

Reconstruct-Anything Model

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Joint work with

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Linear Inverse Problems

Goal: recover signal x from y

$$y = A(x) + \epsilon$$

MEASUREMENTS $\in \mathbb{R}^m$

SIGNAL $\in \mathbb{R}^n$

NOISE $\in \mathbb{R}^m$

Examples

Image denoising

- $A = \text{identity}$

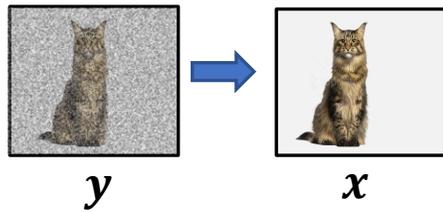
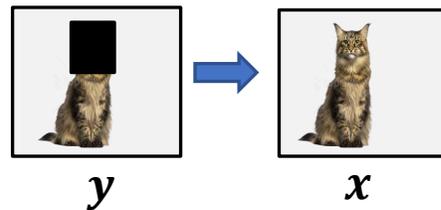
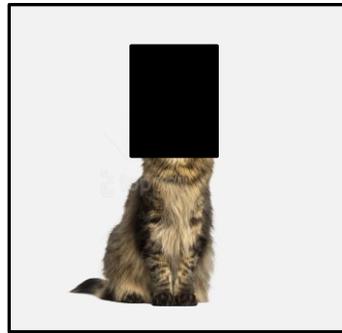


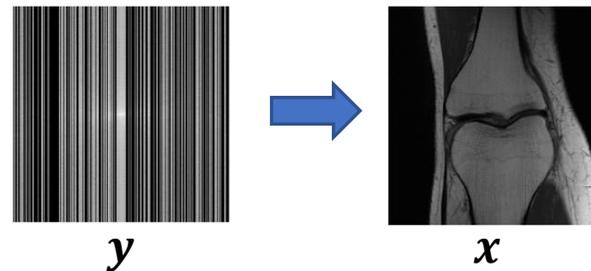
Image inpainting

- $A = \text{diagonal matrix with 1's and 0s.}$



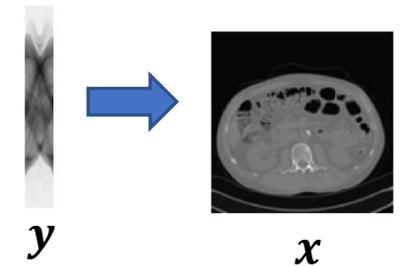
Magnetic resonance imaging

- $A = \text{subset of Fourier modes (} k - \text{space)}$



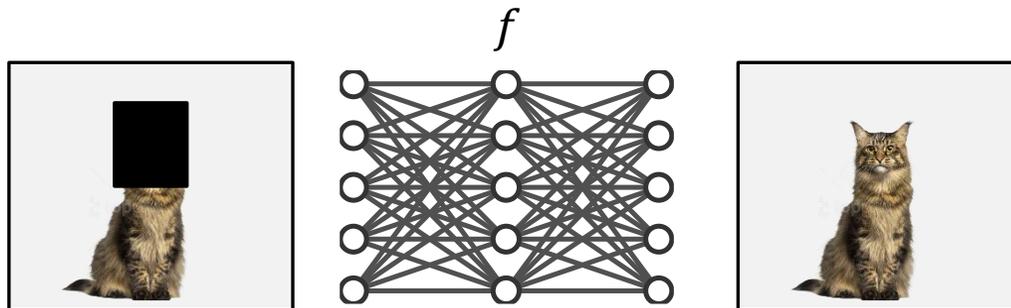
Computed tomography

- $A = \text{1D projections (sinograms)}$



End2End Learning

Idea: use training pairs of signals and measurements to directly learn the inversion function



$$\operatorname{argmin}_f \mathbb{E}_{x,y} \|x - f(y, A)\|^2$$

supervised
dataset



Benefits of End2End

- **Advantages**
 - Outperform plug-and-play/diffusion models on training task
 - Much faster inference (doesn't require multiple iterations as diffusion/PnP)
- **Disadvantages:**
 - Trained for a specific task, doesn't generalize to different tasks
 - Requires a large supervised dataset specific to application
 - It is a black-box, no notion of uncertainty

Taxonomy of End2End

- **Artifact removal networks** (Jin et al., 2017):

$$f(\mathbf{y}, A) = \phi(A^\top \mathbf{y})$$

where ϕ is an image-to-image networks, such as a U-Net.

- **Unrolled neural networks** (Gregor and LeCun, 2010): e.g. proximal gradient descent, we have $f(\mathbf{y}, A) = \mathbf{x}^K$ where

$$\mathbf{x}^{k+1} = \phi_k(\mathbf{x}^k - A^\top A \mathbf{x}^k + A^\top \mathbf{y}) \text{ for } k = 1, \dots, K$$

where ϕ_k are image-to-image networks.

Typically 1 to 3 dB better

How to design f ?

