## Memory AMP and Recent Results on Its Implementation

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  - Problem Formulation
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  - MAMP and GD-MAMP
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    m H}$
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#### Problem Formulation

System model:

$$\Gamma: \quad \boldsymbol{y} = \boldsymbol{A}\boldsymbol{x} + \boldsymbol{n},$$
  
 $\Phi: \quad x_i \sim P_X(x), \ \forall i.$ 

where  $\mathbf{A} \in \mathbb{C}^{M \times N}$ ,  $\mathbf{n} \sim \mathcal{CN}(0, \sigma^2 \mathbf{I}_M)$ , and  $\mathbf{A}, \mathbf{y}, \sigma^2, P_X(\cdot)$  are known.

- Assumptions:
  - (1)  $M, N \to \infty$  with fixed  $\delta = M/N$ .
  - (2) A is right-unitarily-invariant.
  - (3) x is IID. For convenience,  $\mathrm{E}\{x\}=\mathbf{0}$  and  $\frac{1}{N}\mathrm{E}\{\|x\|^2\}=1$ .
- ▶ Goal: Given  $\{y, A, \Gamma, \Phi\}$ , find an MMSE estimate of x:

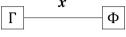
$$MSE \rightarrow \text{mmse}\{\boldsymbol{x}|\boldsymbol{y},\boldsymbol{A},\Gamma,\Phi\}$$

For non-Gaussian x, without the assumptions of  $M, N \to \infty$  and A, finding the optimal solution is generally NP-hard.

# Approximate Message Passing (AMP)

► AMP-type algorithms:

$$\begin{aligned} & \text{linear estimator (LE)}: \quad \boldsymbol{r}_{t} = \gamma_{t}\left(\boldsymbol{x}_{t}\right), \\ & \text{non-linear esitmator (NLE)}: \quad \boldsymbol{x}_{t+1} = \phi_{t}\left(\boldsymbol{r}_{t}\right). \end{aligned}$$



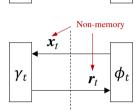
AMP:

$$ext{LE}: \qquad oldsymbol{r}_t = oldsymbol{x}_t + oldsymbol{A}^{ ext{H}}(oldsymbol{y} - oldsymbol{A}oldsymbol{x}_t) + oldsymbol{r}_t^{ ext{Onsager}},$$

 $NLE: \quad \boldsymbol{x}_{t+1} = \phi(\boldsymbol{r}_t) = E\{\boldsymbol{x}|\boldsymbol{r}_t\},\$ 

where 
$$\boldsymbol{r}_t^{\text{Onsager}} = \beta \langle \phi'(\boldsymbol{r}_{t-1}) \rangle (\boldsymbol{r}_{t-1} - \boldsymbol{x}_{t-1}).$$

- √ Bayes optimal
- √ Low-complexity
- imes IID  $oldsymbol{A}$  is required
- □ D. L. Donoho, A. Maleki, and A. Montanari, "Message-passing algorithms for compressed sensing," in *Proc. Nat. Acad. Sci.*, 2009.



## Orthogonal/Vector AMP (OAMP/VAMP)

► OAMP/VAMP:

LE: 
$$\boldsymbol{r}_t = \boldsymbol{x}_t + \frac{1}{\epsilon_t^{\gamma}} \boldsymbol{A}^{\mathrm{H}} (\rho_t \boldsymbol{I} + \boldsymbol{A} \boldsymbol{A}^{\mathrm{H}})^{-1} (\boldsymbol{y} - \boldsymbol{A} \boldsymbol{x}_t),$$
  
NLE:  $\boldsymbol{x}_{t+1} = \frac{1}{\epsilon_{t+1}^{\phi}} [\hat{\phi}_t(\boldsymbol{r}_t) + (1 - \epsilon_{t+1}^{\phi}) \boldsymbol{r}_t],$ 

where  $\epsilon_t^{\gamma}$  and  $\epsilon_t^{\phi}$  are orthogonal parameters.

- √ Bayes optimal (replica)
- $\checkmark$  Unitarily-invariant A
- X High-complexity
- ☐ J. Ma and L. Ping, "Orthogonal AMP," IEEE Access, 2017.
- S. Rangan, P. Schniter, and A. Fletcher, "Vector approximate message passing," *IEEE Trans. Inf. Theory*, 2019.

