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Matrix Completion in Parallel Architectures: A Julia Implementation

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Keegan Lensing,

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University of British Columbia



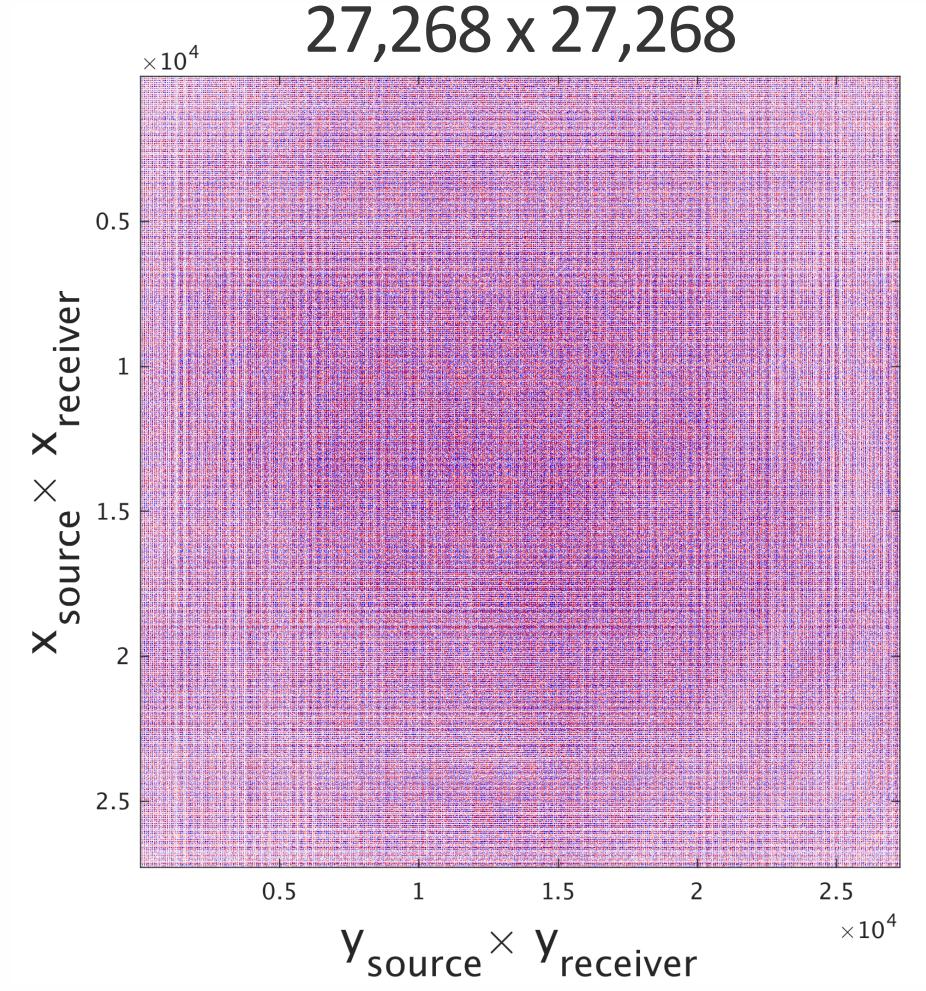
Motivation

- Industry-scale seismic data interpolation
- ▶ Exploit *low-rank* structure of seismic data
 - matrix completion techniques
- Need to improve time complexity
 - design for parallel architectures



Motivation





Seismic data: huge matrices

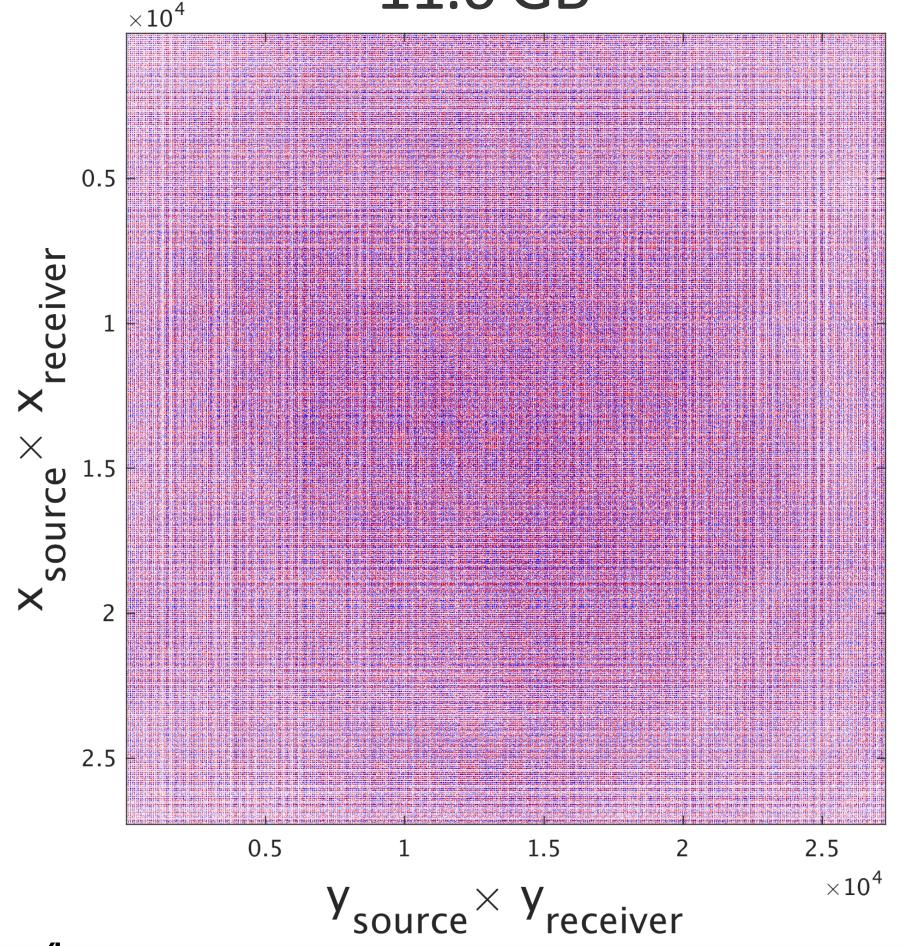
Interpolation quality deteriorates when working on smaller windows

Want to work w/ full matrix

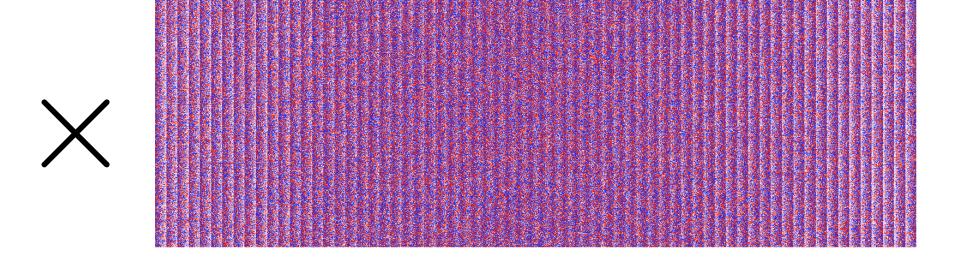


Motivation

11.6 GB



454 MB



No need to store full matrix (96% compression)

Can directly generate gathers



Contributions

Matrix completion: Decoupling method

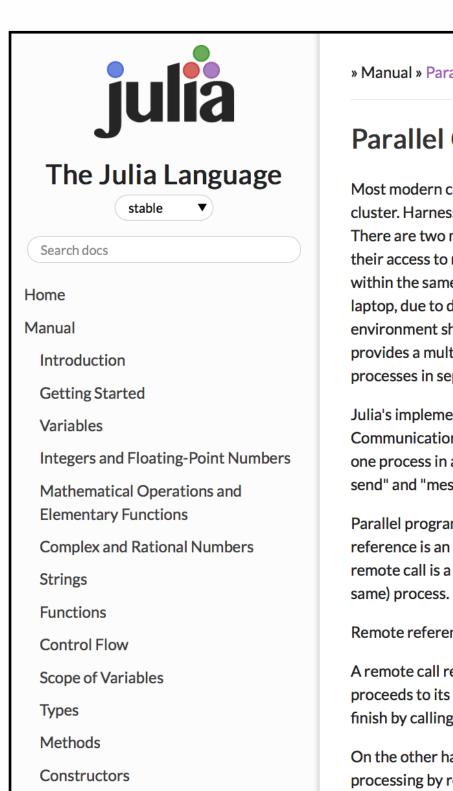
- decompose into independent subproblems
- solve in parallel architectures

Julia implementation:

- efficient multiprocessing environment
- optimize communication time between workers



Contributions



» Manual » Parallel Computing

C Edit on GitHub

Parallel Computing

Most modern computers possess more than one CPU, and several computers can be combined together in a cluster. Harnessing the power of these multiple CPUs allows many computations to be completed more quickly. There are two major factors that influence performance: the speed of the CPUs themselves, and the speed of their access to memory. In a cluster, it's fairly obvious that a given CPU will have fastest access to the RAM within the same computer (node). Perhaps more surprisingly, similar issues are relevant on a typical multicore laptop, due to differences in the speed of main memory and the cache. Consequently, a good multiprocessing environment should allow control over the "ownership" of a chunk of memory by a particular CPU. Julia provides a multiprocessing environment based on message passing to allow programs to run on multiple processes in separate memory domains at once.

Julia's implementation of message passing is different from other environments such as MPI [1]. Communication in Julia is generally "one-sided", meaning that the programmer needs to explicitly manage only one process in a two-process operation. Furthermore, these operations typically do not look like "message send" and "message receive" but rather resemble higher-level operations like calls to user functions.

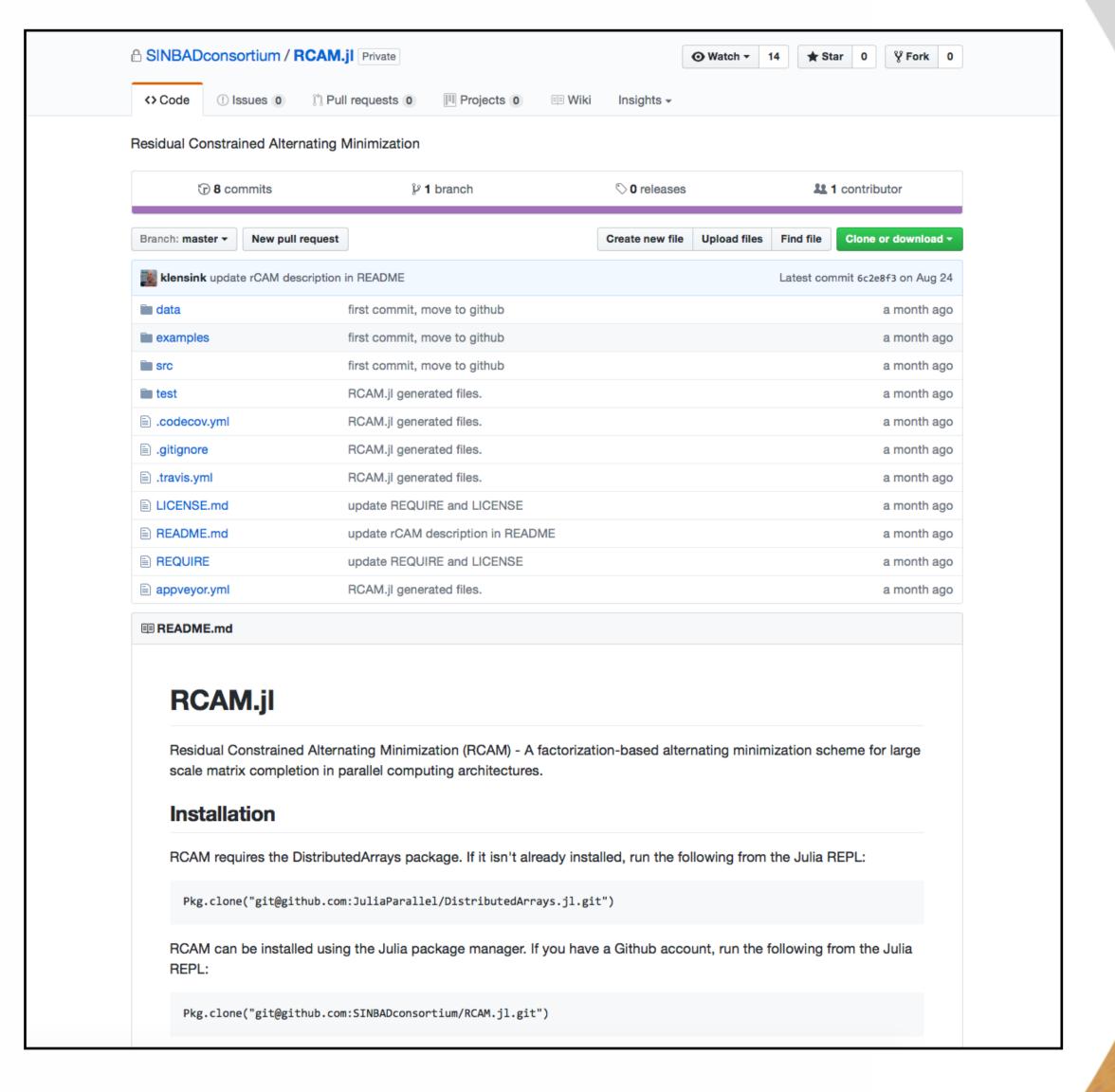
Parallel programming in Julia is built on two primitives: *remote references* and *remote calls*. A remote reference is an object that can be used from any process to refer to an object stored on a particular process. A remote call is a request by one process to call a certain function on certain arguments on another (possibly the same) process.

Remote references come in two flavors: Future and RemoteChannel.

A remote call returns a Future to its result. Remote calls return immediately; the process that made the call proceeds to its next operation while the remote call happens somewhere else. You can wait for a remote call to finish by calling wait() on the returned Future, and you can obtain the full value of the result using fetch().

On the other hand, RemoteChannel s are rewritable. For example, multiple processes can co-ordinate their processing by referencing the same remote Channel.

Software Release Available





Outline

- Matrix completion
 - alternating least squares
 - decoupling method
- Parallel implementation in Julia
- Numerical experiments



Matrix completion

Goal is to approximate $\mathbf{M} \in \mathbb{C}^{n \times m}$, given

observed entries $\Omega \subset \{1,2,...,n\} \times \{1,2,...,m\}$

via

$$\mathbf{b}_{k,\ell} = P_{\Omega}(\mathbf{M})_{k,\ell} = \begin{cases} \mathbf{M}_{k,\ell} & \text{if } (k,\ell) \in \Omega \\ 0 & \text{otherwise} \end{cases}$$



Methodology: Least Squares

If M is approximately rank-r, we solve

$$(\mathbf{L}^{\sharp}, \mathbf{R}^{\sharp}) = \underset{\mathbf{L} \in \mathbb{C}^{n \times r}, \mathbf{R} \in \mathbb{C}^{m \times r}}{\operatorname{arg\,min}} \|P_{\Omega}(\mathbf{L}\mathbf{R}^{*}) - \mathbf{b}\|_{F}$$

and approximate

$$\mathbf{L}^{\sharp}(\mathbf{R}^{\sharp})^{st}pprox \mathbf{M}$$

Prateek Jain, Praneeth Netrapalli, Sujay Sanghavi. "Low-Rank Matrix Completion Using Alternating Minimization".



Methodology: Alternating Least Squares

We alternate the optimization over each factor

$$\mathbf{L}^t = \underset{\mathbf{L} \in \mathbb{C}^{n \times r}}{\operatorname{arg\,min}} \| P_{\Omega}(\mathbf{L}(\mathbf{R}^t)^*) - \mathbf{b} \|_F$$

$$\mathbf{R}^{t+1} = \underset{\mathbf{R} \in \mathbb{C}^{m \times r}}{\operatorname{arg\,min}} \| P_{\Omega}(\mathbf{L}^{t}\mathbf{R}^{*}) - \mathbf{b} \|_{F}$$

starting at initial factor (${f R}^0$) and iteratively obtain

$$\mathbf{L}^T(\mathbf{R}^T)^* pprox \mathbf{M}$$

Prateek Jain, Praneeth Netrapalli, Sujay Sanghavi. "Low-Rank Matrix Completion Using Alternating Minimization".



