# Partially coherent ptychography

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# 1 Models

# 1.1 Coherent model

$$f_j = \left| \mathcal{F} \left( \mathcal{S}_j u \circ \omega \right) \right|^2 \tag{1.1}$$

In a discrete setting,  $u \in \mathbb{C}^n$  is a 2D image with  $\sqrt{n} \times \sqrt{n}$  pixels,  $\omega \in \mathbb{C}^{\bar{m}}$  is a localized 2D probe with  $\sqrt{\bar{m}} \times \sqrt{\bar{m}}$  pixels.

 $f_j \in \mathbb{R}^{\bar{m}}_+(\forall 0 \leq j \leq N-1)$  is a stack of phaseless measurements. Here  $|\cdot|$  represents the element-wise absolute value of a vector, o denotes the elementwise multiplication, and  $\mathcal{F}$  denotes the normalized 2 dimensional discrete Fourier transform. Each  $\mathcal{S}_j \in \mathbb{R}^{\bar{m} \times n}$  is a binary matrix that crops a region j of size  $\bar{m}$  from the image u.

In practice, as the probe is almost never completely known, one has to solve a blind ptychographic phase retrieval (BP-PR) problem:

To find 
$$\omega \in \mathbb{C}^{\bar{m}}$$
 and  $u \in \mathbb{C}^n$  s.t.  $|\mathcal{A}(\omega, u)| = \boldsymbol{a}$ ,

where bilinear operators  $\mathcal{A}: \mathbb{C}^{\bar{m}} \times \mathbb{C}^n \to \mathbb{C}^m$  and  $\mathcal{A}_j: \mathbb{C}^{\bar{m}} \times \mathbb{C}^n \to \mathbb{C}^{\bar{m}} \forall 0 \leq j \leq N-1$  are denoted as follows:

$$\begin{split} \mathcal{A}(\omega, u) &:= \left(\mathcal{A}_0^T(\omega, u), \mathcal{A}_1^T(\omega, u), \dots, \mathcal{A}_{N-1}^T(\omega, u)\right)^T, \mathcal{A}_j(\omega, u) := \mathcal{F}\left(\omega \circ \mathcal{S}_j u\right) \\ \text{and } \boldsymbol{a} &:= \left(\boldsymbol{a}_0^T, \boldsymbol{a}_1^T, \dots, \boldsymbol{a}_{N-1}^T\right)^T \in \mathbb{R}_+^m. \end{split}$$

# 1.2 Specific partially coherent model

## $1.2.1 \quad Model1^{[1]}$

Coherence and vibrations kernels can be combined into one, such that partially coherent ptychography imaging with coherence kernel function  $\kappa$  in a continuous setting:

$$f_{pc,j}(q) = \int |\mathcal{F}_{x \to q} \left( \mathcal{S}_j u(x) \omega(x - y) \right)|^2 \kappa(y) dy$$
 (1.2)

where  $f_{pc}$  is the measured partial coherent intensity and  $\mathcal{F}_{x\to q}$  is the normalized Fourier transform.  $\kappa$  is a function that spikes at 0 like guassians. Setting  $\kappa$  to the Dirac delta function reduces it to the coherent model (1.1).

The partially coherent intensity in a discrete setting is generated as

$$f_{pc,j} = \sum_{i} \kappa_{i} \left| \mathcal{F} \left( \mathcal{S}_{j} u \circ \left( \mathcal{T}_{i} \omega \right) \right) \right|^{2}$$
(1.3)

with translation operator  $\mathcal{T}_i$ , discrete Gaussian weights  $\{\kappa_i\}$ , and periodical boundary condition for the probe.

Generally speaking, solving (1.3) is a non-linear ill-posed problem with an unknown kernel  $\kappa$ , unknown phobe  $\omega$ , and unknown target image u.

#### $1.2.2 \quad Model2^{[9]}$

Another simpler model mentioned here is:

$$f_{nc} = f * \kappa \tag{1.4}$$

where  $f_{pc}$  is the measured partial coherent intensity, f is the coherent intensity in (1.1), \* denotes the convolution operator, and  $\kappa$  is the unknown kernel function (Fourier transform of the complex coherence function).

We remark that (1.3) is quite different from (1.4), since (1.3) illustrates the effects of blurring of images with respect to the probe, while (1.4) can be interpreted as blurring or binning multiple pixels at the detector.

## 1.3 General partially coherent model

This part explains the partially coherent model proposed by physicists. [3]. It is a blind ptychography model based on quantum state tomography<sup>1</sup>. Phobe w is assumed to be in a mixed state to represent a partially coherent effect.

#### 1.3.1 Decompositon model

Find u, r othogonal  $w_k$  s.t.

$$f_{pc,j} = \sum_{k=1}^{r} \left| \mathcal{F} \left( \mathcal{S}_{j} u \circ (\omega_{k}) \right) \right|^{2} \left( 0 \le j \le N - 1 \right)$$

$$\tag{1.5}$$

Denote  $O_j \in C^{\bar{m} \times \bar{m}}$  as a (diagonal) matrix to represent linear transform to w, s.t.  $S_j u \circ \omega = O_j w$ . Denote  $f_q^* \in C^{1 \times \bar{m}}$  as a row vector constructed from Fourier transform  $\mathcal{F}$ , to represent projection on frepuency element. Construct measurement matrix  $\mathcal{I}_{j\mathbf{q}} = O_j^* f_q f_q^* O_j$  and density matrix  $\rho$ , we get another form(actually a natural one in quantum state tomography) of the model:

Find 
$$u, \rho, s.t.$$
  
 $f_{pc,j}(q) = Tr(\mathcal{I}_{j\mathbf{q}}\rho)(0 \le j \le N-1)$  (1.6)  
 $\rho$  is positive semi-definite, with rank  $\le r$ 

Next, we will explain the derivation of this form.

Simple calculation process:

$$f_{pc,j}(q) = |f_q^* O_j w|^2 = (f_q^* O_j w)^* (f_q^* O_j w) = w^* (O_j^* f_q f_q^* O_j) w$$
$$= Tr[w^* (O_j^* f_q f_q^* O_j) w] = Tr[(O_j^* f_q f_q^* O_j) (ww^*)]$$
$$= Tr(\mathcal{I}_{j\mathbf{q}}\rho)$$

It is a bit like the process of phase-lift.

When w is in pure state(a vector in Hilbert space),  $\rho = w^*w$  is a rankone matirx. In partially coherent case, **we use mixed state to model** w. Fow example, with probability 0.5 in state  $\psi_1$  and 0.5 in  $\psi_2$  ( $\psi_1$  and  $\psi_2$  are not necessarily orthogonal here). Now w can no longer be represented by a vector(ps.  $w \neq p_1\psi_1 + p_2\psi_2$ , the latter is still a determined pure state vector).

<sup>&</sup>lt;sup>1</sup>https://homepage.univie.ac.at/reinhold.bertlmann/pdfs/T2\_Skript\_Ch\_9corr.pdfTheorem 9.1. Many symbols in quantum mechanics are included here.

Instead, mixed state is represented by generalizing the density matrix to one with higher rank:

$$\rho = \sum_{k} p_k \psi_k \psi_k^*$$

Easy to find  $\rho$  is a positive semi-definite matrix, we can decompose  $\rho$  using spectral theorem, with  $r(\text{rank of }\rho)$  othogonal state  $w_k$ :

$$\rho = \sum_{k=1}^{r} w_k w_k^* \tag{1.7}$$

$$f_{pc,j}(q) = \operatorname{Tr} \mathcal{I}_{j\mathbf{q}} \rho = \operatorname{Tr} [\mathcal{I}_{j\mathbf{q}} \sum_{k=1}^{r} w_k w_k^*]$$

$$= \sum_{k=1}^{r} w_k^* \mathcal{I}_{j\mathbf{q}} w_k = \sum_{k=1}^{r} |f_q^* O_j w_k|^2$$

And that is exactly  $(1.5)f_{pc,j} = \sum_{k=1}^{r} |\mathcal{F}(\mathcal{S}_{j}u \circ (\omega_{k}))|^{2}$ .  $(f_{pc,j}(q) \text{ is a single value at frequency } q \text{ when } f_{pc,j} \text{ is the whole diffraction image })$ 

We can write it in another quadratic form:

$$f_{pc,j}(q) = \operatorname{Tr} \mathcal{I}_{j\mathbf{q}} \rho = \operatorname{Tr} [(O_j^* f_q f_q^* O_j) \rho] = \operatorname{Tr} [(O_j^* f_q)^* \rho (O_j^* f_q)] = (O_j^* f_q)^* \rho (O_j^* f_q)$$

$$= g_q^* \rho g_q = \sum_{x_1} \sum_{x_2} \overline{g_q(x_1)} \rho(x_1, x_2) g_q(x_2)$$
(1.8)

where 
$$g_q = O_j^* f_q = \overline{S_j u} \circ f_q$$
,  $\overline{g_q} = S_j u \circ \overline{f_q}$ 

Then it is a discrete version of the model in [9].