## Convolutional AMP (CAMP)

► CAMP:

$$egin{aligned} ext{LE}: & m{r}_t = m{x}_t + m{A}^H(m{y} - m{A}m{x}_t) + m{r}_t^{ ext{Onsager}}, \ ext{NLE}: & m{x}_{t+1} = \phi(m{r}_t) = ext{E}\{m{x}|m{r}_t\}, \end{aligned}$$

where 
$$\boldsymbol{r}_t^{\mathrm{Onsager}} = \sum_{\tau=0}^{t-1} \left[ \prod_{t'=\tau}^{t-1} \langle \phi'(\boldsymbol{r}_{t'}) \rangle \right] (\theta_{t-\tau} \boldsymbol{A}^{\mathrm{T}} \boldsymbol{A} - g_{t-\tau}) (\boldsymbol{r}_{\tau} - \boldsymbol{x}_{\tau}).$$

- √ Bayes optimal (replica), if converges
- $\checkmark$  Unitarily-invariant A
- √ Low-complexity
- $\times$  Fails to converge for A with high condition numbers
- ☐ K. Takeuchi, "Bayes-optimal convolutional AMP," *IEEE Trans. Inf. Theory*, 2021.

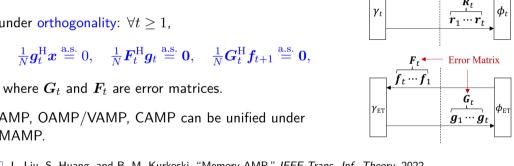
## Memory AMP

► Memory AMP (MAMP):

LE: 
$$r_t = \gamma_t(\boldsymbol{X}_t) = \boldsymbol{Q}_t \boldsymbol{y} + \sum_{i=1}^t \boldsymbol{P}_{t,i} \boldsymbol{x}_i,$$
  
NLE:  $\boldsymbol{x}_{t+1} = \phi_t(\boldsymbol{R}_t),$ 

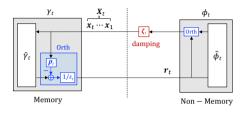
under orthogonality:  $\forall t > 1$ .

► AMP, OAMP/VAMP, CAMP can be unified under MAMP.



L. Liu, S. Huang, and B. M. Kurkoski, "Memory AMP," IEEE Trans. Inf. Theory, 2022.

## Bayes-Optimal MAMP (BO-MAMP) — Principle



- lacktriangle The LE has memory and consists of a local estimator  $\hat{\gamma}_t$  and orthogonalization.
  - 1.  $\hat{\gamma}_t$  approaches  $(\rho_t \boldsymbol{I} + \boldsymbol{A} \boldsymbol{A}^{\mathrm{H}})^{-1} (\boldsymbol{y} \boldsymbol{A} \boldsymbol{x}_t)$  in OAMP/VAMP.
  - 2. The error of  $r_t$  is orthogonal to x and the errors of  $x_1, \dots, x_t$ .
- ▶ Damping  $\zeta$  is added.  $\zeta$  is analytically optimized to guarantee convergence. Also improves convergence speed.
- ► NLE: Same as OAMP/VAMP and has no memory.

#### MAMP with Gradient Descent

#### GD-MAMP:

LE: 
$$\boldsymbol{u}_t = \theta_t \boldsymbol{B} \boldsymbol{u}_{t-1} + \xi_t (\boldsymbol{y} - \boldsymbol{A} \boldsymbol{x}_t),$$

$$\boldsymbol{r}_t = \gamma_t (\boldsymbol{X}_t) = \frac{1}{\varepsilon_t^{\gamma}} (\boldsymbol{A}^{\mathrm{H}} \boldsymbol{u}_t + \sum_{i=1}^t p_{t,i} \boldsymbol{x}_i),$$
NLE:  $\boldsymbol{x}_{t+1} = [\boldsymbol{x}_1 \cdots \boldsymbol{x}_t \ \phi_t(\boldsymbol{r}_t)] \cdot \boldsymbol{\zeta}_{t+1},$ 

- $lackbox{lack} oldsymbol{u}_t$  is an estimate of  $(
  ho_t oldsymbol{I} + oldsymbol{A} oldsymbol{A}^{\mathrm{H}})^{-1} (oldsymbol{y} oldsymbol{A} oldsymbol{x}_t)$ .
- The parameters  $\theta_t$  and  $\xi_t$  are optimized,  $p_{t,i}$  and  $\varepsilon_t^{\gamma}$  are chosen to ensure orthogonality. Computation not shown, but we'll more talk about these later.
- $\triangleright \zeta_{t+1}$  is the optimized damping vector
- $ightharpoonup \phi_t(\cdot)$  is the same as that in OAMP/VAMP,

Gradient descent (GD) is used to approximate  $\frac{\xi_t}{\theta_t}(\rho_t \boldsymbol{I} + \boldsymbol{A} \boldsymbol{A}^{\mathrm{H}})^{-1}(\boldsymbol{y} - \boldsymbol{A} \boldsymbol{x}_t)$  by  $\boldsymbol{u}_t$ .