While unrolled typically obtains SOTA results

- Not enough iterates to ‘converge’
- Convergence guarantees lost when using expressive $\phi_1 \dots \phi_k$

Kimmel: “*DL is a dark monster covered with mirrors. Everyone sees his reflection in it ...*”, Donoho: “*... these mirrors are taken from Cinderella's story, telling each that they are the most beautiful*”

- If we forget optimization perspective:

*main difference between unrolling and artifact removal
is **backpropagation through $A^T A$***

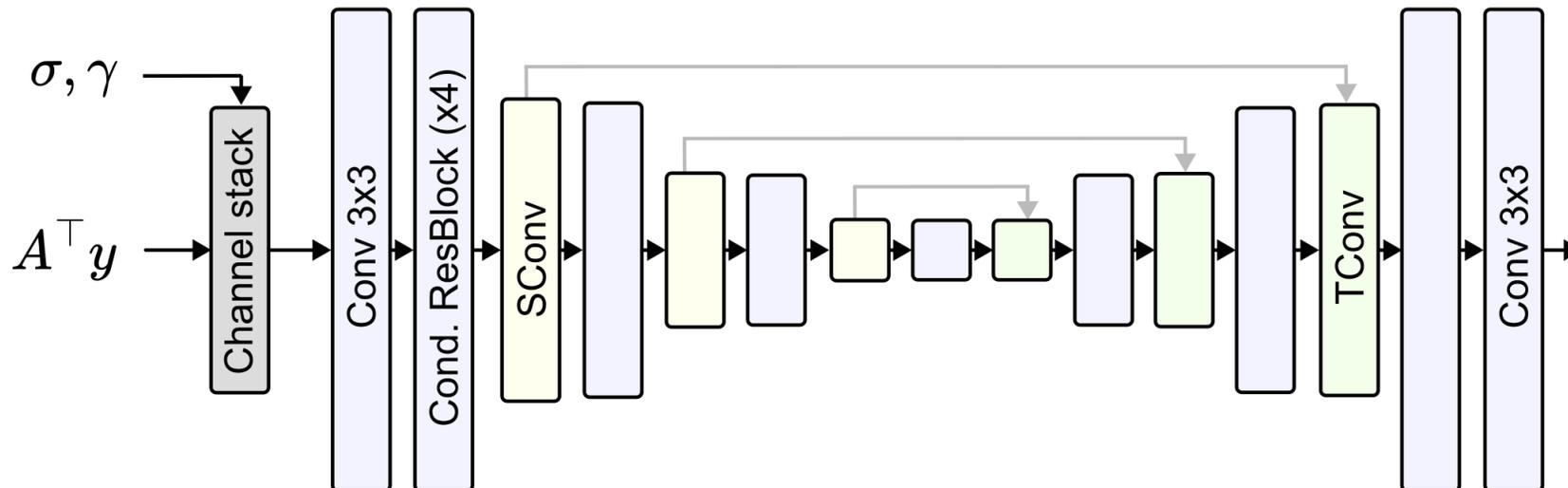
Towards a new model

How can we come up with a better end-to-end model?

- **Hypothesis 1:** A standard U-Net architecture should be enough
- **Hypothesis 2:** Backpropagating through $A^T A$ boosts performance
- **Hypothesis 3:** The network should generalize to multiple A if trained on a family [Gossard & Weiss, 2022]

Towards a new model

- **Hypothesis 1:** A standard U-Net architecture should be enough
- State-of-the-art architecture used for denoising, super-resolution, segmentation, etc.
- We take the **DRUNet architecture** (Zhang et al., 2021) as backbone
- We can start with pretrained denoising weights



Towards a new model

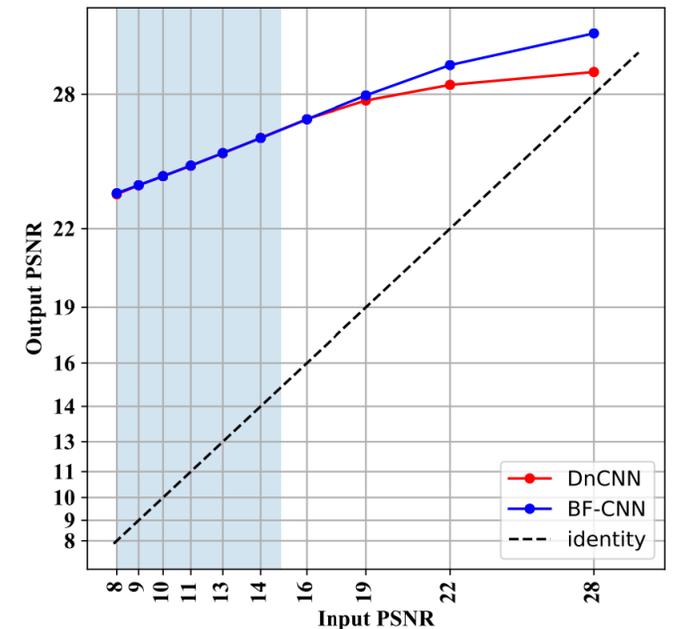
- **Hypothesis 1:** A standard U-Net architecture should be enough
- The network is **scale equivariant** (Mohan et al., 2020), (Herbreteau et al., 2023)

- **Poisson-Gaussian** noise model: $\mathbf{y} = \gamma \mathcal{P}\left(\frac{A\mathbf{x}}{\gamma}\right) + \sigma \epsilon$
- Scaling by $\alpha > 0$ gives $\alpha \mathbf{y} = \alpha \gamma \mathcal{P}\left(\frac{A\alpha \mathbf{x}}{\alpha \gamma}\right) + \alpha \sigma \epsilon$

