Vector Approximate Message Passing

Phil Schniter



Collaborators: Sundeep Rangan (NYU), Alyson Fletcher (UCLA)

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Standard Linear Regression

Goal: Recover $\boldsymbol{x}_o \in \mathbb{R}^N$ from observations $\boldsymbol{y} = \boldsymbol{A} \boldsymbol{x}_o + \boldsymbol{w} \in \mathbb{R}^M$

Examples:

- Compressive Sensing / Medical Imaging:
 - y = measurements $x_o =$ sparse image/signal representation w = sensor noise $A = \Phi \Psi$, Φ measurement operator, Ψ basis
- Wireless communications:
 - y = received samples $x_o =$ finite-alphabet symbols
 - w = noise & interference A = channel operator

Statistics / Machine Learning:

y = experimental outcomesw = model error

 $x_o =$ prediction coefficients A = feature data

Implicit assumptions used in most of this talk

Standard linear regression:

Recover $oldsymbol{x}_o \in \mathbb{R}^N$ from $oldsymbol{y} = oldsymbol{A} oldsymbol{x}_o + oldsymbol{w} \in \mathbb{R}^M$

- A is a known and high dimensional (e.g., $M, N \gtrsim 100$)
- often $N \gg M$ (more unknowns than observations)
- $oldsymbol{w} \sim \mathcal{N}(oldsymbol{0}, au_w oldsymbol{I})$ (additive white Gaussian noise)
- x_o is "structured" (e.g., sparse, natural image, etc.)
- quantities are real-valued (but can be easily extended to complex-valued)

Later will describe extension to generalized linear model: Recover \boldsymbol{x}_o from $\boldsymbol{y} \sim p(\boldsymbol{y}|\boldsymbol{z})$ with hidden $\boldsymbol{z} = \boldsymbol{A}\boldsymbol{x}_o$.

Regularized loss minimization

One way to approach this problem is

$$\widehat{\boldsymbol{x}} = \arg\min_{\boldsymbol{x}} \frac{1}{2} \|\boldsymbol{y} - \boldsymbol{A}\boldsymbol{x}\|^2 + \lambda f(\boldsymbol{x})$$

where

- $\frac{1}{2} \| \boldsymbol{y} \boldsymbol{A} \boldsymbol{x} \|^2$ is the quadratic loss function
- f(x) is a suitably chosen regularizer
 - convex $f(\cdot)$ leads to a convex optimization problem
 - choosing $f(oldsymbol{x}) = \|oldsymbol{x}\|_1$ yields sparse $\widehat{oldsymbol{x}}$
- $\lambda > 0$ is a tuning parameter

Bayesian interpretation:

$$\widehat{oldsymbol{x}} = \mathsf{MAP}$$
 estimate of $oldsymbol{x}$ under

$$\begin{cases} \text{likelihood } p(\boldsymbol{y}|\boldsymbol{x}) = \mathcal{N}(\boldsymbol{y}; \boldsymbol{A}\boldsymbol{x}, \tau_w \boldsymbol{I}) \\ \text{prior } p(\boldsymbol{x}) \propto \exp\left(-\lambda f(\boldsymbol{x})/\tau_w\right) \end{cases}$$

Iterative thresholding

One approach to regularized loss minimization:

initialize $\hat{x}^0 = \mathbf{0}$ for t = 0, 1, 2, ... $v^t = y - A \hat{x}^t$ compute residual $\hat{x}^{t+1} = g(\hat{x}^t + A^{\mathsf{T}} v^t)$ thresholding

where

$$\begin{split} \boldsymbol{g}(\boldsymbol{r}) &= \arg\min_{\boldsymbol{x}} \frac{1}{2} \|\boldsymbol{r} - \boldsymbol{x}\|_{2}^{2} + \lambda f(\boldsymbol{x}) \triangleq \operatorname{prox}_{\lambda f}(\boldsymbol{r}) \\ \|\boldsymbol{A}\|_{2}^{2} < 1 \quad \text{ensures convergence}^{1} \text{ with convex } f(\cdot). \end{split}$$