#### 1.3.2 The relation between models

In this section we explain how the general model connects with specific models (1.3) and (1.4).

First we consider (1.3). We put the  $\kappa_i$  inside:

$$f_{pc,j} = \sum_{i} \left| \mathcal{F} \left( \mathcal{S}_{j} u \circ \left( \sqrt{\kappa_{i}} \mathcal{T}_{i} \omega \right) \right) \right|^{2} = \sum_{i} \left| \mathcal{F} \left( \mathcal{S}_{j} u \circ \left( \hat{\omega}_{i} \right) \right) \right|^{2}$$
(1.9)

Multiple modes  $\hat{w}_i$  are produced by shifted w. Then we can construct density matrix and use truncated SVD to get a low-rank approximation.

$$\rho = \sum_{i} \hat{w}_i \hat{w}_i^* \approx \sum_{k=1}^r w_k w_k^*$$

As for (1.4), we introduce the definition for coherence function:

$$\gamma(x_1, x_2) = \frac{\rho(x_1, x_2)}{w(x_1)\overline{w(x_2)}}$$

In other word:

$$\gamma = \rho \cdot / (ww^*), \rho = \gamma \circ (ww^*) \tag{1.10}$$

where ./ means pairwise division.

In case we assume that coherence function  $\gamma(\mathbf{x}_1, \mathbf{x}_2)$  only depends on the difference between the two points  $\mathbf{x}_1, \mathbf{x}_2$ , i.e.  $\gamma(\mathbf{x}_1, \mathbf{x}_2) = \gamma(0, \mathbf{x}_2 - \mathbf{x}_1)$ , we can write the far-field intensity as a convolution [9]:

 $f_{pc} = \kappa * f$ . If we know the point spread function  $\kappa(q)$ , then we get its inverse Fourier transform  $\gamma(0,x) = (\mathcal{F}^{-1}\kappa)(x)$ . And we get  $\gamma(x_1,x_2)$  based on the assumption above. If we also know the phobe w, then we can get  $\rho$ , again SVD helps us find the main modes  $w_k$  in model (1.5)

With the connection above, we can get the "standard mode decomposition" for simulation experiment which is a good reference.

# 1.4 Gradient Decomposition

Consider specific model 1.3.

$$f_{pc,j}(q) = \int \left| \mathcal{F}_{x \to q} \left( \mathcal{S}_j u(x) \omega(x - y) \right) \right|^2 \kappa(y) dy$$
 (1.11)

Remember that it can connect with the general model through (1.9). In coherent setting:

$$\rho_{cor}(x_1, x_2) = \omega(x_1)\overline{\omega(x_2)}, \rho_{cor, y}(x_1, x_2) = \omega(x_1 - y)\overline{\omega(x_2 - y)}$$

And in partially coherent case:

$$\rho(x_1, x_2) = \int \omega(x_1 - y) \overline{\omega(x_2 - y)} \kappa(y) dy \approx \sum_i \omega_i(x_1) \overline{\omega_i(x_2)}$$
 (1.12)

Consider the space of functions on  $\mathbb{R}^2$  equipped with inner product:

 $< f,g> = \int f(y)g(y)\kappa(y)dy$ . And we may generate an orthogonal sequence which depends on the characteristics of  $\kappa$ . Here we assume that  $\kappa$  is symmetric and separable (like some guassians). That is:

$$\int \kappa(y_1, y_2) y_1^n y_2^m dy = 0$$

for n is odd or m is odd.

Then we can get orthogonal polynomial function sequence:

$$f_n(y) = 1, y_1, y_2, y_1y_2, \dots$$

Next we want to decompose  $\omega(x-y)$  into the part related to the main mode  $\omega(x)$  and that about the shift y. Taylor expansion is used here:

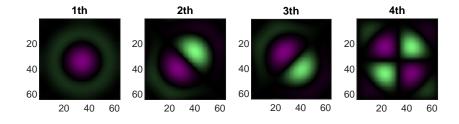
$$\omega(x_1 - y) \approx \omega(x_1) + \nabla_1 \omega(x_1) y_1 + \nabla_2 \omega(x_1) y_2 + \nabla_{12} \omega(x_1) y_1 y_2$$

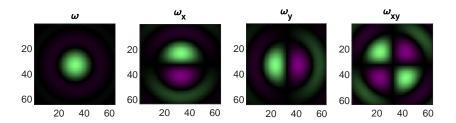
**Remark.** More items of expansion can be considered for a better approximation for  $\omega$ . However, the next terms like  $\nabla_{11}\omega(x_1)*y_1^2$  may not be orthogonal to the latter items  $\{1, y_1, y_2, y_1y_2\}$ , and process like Gram-Smith orthogonalization is needed, which makes coefficients in front of  $f_n(y)$  a linear combination of the gradients of  $\omega$ .

Then we have:

$$\begin{split} \rho(x_1,x_2) &= \int (\omega(x_1) + \nabla_1 \omega(x_1) y_1 + \nabla_2 \omega(x_1) y_2 + \nabla_{12} \omega(x_1) y_1 y_2) \\ \overline{(\omega(x_2) + \nabla_1 \omega(x_2) y_1 + \nabla_2 \omega(x_2) y_2 + \nabla_{12} \omega(x_2) y_1 y_2)} \kappa(y) dy & \xrightarrow{\underline{Orthogonalization}} \\ c_1 \omega(x_1) \overline{\omega(x_2)} + c_2 \nabla_1 \omega(x_1) \overline{\nabla_1 \omega(x_2)} + c_3 \nabla_2 \omega(x_1) \overline{\nabla_2 \omega(x_2)} + c_4 \nabla_{12} \omega(x_1) \overline{\nabla_{12} \omega(x_2)} \\ \text{where } c_n &= \int |f_n(y)|^2 \kappa(y) dy \end{split}$$

This mostly accomplish and since when the small anto of the mostly mode .......





# 2 ADMM-based numerical algorithm

The AP algorithm above can be rewritten into ADMM form, which is more stable and faster. We generalize the ADMM form in [8] to mixed states.

Now  $w \in \mathbb{C}^{(px \times py) \times r}$  is a phobe with r mixed states.  $u \in \mathbb{C}^{Nx \times Ny}$  is an image.  $f \in \mathbb{C}^{(px \times py) \times N}$  is the true(observed) diffraction image stacks. Let  $Y = \sqrt{f}$  be the amplitute of stacks.

An auxiliary variable  $z=\mathcal{A}(\omega,u)\in\mathbb{C}^{(px\times py)\times N\times r}$  is introduced.  $\mathcal{A}$  is an operator generating diffraction image stacks (N) frames for each phobe state) from image u and r different states  $w_k:=w(:,:,k)\in\mathbb{C}^{px\times py}$ . For multi-dimensional vectors, symbol: denotes the free dimensions, and we can fix some indexes to extract particular dimensions from original vectors.