## Why Memory? Intuition 1: Gradient Descent Avoids Matrix Inverse

Want to eliminate matrix inverse in OAMP:  $\frac{\xi_t}{\theta_t}(\rho_t \boldsymbol{I} + \boldsymbol{A}\boldsymbol{A}^{\mathrm{T}})^{-1}(\boldsymbol{y} - \boldsymbol{A}\boldsymbol{x}_t)$ . Solve  $\boldsymbol{W}\boldsymbol{u} = \boldsymbol{b}$  without finding  $\boldsymbol{W}^{-1}$ .

**Intuition 1** Solving  $oldsymbol{u} = oldsymbol{W}^{-1}oldsymbol{b}$  is equivalent to

$$rg \min f(\boldsymbol{u}) = \frac{1}{2} \boldsymbol{u}^{\mathrm{T}} \boldsymbol{W} \boldsymbol{u} - \boldsymbol{b}^{\mathrm{T}} \boldsymbol{u}$$

when  $oldsymbol{W}$  is positive definite. Find solution using gradient descent with step:

$$u_i = u_{i-1} - \alpha \nabla f(u_{i-1})$$
  
=  $u_{i-1} + \alpha (b - Wu_{i-1}).$ 

Then  $u_i$  approaches the correct value u.

Choosing  $\alpha = \frac{2}{\lambda_{\max} + \lambda_{\min}}$  is close to optimal. Shown for real-valued case.

## ChatGPT: Is it possible to find a matrix inverse using gradient descent?



While there might be unconventional methods or iterative approaches that can approximate a matrix inverse, gradient descent is not the typical choice for this particular problem. If you need to find the inverse of a matrix, it's recommended to use established linear algebra techniques for accuracy and efficiency.

## Why Memory? Intuition 2: Neumann Series for Matrix Inverse

Let  $ho(m{C})$  denote the spectral radius of  $m{C}$ . If  $ho(m{C}) < 1$ , then

$$(\boldsymbol{I} - \boldsymbol{C})^{-1} = \sum_{i=0}^{\infty} \boldsymbol{C}^i.$$

Choose C = I - W. When  $\rho(C) \ge 1$ , let  $C' = I - \theta(I - C)$ , where  $\theta$  ensures  $\rho(C') < 1$ . With  $u_0 = 0$ :

$$u_i = C'u_{i-1} + \theta b.$$

Then  $u_i$  approaches  $u = W^{-1}b$ . This iteration is identical to gradient descent.

To accelerate convergence, we can minimize  $\rho(C')$  by:

$$\theta = \frac{2}{\lambda_1 + \lambda_2},$$

where  $\lambda_1$  and  $\lambda_2$  denote the maximum and minimum eigenvalues of I-C.

## Overview of AMP-Type Algorithms

Table 1: Overview of AMP-Type Algorithms

Algorithm	Matrix $m{A}$	Convergence	Time complexity	Optimality
AMP	IID	Converges	Low: $\mathcal{O}(MN)$	Bayes-optimal
OAMP/VAMP	Right unitarily invariant	Converges	High: $\mathcal{O}(M^2N)$	Bayes-optimal
CAMP	Right unitarily invariant	Diverges in high condition numbers	Low: $\mathcal{O}(MN)$	Bayes-optimal
GD-MAMP	Right unitarily invariant	Converges	Low: $\mathcal{O}(MN)$	Bayes-optimal

- D. L. Donoho, A. Maleki, and A. Montanari, "Message-passing algorithms for compressed sensing," in Proc. Nat. Acad. Sci., 2009.
- J. Ma and L. Ping, "Orthogonal AMP," IEEE Access, 2017.
- S. Rangan, P. Schniter, and A. Fletcher, "Vector approximate message passing," *IEEE Trans. Inf. Theory*, 2019.
- ☐ K. Takeuchi, "Bayes-optimal convolutional AMP," *IEEE Trans. Inf. Theory*, 2021.
- L. Liu, S. Huang, and B. M. Kurkoski, "Memory AMP," *IEEE Trans. Inf. Theory*, 2022.