Thus we want:

$$f(\alpha \mathbf{y}, \alpha \sigma, \alpha \gamma) = \alpha f(\mathbf{y}, \sigma, \gamma)$$

which is automatically verified by the DRUNet noise conditioning

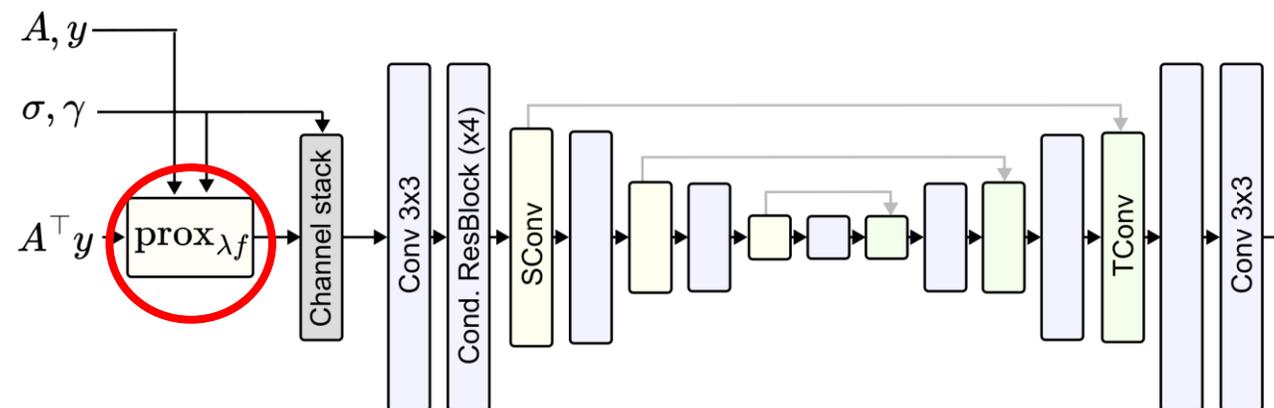


Towards a new model

- **Hypothesis 1:** A standard U-Net architecture should be enough
- We map measurements y to an input of DRUNet as:

$$\text{input} = \operatorname{argmin} \|\mathbf{y} - A\mathbf{x}\|^2 + \gamma \|\mathbf{x} - A^\top \mathbf{y}\|^2$$

where $\gamma > 0$ is a (learnable) parameter which interpolates between $A^\dagger \mathbf{y}$ (variance) and $A^\top \mathbf{y}$ (bias)



Towards a new model

- **Hypothesis 1:** A standard U-Net architecture should be enough
- We validate empirically that a prox input improves performance, with *almost no change in the number of learnable parameters*

Configuration	PSNR (dB)	#Params
base (DRUNet backbone)	25.83	32.6M
base + prox	26.64	32.6M

Towards a new model

- **Hypothesis 2:** Backpropagating through $A^T A$ boosts performance

Unrolled networks incorporate A in two ways

- 1) *Gradient step:* $\mathbf{x}^{k+1} = \mathbf{x}^k - \eta \mathbf{A}^T \mathbf{A} \mathbf{x}^k + \mathbf{A}^T \mathbf{y}$
- 2) *Proximal step:* $\mathbf{x}^{k+1} = (\eta \mathbf{A}^T \mathbf{A} + I)^{-1} (\eta \mathbf{A}^T \mathbf{y} + \mathbf{x}^k)$

Prox. is typically computed via conjugate gradient, a Krylov subspace method,

$$\begin{aligned} \text{where } \mathbf{x}^{k+1} &= \text{span} \left\{ (\eta \mathbf{A}^T \mathbf{A} + I)^i (\eta \mathbf{A}^T \mathbf{y} + \mathbf{x}^k) \right\}_{i=1:n} \\ &= \sum_i^N \alpha_i (\mathbf{A}^T \mathbf{A})^i \mathbf{x}^k + \beta_i (\mathbf{A}^T \mathbf{A})^i \mathbf{A}^T \mathbf{y} \end{aligned}$$