Based on the general model 1.3, the problem is:

Find 
$$\omega, u \ s.t.$$
  
 $\bar{\mathcal{A}}(\omega, u) = Y$ 

Where  $\mathcal{A}: \mathbb{C}^{(px \times py) \times r} \times \mathbb{C}^{Nx \times Ny} \to {}^{(px \times py) \times N \times r}, \mathcal{A}_j: \mathbb{C}^{px \times py} \times \mathbb{C}^{Nx \times Ny} \to \mathbb{C}^{px \times py}, \ \bar{\mathcal{A}}_j: \mathbb{C}^{(px \times py) \times r} \times \mathbb{C}^{Nx \times Ny} \to \mathbb{R}^{(px \times py)}_+, \ \text{and} \ \bar{\mathcal{A}}: \mathbb{C}^{(px \times py) \times r} \times \mathbb{C}^{Nx \times Ny} \to \mathbb{R}^{(px \times py) \times r}_+ (\forall 0 \leq j \leq N-1) \ \text{are denoted as follows:}$ 

$$z(:,:,j,k) = \mathcal{A}_{j}(\omega_{k},u) := \mathcal{F}(\omega_{k} \circ \mathcal{S}_{j}u) \in \mathbb{C}^{px \times py},$$

$$z(:,:,:,k) = \left(\mathcal{A}_{0}^{T}(\omega_{k},u), \mathcal{A}_{1}^{T}(\omega_{k},u), \dots, \mathcal{A}_{N-1}^{T}(\omega_{k},u)\right)^{T} \in \mathbb{C}^{(px \times py) \times N},$$

$$z(:,:,j,:) = \left(\mathcal{A}_{j}^{T}(\omega_{1},u), \mathcal{A}_{j}^{T}(\omega_{2},u), \dots, \mathcal{A}_{j}^{T}(\omega_{r},u)\right)^{T} \in \mathbb{C}^{(px \times py) \times r}.$$
,

$$\begin{split} \bar{\mathcal{A}}_{j}(\omega,u) &:= \sum_{k=1}^{r} |\mathcal{A}_{j}(\omega_{k},u)|^{2} \in \mathbb{R}_{+}^{px \times py}, \\ \bar{\mathcal{A}}(\omega,u) &:= \left(\bar{\mathcal{A}}_{0}^{T}(\omega,u), \bar{\mathcal{A}}_{1}^{T}(\omega,u), \dots, \bar{\mathcal{A}}_{N-1}^{T}(\omega,u)\right)^{T} \in \mathbb{R}_{+}^{(px \times py) \times N}, \\ \text{and } Y &= \left(\boldsymbol{a}_{0}^{T}, \boldsymbol{a}_{1}^{T}, \dots, \boldsymbol{a}_{N-1}^{T}\right)^{T} \in \mathbb{R}_{+}^{(px \times py) \times N}. \end{split}$$

Then  $\mathcal{G}(z) = ||\sqrt{\sum_{k=1}^r |z(:,:,k)|^2} - Y||^2$  measures the difference between values computed by my our model and the groundtruth.

Let  $\mathcal{X}_1$  and  $\mathcal{X}_2$  be the prior range for w and u. Let  $l = px \times py$ ,  $\mathcal{X}_3$  be the index function for orthogonal  $D\alpha \in \mathbb{C}^{l \times r}$  (Orthonormal  $D \in \mathbb{C}^{l \times r}$  s.t.  $D^*D = I$ , and amplitude factors  $\alpha \in \mathbb{R}^{r \times r}$ ).  $\Omega$  is a reformulation operator, and  $\Omega(D\alpha) := reshape(D\alpha, [px, py, r]) \in \mathbb{C}^{(px \times py) \times r}$ . And sometimes we simplify  $\Omega(D^k\alpha^k)$  to  $\Omega^k$  and  $\Omega(D\alpha)$  to  $\Omega$ .

Then we get the following:

$$\min_{\omega, u, z} \mathcal{G}(z) + \mathbb{I}_{\mathcal{X}_1}(\omega) + \mathbb{I}_{\mathcal{X}_2}(u) + \mathbb{I}_{\mathcal{X}_3}(D\alpha) 
s.t. \quad z - \mathcal{A}(\omega, u) = 0, \quad \Omega(D\alpha) - w = 0.$$
(2.1)

The corresponding augmented Lagrangian reads

$$\Upsilon_{\beta}(\omega, u, z, \Lambda) := \mathcal{G}(z) + \mathbb{I}_{\mathcal{X}_{1}}(\omega) + \mathbb{I}_{\mathcal{X}_{2}}(u) + \Re(\langle z - \mathcal{A}(\omega, u), \Lambda \rangle) + \frac{\beta}{2} \|z - \mathcal{A}(\omega, u)\|^{2} + \Re(\langle \Omega(D\alpha) - w, \Lambda_{2} \rangle) + \frac{\beta_{2}}{2} \|\Omega(D\alpha) - \omega\|^{2}$$

where  $\Lambda, \Lambda_2 \in \mathbb{C}^{(px \times py) \times N \times r}$  is the multiplier. Let  $\Lambda = \Lambda/\beta, \Lambda_2 = \Lambda_2/\beta_2$ , we can combine the  $\Re$  part and the augmented part to get:

$$\begin{split} \Upsilon_{\beta}(\omega,u,z,\Lambda) := & \mathcal{G}(z) + \mathbb{I}_{\mathcal{X}_{1}}(\omega) + \mathbb{I}_{\mathcal{X}_{2}}(u) + \mathbb{I}_{\mathcal{X}_{3}}(D\alpha) \\ & + \frac{\beta}{2} \|z - \mathcal{A}(\omega,u) + \Lambda\|^{2} - \frac{\beta}{2} ||\Lambda||^{2} + \frac{\beta_{2}}{2} \|\Omega(D\alpha) - w + \Lambda_{2}\|^{2} - \frac{\beta_{2}}{2} ||\Lambda_{2}||^{2} \end{split}$$

In ADMM, one seeks a saddle point of the following problem:

$$\max_{\Lambda,\Lambda_2} \min_{\omega,u,z,D,\alpha} \Upsilon_{\beta}(\omega,u,z,\Lambda,\Lambda_2,D,\alpha)$$

A natural scheme to solve the above saddle point problem is to split them, which consists of four-step iterations for the generalized ADMM (only the subproblems w.r.t.  $\omega$  or u have proximal terms), as follows:

Step 1: 
$$\omega^{k+1} = \arg\min_{\omega} \frac{\beta}{2} \|z - \mathcal{A}(\omega, u) + \Lambda\|^2 + \frac{\beta_2}{2} \|\omega - (\Omega(D^k \alpha^k) + \Lambda_2)\|^2 + \frac{\alpha_1}{2} \|\omega - \omega^k\|_{M_1^k}^2$$
,  
Step 2:  $u^{k+1} = \arg\min_{u} \frac{\beta}{2} \|z - \mathcal{A}(\omega, u) + \Lambda\|^2 + \frac{\alpha_2}{2} \|u - u^k\|_{M_2^k}^2$ ,  
Step 3:  $z^{k+1} = \arg\min_{z} \frac{\beta}{2} \|z - \mathcal{A}(\omega, u) + \Lambda\|^2$ ,  
Step 4:  $D^{k+1} = \arg\min_{D} \mathbb{I}_{\mathcal{X}_3}(D\alpha) + \frac{\beta_2}{2} \|\Omega(D^k \alpha^k) - \omega^{k+1}\|^2$   
Step 5:  $\alpha^{k+1} = \arg\min_{\alpha} \frac{\beta_2}{2} \|\Omega(D^{k+1} \alpha^k) - \omega^{k+1}\|^2$   
Step 6:  $\Lambda^{k+1} = \Lambda^k + (z^{k+1} - \mathcal{A}(\omega^{k+1}, u^{k+1}))$  (2.3)  
Step 7:  $\Lambda_2^{k+1} = \Lambda_2^k + (\Omega(D^{k+1} \alpha^{k+1}) - \omega^{k+1})$ 

For simplicity, we ignore the stable quadratic terms in Step1 and Step2 in the following analysis.