For example, $f(x) = \|x\|_1$ gives "soft thresholding" $[g(r)]_j = \text{sgn}(r_j) \max\{0, |r_j| - \lambda\}$



¹Daubechies, Defrise, DeMol–CPAM'04

Approximate Message Passing (AMP)

A modification of iterative thresholding:

$$\begin{array}{l} \text{initialize } \widehat{\boldsymbol{x}}^0 = \boldsymbol{0}, \ \boldsymbol{v}^{-1} = \boldsymbol{0} \\ \text{for } t = 0, 1, 2, \dots \\ \boldsymbol{v}^t = \boldsymbol{y} - \boldsymbol{A} \widehat{\boldsymbol{x}}^t + \frac{N}{M} \boldsymbol{v}^{t-1} \big\langle \boldsymbol{g}^{t-1'} (\widehat{\boldsymbol{x}}^{t-1} + \boldsymbol{A}^\mathsf{T} \widehat{\boldsymbol{v}}^{t-1}) \big\rangle & \text{corrected residual} \\ \widehat{\boldsymbol{x}}^{t+1} = \boldsymbol{g}^t (\widehat{\boldsymbol{x}}^t + \boldsymbol{A}^\mathsf{T} \boldsymbol{v}^t) & \text{thresholding} \end{array}$$

where

$$\left\langle \boldsymbol{g}'(\boldsymbol{r}) \right\rangle \triangleq rac{1}{N} \sum_{j=1}^{N} rac{\partial g_j(\boldsymbol{r})}{\partial r_j}$$
 "divergence."

Note:

- The residual v^t now includes an "Onsager correction."
- The thresholding $g^t(\cdot)$ can vary with iteration t.
- Can be derived using Gaussian & Taylor-series approximations of min-sum belief-propagation / message passing.

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Vector Approximate Message Passing

AMP vs ISTA (and FISTA)

Typical convergence behavior with i.i.d. Gaussian A:



Experiment:

- M = 250, N = 500
- Pr $\{x_n \neq 0\} = 0.1$
- SNR= 40 dB
- ISTA, FISTA², AMP all reach same solution: NMSE= -36.8dB
- Convergence to -35dB:
 - ISTA: 2407 iterations
 - FISTA: 174 iterations
 - AMP: 25 iterations

²Beck, Teboulle–JIS'09

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AMP's denoising property

Assumption 1 • $A \in \mathbb{R}^{M \times N}$ is i.i.d. Gaussian • $M, N \to \infty$ s.t. $\frac{M}{N} = \delta \in (0, \infty)$ • $f(x) = \sum_{j=1}^{N} f(x_j)$ with Lipschitz f

Under Assumption 1, something remarkable happens to the input to the thresholder: $^{\rm 3}$

$$oldsymbol{r}^t \triangleq \widehat{oldsymbol{x}}^t + oldsymbol{A}^{\mathsf{T}} oldsymbol{v}^t = oldsymbol{x}_o + \mathcal{N}(oldsymbol{0}, au_r^t oldsymbol{I})$$

with $au_r^t = rac{1}{M} \|oldsymbol{v}^t\|^2 \triangleq \widehat{ au}_r^t$

In other words, r^t is a noisy version of the true signal x_o , where the noise is Gaussian with known variance.

³Bayati,Montanari–TranslT'11

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AMP's state evolution

Define the iteration-t mean-squared error (MSE)

$$\mathcal{E}^t = \frac{1}{N} \operatorname{E} \left\{ \| \widehat{\boldsymbol{x}}^t - \boldsymbol{x}_o \|^2 \right\}.$$

Under Assumption 1, AMP has the following scalar state evolution (SE):

for t = 0, 1, 2, ... $\tau_r^t = \tau_w + \frac{N}{M} \mathcal{E}^t$ $\mathcal{E}^{t+1} = \frac{1}{N} \operatorname{E} \left\{ \left\| \boldsymbol{g}^t \left(\boldsymbol{x}_o + \mathcal{N}(\boldsymbol{0}, \tau_r^t \boldsymbol{I}) \right) - \boldsymbol{x}_o \right\|^2 \right\}$

The rigorous proof⁴ of the SE uses Bolthausen's conditioning trick from the statistical physics literature.