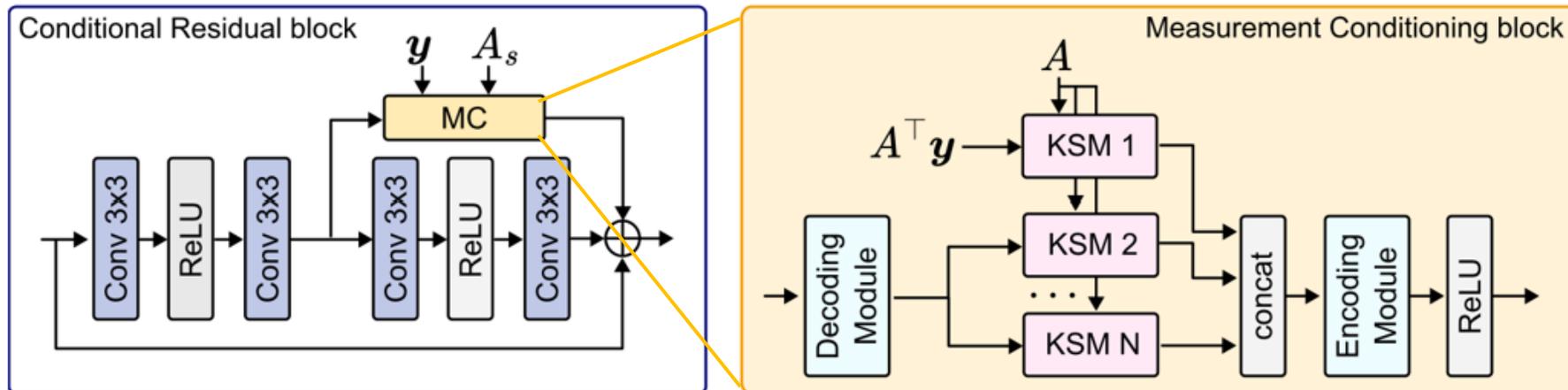
for some scalars $\{(\alpha_i, \beta_i)\}_{i=1:n}$

Towards a new model

- **Hypothesis 2:** Backpropagating through $A^T A$ boosts performance

We generalize the grad and prox steps via a Krylov embedding layer

of order N , written as $\phi(\text{concat} \left[(A^T A)^i \mathbf{x}^k, (A^T A)^i A^T \mathbf{y} \right]_{i=1:N})$ with ϕ being some conv layers.

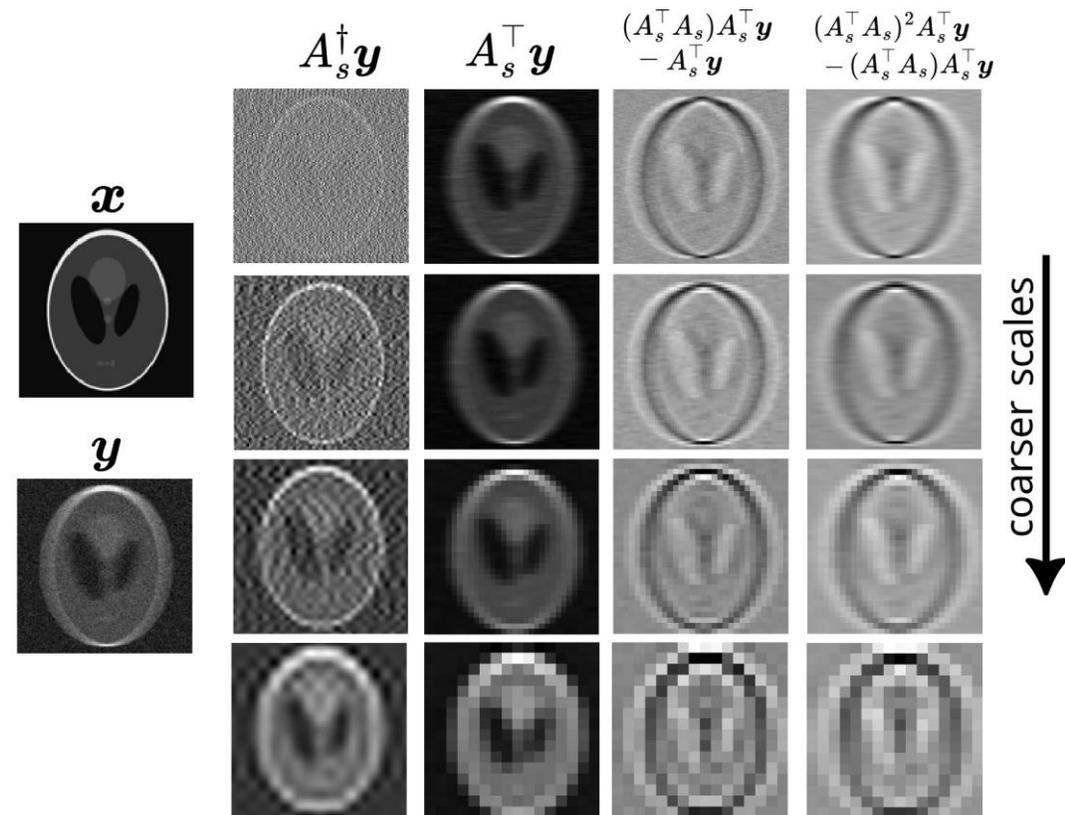


Towards a new model

- **Hypothesis 2:** Backpropagating through $A^T A$ boosts performance
- We can incorporate the operator at every scale of the UNet, as

$A_s = AU_s$ where U_s is an upsampling operator

- The operator becomes better conditioned at coarse scales
- We can often compute $A_s^T A_s$ efficiently without passing through the fine scale



Towards a new model

- **Hypothesis 2:** Backpropagating through $A^T A$ boosts performance
- Again, we see that incorporating the operator improves performance, at additional little parameter cost:

Configuration	PSNR (dB)	#Params
base (DRUNet backbone)	25.83	32.6M
base + prox	26.64	32.6M
base + prox + embed y	26.71	33.4M
base + prox + embed y + Krylov (RAM)	27.61	35.6M

Towards a new model

- **Hypothesis 3:** The network should generalize to multiple A if trained on a family
- We evaluate this claim by training on a single or various tasks:

training on all the tasks performs the best or very close to.