# 2.1 Subproblems w and u

w.r.t. the probe  $\omega$ :

$$\begin{split} & \omega^{k+1} = \arg\min_{\omega \in \mathcal{X}_1} \frac{1}{2} \left\| z^k + \Lambda^k - \mathcal{A} \left( \omega, u^k \right) \right\|^2 + \frac{\beta_2}{2} ||\omega - (\Omega^k + \Lambda_2^k)||^2 \\ & = \arg\min_{\omega \in \mathcal{X}_1} \frac{1}{2} \left\| \hat{z}^k - \mathcal{A} \left( \omega, u^k \right) \right\|^2 + \frac{\beta_2}{2} ||\omega - \hat{\Omega}^k||^2 \\ & = \arg\min_{\omega \in \mathcal{X}_1} \frac{1}{2} \sum_{j,i} \left\| \mathcal{F}^{-1} \hat{z}(:,:,j,i)^k - \omega(:,:,i) \circ \mathcal{S}_j u^k \right\|^2 + \frac{\beta_2}{2} \sum_i ||\omega(:,:,i) - \hat{\Omega}^k(:,:,i)||^2 \\ & \text{with } \hat{z}^k := z^k + \Lambda^k, \hat{\Omega}^k := \Omega^k + \Lambda_2^k \end{split}$$

The close form solution of Step 1 is given as(details are in the Appendix 6.1)

$$\omega^{k+1} = \operatorname{Proj}\left(\frac{\beta \sum_{j} (\mathcal{S}_{j} u^{k})^{*} \circ [(\mathcal{F}^{-1} \hat{z}^{k}) (:,:,j,:)] + \beta_{2} \hat{\Omega}^{k}}{\beta \sum_{j} |\mathcal{S}_{j} u^{k}|^{2} + \beta_{2}}; \mathcal{X}_{1}\right)$$
(2.5)

with the projection operator onto  $\mathcal{X}_1$  defined as  $\operatorname{Proj}(\omega; \mathcal{X}_1) := \mathcal{F}^{-1}(\mathcal{F}(\omega) \circ C_w)$ , where  $\mathcal{X}_1 = \{\omega : \mathcal{F}(\omega) \text{ supports on the index function } C_w\}$ .  $\mathcal{F}^{-1}$  acts on the first two dimensions of  $\hat{z}$  (i.e.  $\hat{z}_{j,i} := \hat{z}(:,:,j,i)$ ) and  $\omega$  (i.e.  $\omega_i := \omega(:,:,i)$ ).

Similarly we have:

$$u^{k+1} = \operatorname{Proj}\left(\frac{\sum_{j,i} \mathcal{S}_{j}^{T} \left(\left(\omega_{i}^{k+1}\right)^{*} \circ \mathcal{F}^{-1} \hat{z}_{j,i}^{k}\right)}{\sum_{j,i} \left(\mathcal{S}_{j}^{T} \left|\omega_{i}^{k+1}\right|^{2}\right)}; \mathcal{X}_{2}\right) . \tag{2.6}$$

Here  $S_i^T$  is an operator mapping its augment to target position j in image u.

#### 2.2 Subproblem z

$$\begin{split} z^{k+1} &= \arg\min_{z} \mathcal{G}(z) + \frac{\beta}{2} \left\| z - \mathcal{A} \left( \omega^{k+1}, u^{k+1} \right) + \Lambda^{k} \right\|^{2} \\ &= \arg\min_{z} \frac{1}{2} || \sqrt{\sum_{i=1}^{r} |z(:,:,:,i)|^{2}} - Y ||^{2} + \frac{\beta}{2} \left\| z - z^{+} \right\|^{2} \\ &= \arg\min_{z} \sum_{x,y,j} [\frac{1}{2} (\sqrt{\sum_{i=1}^{r} |z(x,y,j,i)|^{2}} - Y(x,y,j))^{2} + \frac{\beta}{2} ||z(x,y,j,:) - z^{+}(x,y,j,:)||^{2}] \\ &\text{where } z^{+} = \mathcal{A} \left( \omega^{k+1}, u^{k+1} \right) - \Lambda^{k} \end{split}$$

The close form solution of Step 3 is given as(details are in the Appendix 6.1):

$$z_i^{k+1} = \frac{z_i^+ \frac{Y}{M^k} + \beta z_i^+}{1+\beta}, 1 \le i \le r$$
 where  $z_i := z(:,:,:,i)$  and  $M^k = \sqrt{\sum_i |z_i^+|^2} \in \mathbb{C}^{px \times py \times N}$ 

# 2.3 Subproblem D and $\alpha$

$$\begin{split} D^{k+1} &= \arg \min_{D} \|\Omega - w^{k+1} + \Lambda_2^k\|^2 \\ &= \arg \min_{D} \|D\alpha^k - \hat{w}^{k+1}\|^2 \end{split}$$

where  $\hat{w}^{k+1} = reshape(\omega^{k+1} - \Lambda_2^k, [l, r]), D^*D = I$ 

The close form solution for  $D^{k+1}$  is (details are in the Appendix 6.1):

$$D^{k+1} = UV^* (2.8)$$

The updation of  $\alpha$  is easier:

$$\alpha^{k+1} = \arg\min_{\alpha} \|D^{k+1}\alpha - \hat{w}^{k+1}\|^2 = \arg\min_{\alpha} \sum_{i} ||\alpha_{i}D^{k+1}(:,i) - \hat{w}^{k+1}(:,i)||^2$$

Notice that each  $\alpha_i \in \mathbb{R}$  can be solved independently, the first optimality condition gives:

$$\alpha_i^{k+1} = \sum_{i_0} \Re[\overline{D^{k+1}(i_0, i)}\hat{w}^{k+1}(i_0, i)] (1 \le i \le r)$$
 (2.9)

# 3 Numerical experiments

The codes are implemented in MATLAB. First, we introduce the setting in experiments, then we conduct experiments on simulation data generated from specific models in 1.2. In each experiment, we compare the reconstructed images and modes in a different setting. We also generate ideal results as in 1.3.2 and compare ours with them.

#### **Algorithm 1:** ADMM for general mixed-state model(1.5)

**Initialization**: Set the number of states r,

$$\omega^{0}, u^{0}, z^{0} = \mathcal{A}\left(\omega^{0}, u^{0}\right), \Lambda^{0}, \Lambda^{0}_{2} = 0;$$
 $D^{0}$  and  $\alpha^{0}$  from SVD on  $\omega^{0}$ 

maximum iteration number Iter $_{\rm Max}$  , and parameter  $\beta,\!\beta_2$ 

**Output**:  $u^{\star} := u^{Iter_{Max}}$  and  $\omega^{\star} := \omega^{Iter_{Max}}$ 

1 for ii = 0 to  $Iter_{Max} - 1$  do

- Compute  $\omega^{k+1}$  by (2.5) with  $\hat{z}^k := z^k + \Lambda^k$ ;
- **3** Compute  $u^{k+1}$  (2.6). with  $\hat{z}^k$  the same as above;
- 4 Compute  $z_i^{k+1}, 1 \le i \le r$  by (6.1). with  $z^+ = \mathcal{A}\left(\omega^{k+1}, u^{k+1}\right) \Lambda^k$ ;
- 5 Compute  $D^{k+1}$  (6.2).;
- 6 Compute  $\alpha_i, 1 \le i \le r$  (2.9).;
- 7 Update the multiplier as Step 6 and Step 7 of (2.3) and (2.4);
- 8 end

## 3.1 Experiment setting

#### 3.1.1 Parameters

Parameters	Illustration	Values
$N_x, N_y$	size of image $u$	128,128
$p_x, p_y$	size of phobe $w$	64,64
Dist	scan distance between neighborhood frames	4,8,16
N	number of frames in diffusion image stacks	
$\overline{r}$	number of states(modes)	1(coherent) to 15
gridFlag	types of scan methods	1(rectangular lattice),
		$2 ({\rm hexagonal}), 3 ({\rm randomly\ disturb\ on\ } 2)$
blurFlag	types of partially coherent effect	1(1.4),2(1.3)

In order to deal with nontrivial ambiguities, non-periodical lattice-based scanning can be considered experimentally to remove the periodicity of the scanning geometry, e.g., adding a small number of random offsets to a set of the lattice. (gridFlag = 3). Or we can add prior knowledge about masks as an additional constraint, e.g. restricting the masks(frequency domain) in a circle.

Dist is an important parameter for successful reconstruction. Generally speaking, the smaller the Dist, the more the overlapping area and redundancy in data, we can get reconstruction images with higher qualities. Here we find

Dist = 8 is enough while Dist = 16 always fails.

#### 3.1.2 Performance metrics

In order to evaluate the performances of algorithms, we introduce 3 metrics.

1. Relative error err and signal-to-noise ratio snr

$$err^{k} = \frac{||cu^{k} - u_{true}||_{F}}{||cu^{k}||_{F}}, c = \frac{sum(u_{true} \circ \overline{u^{k}})}{||u^{k}||_{F}^{2}}$$
  
 $snr^{k} = -20\log_{10}(err^{k})$ 

 $||\cdot||_F$  is the Frobenius norm. err measures the difference between a reconstructed image and the groundtruth image. c is an estimated scale factor, and sum means the sum of all elements in the target matrix.

2. R-factor R

Let 
$$zz = \mathcal{A}_i \left( \omega^k, u^k \right)$$

$$R^k := \frac{\left\| \sqrt{\sum_{i=1}^r |zz(:,:,:,i)|^2} - Y \right\|_1}{\|Y\|_1}$$

R measures the difference between the reconstruction diffraction stacks and groundtruth stacks Y. We don't always know  $u_{true}$ , Y is the only input data for our algorithm, and R-factor can be used to verify the convergence.