⁴Bayati,Montanari–TranslT'11

Choice of denoiser in AMP

1) LASSO/BPDN

- Goal: compute " $\widehat{\boldsymbol{x}} = rg \max_{\boldsymbol{x}} \frac{1}{2} \| \boldsymbol{y} \boldsymbol{A} \boldsymbol{x} \|^2 + \lambda \| \boldsymbol{x} \|_1$."
- Use $g^t(r) = \text{soft}(r; \alpha \sqrt{\widehat{\tau}_r^t})$, where α has a one-to-one map to λ .

2) Bayesian MMSE

- Goal: compute/approximate MMSE estimate $\widehat{x} = \mathrm{E}\{x|y\}$.
- Suppose $x_o \sim \text{i.i.d.} p(x_j)$ with known $p(x_j)$.
- Use $[\boldsymbol{g}^t(\boldsymbol{r})]_j = \mathrm{E}\left\{x_j \middle| r_j = x_{o,j} + \mathcal{N}(0, \hat{\tau}_r^t)\right\} \dots$ scalar denoising!
- MMSE is achieved when the SE has a unique fixed point!

The choice of denoiser determines the problem solved by AMP.

Choice of denoiser in AMP (cont.)

3) Non-parametric (or model free) estimation

- Goal: compute MMSE estimate without knowing i.i.d. prior $p(x_j)$.
- Assume scalar $GMM(\theta)$ with unknown parameters θ .
- Use MMSE scalar estimator for $GMM(\theta^t)$ at iteration t.
- Use EM algorithm to update θ^t . Details given later...

4) Black-Box Denoisers⁵

- Goal: leverage sophisticated off-the-shelf denoisers like BM3D for natural images or BM4D for image sequences.
- Use $g^t(r) = BM3D(r; \tau_r^t)$.

• Approximate divergence as $\langle g^{t'}(\boldsymbol{r}) \rangle \approx \frac{1}{N} \sum_{j=1}^{N} \frac{g_{j}^{t}(\boldsymbol{r} + \epsilon \boldsymbol{s}) - s_{j}g_{j}^{t}(\boldsymbol{r})}{\epsilon}$ where $\{s_{j}\} \sim \text{i.i.d. uniform } \pm 1$.

⁵Metzler, Maleki, Baraniuk–TIT'16

The limitations of AMP

The good:

- For large i.i.d. sub-Gaussian A, AMP performs provably well.⁶
- Finite-sample analysis shows mild degradation with not-so-large i.i.d. Gaussian A.⁷
- Empirical evidence shows good performance in some other cases (e.g., randomly sub-sampled Fourier A & i.i.d. sparse x)

The bad:

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For general A, AMP can perform poorly
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The ugly:

For general A, AMP may fail to converge!

- lacksim ill-conditioned A
- lacksim non-zero mean A

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<sup>7</sup>Rush, Venkataraman–ISIT'16
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⁶Bayati,Lelarge,Montanari–AAP'15