Training tasks	Inpainting	Deblurring	SR ($\times 2$)
Inpainting only	30.84	23.03	26.17
Deblurring only	14.06	25.62	28.02
SR $\times 2$, $\sigma = 0$ only	n/a	21.25	28.60
All three tasks	30.73	25.58	29.79

Towards a new model

- **Hypothesis 3:** The network should generalize to multiple A if trained on a family
- We train the final model on a family of inverse problems:
 - *Deblurring: random motion, confocal, diffraction, gaussian blurs*
 - *Super-resolution: bicubic, bilinear, sinc*
 - *Tomography: random angles*
 - *Single-coil MRI: random masks (uniform, gaussian, etc)*
 - *Inpainting: random masks (missing lines, Bernoulli)*
 - *Demosaicing: Bayer*
 - *Denoising*
- All tasks are corrupted with Gaussian, Poisson and Poisson-Gaussian noise with random noise parameters



Section Navigation

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- Optimization ▾
- Plug-and-Play ▾
- Sampling ▾
- Unfolded ▾
- Patch Priors ▾
- Self-Supervised Learning ▾
- Adversarial Learning ▾
- Advanced ▾

🏠 > Examples

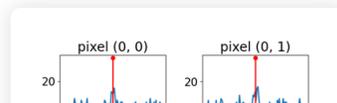
Examples

All the examples have a download link at the end. You can load the example's notebook on [Google Colab](#) and run them by adding the line

```
pip install git+https://github.com/deepinv/deepinv.git#egg=deepinv
```

to the top of the notebook (e.g., [as in here](#)).

Basics





Deep Inverse



Towards a new model

- **Hypothesis 3:** The network should generalize to multiple A if trained on a family
- We consider **multi-channel data:**
 - Single channel images – LIDC-IDRI 1k samples
 - Complex images – FastMRI 2k samples
 - RGB images – LSDIR 90k samples

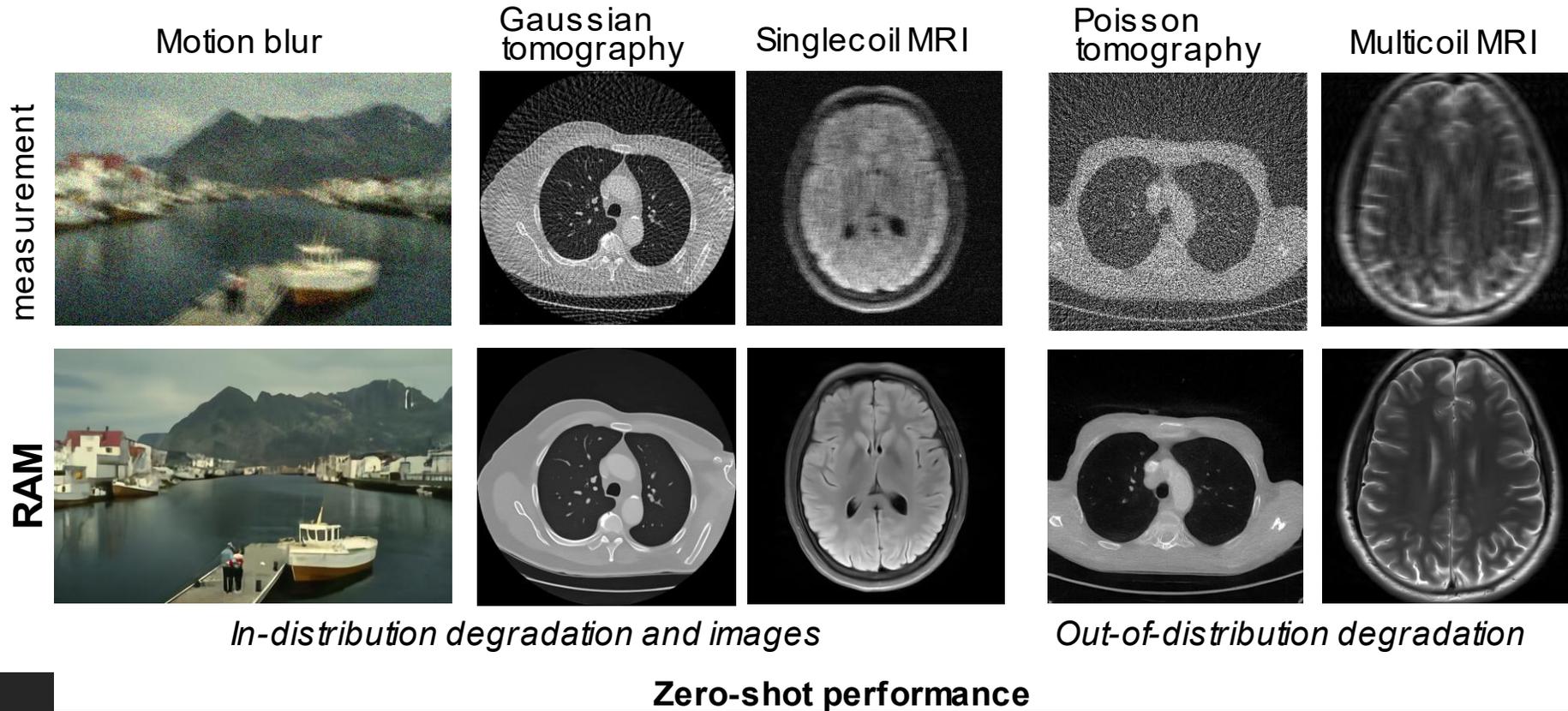
The resulting supervised loss is:

$$\operatorname{argmin}_f \sum_{i=1}^N \mathbb{E}_A \mathbb{E}_{\mathbf{y}_i | x_i, A} w_i \|\mathbf{x}_i - f(\mathbf{y}_i, A)\|_1 \quad \text{with } w_i \propto \frac{\|\mathbf{x}_i\|}{\|A^\top \mathbf{y}_i - \mathbf{x}_i\|}$$

- The network is trained for 200k gradient steps with the Adam optimizer.

Reconstruct Anything Model

- The model obtains very good results on in-distribution tasks, and other out-of-distribution tasks which are close to the training ones.



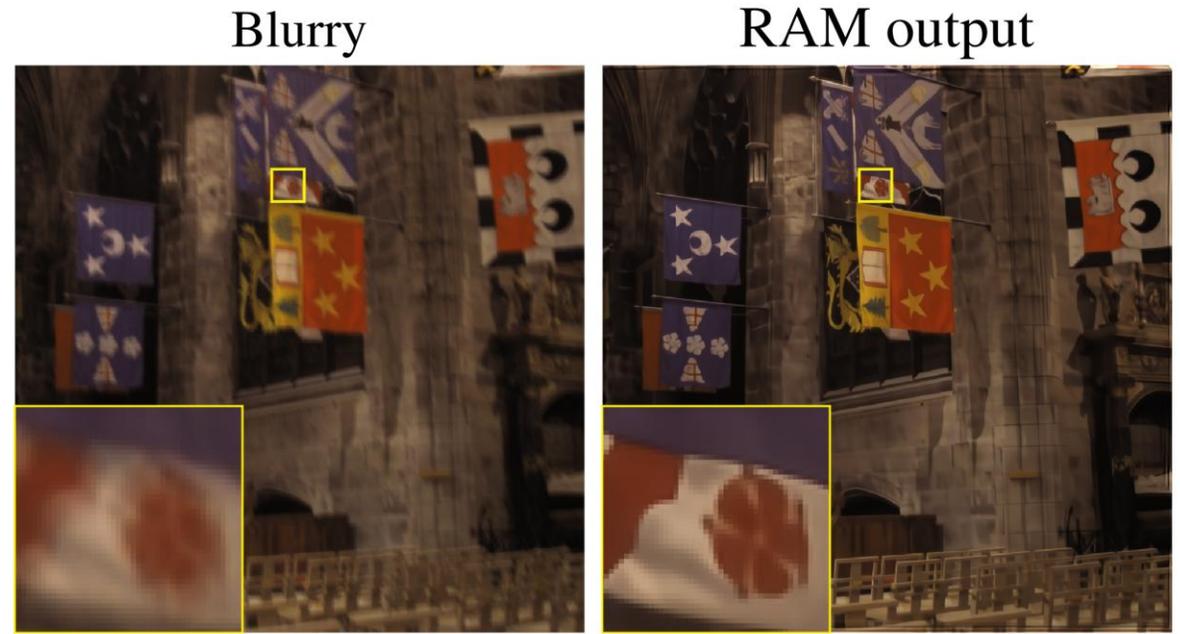
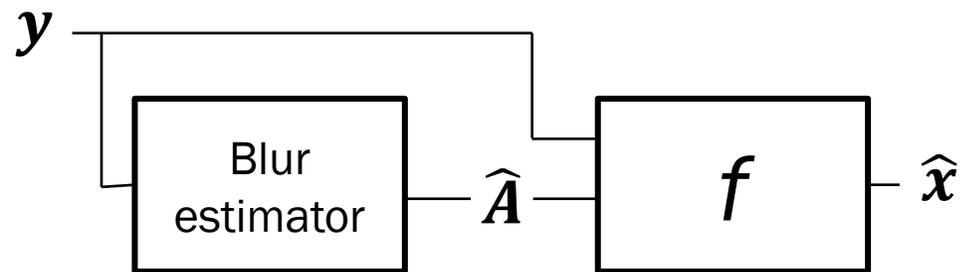
Reconstruct Anything Model

- We perform equally or better than PnP, diffusion and unrolled models, *all of them with many more parameters, and inference time*

Method	Params	CBSD68			Urban100			Div2K			CBSD68			Urban100			Div2K		
		Easy	Med.	Hard	Easy	Med.	Hard	Easy	Med.	Hard	Easy	Med.	Hard	Easy	Med.	Hard	Easy	Med.	Hard
Iterative DPIR	32M	33.45	26.53	24.42	34.39	28.12	24.38	36.32	29.97	27.16	32.10	25.70	22.73	32.53	23.78	20.38	35.23	28.14	24.71
DiffPIR	552M	31.83	25.08	23.80	32.03	27.03	23.60				20.02	24.06	22.51	21.17	22.62	20.27	-	-	-
DDRM*	552M	33.10															34.46	28.25	24.89
PDNet	456K	30.54															32.01	27.29	24.03
Restormer	26M	30.96															30.01	27.20	25.00
uDPIR/t	32M	33.78															34.63	28.45	25.20
uDPIR/u	256M	33.64															34.47	28.37	25.23
RAM	36M	34.04															35.23	28.56	25.16
		GFLOPS			60			2234			2234			360					
		Test memory (BS=1)			52			1213			298			354					
		Train memory (BS=8)			5815			37288			36374			9670					