3. Masks approximation error  $err_M$ 

$$err_{M}^{k} = \frac{||c\rho^{k} - \rho_{true}||_{F}}{||\rho_{true}||_{F}}, c = \frac{sum(\rho_{true} \circ \overline{\rho^{k}})}{||\rho^{k}||_{F}^{2}}$$

 $err_M$  measures the difference between the reconstructed density matrix  $\rho^k$  from masks and the standard density matrix  $\rho_{true}$  from the theoretical model.

#### 3.1.3 Operations on modes

1. Initialization

In the following tests, we set  $u^0 = \mathbf{1}_{N_x \times N_y}$  and initial phobe  $m = \frac{1}{N} \mathcal{F}^{-1} \left( \sum_j Y(:,:,j) \right)$ . We generate other modes by randomly disturb intial mask m. r initial probes were created by multiplying the initial mask by different arrays of random complex values (modulus part varies within the range of [0,1] and phase part  $[0,2\pi]$ ). Then we get  $w^0$ .

#### 2. Orthogonalization

Modes  $w_k$  computed in our algorithm are not always orthogonal. However, we can easily orthogonalize them by generating a density matrix  $\rho$  and perform spectral decomposition on  $\rho$  like in (1.7). The orthogonal representation of modes is always unique(Specifically, when the eigenvalues of  $\rho$  are all different, and this always happens in real-world data).

Notice that the number of modes is always small i.e.  $r \ll l := p_x \times p_y$ , we perform SVD on the  $l \times r$  phobe matrix directly instead of  $l \times l$  density matrix  $\rho$ . And we can also select the first  $r' \leq r$  modes for an approximation.

## Algorithm 2: SVD-based orthogonalization for phobes(Matlab)

```
Input: w \in \mathbb{C}^{px \times py \times r}, the number of modes needed r'
Output: w_{ort} \in \mathbb{C}^{px \times py \times r'}

1 phobe matrix ss = reshape(w, [l\ r]);

2 [U, S, V] = svd(ss, 'econ');

3 q = U(:, 1:r') * S(1:r', 1:r');

4 w_{ort} = reshape(q, [px\ py\ r']);
```

Orthogonalization operation is always performed before we display final modes to get a clearer representation. We also want to figure out whether orthogonalization can be added to improve our Algorithm1.

#### 3. Compression

In noisy case, we consider extracting fewer main modes to avoid the distortion from noise.

#### **Algorithm 3:** SVD-based compression for phobes(Matlab)

```
Input: w \in \mathbb{C}^{px \times py \times r}, the number of main modes kept r'

Output: w_{com} \in \mathbb{C}^{px \times py \times r}

1 phobe matrix ss = reshape(w, [l\ r]);

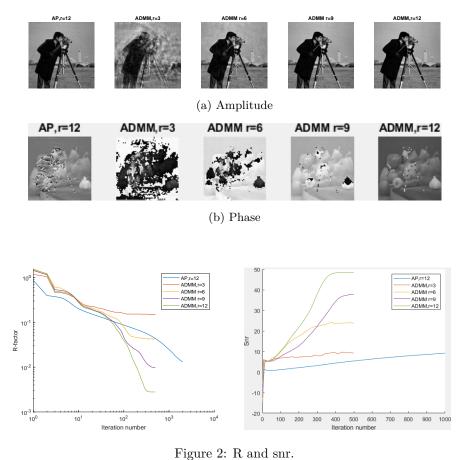
2 [U, S, V] = svd(ss, econ');

3 q = U(:, 1:r') * S(1:r', 1:r') * V(:, 1:r');

4 w_{com} = reshape(q, [px\ py\ r]);
```

# 3.2 Approximation by modes

 $Dist=8,\ gridFlag=1,\ blurFlag=2$  and  $\kappa$  is a guassian kernel with



rigure 2. It and sin.

We first compare snr and R for reconstructed images using a different number of modes. When the number of modes increases, the R-factor decreases, and the snr increases. That indicates that the quality of the reconstructed image increases.

The R for 3,6,9,12 modes using ADMM algorithm are stable after 500 iterations at 0.15, 0.042, 0.0099, 0.0028, respectively, when R for AP does not

converge after 2000 iterations.

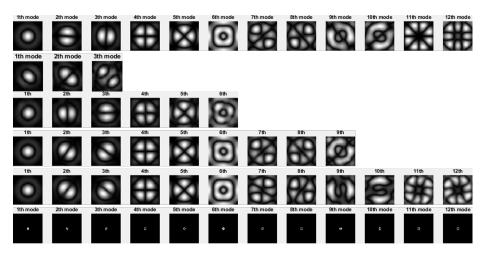


Figure 3: Mode pattern. The first row represents the standard mode pattern. And the last two rows represent the mode pattern for 12 modes. Mode patterns are in the time domain except that the last row is in the frequency domain.

Standard mode pattern is obtained through performing SVD on standard density matrix generated from the model and extracting the first 12 modes. As shown in Figure 3, our algorithm can generally catch the main modes and get an optimal approximation.

Further, denote  $err_M$  for r modes as  $err_M^r$ . Optimal  $err_M^{r,*}$  can be calculated from the theory of low-rank approximation, and we compare our  $err_M^r$  with it. Denote the singular values of standard density matrix  $\rho_{true}$  as  $s_i, i = 1...d = rank(\rho_{true})$ 

$$err_{M}^{r,*} = \min_{\rho, rank(\rho) = r} \frac{||\rho_{true} - \rho||_{F}}{||\rho_{true}||_{F}} = \sqrt{\frac{\sum_{i=r+1}^{d} s_{i}^{2}}{\sum_{i} s_{i}^{2}}} = \sqrt{1 - s_{cum}(r)}$$

$$where \ s_{cum}(r) = \frac{\sum_{i=1}^{r} s_{i}^{2}}{\sum_{i} s_{i}^{2}}$$

As shown in Figure 5, the  $err_M^*$  reaches around 0.01, and  $err_M$  is close to  $err_M^*$  with 12 modes, which indicates that our algorithm does well in low-rank approximation to standard density matrix.

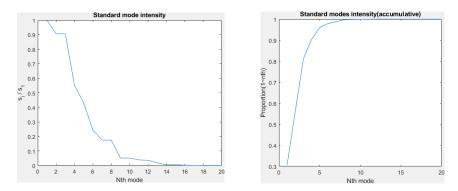
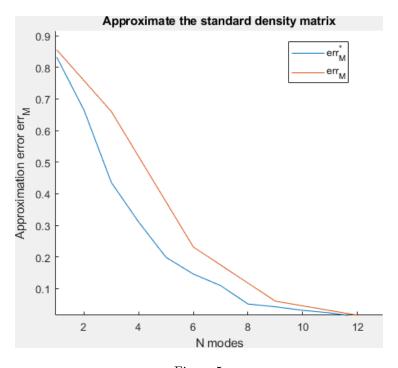


Figure 4: The distribution of singular values of the standard density matrix. The vertical axis in the left subfigure represents the ratio of  $i^{th}$  largest singular to the first one  $s_i/s_1$ , and that in the right one represents  $S_{cum}(i)$ . The singular value decreases exponentially and the matrix is approximately low-rank.



 $Figure\ 5$ 

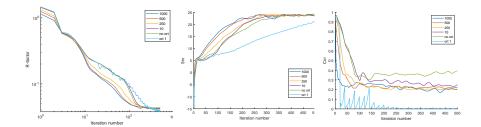


Figure 6: The vertical axis and horizontal axis are set in log-scale in the first subfigure for R-factor. The blue line represents  $\beta_2/\beta=1000$  and works the best in this case. The green line represents the result without orthogonalization. 'ort 1' means performing orthogonalization every 20 iterations.

# 3.3 Add orthogonalization constraint

The experiment setting is the same as above except that we change  $ratio = \beta_2\beta$  to introduce orthogonalization constraint in ADMM. The larger the  $\beta_2$ , the stricter the orthogonalization constraint. As a reference, we also tried performing orthogonalization as in Algorithm 2 every 20 iterations.

The reconstructed images are similar after 500 iterations, while snr and Rfactor using the ADMM algorithm with orthogonalization constraint improve faster. And the degree of correlation coherence between modes also decreases faster with larger  $\beta_2$ .

