{1}{2} \left(\|\mathbf{A}^T \mathbf{E}\|_F^2 + \|\mathbf{E}^T \mathbf{E} - \mathbf{I}\|_F^2 \right)$$

↑ promote

- Enough to represent various **SV**
- Avoid the trivial solution

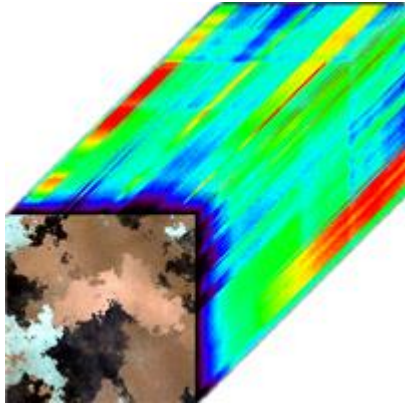


Roadmap

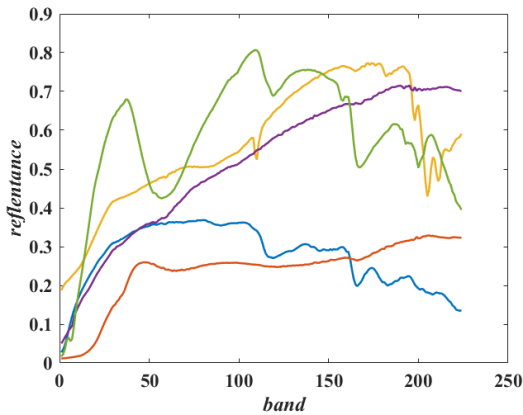
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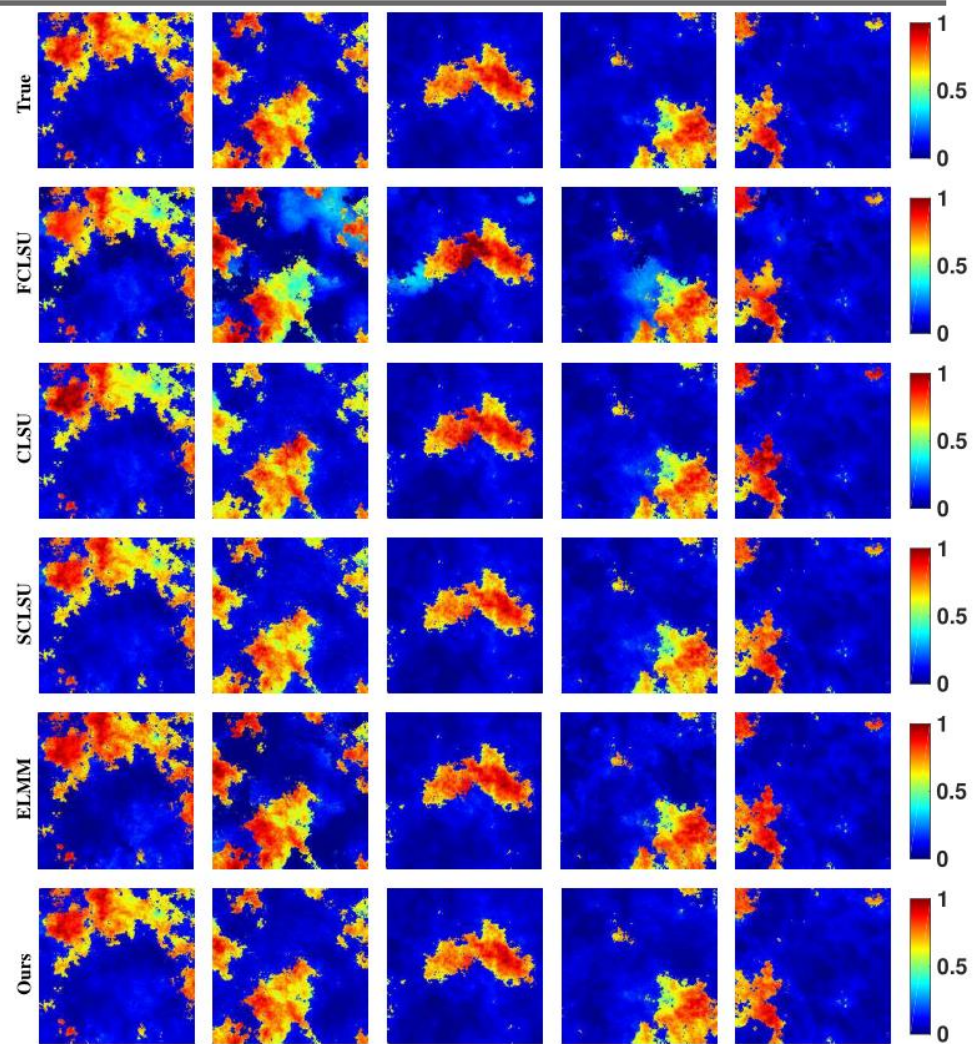
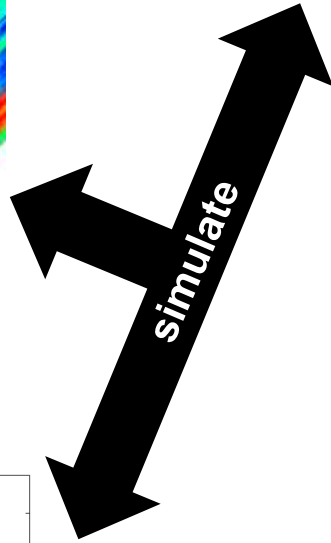
Abundance Map for Simulated Data