## Comparison of AMP-Type Algorithms

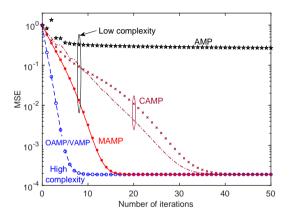


Figure 1:  $M=2^{13}, N=2^{14}, \kappa(\boldsymbol{A})=10, \mathrm{SNR}=30\mathrm{dB}$ 

- ► AMP: low complexity poor MSE
- OAMP/VAMP: fastest convergence high complexity
- ► CAMP: low complexity slow convergence incorrect state evolution
- ► GD-MAMP: low complexity fast convergence correct state evolution

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#### Overflow Problem in GD-MAMP

- ▶ Representation of floating point numbers from the IEEE 754 technical standard:
  - binary32 (single precision):  $\pm 1.18 \times 10^{-38} \sim \pm 3.4 \times 10^{38}$
  - binary64 (double precision):  $\pm 2.23 \times 10^{-308} \sim \pm 1.8 \times 10^{308}$

Double-precision is widely used, including in Matlab and Python

- The dreaded NaN will appear for values that are too large.
- Overflow problem: In GD-MAMP, some intermediate variables may increase exponentially, and exceed the maximum value of double precision.

## Which intermediate variables cause overflow in GD-MAMP?

- In iteration t, the parameters (1)  $\xi_t$  (2)  $p_{t,i}$  and (3)  $v_{t,t}^{\gamma}$  (variance of  $r_t$ ) require  $w_t$ .
- lacksquare  $\lambda^\dagger = (\lambda_{\max} + \lambda_{\min})/2$  and  $m{B} = \lambda^\dagger m{I} m{A} m{A}^{\mathrm{H}}$ , where  $\lambda_{\max}, \lambda_{\min}$  denotes the maximum and minimum eigenvalues of  $m{A} m{A}^{\mathrm{H}}$ .

$$b_k \equiv \frac{1}{N} \operatorname{tr} \{ \boldsymbol{B}^k \},$$
  

$$w_k \equiv \frac{1}{N} \operatorname{tr} \{ \boldsymbol{A}^{\mathrm{H}} \boldsymbol{B}^k \boldsymbol{A} \} = \lambda^{\dagger} b_k - b_{k+1}.$$

- ▶  $b_k$  can be computed if the eigenvalues of  $AA^H$  are known. Otherwise, there are simple methods to approximate  $\lambda_{\max}, \lambda_{\min}$  and  $b_k$ .
- ▶ If  $\lambda_{\max} > \lambda_{\min} + 2$ , the spectral radius  $\rho(B) > 1$ ,  $b_{2k}$  increases exponentially.
- **Even**  $w_k$  may increase exponentially.

## Overflow of $w_k$

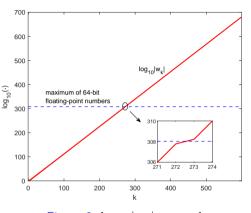


Figure 2:  $\log_{10}|w_k|$  versus k

- $\blacktriangleright w_1, \cdots, w_{2T}$  are required, where T is the maximum number of iterations.
- ▶ As shown in Figure 2,  $|w_k|$  increases exponentially as k increases.

#### Overflow Problem in GD-MAMP

▶ To compute  $\xi_t$ ,  $p_{t,i}$  and  $v_{t,t}^{\gamma}$ , we need  $w_k$ :

$$b_k \equiv \frac{1}{N} \operatorname{tr} \{ \boldsymbol{B}^k \},$$
  

$$w_k \equiv \frac{1}{N} \operatorname{tr} \{ \boldsymbol{A}^{\mathrm{H}} \boldsymbol{B}^k \boldsymbol{A} \} = \lambda^{\dagger} b_k - b_{k+1}.$$