#### Remark.

- 1. We suppose that orthogonalization constraints can improve robustness on different initial values.
- 2. When the number of modes is large enough(like 12 here), the effect of the orthogonalization constraint seems not noticeable.

### 3.4 Noisy case

 $Dist=8,\ gridFlag=3,\ blurFlag=2$  and  $\kappa$  is a guassian kernel with  $\sigma=(15,15),\ \beta=0.05.$ 

Poisson noise is added on diffraction images Y through MATLAB command:

$$Y_{noise} = poissrnd(Y * (\eta))/\eta;$$

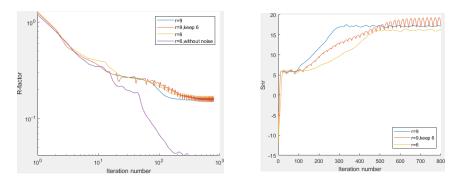


Figure 8: R and snr. The red line represents the result with compression.

where  $\eta = 0.0675$  is used here.

Four different settings are compared: 9 modes with noise, 9 modes with



Figure 7: Reconstructed images in the noisy case

The result periodically fluctuates with compression operation, while it can finally behave the best in the noisy case. Although the difference is not noticeable, we can consider adding compression in constraint instead of performing the direct truncated operation.

# 4 进一步地讨论

虽然我们用一般化模型也较好地求解了偏向干的相位恢复问题,但没用用到特殊模型 1.9 中特殊的结构。我们要恢复的不是一个一般的密度矩阵,而是有一定结构的,下面将进行推倒阐述:

## 4.1 Density matrix in Fourier space

$$f_{pc,j}(q) = \int |\mathcal{F}_{x \to q} (\mathcal{S}_{j} u(x) \omega(x - y))|^{2} \kappa(y) dy$$

$$= \int \mathcal{F}_{x \to q} (\mathcal{S}_{j} u(x) \omega(x - y)) \overline{\mathcal{F}_{x \to q} (\mathcal{S}_{j} u(x) \omega(x - y))} \kappa(y) dy$$

$$= \int \int \mathcal{S}_{j} u(x_{1}) \omega(x_{1} - y) e^{-iqx_{1}} dx_{1} \overline{\int} \mathcal{S}_{j} u(x_{2}) \omega(x_{2} - y) e^{-iqx_{2}} dx_{2} \kappa(y) dy$$

$$= \int \int \mathcal{S}_{j} u(x_{1}) e^{-iqx_{1}} \omega(x_{1} - y) dx_{1} \overline{\int} \mathcal{S}_{j} u(x_{2}) e^{-iqx_{2}} \omega(x_{2} - y) dx_{2} \kappa(y) dy$$

$$= \int \int \mathcal{S}_{j} u(x_{1}) e^{-iqx_{1}} \omega(x_{1} - y) dx_{1}' \overline{\int} (\mathcal{S}_{j} u) \widehat{(x_{2})} e^{-iqx_{2}} \omega(x_{2} - y) dx_{2}' \kappa(y) dy$$

$$= \int \int \mathcal{S}_{j} u(x_{1}) e^{-iqx_{1}} \widehat{\omega}(x_{1}') e^{-iyx_{1}'} \overline{\mathcal{S}_{j} u(x_{2})} e^{-iqx_{2}} \widehat{\omega}(x_{2}') e^{-iyx_{2}'} \int \kappa(y) dy dx_{1}' dx_{2}'$$

$$= \int \int \widehat{(\mathcal{S}_{j} u)} (x_{1}' + q) \widehat{\omega}(x_{1}') \overline{\widehat{(\mathcal{S}_{j} u)} (x_{2}' + q) \widehat{\omega}(x_{2}')} \widehat{(\int} \kappa(y) e^{-iy(x_{1}' - x_{2}')} dy)$$

$$= \int \int \widehat{(\mathcal{S}_{j} u)} (x_{1}' + q) \widehat{\omega}(x_{1}') \overline{\widehat{(\mathcal{S}_{j} u)} (x_{2}' + q) \widehat{\omega}(x_{2}')} \widehat{\kappa}(x_{1}' - x_{2}') dx_{1}' dx_{2}'$$

$$(4.1)$$

If we can decompose  $\hat{\kappa}(x_1', x_2') = \sum_{i=1}^r a_i(x_1') \overline{a(x_2')}$ , then we can get the density matrix in Fourier space:

$$\hat{\rho}(x_1', x_2') = \sum_{i=1}^r \hat{\omega}(x_1') a_i(x_1') \overline{\hat{\omega}(x_2') a_i(x_2')}$$

which is similar to 1.10. We have:

$$\hat{\rho} = \hat{\omega}\hat{\omega}'. * \hat{\kappa}$$

(This has been verified through experiment)

#### 4.2 Ideas

Notice that if  $x_1', x_2' \in R^1$ ,  $\hat{\kappa}(x_1' - x_2')$  can be represented by a Toeplitz matrix in discrete case. And when  $\kappa(-x) = \kappa(x)$ , it is a symmetric Toeplitz matrix. Under certain condition, it may be a semi-positive definite one. This gives a special structure of the density matrix that we want to decompose and approximate. Further, refer to the vandemonde decomposition<sup>[11]</sup>, a Toeplitz matrix can be decomposed as followed: Suppose that T is an  $n \times n$  positive

semidefinite (PSD) Toeplitz matrix of rank r < n. The result states that T can be uniquely decomposed as

$$T = APA^H$$

where  $\boldsymbol{P}$  is an  $r \times r$  positive definite diagonal matrix and  $\boldsymbol{A}$  is an  $n \times r$  Vandermonde matrix whose columns correspond to uniformly sampled complex sinusoids with different frequencies. That is,

$$\boldsymbol{A}_{n}(\boldsymbol{f}) := \left[\boldsymbol{a}_{n}\left(f_{1}\right), \ldots, \boldsymbol{a}_{n}\left(f_{r}\right)\right]$$

specific,  $\mathbf{a}_n(f) = n^{-\frac{1}{2}} \left[ 1, e^{i2\pi f}, \dots, e^{i2\pi(n-1)f} \right]^T \in \mathbb{C}^n$  denote a uniformly sampled complex sinusoid with frequency  $f \in \mathbb{T}([0,1])$  and unit power, where  $i = \sqrt{-1}$ . It follows that:  $\mathbf{P} = \operatorname{diag}\left( |c_1|^2, \dots, |c_r|^2 \right)$ . Then, the  $\hat{\kappa}(x_1', x_2')$  can be decomposed as:

$$\hat{\kappa} = \boldsymbol{A}_n(\boldsymbol{f}) \boldsymbol{P} \boldsymbol{A}_n^H(\boldsymbol{f}) = \sum_{j=1}^r |c_j|^2 a_n(f_j) \boldsymbol{a}_n^H(f_j)$$

And the density matrix in Fourier space can be decomposed as:

$$\hat{\rho} = \sum_{j=1}^{r} |c_j|^2 \left[ \hat{\omega} a_n \left( f_j \right) \right] \left[ \hat{\omega} \boldsymbol{a}_n \left( f_j \right) \right]^H$$

Remark: more investigation needed below. However, in our case,  $x'_1, x'_2 \in \mathbb{R}^2$ . In [11], the author extends it to higher dimensional cases If we can write

$$\hat{\rho} = \sum_{j=1}^{r} |c_j|^2 \left[ \hat{\omega} a_n \left( \overrightarrow{f_j} \right) \right] \left[ \hat{\omega} \boldsymbol{a}_n \left( \overrightarrow{f_j} \right) \right]^H$$

where

$$\boldsymbol{a}_n(\overrightarrow{f}) = n^{-\frac{1}{2}} \left[ 1, e^{i2\pi \overrightarrow{f} \cdot \overrightarrow{x'}}, \dots, e^{i2\pi(n-1)\overrightarrow{f} \cdot \overrightarrow{x'}} \right]^T \in \mathbb{C}^n$$

density matrix in time space will be amazingly written like:

$$\hat{\rho} = \sum_{j=1}^{r} |c_j|^2 \, \omega(\overrightarrow{x_1} - \overrightarrow{f}) \overline{\omega(\overrightarrow{x_2} - \overrightarrow{f})}$$

# 5 Convergence analysis

#### 5.1 Framework of convergence analysis

#### 5.1.1 Lipchitz property

Definition 1 (Lipschitz property). The function  $\Phi: \mathbb{C}^m \to \mathbb{R}$  has the Lipschitz property in  $\mathbb{C}^m$  with a Lipschitz constant L if

(1) it is Lipschitz differentiable (its gradient is Lipschitz continuous), i.e.,

$$\|\nabla\Phi(v_1) - \nabla\Phi(v_2)\| \le L \|v_1 - v_2\| \, \forall v_1, v_2 \in \mathbb{C}^m$$
 (5.1)

(2) it has the descent property as

$$(2.5) \Phi(v_2) - \Phi(v_1) - \Re(\langle \nabla \Phi(v_1), v_2 - v_1 \rangle) \le \frac{L}{2} \|v_2 - v_1\|^2 \, \forall v_1, v_2 \in \mathbb{C}^m.$$
(5.2)

Show the function  $\mathcal{G}(\cdot)$  has the Lipschitz property.