A false color simulated data



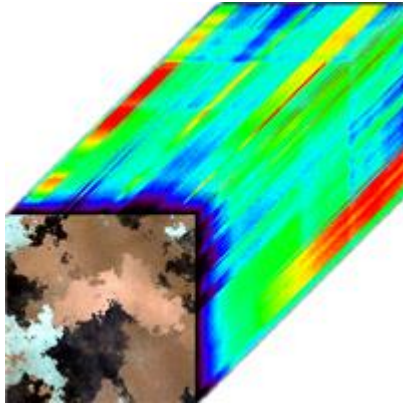
Endmembers used for generating the simulated data



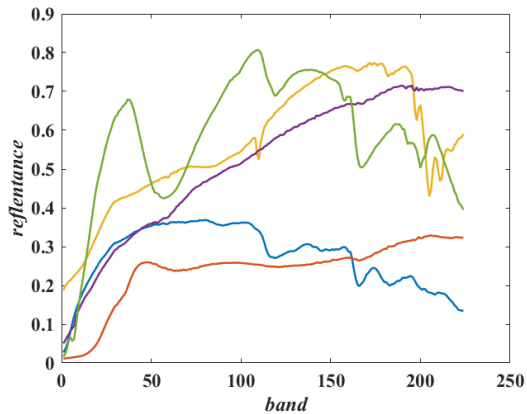
The abundance maps (each column corresponds to one endmember extracted by VCA) and the first row shows the ground truth