This talk: Vector AMP

For SLR y = Ax + w, the vector AMP algorithm is⁸

$$\begin{split} & \text{for } t = 0, 1, 2, \dots \\ & \widehat{\boldsymbol{x}}_{1}^{t} = \boldsymbol{g}(\boldsymbol{r}_{1}^{t}; \gamma_{1}^{t}) & \text{denoising} \\ & \alpha_{1}^{t} = \left\langle \boldsymbol{g}'(\boldsymbol{r}_{1}^{t}; \gamma_{1}^{t}) \right\rangle & \text{divergence} \\ & \boldsymbol{r}_{2}^{t} = \frac{1}{1-\alpha_{1}^{t}} \left(\widehat{\boldsymbol{x}}_{1}^{t} - \alpha_{1}^{t} \boldsymbol{r}_{1}^{t} \right) & \text{Onsager correction} \\ & \gamma_{2}^{t} = \gamma_{1}^{t} \frac{1-\alpha_{1}^{t}}{\alpha_{1}^{t}} & \text{precision of } \boldsymbol{r}_{2}^{t} \\ \hline & \widehat{\boldsymbol{x}}_{2}^{t} = \left(\boldsymbol{A}^{\mathsf{T}} \boldsymbol{A} / \widehat{\boldsymbol{\tau}}_{w} + \gamma_{2}^{t} \boldsymbol{I} \right)^{-1} \left(\boldsymbol{A}^{\mathsf{T}} \boldsymbol{y} / \widehat{\boldsymbol{\tau}}_{w} + \gamma_{2}^{t} \boldsymbol{r}_{2}^{t} \right) & \text{LMMSE} \\ & \alpha_{2}^{t} = \frac{\gamma_{2}^{t}}{N} \operatorname{Tr} \left[\left(\boldsymbol{A}^{\mathsf{T}} \boldsymbol{A} / \widehat{\boldsymbol{\tau}}_{w} + \gamma_{2}^{t} \boldsymbol{I} \right)^{-1} \right] & \text{divergence} \\ & \boldsymbol{r}_{1}^{t+1} = \frac{1}{1-\alpha_{2}^{t}} \left(\widehat{\boldsymbol{x}}_{2}^{t} - \alpha_{2}^{t} \boldsymbol{r}_{2}^{t} \right) & \text{Onsager correction} \\ & \gamma_{1}^{t+1} = \gamma_{2}^{t} \frac{1-\alpha_{2}^{t}}{\alpha_{2}^{t}} & \text{precision of } \boldsymbol{r}_{1}^{t+1} \end{split}$$

Note similarities with standard AMP.

⁸Rangan,Schniter,Fletcher-arXiv:1610.03082.

Vector AMP without matrix inverses

Can avoid matrix inverses using an "economy" SVD $A = USV^{T}$:

$$\begin{aligned} & \text{for } t = 0, 1, 2, \dots \\ & \widehat{\boldsymbol{x}}^t = \boldsymbol{g}(\boldsymbol{r}_1^t; \gamma_1^t) & \text{denoising} \\ & \alpha_1^t = \left\langle \boldsymbol{g}'(\boldsymbol{r}_1^t; \gamma_1^t) \right\rangle & \text{divergence} \\ & \boldsymbol{r}_2^t = \frac{1}{1-\alpha_1^t} \left(\widehat{\boldsymbol{x}}^t - \alpha_1^t \boldsymbol{r}_1^t \right) & \text{Onsager} \\ & \frac{\gamma_2^t = \gamma_1^t \frac{1-\alpha_1^t}{\alpha_1^t}}{\alpha_1^t} & \text{precision} \\ & \frac{\alpha_2^t = \frac{1}{N} \sum_j \gamma_2^t / (s_j^2 / \widehat{\tau}_w + \gamma_2^t)}{\mathbf{t}_1^{t-\alpha_2^t} \mathbf{V} \left(\mathbf{S}^2 + \widehat{\tau}_w \gamma_2^t \mathbf{I} \right)^{-1} \mathbf{S} \left(\mathbf{U}^\mathsf{T} \boldsymbol{y} - \mathbf{S} \mathbf{V}^\mathsf{T} \boldsymbol{r}_2^t \right) & 2 \text{ matvec} \\ & \gamma_1^{t+1} = \gamma_2^t \frac{1-\alpha_2^t}{\alpha_2^t} & \text{precision} \end{aligned}$$

Note economy SVD computable with $O(M^3 + M^2N)$ operations.

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1) Can be derived using an approximation of message passing on a factor graph, now with vector-valued variable nodes.