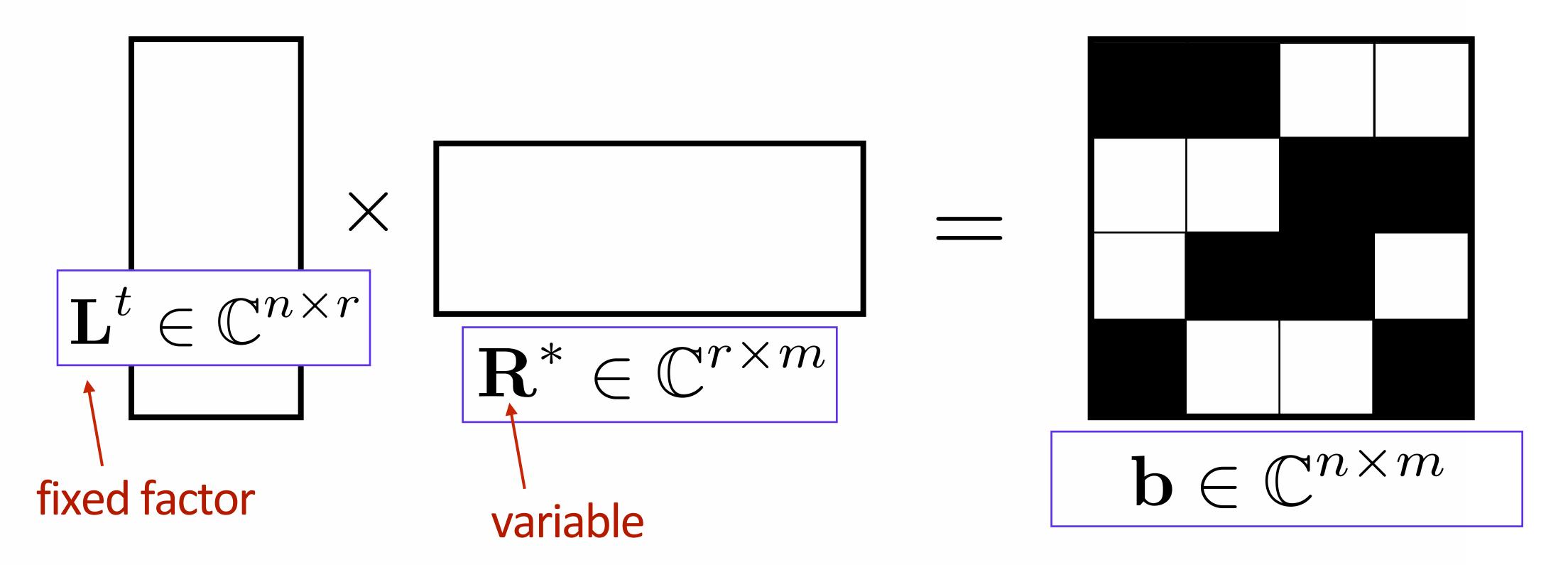
Each subproblem, e.g.,

$$\mathbf{R}^{t+1} = \underset{\mathbf{R} \in \mathbb{C}^{m \times r}}{\operatorname{arg\,min}} \| P_{\Omega}(\mathbf{L}^{t}\mathbf{R}^{*}) - \mathbf{b} \|_{F}$$

can be decoupled to solve each row of \mathbf{R}^{t+1} independently.

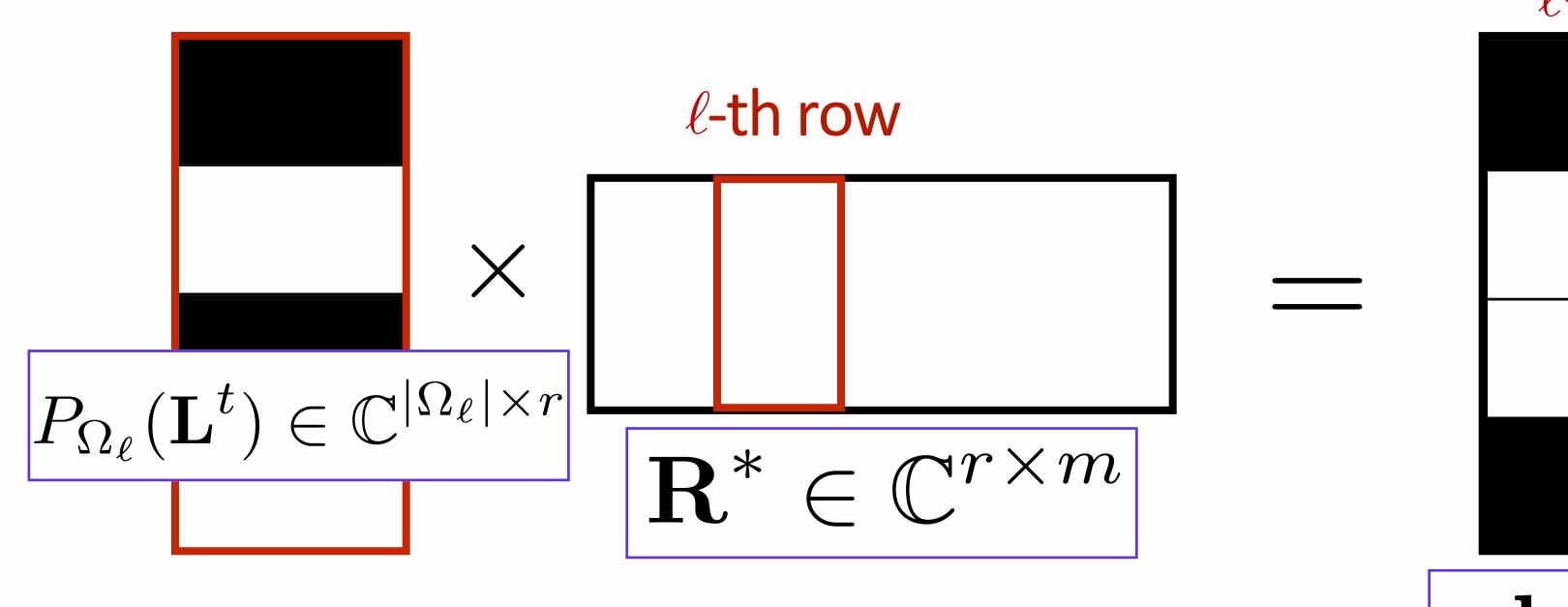


Decoupling method: Visualization

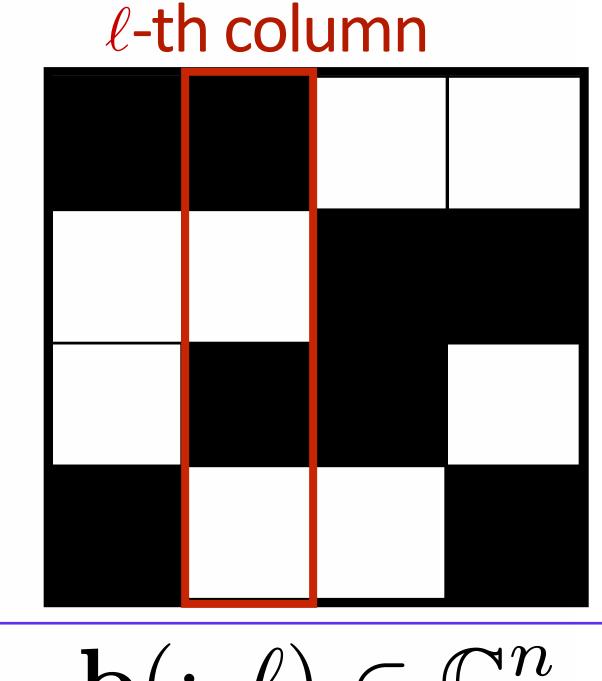




Decoupling method: Visualization



$$\mathbf{R}^{t+1}(\ell,:) = \underset{v \in \mathbb{C}^r}{\arg\min} \|P_{\Omega_{\ell}}(\mathbf{L}^t)v - \mathbf{b}(:,\ell)\|_2$$



$$\mathbf{b}(:,\ell) \in \mathbb{C}^n$$



So each row can be solved independently as

$$\mathbf{R}^{t+1}(\ell,:) = \underset{v \in \mathbb{C}^r}{\arg\min} \|P_{\Omega_{\ell}}(\mathbf{L}^t)v - \mathbf{b}(:,\ell)\|_2$$

 $P_{\Omega_\ell}(\mathbf{L}^t)$ is \mathbf{L}^t restricted to the entries observed in the ℓ -th column



$$\mathbf{R}^{t+1}(\ell,:) = \underset{v \in \mathbb{C}^r}{\arg\min} \|P_{\Omega_{\ell}}(\mathbf{L}^t)v - \mathbf{b}(:,\ell)\|_2$$

closed form solution is

$$\mathbf{R}^{t+1}(\ell,:) = \left(P_{\Omega_{\ell}}(\mathbf{L}^t)^* P_{\Omega_{\ell}}(\mathbf{L}^t)\right)^{-1} P_{\Omega_{\ell}}(\mathbf{L}^t)^* \mathbf{b}(:,\ell)$$



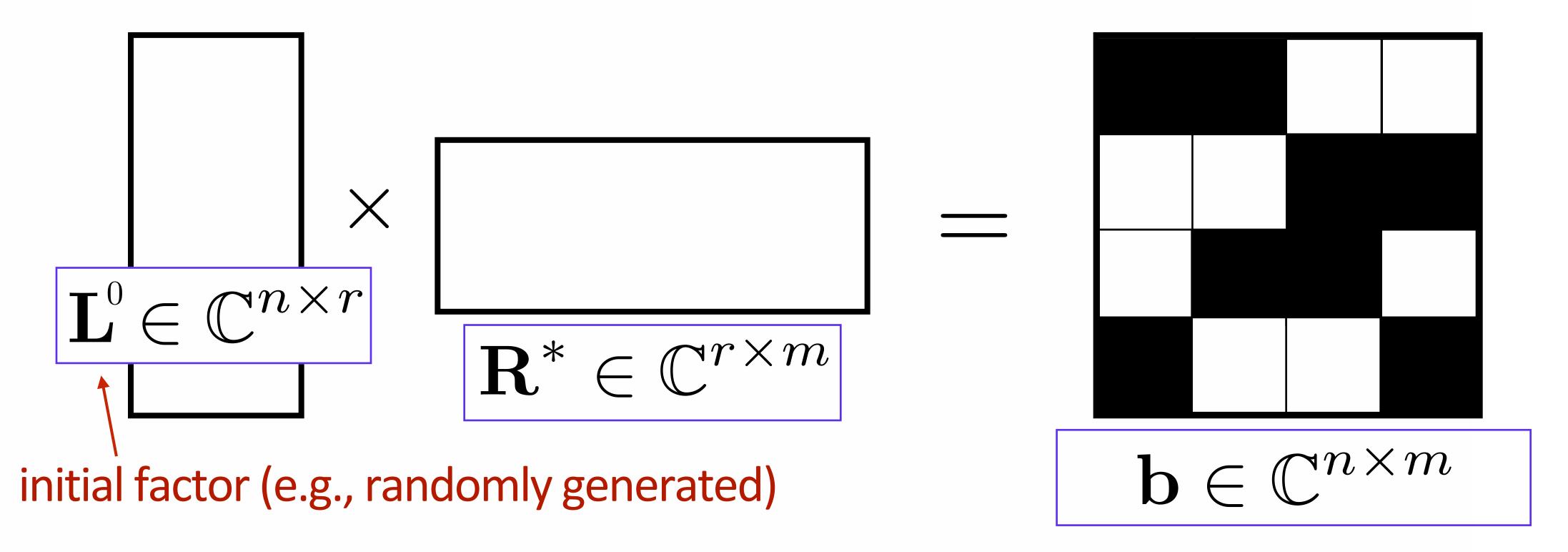
$$\mathbf{R}^{t+1}(\ell,:) = \left(P_{\Omega_{\ell}}(\mathbf{L}^t)^* P_{\Omega_{\ell}}(\mathbf{L}^t)\right)^{-1} P_{\Omega_{\ell}}(\mathbf{L}^t)^* \mathbf{b}(:,\ell)$$

 $\mathrm{cheap}\,\mathrm{if}\;r\ll\min\{n,m\}$

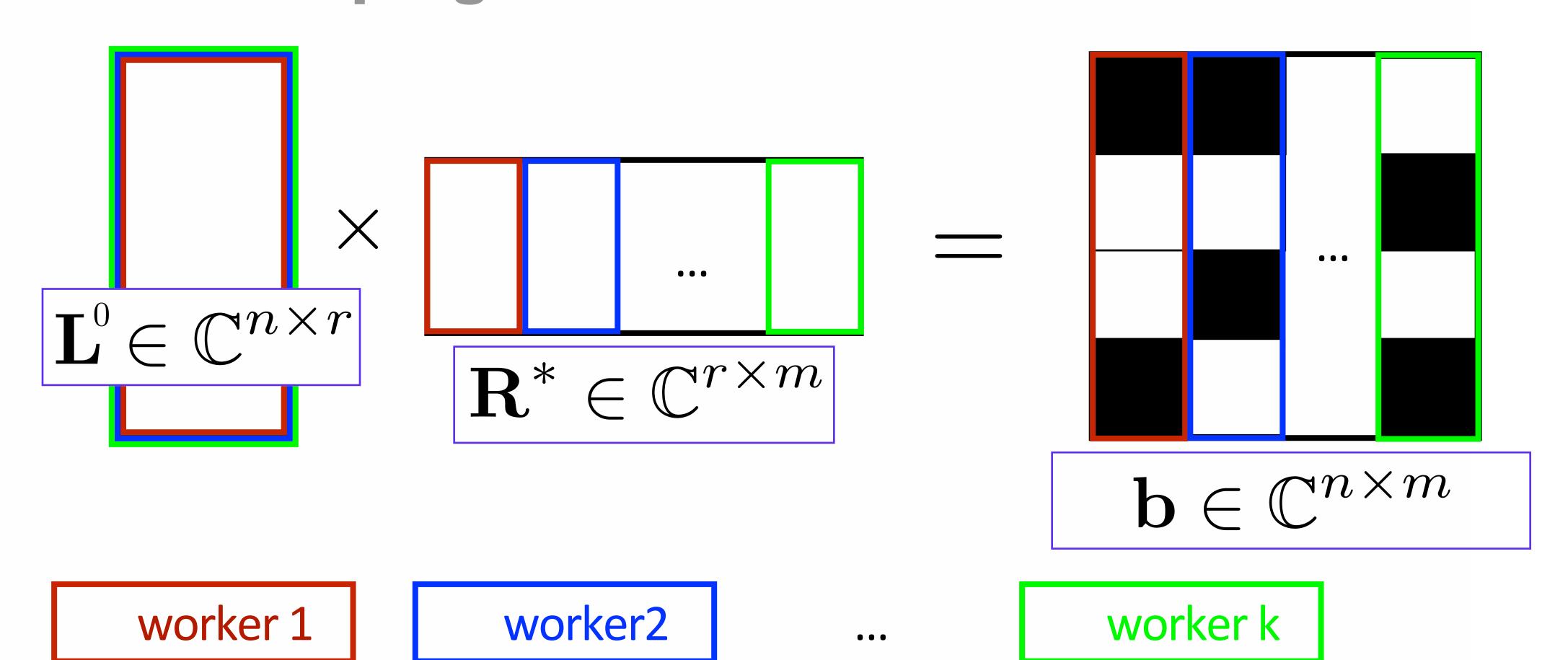
since $P_{\Omega_{\ell}}(\mathbf{L}^t)^*P_{\Omega_{\ell}}(\mathbf{L}^t) \in \mathbb{C}^{r \times r}$

(e.g., via Cholesky factorization)