Non-linear problems

- The model can be extended to do **phase retrieval** and **blind deblurring** *without any retraining*
- For **blind deblurring**, we first estimate a space-varying blur (Carbajal et al., 2023), and then use RAM

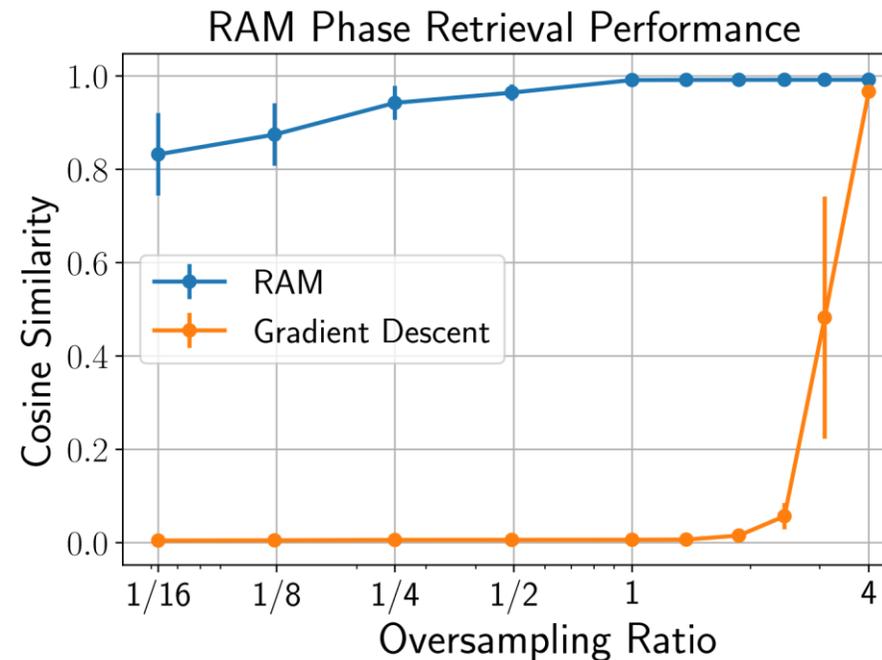


Non-linear problems

- For **phase retrieval** $y = |Ax|^2$, we use RAM inside the Gerchberg Saxton algorithm

for $k = 1, \dots, K$

- $\phi^k = \text{phase}(Ax^k)$
- $x^{k+1} = f(y \circ e^{i\phi^k}, A)$



Self-Supervised Finetuning

- If the inverse problem is ‘far’ from the training distribution, the performance may degrade
- To alleviate this, we can finetune RAM with **measurement data only** by minimizing a self-supervised loss, such as SURE for gaussian noise:

$$\operatorname{argmin}_f \sum_i \|\mathbf{y}_i - Af(\mathbf{y}_i)\|^2 + 2\sigma^2 \operatorname{div}(A \circ f)(\mathbf{y}_i)$$

- SURE can be replaced by UNSURE, Recorrup2Recorrup, etc., according to knowledge of noise
- if A is incomplete/ill-posed, we add the EI, MOI or splitting loss
- See Tachella and Davies “Learning from Noisy and Incomplete Data”, Foundations & Trends in SP, 2026

Self-Supervised Finetuning

- The model can be finetuned with self-supervised losses on up to a single y ($N = 1$) **in minutes**

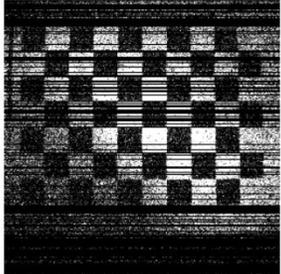
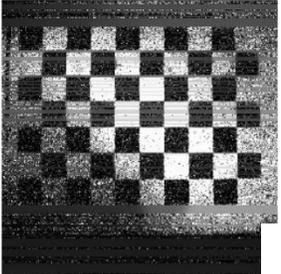
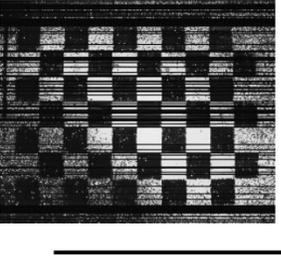
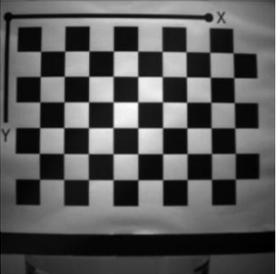
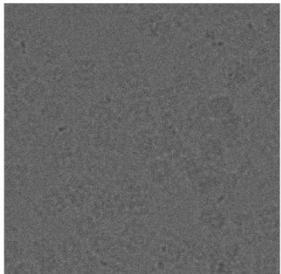
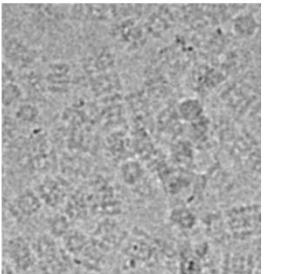
	$A^T y$	Baseline	DRUNet zero-shot		Reference	
SPAD						
				Compressed Sensing	Demosaicing	
Cryo-EM				$N = 1$	43	60
				$N = 10$	80	107
				$N = 100$	228	267
Comp. Sens.						
						