#### 5.1.2 Properties of iterative sequence

First, we briefly review the framework for the convergence analysis of an iterative algorithm for nonconvex optimization problems. Consider an optimization problem minv  $\Phi(\mathbf{v})$ , with the functional  $\Phi$  being proper, lower semicontinuous, and bounded below. Let an iterative sequence  $v_{k_{k=0}}^{\infty}$  be generated by some algorithm to solve the above minimization problem.

(i) sufficient decrease condition as

$$\Phi\left(v^{k}\right) - \Phi\left(v^{k+1}\right) \ge c_1 \left\|v^{k+1} - v^{k}\right\|^2, \text{ with a positive constant } c_1 \qquad (5.3)$$

(ii) relative error condition as

$$(4.2) \|p^{k+1}\| \le c_2 \|v^{k+1} - v^k\|, \text{ with some } p^{k+1} \in \partial \Phi\left(v^{k+1}\right), \text{ and a positive constant } c_2,$$

$$(5.4)$$

(iii) Kurdyka Łojasiewicz (KL) property [1].

The first two conditions can guarantee that each limiting point of the iterative sequence is a stationary point of  $\Phi$ , which demonstrates the subsequence convergence. Furthermore, by the KL property of  $\Phi$ , the iterative sequence can be proved to be a Cauchy sequence and therefore convergence can be reached. Along this line, the convergence for ADMM was established for noncovex problems [22, 40, 24, 31, 45].

# 5.2 Convergence analysis for algorithm without the orthogonal constraint

$$\mathcal{G}(z) = \frac{1}{2} ||\sqrt{|z|^2 + \epsilon 1_m} - \sqrt{f + \epsilon 1_m}||^2 = \frac{1}{2} ||\sqrt{\sum_i |z_i|^2 + \epsilon 1_m} - \sqrt{f + \epsilon 1_m}||^2$$

Compute  $\nabla g(z) = \langle \nabla_1 g(z_1), \nabla_2 g(z_2), \cdots, \nabla_r g(z_r) \rangle$  for a single  $\nabla_i g(z_i)$  we get

$$||\nabla_i g(z_{i,1}) - \nabla_i g(z_{i,2})|| \le (1 + \frac{2}{\sqrt{\epsilon}})||z_{i,1} - z_{i,2}||$$

Combine i and get (5.1), then from smoothness we have (5.2)

We can almost follow the framework of proof in coherent case. We only need to split z to  $z_i$  and  $\omega$  to  $\omega_i$  for different modes, verify them separately.

Question: I don't know how to verify KL property.

# 5.3 Convergence analysis for algorithm without the orthogonal constraint

 $\mathcal{G}(z)$  is the same as above, so the LP property has been verified.

Next we try to verify three properties of the iterative sequence.

The first order conditions give: 4.1. Convergence of Algorithm 1. We show the first-order optimality conditions for subproblems of (3.4). Letting  $X^k := \left(\omega^k, u^k, z^k, \Lambda^k, D^k, \alpha^k, \Lambda_2^k\right)$  be generated by Algorithm 1, the following relations hold:

$$0 \in \omega^{k+1} \circ \sum_{j} \left| \mathcal{S}_{j} u^{k} \right|^{2} - \sum_{j} \left( \mathcal{S}_{j} u^{k} \right)^{*} \circ \mathcal{F}^{-1} \left( z_{j}^{k} + \frac{1}{\beta_{1}} \Lambda_{j}^{k} \right)$$

$$+ \frac{1}{\beta_{1}} \partial \mathbb{I}_{\mathcal{X}_{1}} \left( \omega^{k+1} \right) + \frac{\alpha_{1}}{\beta_{1}} \operatorname{diag} \left( M_{1}^{k} \right) \circ E_{\omega}^{k+1} + \frac{\beta_{2}}{\beta_{1}} (\omega^{k+1} - \hat{\Omega}^{k}), \hat{\Omega} = \Omega + \Lambda_{2};$$

$$0 \in u^{k+1} \circ \sum_{j} \mathcal{S}_{j}^{T} \left| \omega^{k+1} \right|^{2} - \sum_{j} \mathcal{S}_{j}^{T} \left( \left( \omega^{k+1} \right)^{*} \circ \mathcal{F}^{-1} \left( z_{j}^{k} + \frac{1}{\beta} \Lambda_{j}^{k} \right) \right)$$

$$+ \frac{1}{\beta} \partial \mathbb{I}_{\mathcal{X}_{2}} \left( u^{k+1} \right) + \frac{\alpha_{2}}{\beta} \operatorname{diag} \left( M_{2}^{k} \right) \circ E_{u}^{k+1}$$

$$0 = \nabla \mathcal{G} \left( z^{k+1} \right) + \beta \left( z^{k+1} - \mathcal{A} \left( \omega^{k+1}, u^{k+1} \right) \right) + \Lambda^{k}.$$

Using (4.6) and the Step 4 of (3.4), one can derive

$$0 = \nabla \mathcal{G}(z^{k+1}) + \Lambda^{k+1} \forall k \ge 0.$$

The optimality condition for D is unclear because of the orthogonal constraint.

$$(?)0 \in \frac{1}{\beta_2} \partial \mathbb{I}_{\mathcal{X}_3} \left( D^{k+1} \right) + (\Omega(D^{k+1} \alpha^k) - \omega^{k+1} + \Lambda_2^k) \alpha_k^k$$
$$\alpha_i^{k+1} = \sum_{i_0} \Re[\overline{D^{k+1}(i_0, i)} \hat{w}^{k+1}(i_0, i)] (1 \le i \le r)$$

$$(?)\Lambda_2^{k+1} = \Lambda_2^k + (\Omega(D^{k+1}\alpha^{k+1}) - \omega^{k+1})$$

$$\begin{split} \Upsilon_{\beta}(\omega, u, z, \Lambda) := & \mathcal{G}(z) + \mathbb{I}_{\mathcal{X}_{1}}(\omega) + \mathbb{I}_{\mathcal{X}_{2}}(u) + \mathbb{I}_{\mathcal{X}_{3}}(D\alpha) \\ & + \frac{\beta}{2} \|z - \mathcal{A}(\omega, u) + \Lambda\|^{2} - \frac{\beta}{2} ||\Lambda||^{2} + \frac{\beta_{2}}{2} \|\Omega(D\alpha) - w + \Lambda_{2}\|^{2} - \frac{\beta_{2}}{2} ||\Lambda_{2}||^{2} \\ & (5.5) \end{split}$$

(5.3)  $u, z, \Lambda$  are the same as previous.

Ps: the update for z is not easy. We need the LP property of  $\mathcal{G}(z)$ , Cauchy's  $\Re\left(\langle z_1, z_2 \rangle\right) \geq -\frac{L}{2} \|z_1\|^2 - \frac{1}{2L} \|z_2\|^2 \,\forall \, z_1, z_2 \in \mathbb{C}^m$ , and carefully estimate.

As for  $\omega$ , it is like another stable term  $\|\Omega(D\alpha) - w + \Lambda_2\|^2$  is added, and can be derived by calculating the Gateaux derivative.

 $\alpha$  is OK.

from  $D^k$  to  $D^{k+1}$ , the decrease is unclear since the update step is complex. However, we can prove it if add the assumption that  $\omega \alpha = USV^*$ ,  $s_i$  is bounded below.

 $\Lambda_2$  is unclear because it has no relationship with  $\nabla \mathcal{G}(z)$ , we have to bound below  $(\Omega(D^{k+1}\alpha^{k+1}) - \omega^{k+1})$ , but it seems unreasonable.