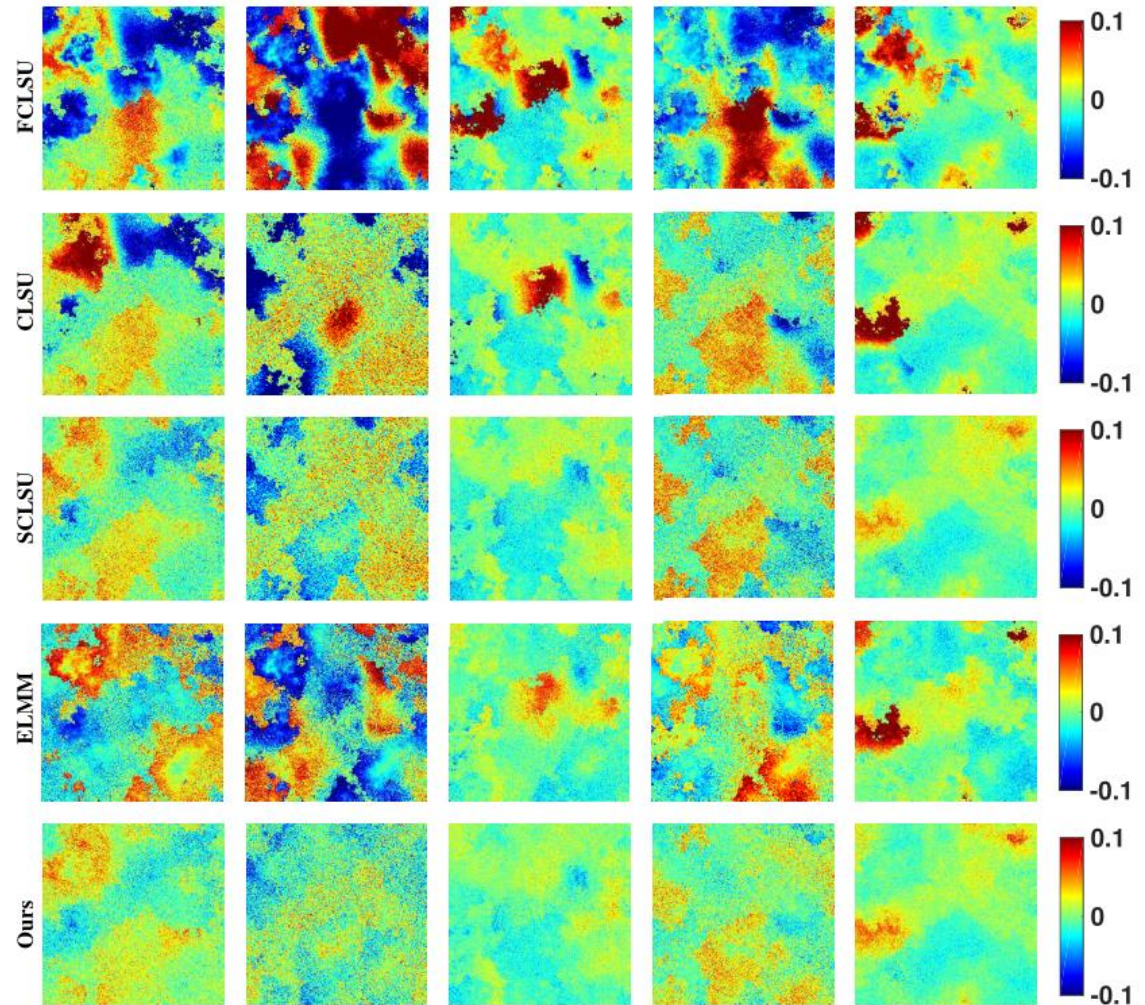
Abundance Difference Map for Simulated Data



A false color simulated data



Endmembers used for generating the simulated data



The abundance difference maps obtained between estimated abundance map and true map (each column corresponds to one endmember extracted by VCA)



Quantity Experiments (Simulated Data)

$$xRMSE = \frac{1}{N} \sum_{k=1}^N \sqrt{\frac{1}{P} \sum_{p=1}^P \left(\mathbf{x}_{pk} - \hat{\mathbf{x}}_{pk} \right)^2}$$

$$yRMSE = \frac{1}{N} \sum_{k=1}^N \sqrt{\frac{1}{L} \sum_{l=1}^L \left(\mathbf{y}_{pk} - \hat{\mathbf{y}}_{pk} \right)^2}$$

$$ySAM = \frac{1}{N} \sum_{k=1}^N \arccos \left(\frac{\mathbf{y}_k \bullet \hat{\mathbf{y}}_k}{\|\mathbf{y}_k\| \times \|\hat{\mathbf{y}}_k\|} \right)$$

N : the number of pixel

$xRMSE$: Abundance Root Mean Square Error

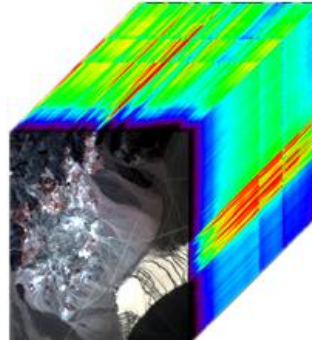
$yRMSE$: Reconstruction Root Mean Square Error

$ySAM$: Average Spectral Angle Mapper

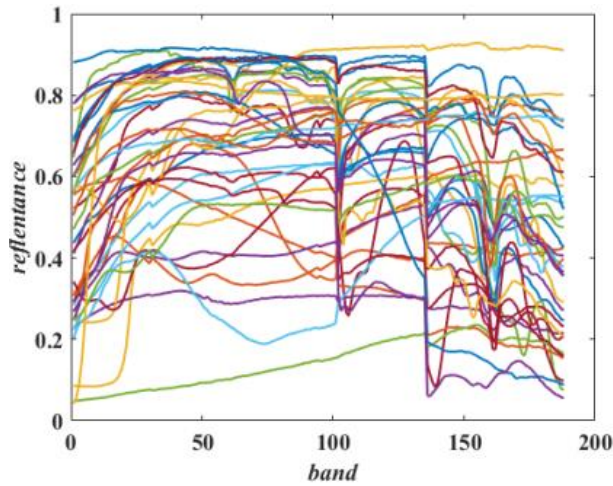
Algorithm	FCLSU	CLSU	CLSU+Sparse	SCLSU	SCLSU+Sparse	ELMM	Ours
$xRMSE$	0.0524	0.0380	0.0379	0.0251	0.0248	0.0337	0.0206
$yRMSE$	0.0151	0.0123	0.0127	0.0123	0.0127	0.0088	0.000006
$ySAM$	1.9600	1.7713	1.7715	1.7713	1.7715	1.2998	0.0007