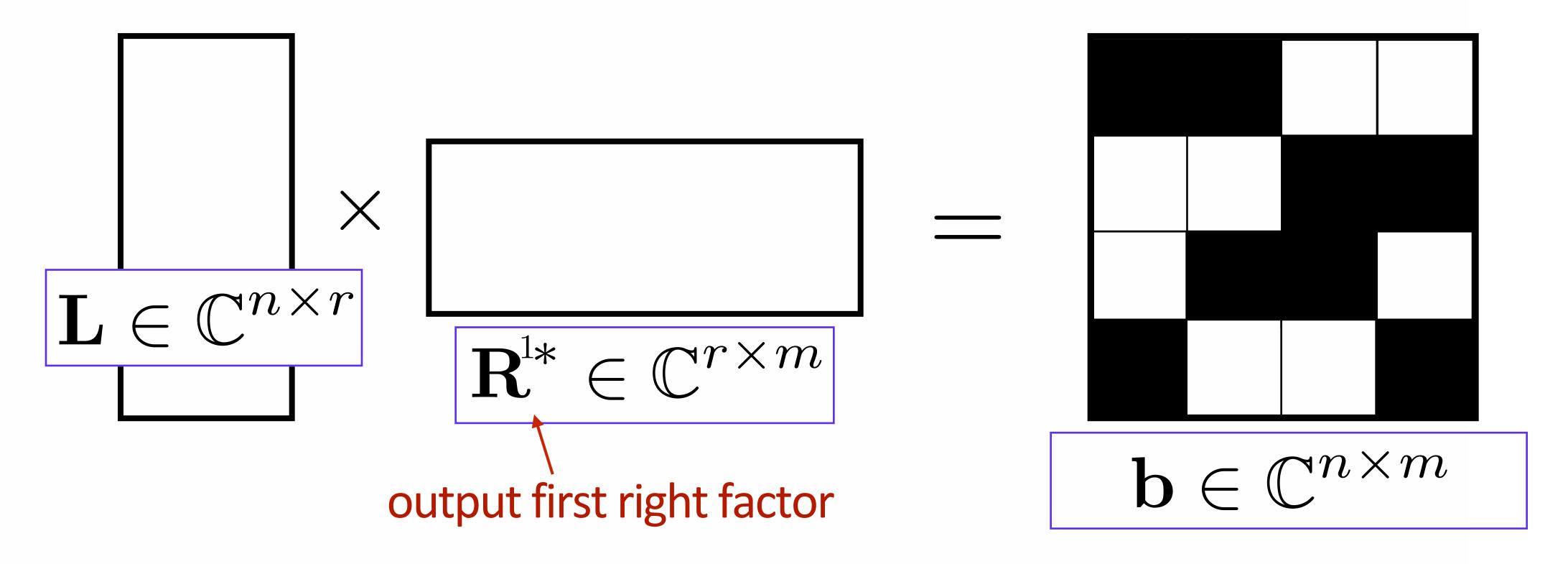




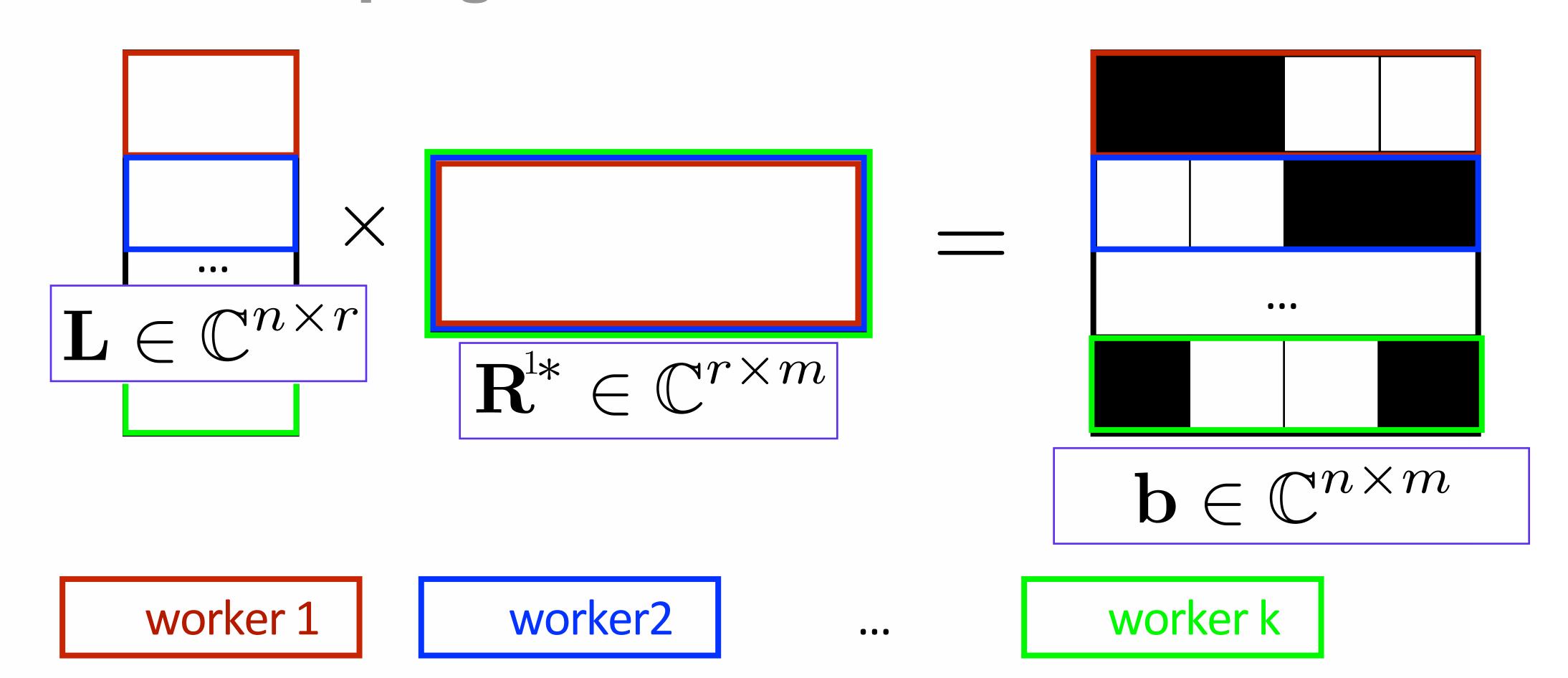


18

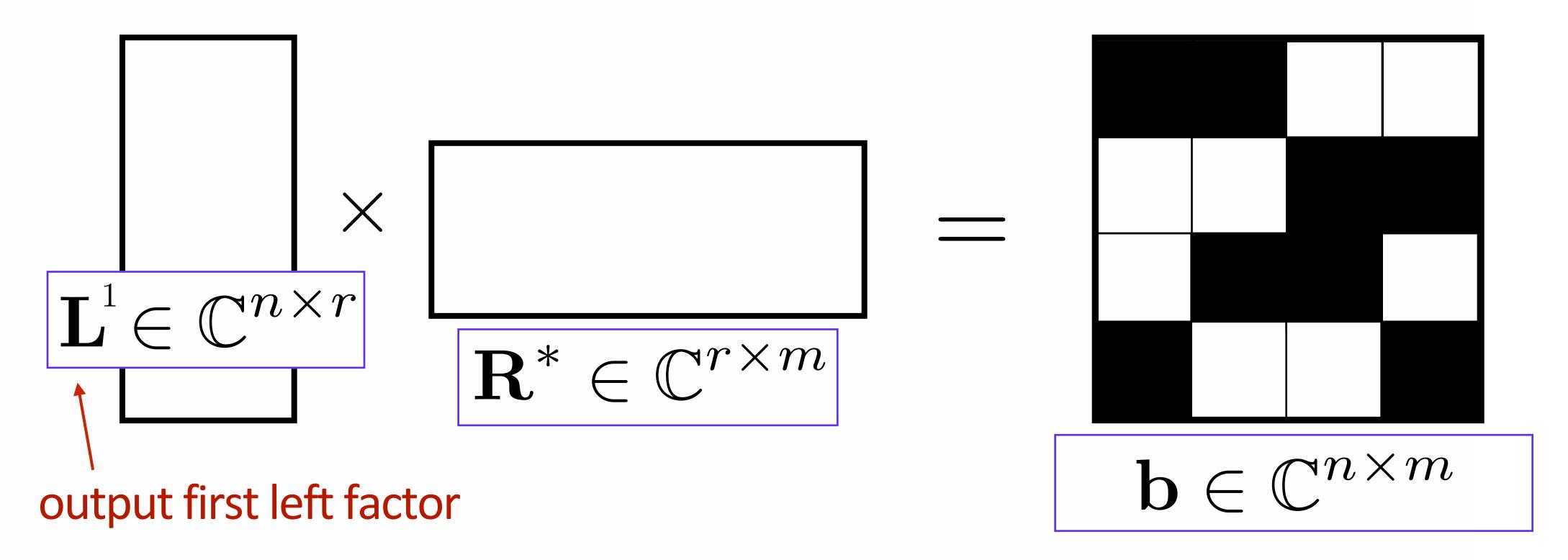




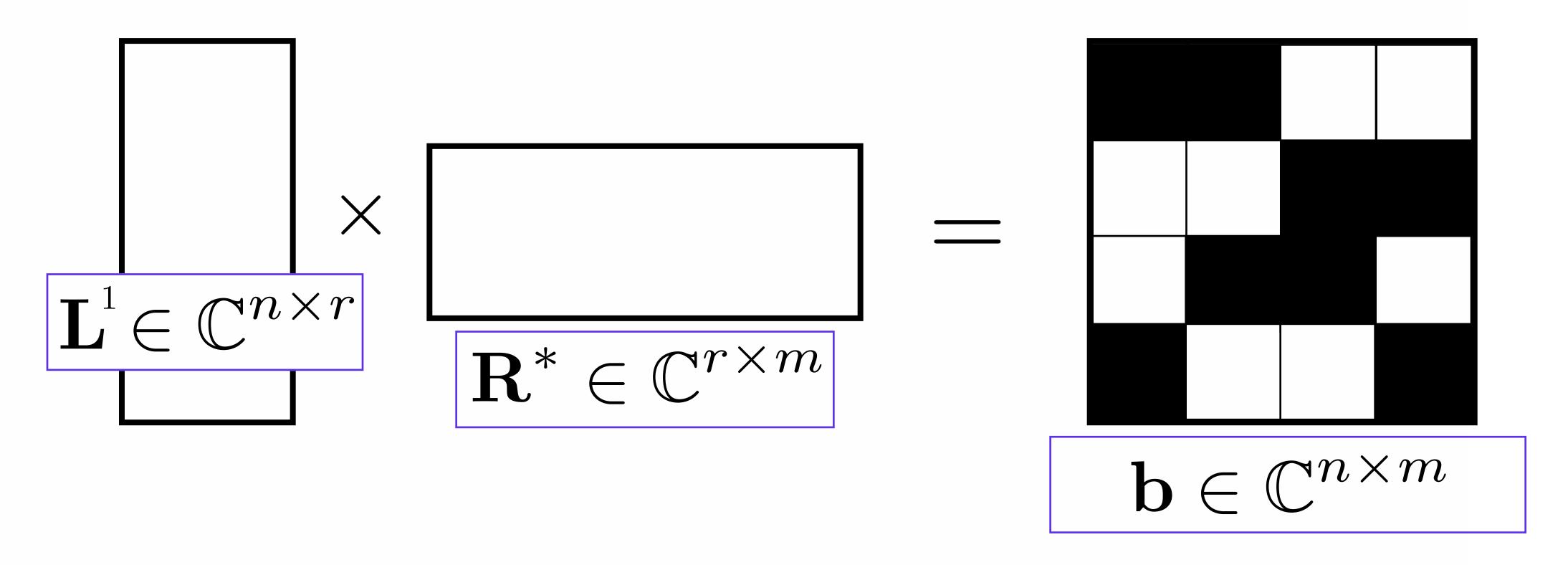












...and so on solve for $(\mathbf{L}^2, \mathbf{R}^2), \ (\mathbf{L}^3, \mathbf{R}^3), \ \dots, \ (\mathbf{L}^T, \mathbf{R}^T)$



Outline

- Matrix completion
 - alternating least squares
 - decoupling method
- Parallel implementation in Julia
- Numerical experiments

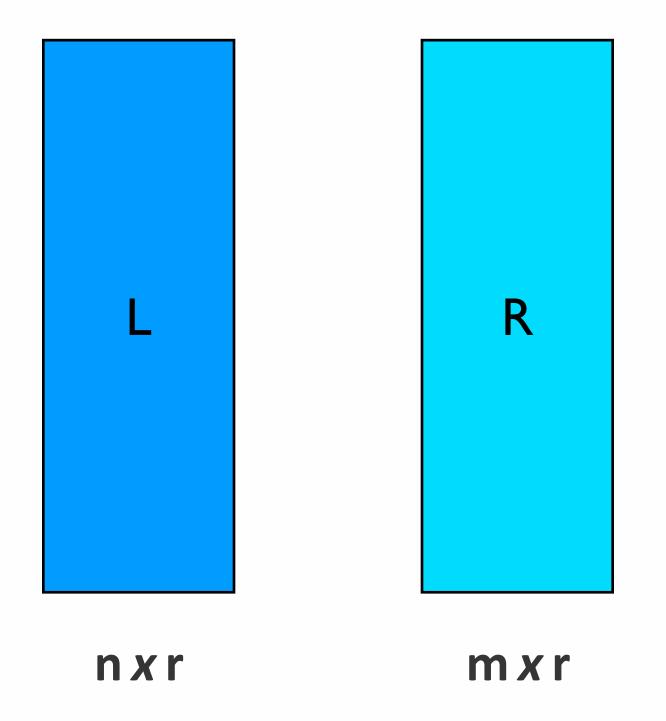


Dedicate a worker to store each L and R factor, and handle all messaging related to factor updates.

Remaining workers will store distributed data, **b**, and solve **L** & **R** on local portion of **b**.

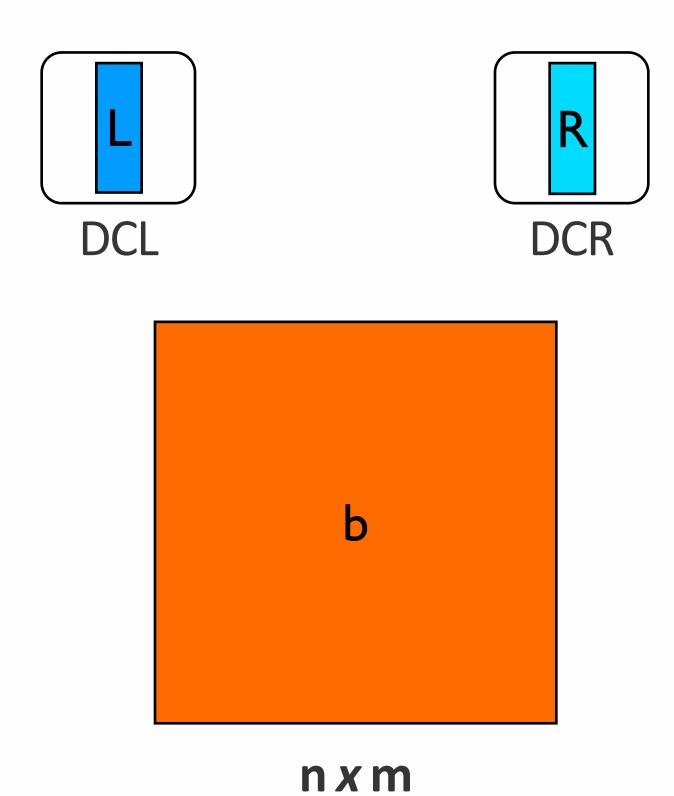


1) Dedicate a worker to store each L & R factor, and handle all messaging related to factor updates.



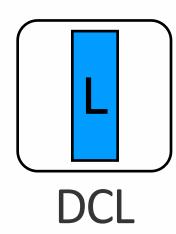


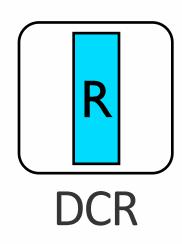
2) Remaining workers will store distributed data, b, and solve L & R on only local portion of **b**.

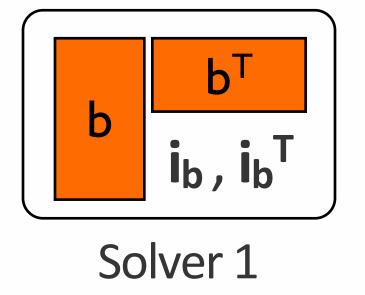


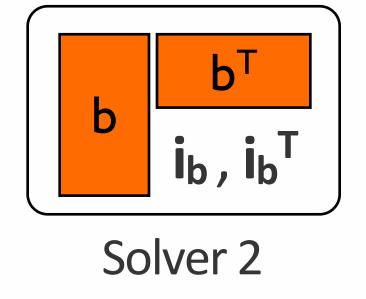


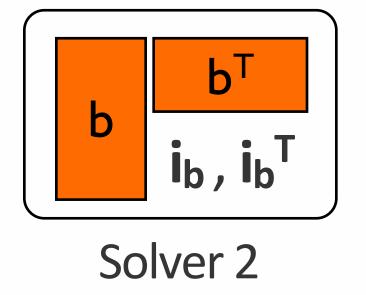
3) Distribute non-zero elements of each column of **b** & **b**^T, and corresponding indexes for non-zero values in each column, \mathbf{i}_b & \mathbf{i}_b ^T