2) Performance can be rigorously characterized by a state-evolution in the high-dimensional limit of certain random A:

 $SVD A = USV^{\mathsf{T}}$

- **U** is deterministic
- S is deterministic
- lacksquare V is uniformly distributed on the group of orthogonal matrices

"A is right-rotationally invariant"

Message-passing derivation of VAMP

• Write joint density as $p(\boldsymbol{x}, \boldsymbol{y}) = p(\boldsymbol{x})p(\boldsymbol{y}|\boldsymbol{x}) = p(\boldsymbol{x})\mathcal{N}(\boldsymbol{y}; \boldsymbol{A}\boldsymbol{x}, \tau_w \boldsymbol{I})$ $p(\boldsymbol{x}) = \mathcal{O}^{\boldsymbol{x}} \mathcal{N}(\boldsymbol{y}; \boldsymbol{A}\boldsymbol{x}, \tau_w \boldsymbol{I})$

• Variable splitting: $p(\boldsymbol{x}_1, \boldsymbol{x}_2, \boldsymbol{y}) = p(\boldsymbol{x}_1)\delta(\boldsymbol{x}_1 - \boldsymbol{x}_2)\mathcal{N}(\boldsymbol{y}; \boldsymbol{A}\boldsymbol{x}_2, \tau_w \boldsymbol{I})$

$$p(\boldsymbol{x}_1) = \underbrace{\boldsymbol{x}_1 \quad \boldsymbol{x}_2}_{\delta(\boldsymbol{x}_1 - \boldsymbol{x}_2)} \quad \boldsymbol{\mathcal{N}}(\boldsymbol{y}; \boldsymbol{A}\boldsymbol{x}_2, \tau_w \boldsymbol{I})$$

Perform⁹ message-passing with messages approximated as $\mathcal{N}(\boldsymbol{\mu}, \sigma^2 \boldsymbol{I})$. An instance of expectation-propagation¹⁰ (EP).

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⁹Rangan,Schniter,Fletcher–arXiv:1610.03082.

¹⁰Minka–Dissertation'01

Free-energy derivation of VAMP

Want to compute posterior density:

$$p(\boldsymbol{x}|\boldsymbol{y}) = \frac{p(\boldsymbol{x})\ell(\boldsymbol{x})}{Z} \text{ with } \begin{cases} p(\boldsymbol{x}) = \text{prior} \\ \ell(\boldsymbol{x}) = N(\boldsymbol{y}; \boldsymbol{A}\boldsymbol{x}, \tau_w \boldsymbol{I}), \text{ likelihood} \\ Z = \int p(\boldsymbol{x})\ell(\boldsymbol{x}) \mathrm{d}\boldsymbol{x}, \text{ partition fxn} \end{cases}$$

but difficult due to high-dimensional integral.

What if we compute the density via

$$rg\min_{b(\boldsymbol{x})} D(b(\boldsymbol{x}) \| p(\boldsymbol{x}|\boldsymbol{y}))$$

where the KL divergence can be written as

$$D(b||p) = \underbrace{D(b||p) + D(b||\ell) + H(b)}_{\text{const}} + \text{const},$$

Gibbs free energy

thus avoiding the partition function Z. Still difficult...

Free-energy derivation of VAMP (cont.)

• What about splitting the belief b(x):

 $\arg\min_{b_1,b_2} \max_q J(b_1,b_2,q) \text{ s.t. } b_1 = b_2 = q$ $J(b_1,b_2,q) = D(b_1 || p) + D(b_2 || \ell) + H(q)$ noting that $D(\cdot || p)$ is convex and $H(\cdot)$ is concave? Still difficult due to the pdf constraint...

So, relax the pdf constraint to moment-matching constraints:

 $b_1 = b_2 = q \quad \longrightarrow \quad \begin{cases} \mathrm{E}\{\boldsymbol{x}|b_1\} = \mathrm{E}\{\boldsymbol{x}|b_2\} = \mathrm{E}\{\boldsymbol{x}|q\} \\ \mathrm{Tr}[\mathrm{Cov}\{\boldsymbol{x}|b_1\}] = \mathrm{Tr}[\mathrm{Cov}\{\boldsymbol{x}|b_2\}] = \mathrm{Tr}[\mathrm{Cov}\{\boldsymbol{x}|q\}] \end{cases}$

An instance of expectation-consistent approximation¹¹ (EC).