Table 5. Self-supervised finetuning time in seconds.

Uncertainty Quantification

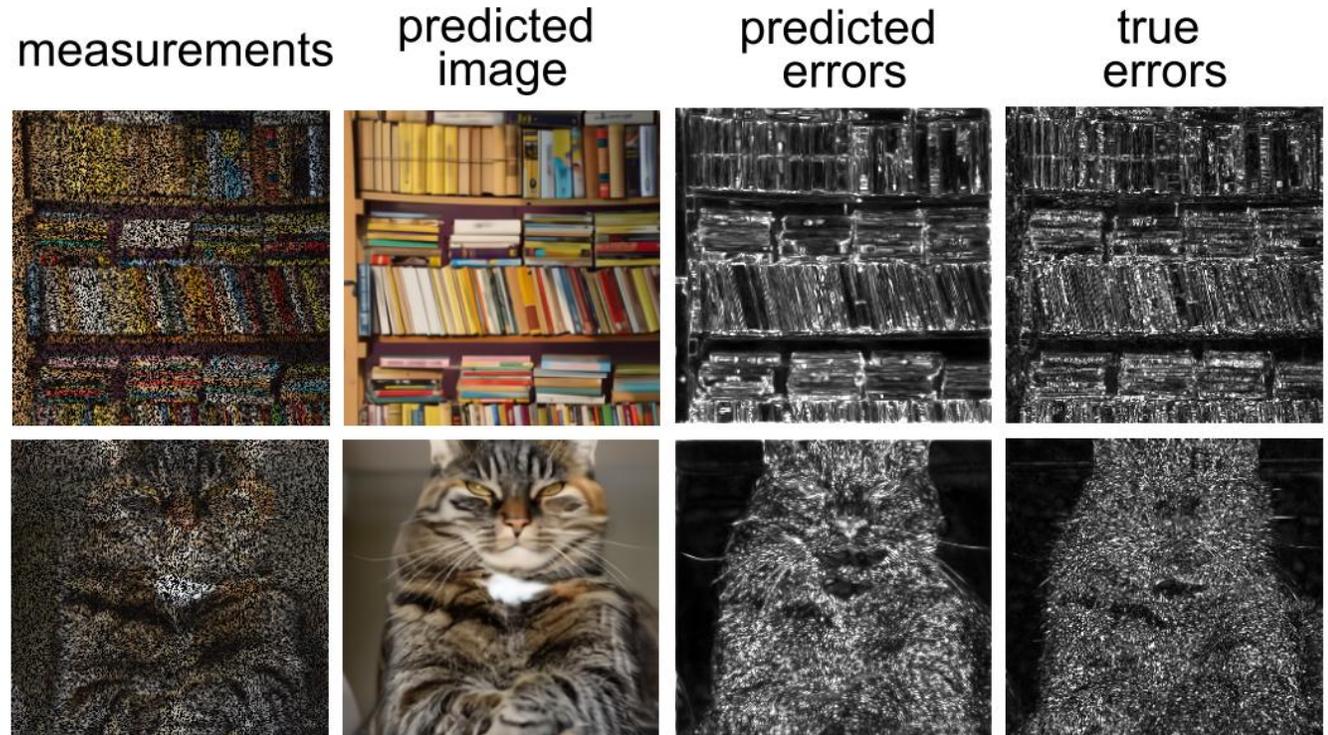
- Since the model is lightweight, we can inference it multiple times to quantify uncertainty
- *Equivariant bootstrap* (Tachella & Pereyra, 2024):
Using $\hat{x} = f(y)$ as 'ground-truth':

For $j = 1, \dots, r$

Sample transform g and noise ϵ_j

Bootstrap $x^j = T_g^{-1} f(AT_g \hat{x} + \epsilon_j)$

Error estimates: $e^j = \hat{x} - x^j$



Conclusions

- **Disadvantages:**

- ~~Trained for a specific task, doesn't generalize to different tasks~~
- ~~Requires a large supervised dataset specific to application~~
- ~~It is a black box, no notion of uncertainty~~

$p(A)$ is low-dimensional

If you want to try out the model:

```
import deepinv as dinv
x = dinv.utils.load_example("butterfly.png")
physics = dinv.physics.Inpainting(...)
y = physics(x)
model = dinv.models.RAM()
x_hat = model(y, physics) # run model
```

Open questions

- Can we iterate it for better perceptual performance (Delbraccio, 2023)?
- What is the **least** number of A evaluations necessary?
- Can we go further by **scaling up** the model?
- Can we express our uncertainty about A ?
- What about 3D or 3D+time?

Stay tuned, RAMv2
coming soon!