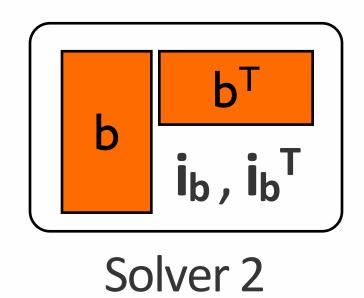






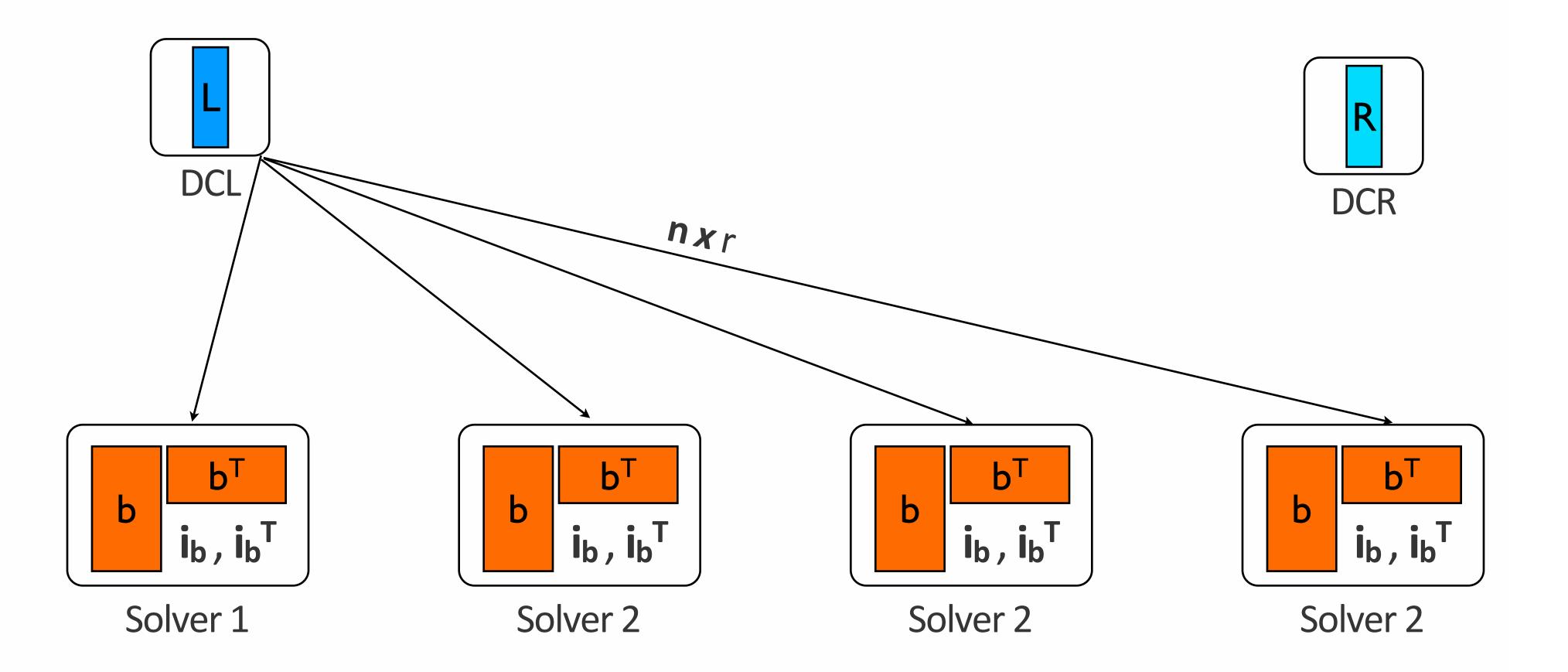






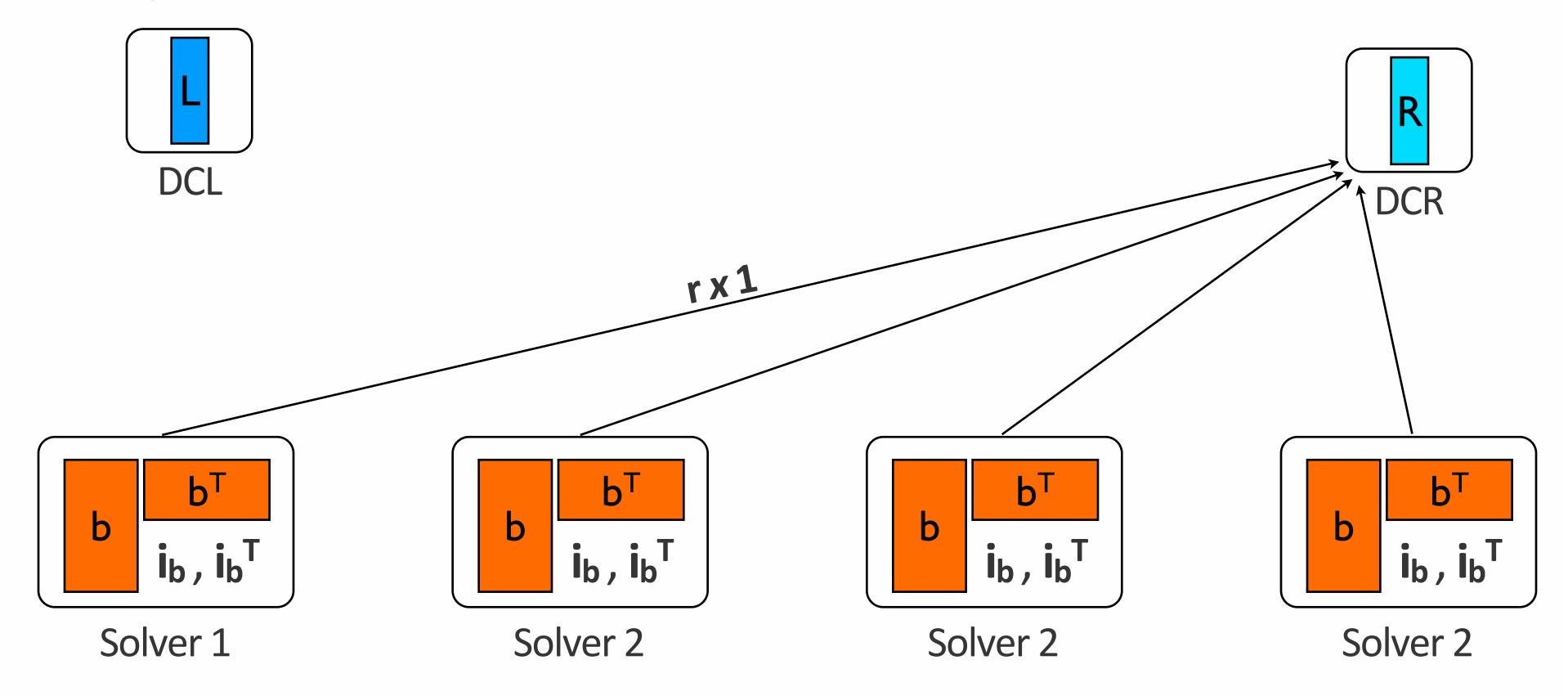


4) Solvers get latest L update from DCL



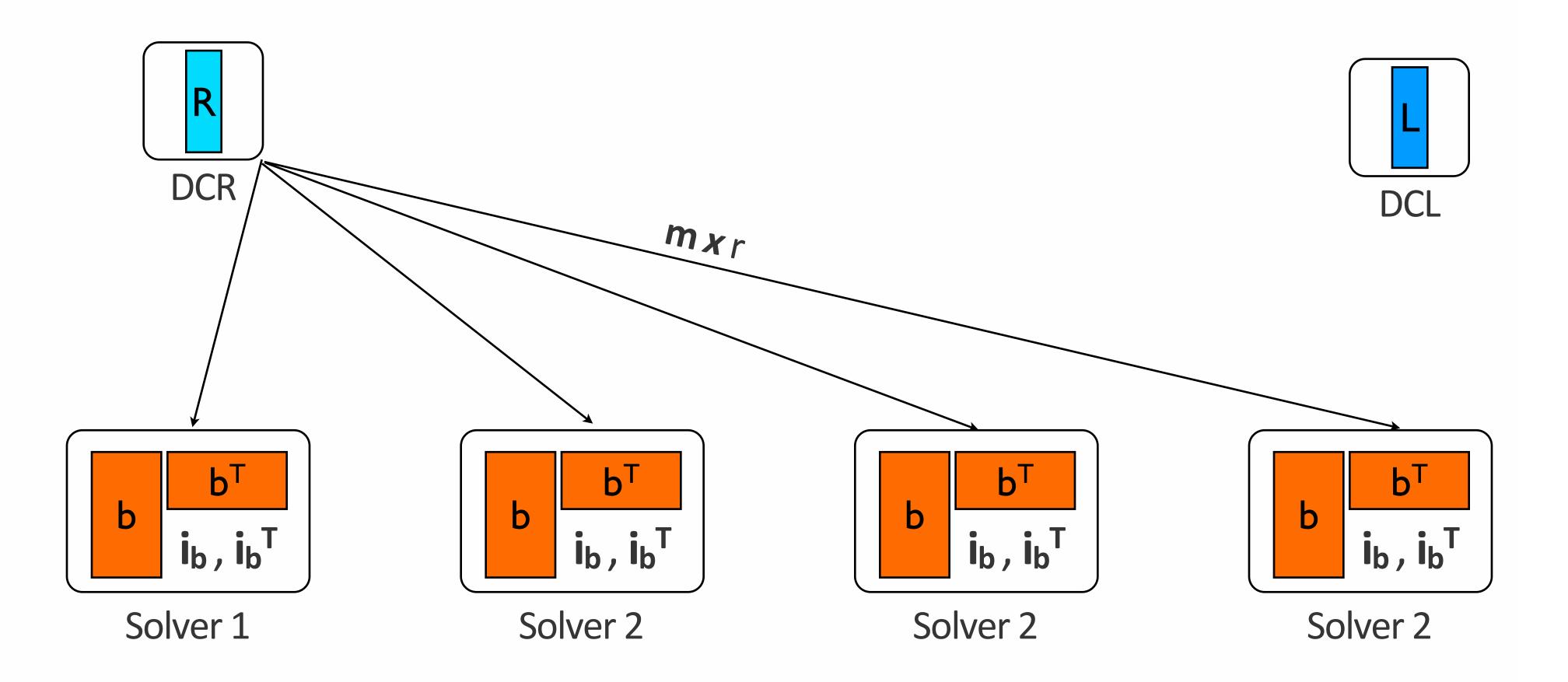


4) Solvers compute a row update of R, using L and first column of local part of **b**





4) Vice versa for R update





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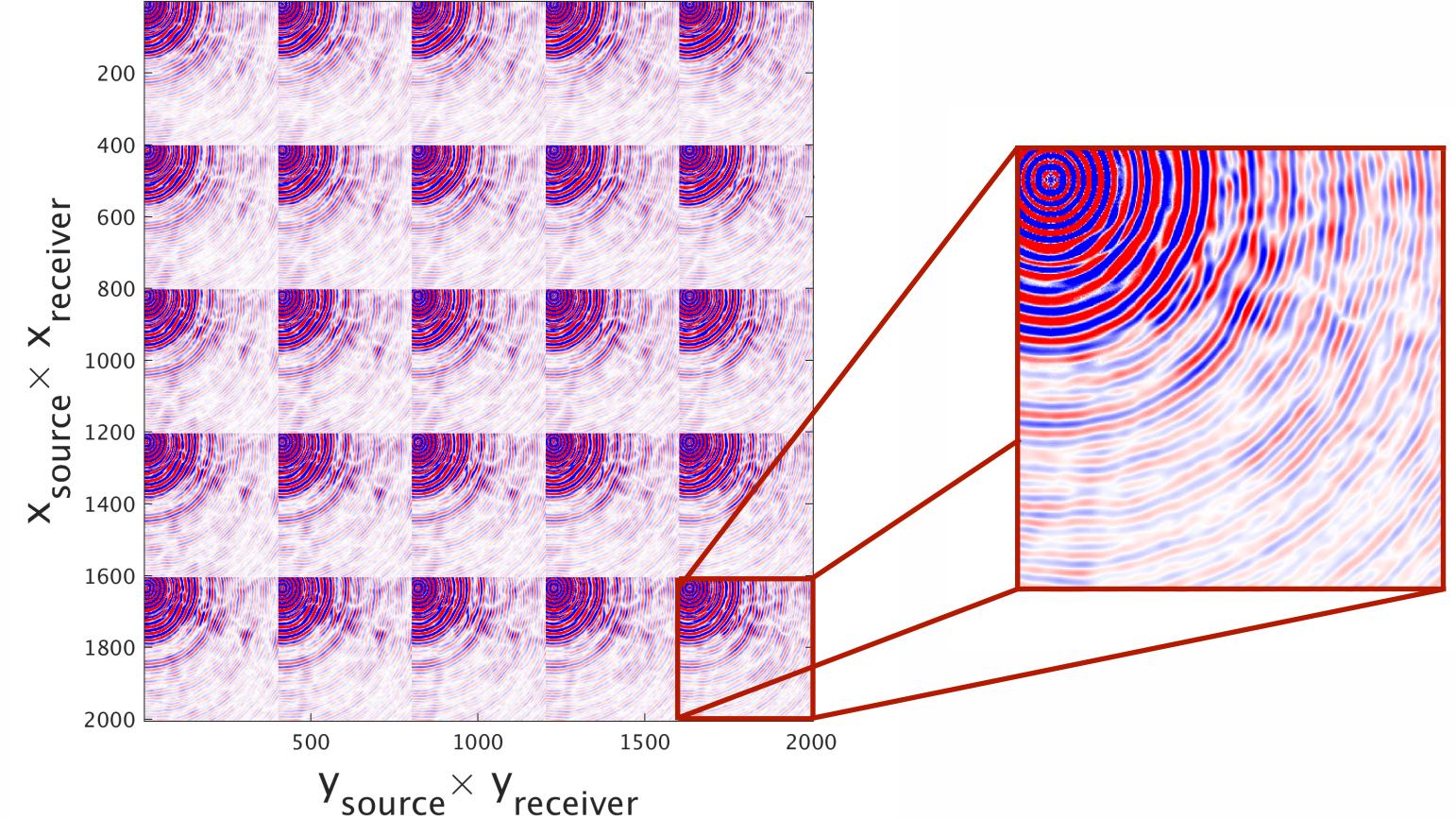
Interpolation: Synthetic BG 3D Model

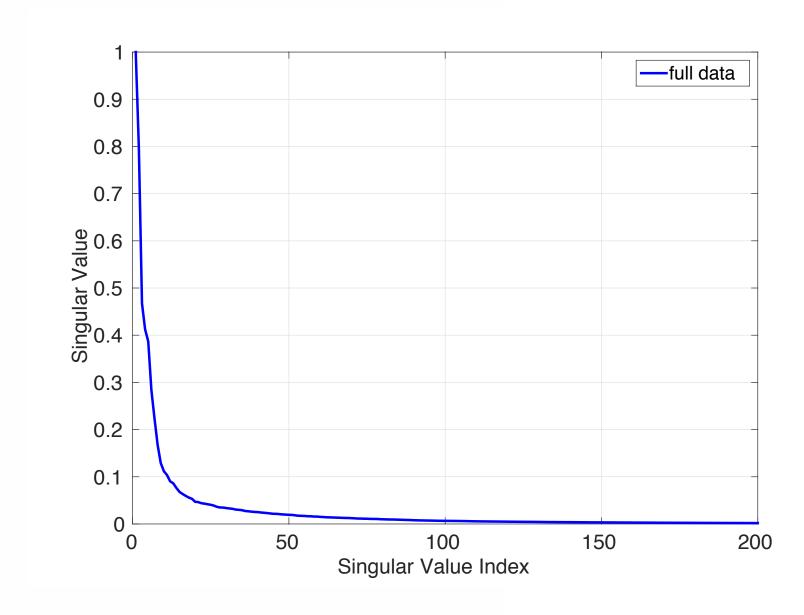
- ▶ 68 x 68 sources with 401 x 401 receivers
- ▶ Data at 7.34 Hz and 12.3 Hz.
- ▶ Matricize in "(rec,rec)"-form