(5.4) From the first order optimality conditions, we can verify the relative error condition for all variables.

We can not easily get the boundedness in:

Lemma 4.2 (boundedness). Letting  $\beta > \frac{3+\sqrt{17}}{2}L$ , the sequence  $\{X^k\}$  generated by Algorithm 1 is bounded. Furthermore, the sequence of augmented Lagrangian  $\{\Upsilon_\beta(X^k)\}$  is bounded and nonincreasing.

 $\Lambda_2$  is unclear because it has no relationship with  $\nabla \mathcal{G}(z)$ , and the boundedness relationship can not be passed from variable to variable.

# 6 Appendix

### 6.1 Subproblems in ADMM

#### 6.1.1 $\omega$

Essentially in this subproblem, each state  $\omega_i = w(:,:,i)$  is independent. Then we can optimize each  $w_i$  separately.

$$\omega_i^{k+1} = \arg\min_{\omega \in \mathcal{X}_1} \frac{1}{2} \sum_j \left\| \mathcal{F}^{-1} \hat{z}(:,:,j,i)^k - \omega(:,:,i) \circ \mathcal{S}_j u^k \right\|^2$$

Essentially in this subsubproblem, each element in  $w_i$  is independent. such that one just needs to solve the following 1D constraint quadratic problem:

$$\omega_i^{k+1}(t) = \arg\min_{|x| < C_{\omega}} \rho_t^k(x).$$

where  $\rho_t^k(x) := \frac{1}{2} \sum_j \left| \left( \mathcal{F}^{-1} \hat{z}(:,:,j,i)^k \right)(t) - x \times \left( \mathcal{S}_j u^k \right)(t) \right|^2 \forall x \in \mathbb{C}$ The derivative of  $\rho_t^k(x)$  is calculated <sup>2</sup> as

$$\nabla \rho_t^k(x) = \frac{d\rho_t^k(x)}{dx^*}$$

$$= \sum_j \left( x \times \left| \left( \mathcal{S}_j u^k \right) (t) \right|^2 - \left( \mathcal{S}_j u^k \right)^* (t) \left( \mathcal{F}^{-1} \hat{z}_{j,i}^k \right) (t) \right)$$

$$= x \times \left( \sum_j \left| \left( \mathcal{S}_j u^k \right) (t) \right|^2 \right) - \sum_j \left( \left( \mathcal{S}_j u^k \right)^* (t) \left( \mathcal{F}^{-1} \hat{z}_{j,i}^k \right) (t) \right)$$

The first order optimality condition is  $\nabla \rho_t^k(x) = 0$ . Then the close form solution of  $w_i$  is given as

$$\omega_i^{k+1} = \operatorname{Proj}\left(\frac{\sum_j \left(\mathcal{S}_j u^k\right)^* \circ \left(\mathcal{F}^{-1} \hat{z}_{j,i}^k\right)}{\sum_j \left|\mathcal{S}_j u^k\right|^2}; C_{\omega}\right)$$

#### 6.1.2 z

$$\begin{split} z^{k+1} &= \arg\min_{z} \mathcal{G}(z) + \frac{\beta}{2} \left\| z - \mathcal{A} \left( \omega^{k+1}, u^{k+1} \right) + \Lambda^{k} \right\|^{2} \\ &= \arg\min_{z} \frac{1}{2} || \sqrt{\sum_{i=1}^{r} |z(:,:,:,i)|^{2}} - Y ||^{2} + \frac{\beta}{2} \left\| z - z^{+} \right\|^{2} \\ &= \arg\min_{z} \sum_{x,y,j} [\frac{1}{2} (\sqrt{\sum_{i=1}^{r} |z(x,y,j,i)|^{2}} - Y(x,y,j))^{2} + \frac{\beta}{2} ||z(x,y,j,:) - z^{+}(x,y,j,:)||^{2}] \\ &\text{where } z^{+} = \mathcal{A} \left( \omega^{k+1}, u^{k+1} \right) - \Lambda^{k} \end{split}$$

$$z^*(x,y,j,:) = \arg\min_{z_{x,y,j} \in \mathbb{C}^r} \frac{1}{2} (||z_{x,y,j}|| - Y_{x,y,j})^2 + \frac{\beta}{2} ||z_{x,y,j} - z_{x,y,j}^+||^2$$

For any fixed x, y, j and free i, the problem can be seen as:

<sup>&</sup>lt;sup>2</sup>Notice that here  $\rho$  is a real value function with complex variable, and we use wirtinger derivatives here. More properties and calculation rules are listed in this link: https://blog.csdn.net/weixin\_37872766/article/details/107673096

Notice that for fixed  $||z_{x,y,j}||$ , the first term in expression is fixed. To optimize the second term, we should always choose  $z_{x,y,j}$  with the same direction as  $z_{x,y,j}^+$ . So we have  $||z_{x,y,j}-z_{x,y,j}^+||^2=(||z_{x,y,j}||-||z_{x,y,j}^+||)^2$ 

$$\frac{z(x,y,j,i)}{||z_{x,y,j}||} = \frac{z^+(x,y,j,i)}{||z_{x,y,j}^+||}, z(x,y,j,i) = ||z_{x,y,j}|| \frac{z^+(x,y,j,i)}{||z_{x,y,j}^+||}$$

To determine  $z_{x,y,j}$ , we only need to determine  $||z_{x,y,j}||$ . Denote it as a.

$$||z_{x,y,j}||^* = \arg\min_{a \in \mathbb{R}} \frac{1}{2} (a - Y_{x,y,j})^2 + \frac{\beta}{2} (a - ||z_{x,y,j}^+||)^2$$

The first optimality condition easily gives:

$$a = \frac{Y_{x,y,j} + \beta ||z_{x,y,j}^+||}{1 + \beta}$$

The close form solution of Step 3 is given as:

$$z_i^{k+1} = \frac{z^+ \frac{Y}{M^k} + \beta z_i^+}{1 + \beta}, 1 \le i \le r$$
 (6.1)

where  $M^k = \sqrt{\sum_i |z^+(:,:,:,i)|^2} \in \mathbb{C}^{px \times py \times N}$ 

# 6.1.3 D

$$\begin{split} D^{k+1} &= \arg \min_{D} \|\Omega - w^{k+1} + \Lambda_2^k\|^2 \\ &= \arg \min_{D} \|D\alpha^k - \hat{w}^{k+1}\|^2 \end{split}$$

where  $\hat{w}^{k+1} = reshape(\omega^{k+1} - \Lambda_2^k, [px \times py, r]), D^*D = I$ 

This is a special case in Orthogonal Procrustes problem  $^3$ 

$$\begin{split} \|D\alpha^k - \hat{w}^{k+1}\|^2 = & Tr[(D\alpha^k - \hat{w}^{k+1})^*(D\alpha^k - \hat{w}^{k+1})] \\ = & \|\alpha^k\|_F^2 - Tr[(\alpha^k)^*D^*\hat{\omega}^{k+1}] - Tr[(\hat{\omega}^{k+1})^*D\alpha^k] + \|\hat{\omega}^{k+1}\|_F^2 \\ D^{k+1} = & \arg\max_D Tr[(\alpha^k)^*D^*\hat{\omega}^{k+1}] + Tr[(\hat{\omega}^{k+1})^*D\alpha^k] \\ = & \arg\max_D \Re(Tr[(\alpha^k)^*D^*\hat{\omega}^{k+1}]) \\ = & \arg\max_D \Re(Tr[D^*(\hat{\omega}^{k+1}\alpha^k)]) \end{split}$$

 $<sup>^3 \</sup>verb|https://en.wikipedia.org/wiki/Orthogonal_Procrustes_problem|$ 

Consider the SVD decomposition:  $\hat{\omega}^{k+1}\alpha^k = USV^*$ 

$$\begin{split} D^{k+1} &= \arg\max_{D} \Re(Tr[D^*USV^*]) = \arg\max_{D} \Re(Tr[(V^*D^*U)S]) \\ & \xrightarrow{\hat{D} = V^*D^*U \text{ is orthonormal}} \arg\max_{\hat{D}} \Re(Tr[\hat{D}S]) \end{split}$$

We can easily see  $\hat{D} = I$  is optimal, and:

$$D^{k+1} = UV^* \tag{6.2}$$

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