- ▶ While  $w_k$  increases exponentially, it always appears in the product  $\vartheta w_k$ , which is bounded (i.e.  $\vartheta$  is small).
- ▶ For any  $\vartheta \in \mathbb{R} \setminus \{0\}$ , the following holds:

$$\vartheta w_k = \frac{\operatorname{sgn}(\vartheta)}{N} \mathbf{1}^{\mathrm{T}} \big[ (\lambda^{\dagger} \mathbf{1} - \lambda_B) \circ s_{\lambda}^{\circ k} \circ e^{\circ \log |\vartheta| \mathbf{1} + k \lambda_B^{\log}} \big],$$

where  $\lambda_B$  denotes the eigenvalues of B,  $s_\lambda \equiv \mathrm{sgn}(\lambda_B)$ ,  $\lambda_B^{\log} \equiv \log^\circ |\lambda_B|$  and  $\circ$  is component-wise operation.

# Overflow-Avoiding GD-MAMP with Eigenvalues of ${m A}{m A}^{ m H}$

- Define

$$\chi_k \equiv \theta_0^k w_k,$$

where  $\theta_0 = (\lambda^{\dagger} + \sigma^2)^{-1} > 0$ . We pre-compute  $\chi_1, \dots, \chi_{2T-1}$  before the iterations. Computing  $\vartheta w_k$  can be reduced to a scalar operation:

$$\vartheta w_k = \operatorname{sgn}(\vartheta) e^{\log |\alpha| - k \log \vartheta} \chi_k.$$

Pre-computing  $\chi_1, \dots, \chi_{2T-1}$  costs  $\mathcal{O}(MT)$ , and computing terms involving  $w_k$  in iterations costs  $\mathcal{O}(T^3)$ . The overall complexity is reduced to  $\mathcal{O}(MT + T^3)$ .

# Overflow-Avoiding GD-MAMP with Eigenvalues of ${m A}{m A}^{ m H}$

Theorem (2) For any  $k \ge 0$ ,

$$|\chi_k| \le \delta(\lambda^{\dagger} + \theta_0^{-1}),$$

where  $\delta = M/N$  .

▶ Theorem 2 shows that  $\chi_k$  is bounded. In other words, computing  $\chi_k$  has no risks of overflow.

## Overflow-Avoiding GD-MAMP without Eigenvalues of $m{A}m{A}^{ ext{H}}$

In large-scale systems, computation of eigenvalues of  ${m A}{m A}^{
m H}$  may be impractical.

- ▶ A method to estimate the maximum and minimum eigenvalue  $\lambda_{max}$  and  $\lambda_{min}$  was given in [LHK22].
- ▶ For  $k \ge 0$ ,  $\chi_k$  can be estimated by

$$\chi_k = \bar{\boldsymbol{h}}_i^{\mathrm{H}} \bar{\boldsymbol{h}}_{k-i}$$

where  $i=\lceil k/2 \rceil$  and  $\bar{\boldsymbol{h}}_i$  is given by a recursion

$$\bar{\boldsymbol{h}}_i = \theta_0 (\lambda^{\dagger} \boldsymbol{I} - \boldsymbol{A} \boldsymbol{A}^{\mathrm{H}}) \bar{\boldsymbol{h}}_{i-1}$$

with  $ar{m{h}}_0 = m{A}m{h}_0$ ,  $m{h}_0 \sim \mathcal{N}(m{0}, \frac{1}{N}m{I}_N)$ .

#### Simulation Results

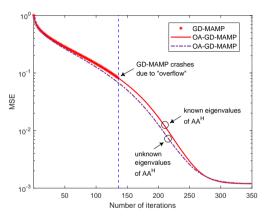


Figure 3:  $M=2^{13}, N=2^{14}, \kappa(\mathbf{A})=1000, \mathrm{SNR}=35\mathrm{dB}$ 

- ▶ When t > 136, the unmodified GD-MAMP does not reach the fixed point since  $b_{137}$  and  $w_{137}$  overflows.
- lacktriangle Both OA-GD-MAMP with and without eigenvalues of  $AA^{
  m H}$  work properly.

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## Complexity Analysis of GD-MAMP

Let T be the number of iterations. The main complexity of GD-MAMP is  $\mathcal{O}(MNT)$ , dominated by the number of matrix-vector products, each with  $\mathcal{O}(MN)$ 

LE: 
$$\boldsymbol{u}_t = \theta_t \boldsymbol{B} \boldsymbol{u}_{t-1} + \xi_t (\boldsymbol{y} - \boldsymbol{A} \boldsymbol{x}_t),$$

$$\boldsymbol{r}_t = \frac{1}{\varepsilon_t^{\gamma}} (\boldsymbol{A}^{\mathrm{H}} \boldsymbol{u}_t + \sum_{i=1}^t p_{t,i} \boldsymbol{x}_i),$$
NLE:  $\boldsymbol{x}_{t+1} = [\boldsymbol{x}_1 \cdots \boldsymbol{x}_t \ \phi_t(\boldsymbol{r}_t)] \cdot \zeta_{t+1}.$ 

GD-MAMP requires 4 matrix-vector products per iteration?