Data Matricized - (rec,rec) form

BG 3D Dataset 7.35 Hz

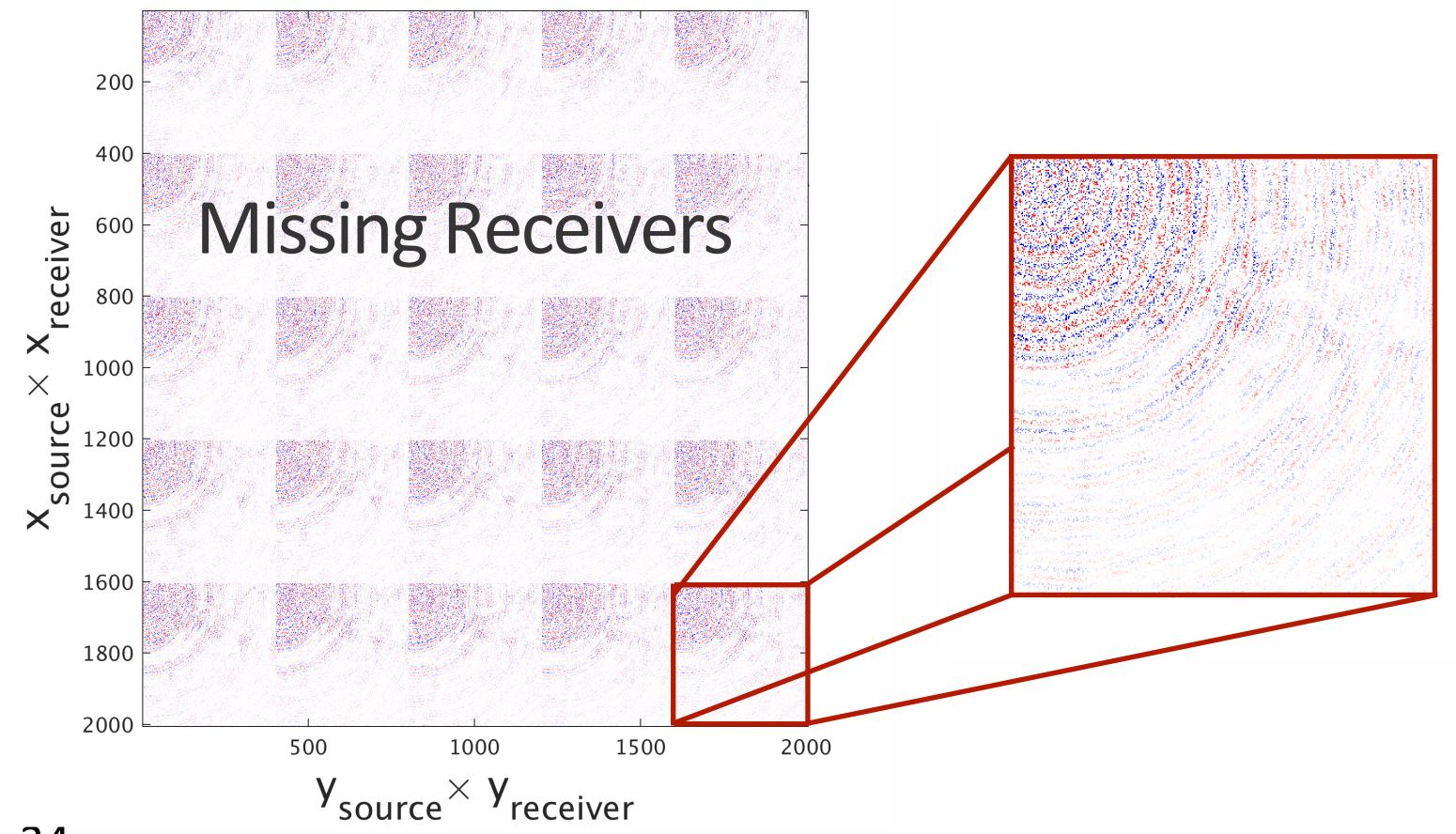


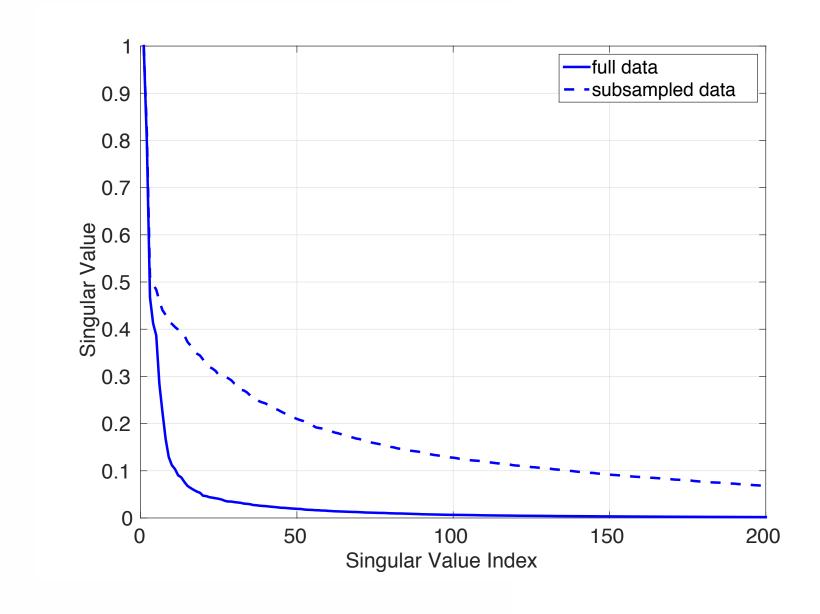




Data Matricized - (rec,rec) form

BG 3D Dataset 7.35 Hz





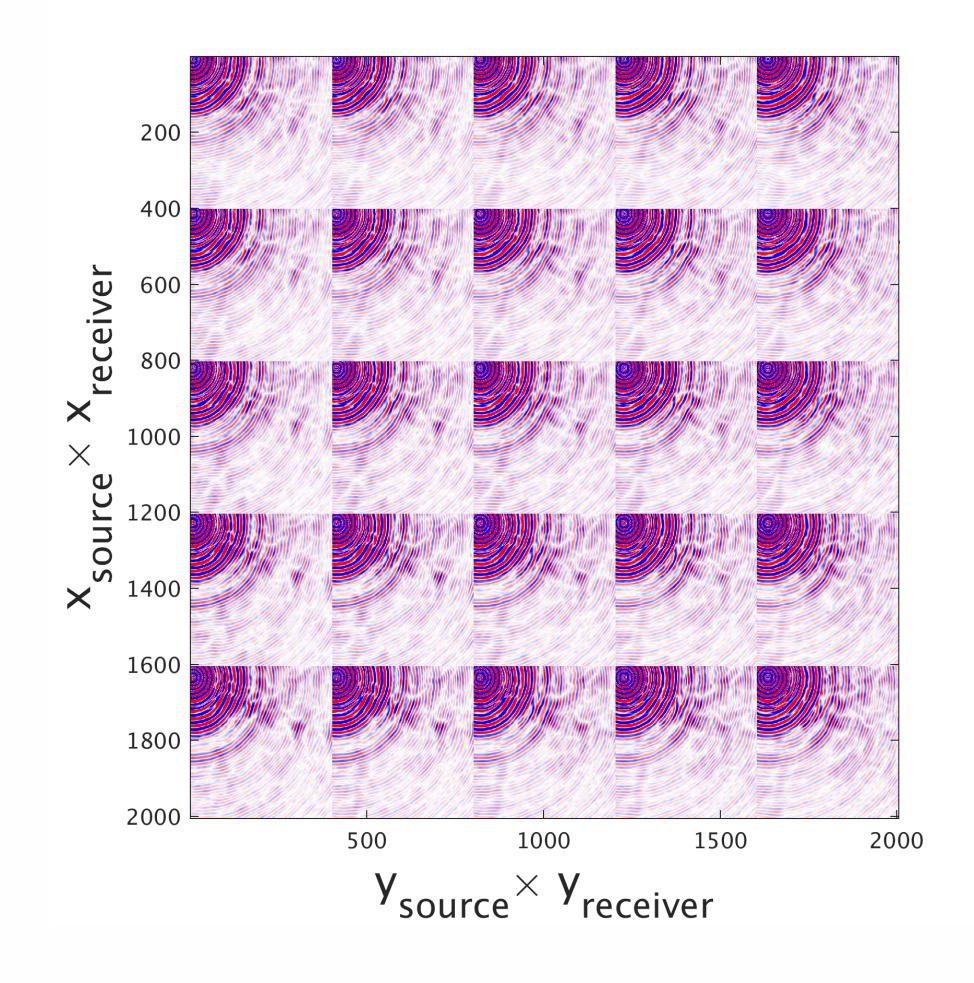
34



3D Interpolation Experiment

BG 3D Dataset

7.35 Hz



Size: 27,268 x 27,268

(full slice, no windowing)

Remove 80 % of Receivers randomly

Compare Interpolation via:

- SPG-LR
- Decoupling method



How to choose the rank parameter?

Issue: need

$$\mathbf{R}^{t+1}(\ell,:) = \left(P_{\Omega_{\ell}}(\mathbf{L}^t)^* P_{\Omega_{\ell}}(\mathbf{L}^t)\right)^{-1} P_{\Omega_{\ell}}(\mathbf{L}^t)^* \mathbf{b}(:,\ell)$$

How do we know $P_{\Omega_{\ell}}(\mathbf{L}^t)^* P_{\Omega_{\ell}}(\mathbf{L}^t) \in \mathbb{C}^{r \times r}$ is invertible?



How to choose the rank parameter?

Theorem:

Let Ω be chosen uniformly at random.

Let $\mathbf{L} \in \mathbb{C}^{n \times r}$ be full rank, define $\tilde{\mathbf{L}} = \operatorname{orth}(\mathbf{L})$ and $\beta := \max_{k,\ell} \left(\tilde{\mathbf{L}}_{k,\ell} \right)^2$.

Then if
$$|\Omega| \geq \alpha \frac{8}{3} \beta n r \log(nr)$$

 $P_{\Omega_\ell}(\mathbf{L}^t)^*P_{\Omega_\ell}(\mathbf{L}^t)$ is invertible for every $\ell \in \{1,2,...,n\}$ with probability $\geq 1-2n^{1-\alpha}$



How to choose the rank parameter?

$$|\Omega| \ge \alpha \frac{8}{3} \beta n r \log(nr)$$

In our case:
$$|\Omega|=.2\cdot nm$$

$$n=m=27,268$$

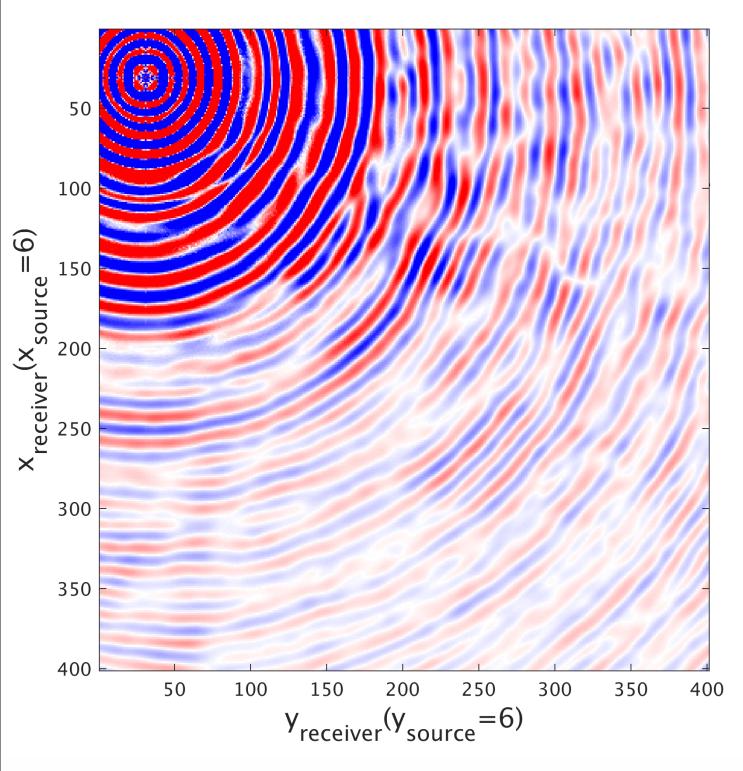
(ignoring constants)
$$\implies r \leq 534$$

choose upper bound as rank, gives well defined procedure.