Abundance Map for Real Data

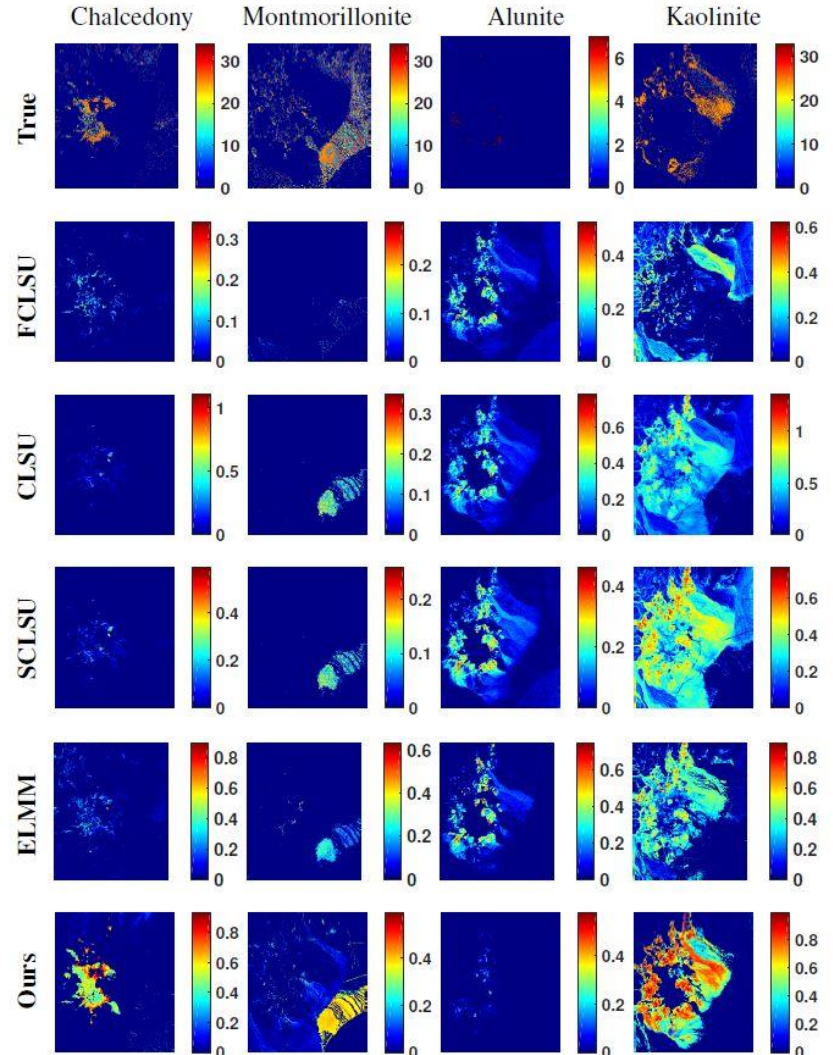


Cuprite Data



Selected Endmembers from Spectral Library *

* Due to the complexity of spectral variability in Cuprite Data, here we get help from spectral library to construct endmember dictionary.



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Conclusion and Future work

Conclusion

1. A novel spectral mixing model is proposed to simultaneously consider the scaling factors (main) and other spectral variabilities.
2. A data-driven dictionary learning method is explored with the low-coherent regularization to design the spectral variability dictionary.
3. An alternating iterative optimization strategy is applied to solve the proposed model using ADMM-based framework.

Future work

1. Spatial regularization should be able to further improve the performance of spectral unmixing.
2. Distributed strategy could be introduced to promote a large-scale spectral unmixing.



Thank you for your attention!

Knowledge for Tomorrow