- ▶ Computing  $A\phi_{t-1}(r_{t-1})$  to estimate  $v_{t,1}^{\phi}, \dots, v_{t,t}^{\phi}$  requires one (hidden in  $\zeta_t$ ).
- lacksquare Computing  $m{B}m{u}_{t-1}=(\lambda^\daggerm{I}-m{A}m{A}^\mathrm{H})m{u}_{t-1}$  requires two.
- ightharpoonup Computing  $m{A}^{\mathrm{H}}m{u}_t$  requires one.

Easily eliminate one matrix-vector product:

lacktriangle Two of the products are  $m{A}^{
m H}m{u}_t$ , these only need to be computed once.

GD-MAMP requires 3 matrix-vector products per iteration!

# Complexity-Reduced GD-MAMP Using Approximate $\xi_t$

Finding the damping vector  $\zeta_t$  nominally requires one matrix-vector product:

- 1. Compute  $z_t = y A\phi_{t-1}(r_{t-1})$ .
- 2. Estimate  $v_{t,1}^{\phi}, \dots, v_{t,t}^{\phi}$  by using  $z_t$ , where  $v_{t,i}^{\phi}$  denotes the covariance of  $\phi_{t-1}(r_{t-1})$  and  $x_i$  for i < t.
- 3. Compute  $\zeta_t$  from the covariance matrix  $m{V}^\phi_t$  of  $m{x}_1,\cdots,m{x}_{t-1}$  and  $\phi_{t-1}(m{r}_{t-1})$ .

#### To remove the above matrix-vector product:

- (1) We move the damping from the NLE to the LE (details omitted).
- (2)  $\xi_t$  nominally depends on  $z_t$  and  $V_t^{\phi}$ . But we found that approximating  $\xi_t$  gave little to no performance loss:

$$\tilde{\xi}_t = 1/(v_{t,t}^{\phi} + \sigma^2).$$

where  $v_{t,t}^{\phi}$  is the variance of  $x_t$ , given as that in OAMP/VAMP.

▶ The resulting complexity-reduced GD-MAMP requires only 2 matrix-vector products per iteration.

#### Simulation Results

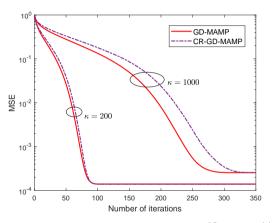


Figure 4: MSE versus number of iterations,  $M=2^{13}, N=2^{14}, \mathrm{SNR}=35\mathrm{dB}$ 

► CR-GD-MAMP requires a few more iterations to converge.

#### Simulation Results

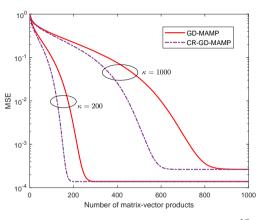


Figure 5: MSE versus number of matrix-vector products,  $M=2^{13}, N=2^{14}, \mathrm{SNR}=35\mathrm{dB}$ 

► CR-GD-MAMP achieves the same MSE as GD-MAMP while requiring about 2/3 matrix-vector products.

#### Conclusion

GD-MAMP is a memory AMP algorithm that:

- Converges for unitarily-invariant matrices
- is low complexity, avoiding matrix inverse
- converges for A of high condition number

overcoming the weakness of AMP, OAMP/VAMP and CAMP.

- (1) To solve the overflow problem, we propose OA-GD-MAMP:
  - ightharpoonup With known eigenvalues of  $AA^{\mathrm{H}}$ , OA-GD-MAMP is equivalent to GD-MAMP.
  - ▶ Otherwise, OA-GD-MAMP can achieve nearly the same performance.
- (2) GD-MAMP requires three matrix-vector products per iteration. To reduce it:
  - We propose CR-GD-MAMP as a variant of GD-MAMP. It requires only two matrix-vector products per iteration with almost the same convergence speed.