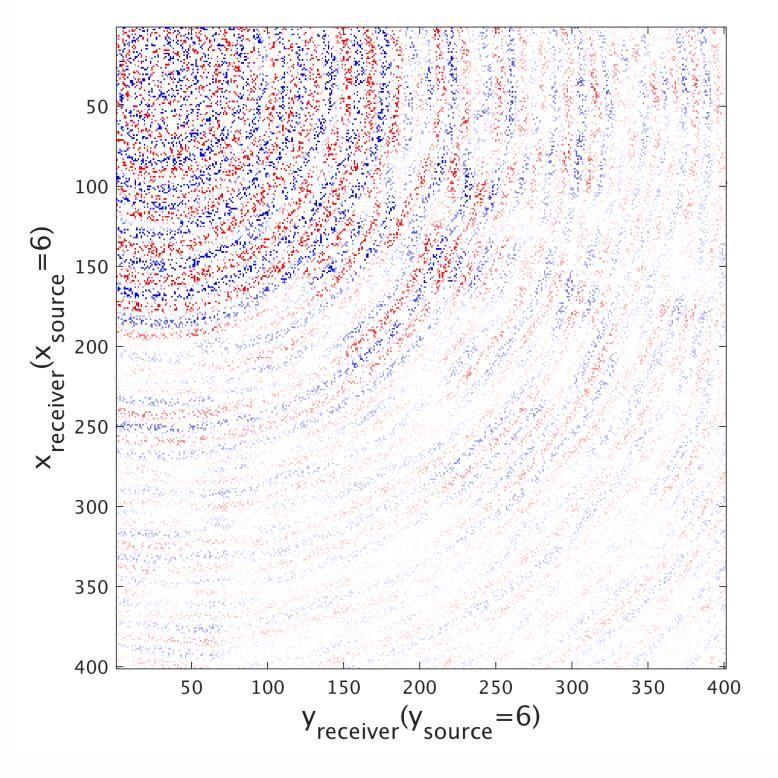


Common Source Gather

True Source Gather



Subsampled Source Gather

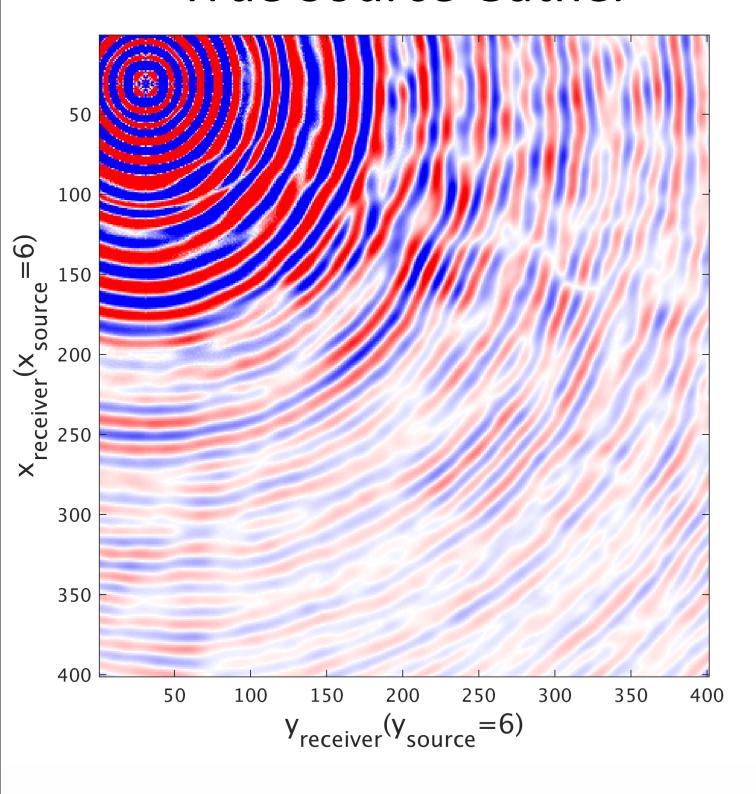


Remove 80 % of Receivers randomly

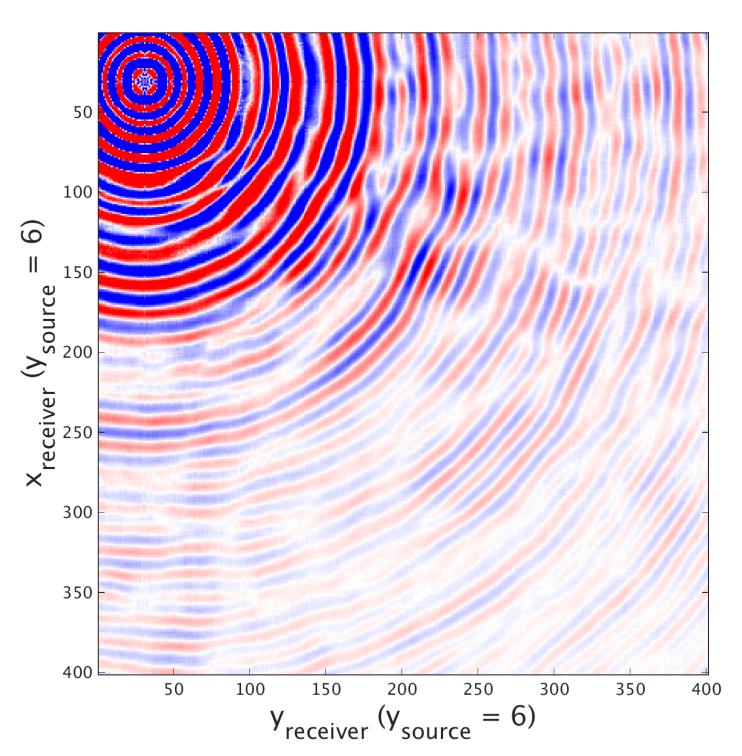


Results: SPG-LR

True Source Gather



Recovered Source Gather



SPG-LR iterations: 400

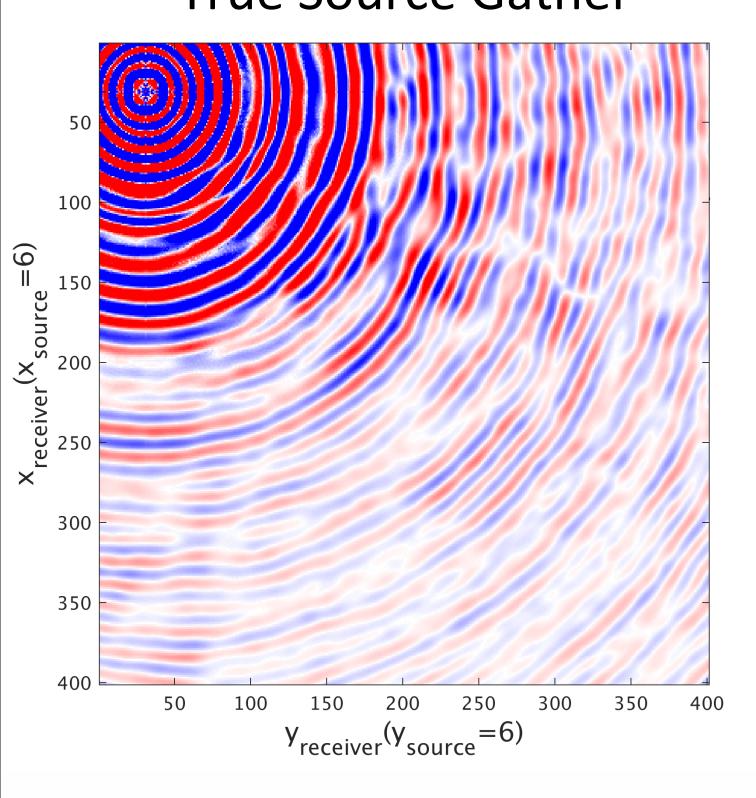
SNR = 26.1 dB

Time = 82 hrs and 40 min

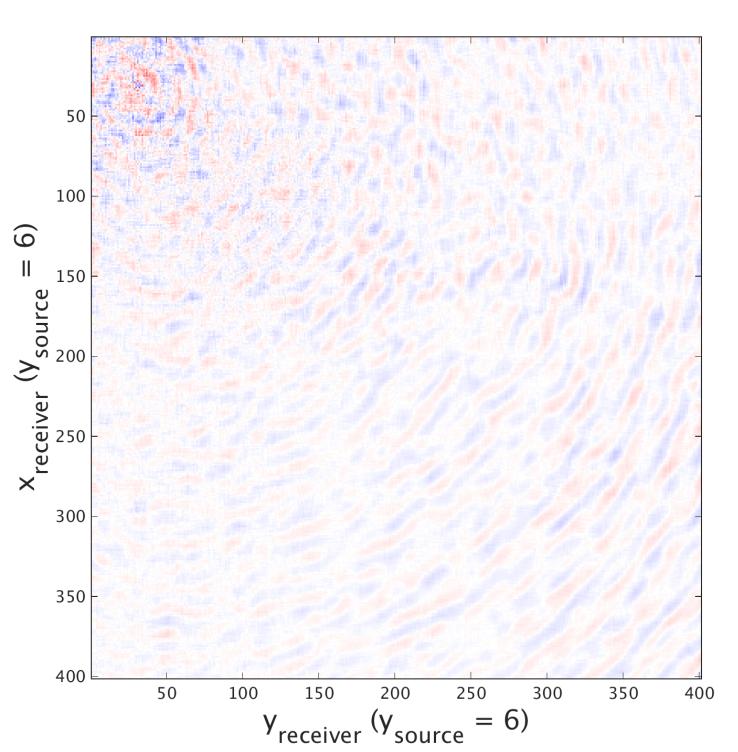


Results: SPG-LR

True Source Gather



Difference Plot



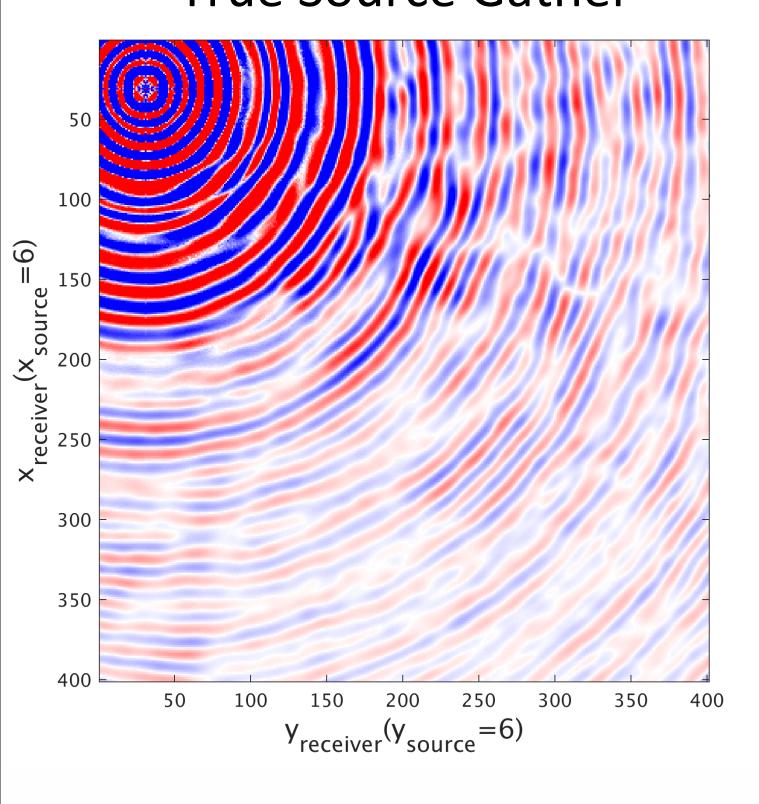
SPG-LR iterations: 400

SNR = 26.1 dB

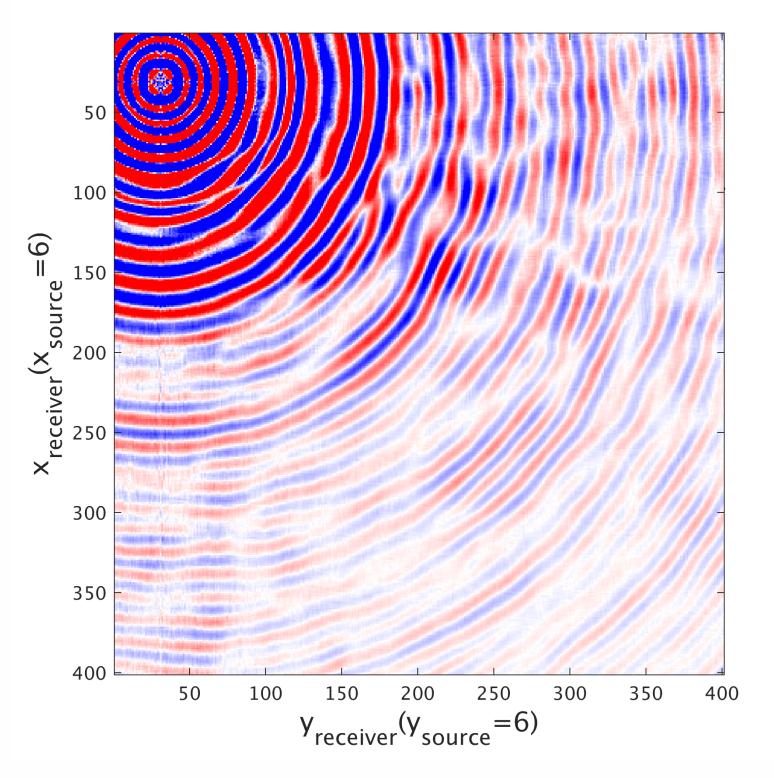
Time = 82 hrs and 40 min



True Source Gather



Recovered Source Gather



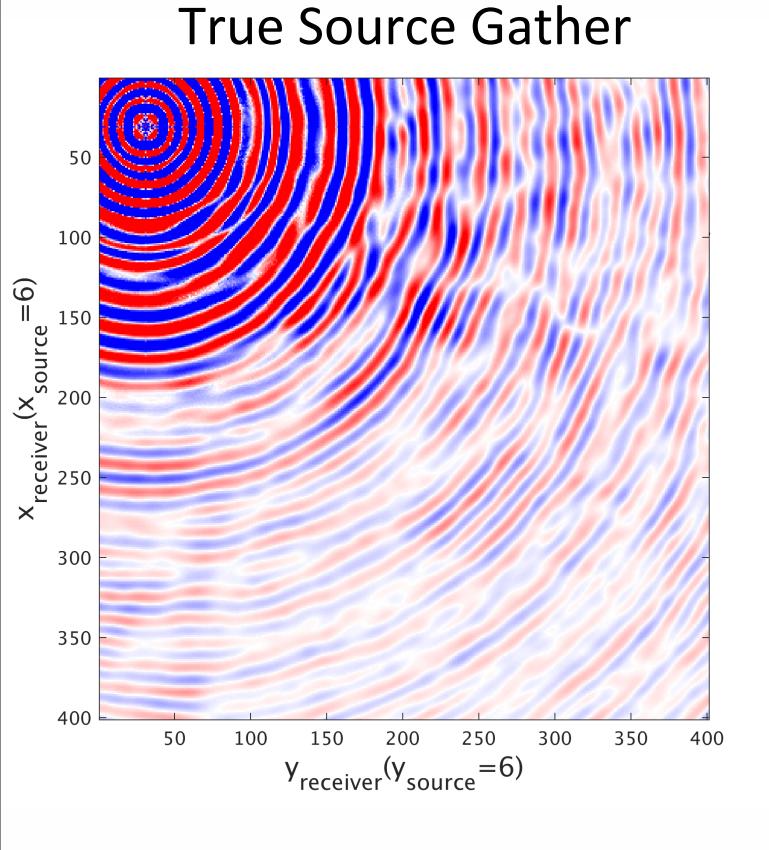
60 workers

Alternations: 5

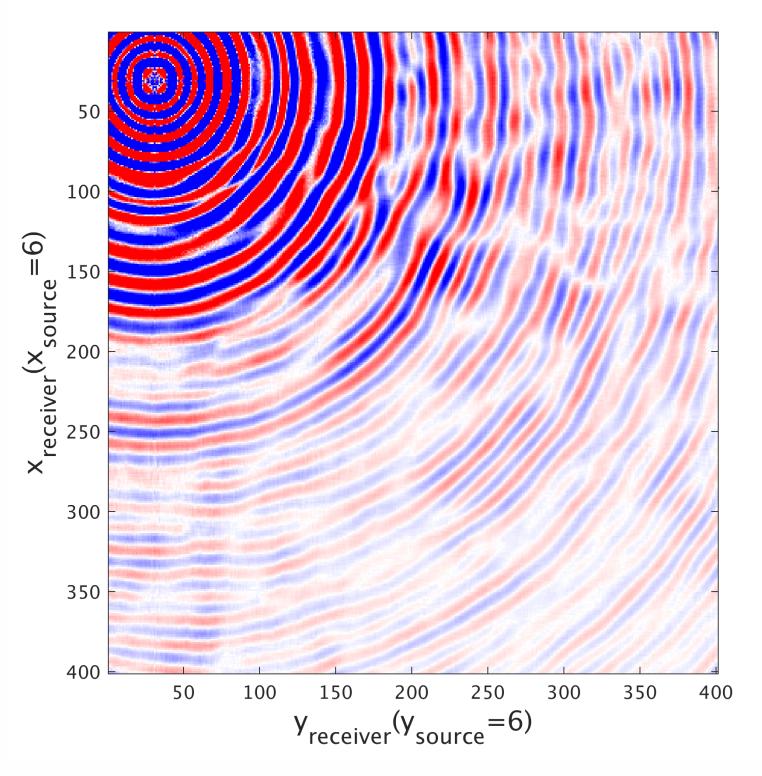
SNR = 24.2 dB

Time = 1 hr and 7 mins





Recovered Source Gather



60 workers

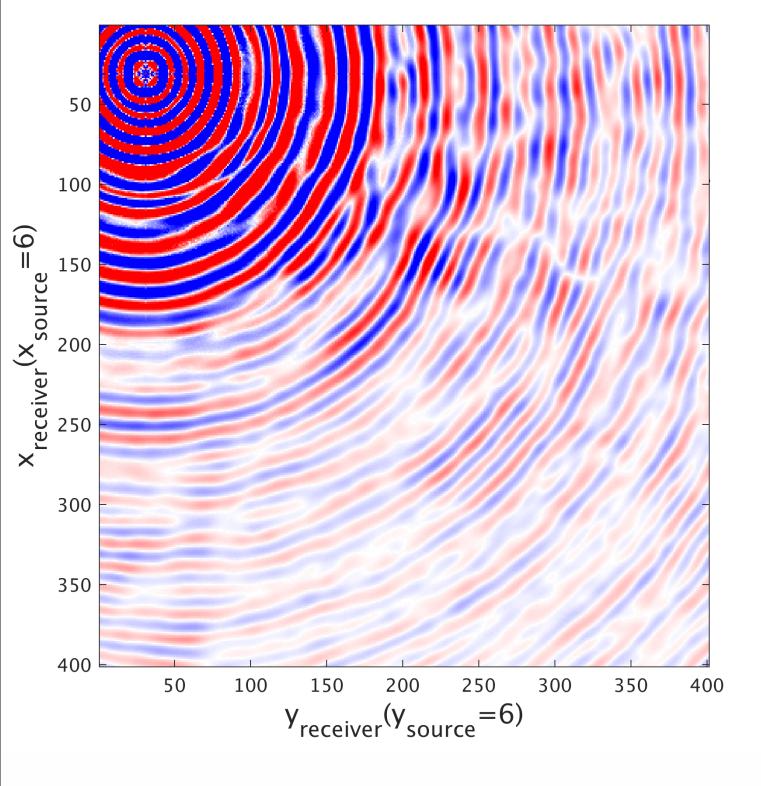
Alternations: 7

SNR = 25.1 dB

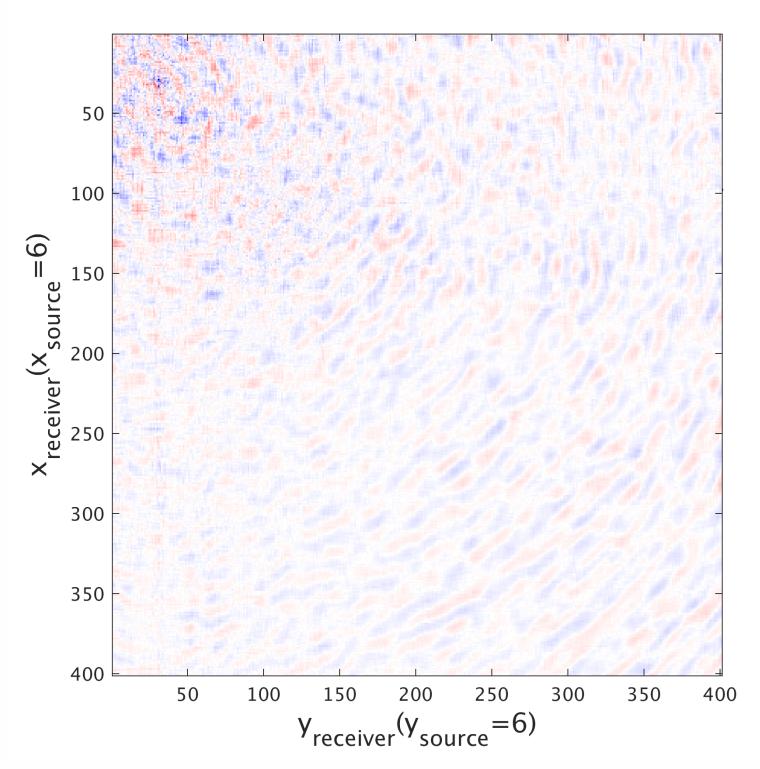
Time = 1 hr and 33 mins







Difference Plot



60 workers

Alternations: 7

SNR = 25.1 dB

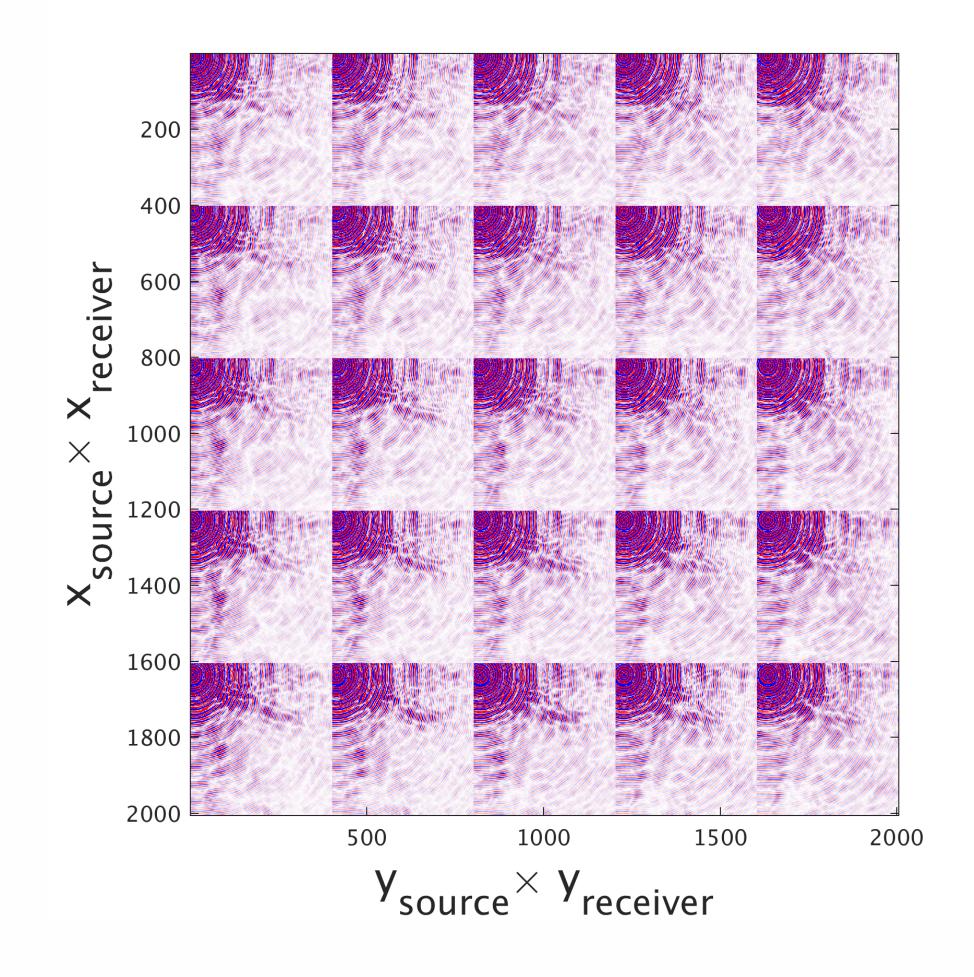
Time = 1 hr and 33 mins



3D Interpolation Experiment

BG 3D Dataset

12.3 Hz



Size: 27,268 x 27,268

(full slice, no windowing)

Remove 80 % of Receivers randomly

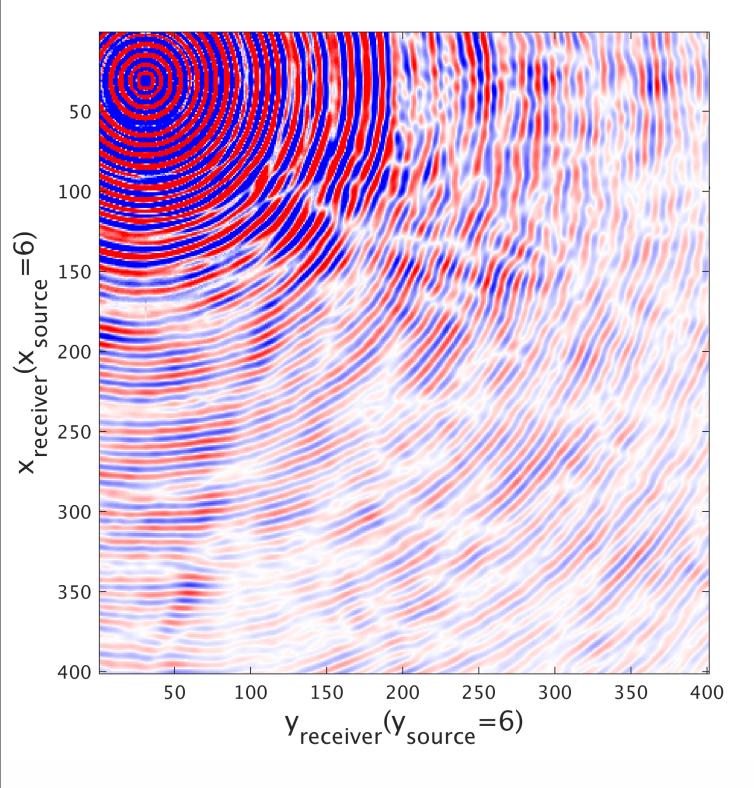
Compare Interpolation via:

- SPG-LR
- Decoupling method

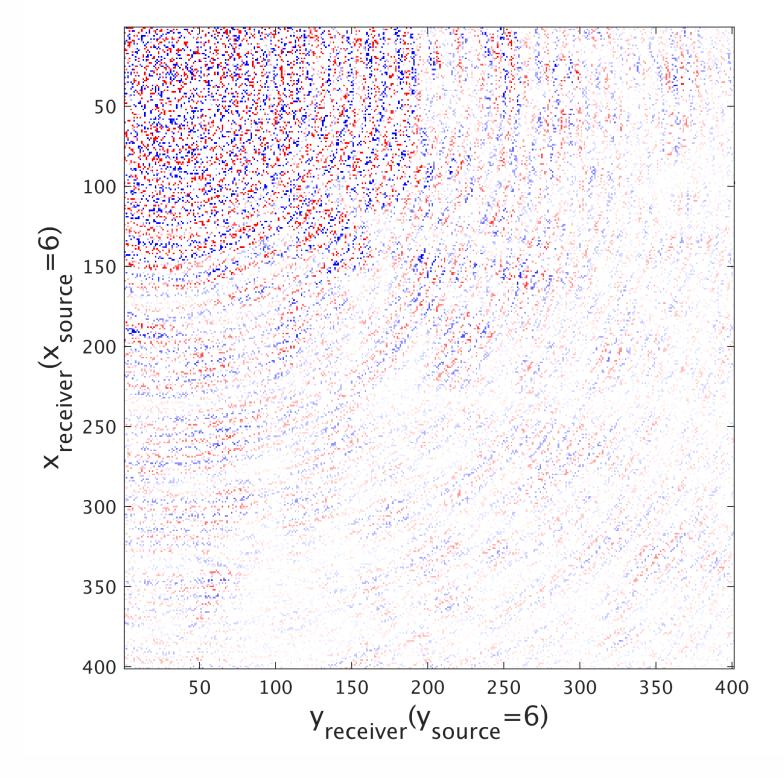


Common Source Gather

True Source Gather



Subsampled Source Gather

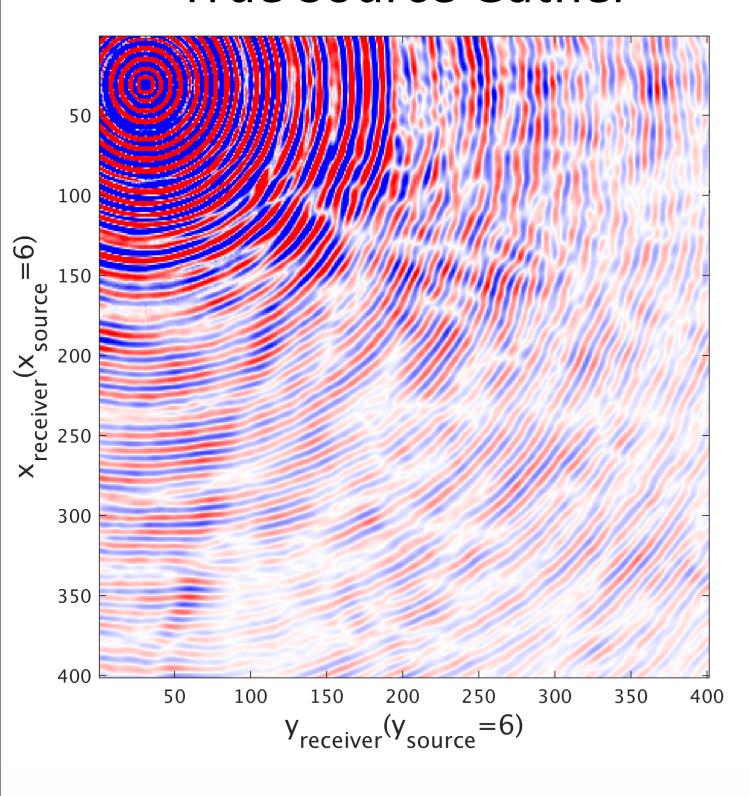


Remove 80 % of Receivers randomly

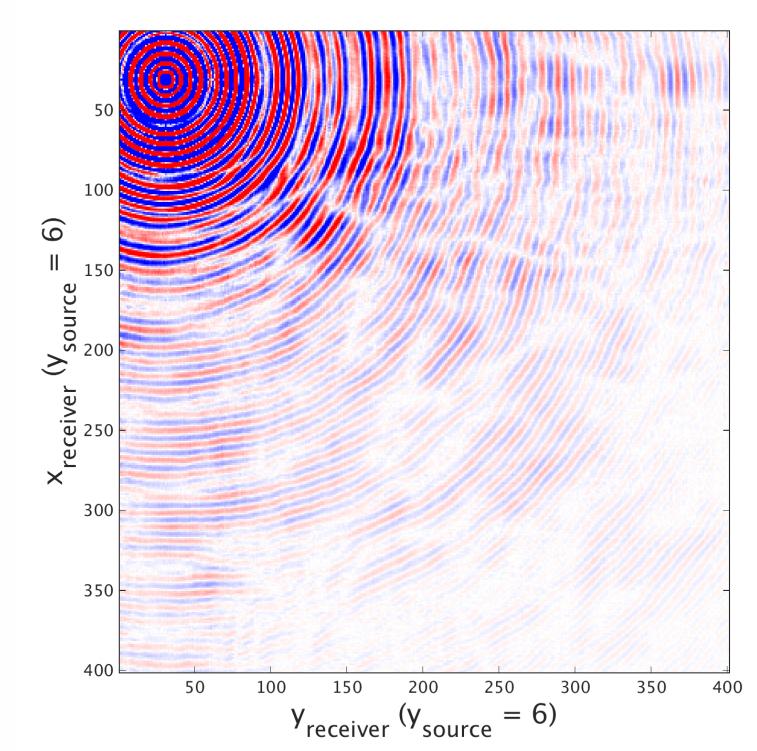


Results: SPG-LR

True Source Gather



Recovered Source Gather



SPG-LR iterations: 400

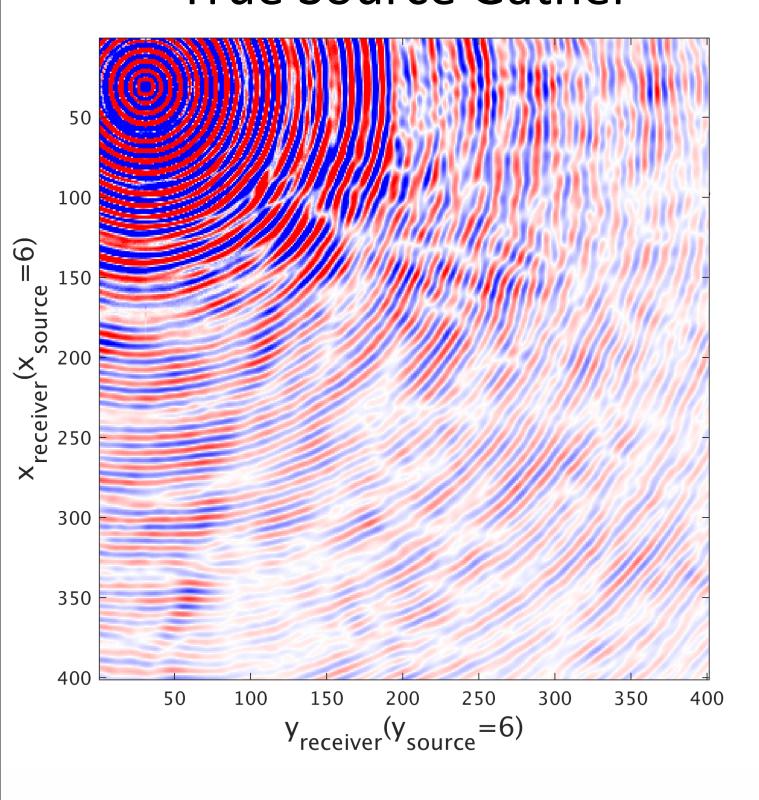
SNR = 20.5 dB

Time = 137 hrs and 20 min

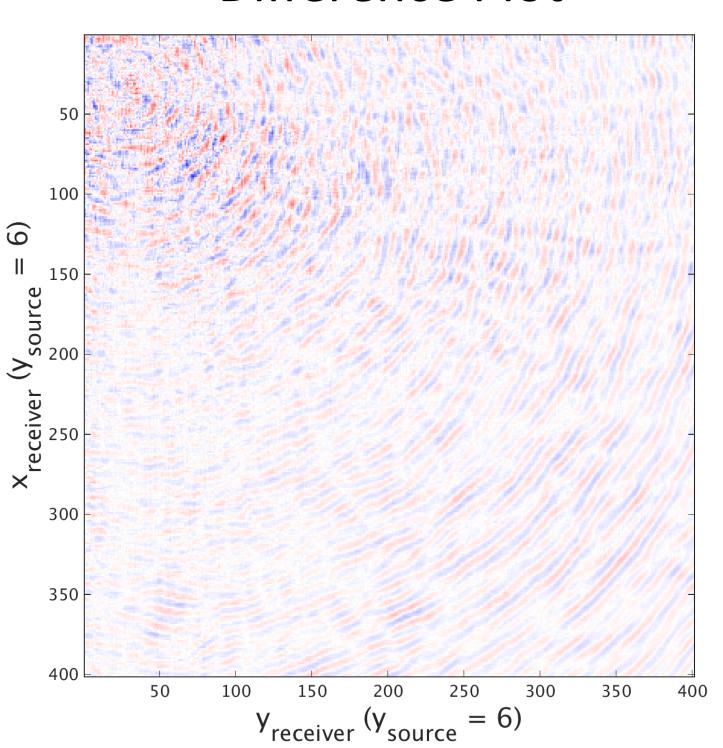


Results: SPG-LR

True Source Gather



Difference Plot



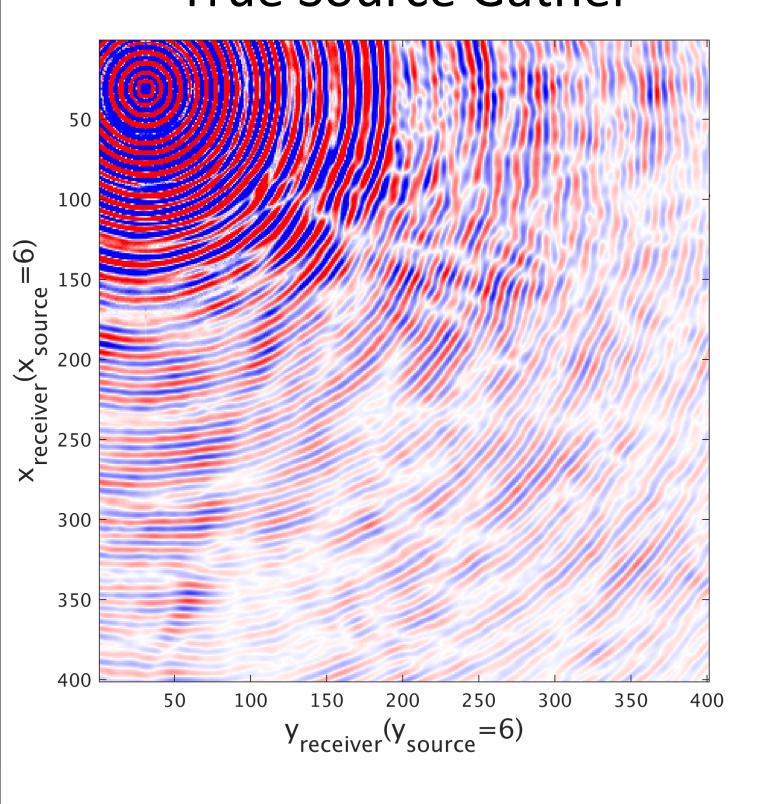
SPG-LR iterations: 400

SNR = 20.5 dB

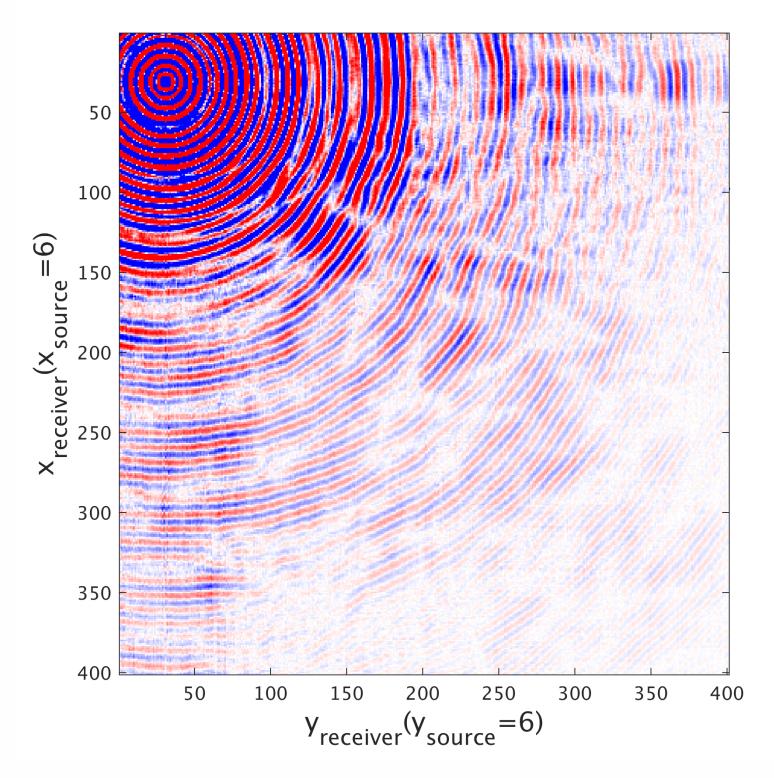
Time = 137 hrs and 20 min



True Source Gather



Recovered Source Gather



60 workers

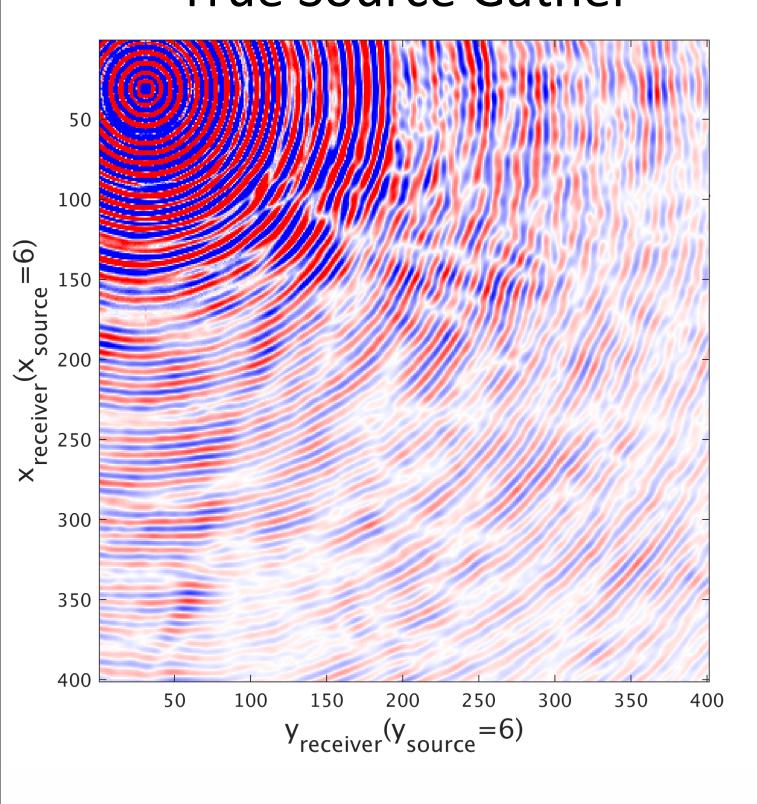
Alternations: 5

SNR = 19 dB

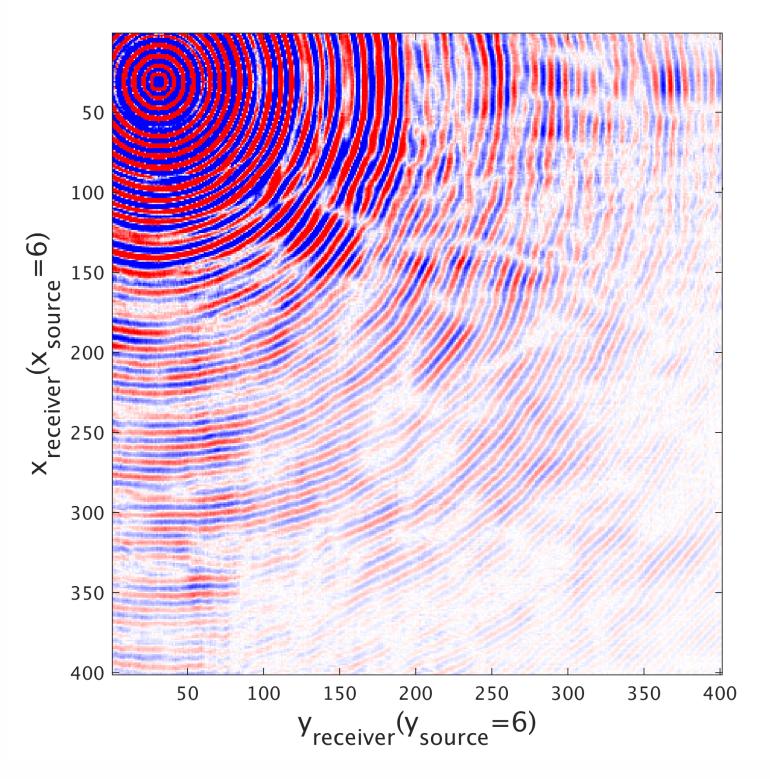
Time = 1 hrs and 7 mins



True Source Gather



Recovered Source Gather



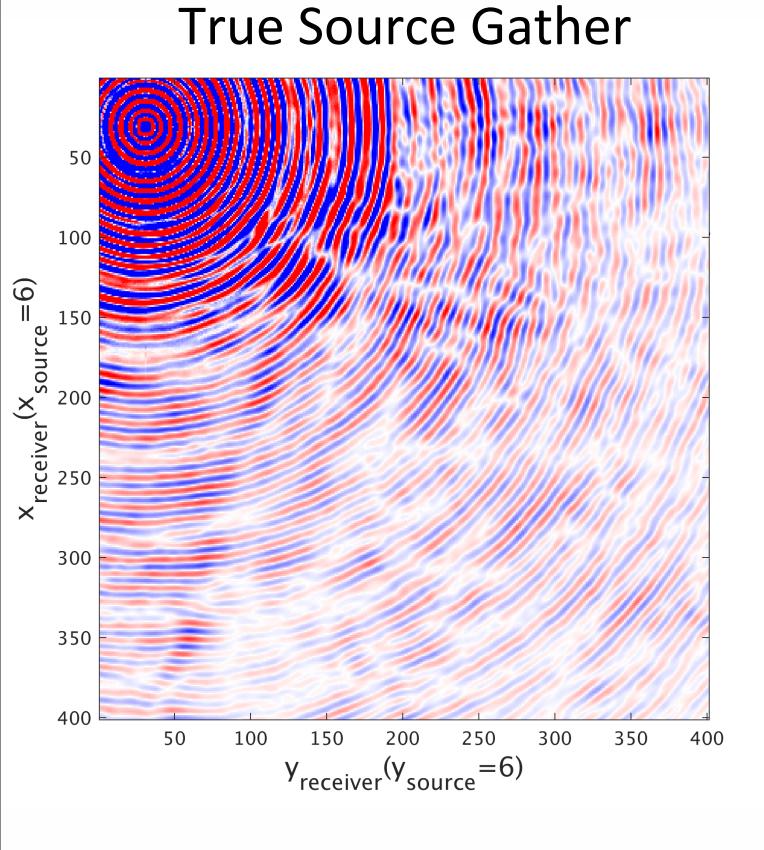
60 workers

Alternations: 7

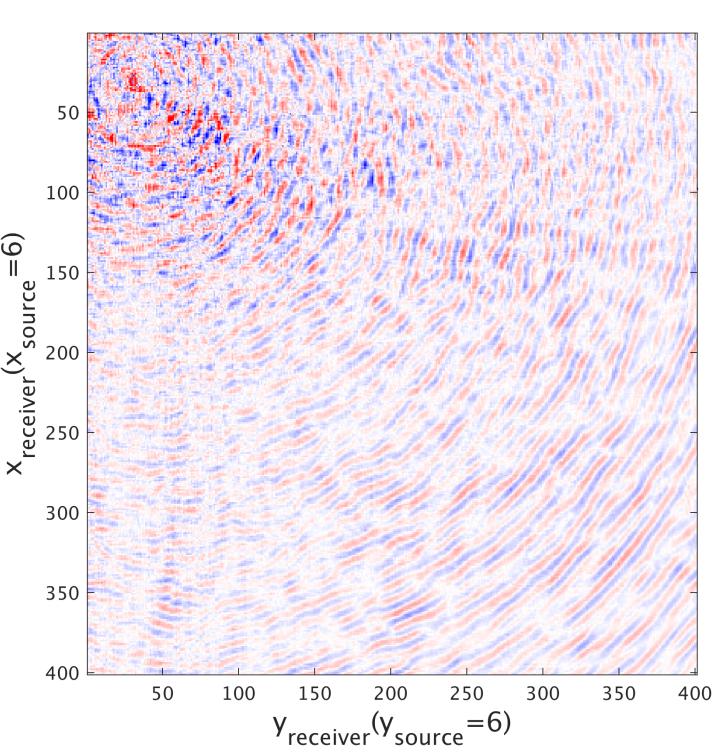
SNR = 20 dB

Time = 1 hrs and 36 mins





Difference Plot



60 workers

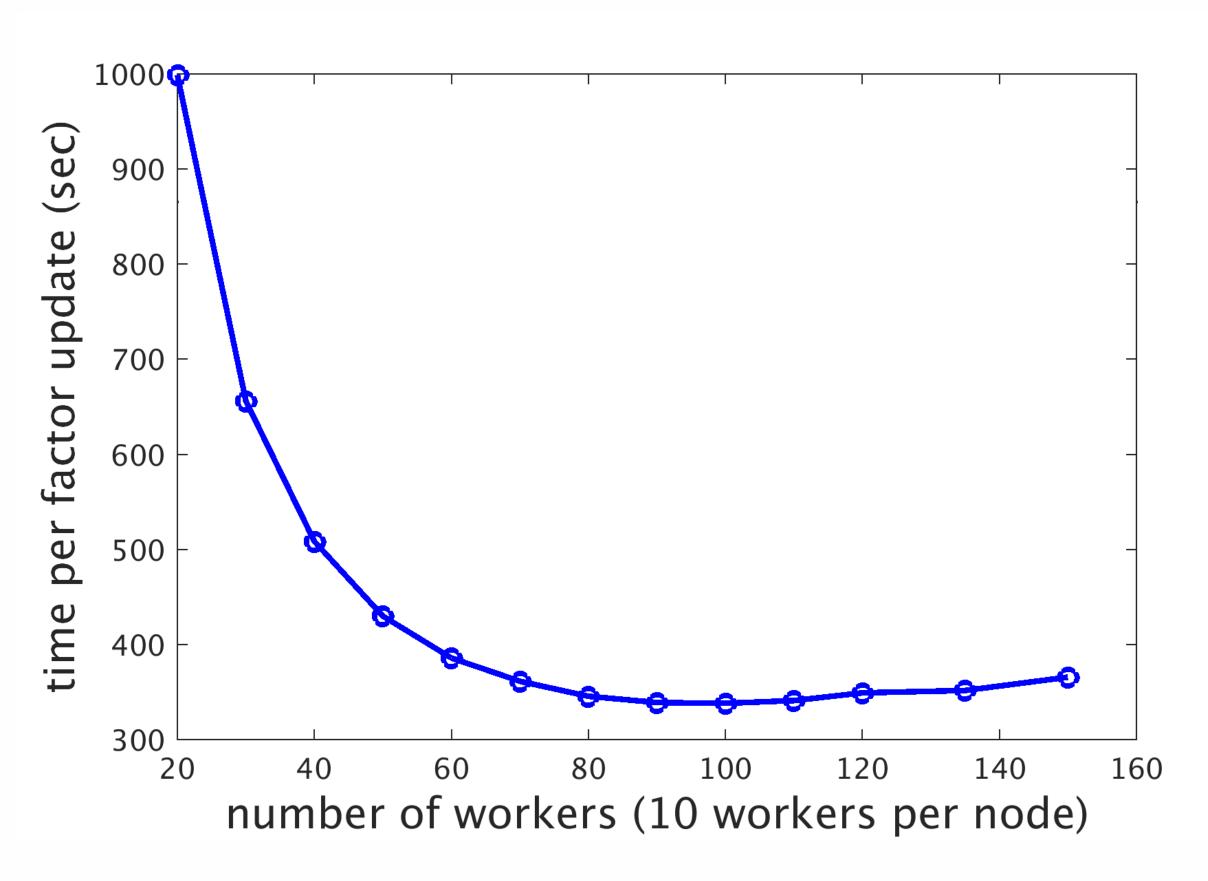
Alternations: 7

SNR = 20 dB

Time = 1 hrs and 36 mins



Scalability: time vs # workers



Matrix Size: 27,268 x 27,268

(full slice, no windowing)

rank = 534

missing 80% receivers



To window or not to window?

Look at recovery with various windows sizes

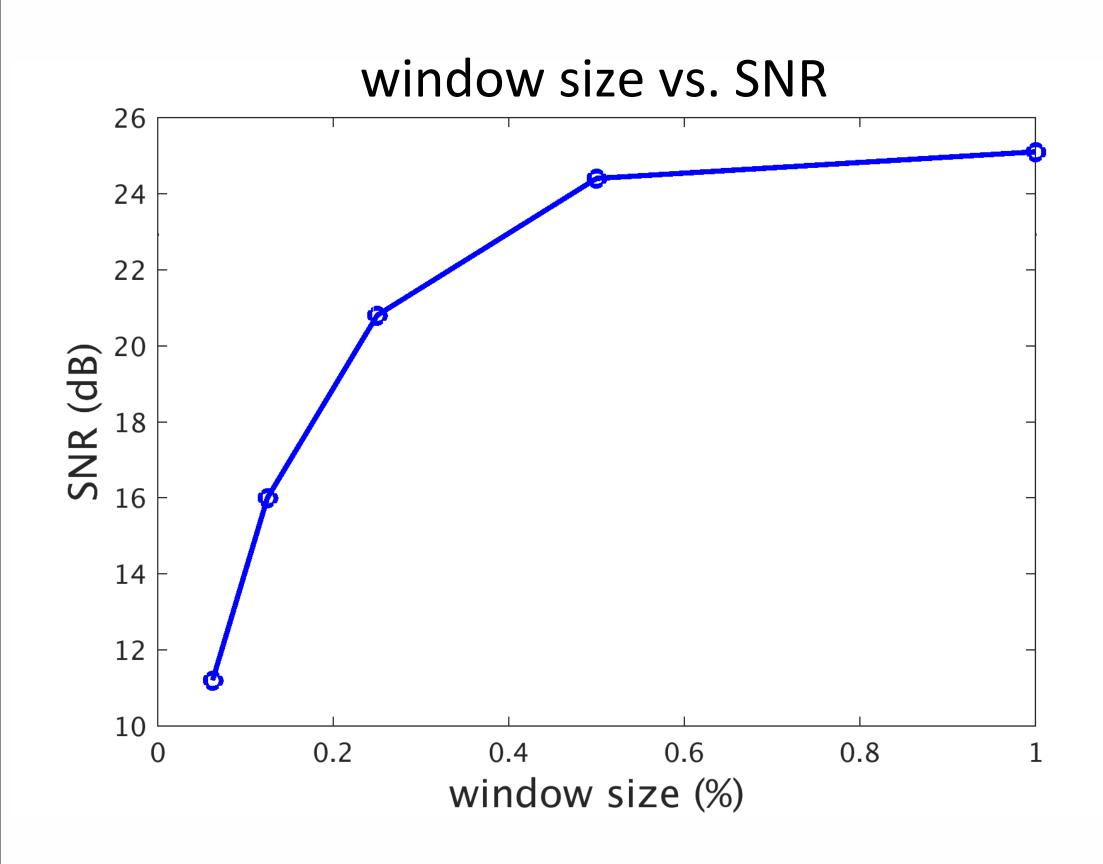
Rank chosen according to window size (as before)

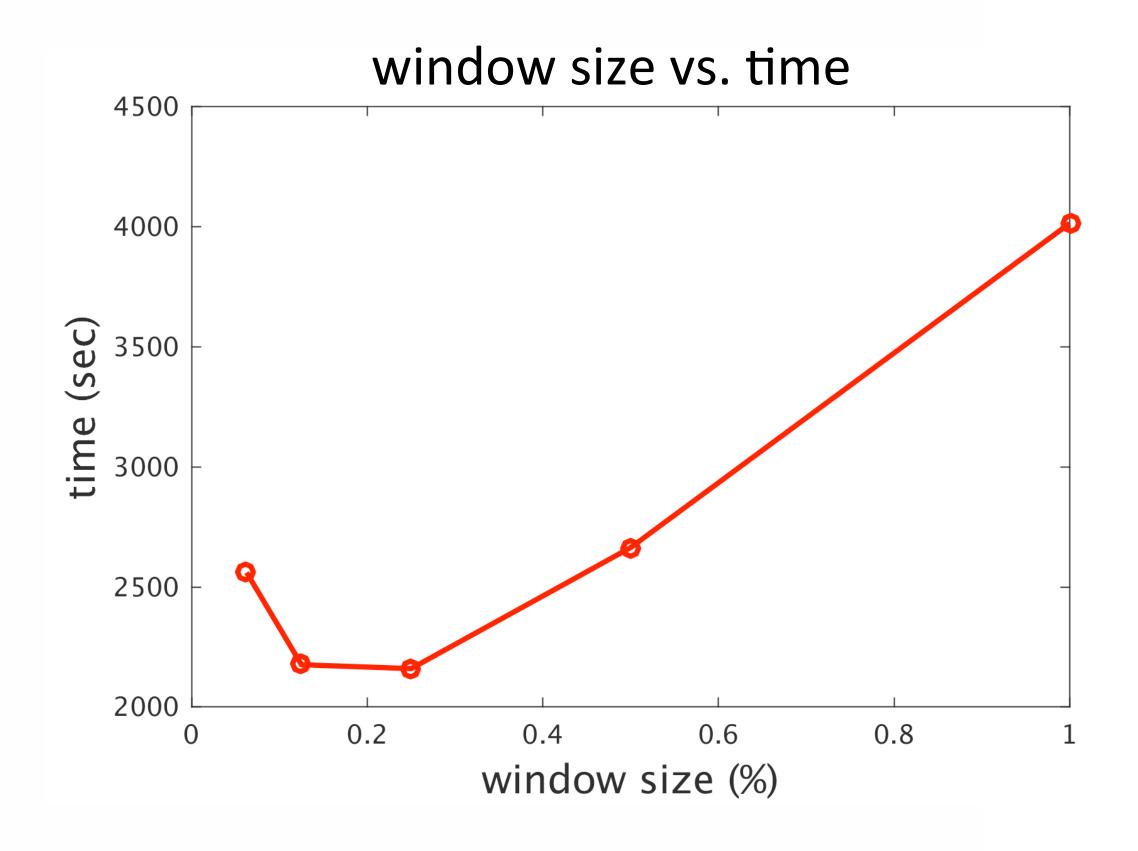
Missing 80% receivers

7 Alternations



To window or not to window?







Conclusions

- Significant improvement in computation time
- Equivalent SNR output
- Optimized communication time between workers
- Parameter free



R. Kumar, O. Lopez, D. Davis, A. Aravkin and F. Herrmann. "Beating-Level Set Methods for 5D Seismic Data Interpolation: A Primal Dual Alternating Approach"

RCAM

(Residual Constrained Alternating Minimization)

Distributed implementation to penalize norm with noise constraint

$$\mathbf{L}^{t} = \underset{\mathbf{L} \in \mathbb{C}^{n \times r}}{\min} \|\mathbf{L}\|_{F}^{2} \quad \text{s.t.} \quad \|P_{\Omega}(\mathbf{L}(\mathbf{R}^{t})^{*}) - \mathbf{b}\|_{F} \leq \eta$$

$$\mathbf{R}^{t+1} = \underset{\mathbf{R} \in \mathbb{C}^{m \times r}}{\min} \|\mathbf{R}\|_F^2 \quad \text{s.t.} \quad \|P_{\Omega}(\mathbf{L}^t \mathbf{R}^*) - \mathbf{b}\|_F \le \eta$$

- Avoid overfitting noise
- Robust (e.g., not sensitive to overshooting rank)
- time(ALS) < time(RCAM) << time(SPG-LR)</p>



Future Work

- \blacktriangleright Design for other measurement operators, \mathcal{A} .
 - incorporate "off-the-grid" measurements
 - source separation



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Software release available https://github.com/SINBADconsortium/RCAM.jl