¹¹Opper,Winther–NIPS'04, Fletcher,Rangan,Schniter–ISIT'16

Free-energy derivation of VAMP (cont.)

The stationary points of the EC optimization are

$$egin{aligned} b_1(oldsymbol{x}) &\propto p(oldsymbol{x}) \mathcal{N}(oldsymbol{x};oldsymbol{r}_1;oldsymbol{I}/\gamma_1) \ b_2(oldsymbol{x}) &\propto \ell(oldsymbol{x}) \mathcal{N}(oldsymbol{x};oldsymbol{r}_2;oldsymbol{I}/\gamma_2) \ q(oldsymbol{x}) &= \mathcal{N}(oldsymbol{x};\widehat{oldsymbol{x}};oldsymbol{I}/\eta) \end{aligned}$$

for parameters $m{r}_1, \gamma_1, m{r}_2, \gamma_2, \widehat{m{x}}, \eta$ that satisfy

$$\widehat{\boldsymbol{x}} = \mathrm{E}\{\boldsymbol{x}|b_1\} = \mathrm{E}\{\boldsymbol{x}|b_2\} = \mathrm{E}\{\boldsymbol{x}|q\}$$
$$1/\eta = \frac{1}{N}\operatorname{Tr}[\operatorname{Cov}\{\boldsymbol{x}|b_1\}] = \frac{1}{N}\operatorname{Tr}[\operatorname{Cov}\{\boldsymbol{x}|b_2\}] = \frac{1}{N}\operatorname{Tr}[\operatorname{Cov}\{\boldsymbol{x}|q\}].$$

 Can then construct algorithms whose fixed points coincide with these stationary points (e.g., EC, ADATAP,¹² S-AMP¹³). But convergence is not guaranteed.

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¹²Opper,Winther–NC'00

¹³Cacmak,Winter,Fleury–ITW'14

Putting things in perspective

- The aforementioned belief-propagation and free-energy derivations are both well known and heuristic (in general).
 - The resulting algorithms may not converge to their fixed points
 - S-AMP diverges with mildly ill-conditioned matrices
 - Even if they do converge, the accuracy of the fixed points is unclear:
 - EP generally suboptimal due to approximation of messages
 - EC generally suboptimal due to approximation of constraint
- The important question is whether/when a given heuristic can be rigorously analyzed and proven to work well.

AMP rigorous analyzed under large i.i.d. Gaussian A and Bayes optimal under certain combinations of $\{p(x), \ell(x)\}$.

VAMP state evolution

VAMP has a rigorous SE¹⁴

Assuming empirical convergence of $\{s_j\} \rightarrow S$ and $\{(r_{1,j}^0, x_{o,j})\} \rightarrow (R_1^0, X_o)$ and Lipschitz continuity of g and g', the VAMP-SE under $\hat{\tau}_w = \tau_w$ is as follows:

$$\begin{aligned} & \text{for } t = 0, 1, 2, \dots \\ & \mathcal{E}_1^t = \mathrm{E}\left\{\left[g\left(X_o + \mathcal{N}(0, \tau_1^t); \overline{\gamma}_1^t\right) - X_o\right]^2\right\} & \text{MSE} \\ & \overline{\alpha}_1^t = \mathrm{E}\left\{g'(X_o + \mathcal{N}(0, \tau_1^t); \overline{\gamma}_1^t)\right\} & \text{divergence} \\ & \overline{\gamma}_2^t = \overline{\gamma}_1^t \frac{1 - \overline{\alpha}_1^t}{\overline{\alpha}_1^t}, \quad \tau_2^t = \frac{1}{(1 - \overline{\alpha}_1^t)^2} \left[\mathcal{E}_1^t - \left(\overline{\alpha}_1^t\right)^2 \tau_1^t\right] \\ & \mathcal{E}_2^t = \mathrm{E}\left\{\left[S^2/\tau_w + \overline{\gamma}_2^t\right]^{-1}\right\} & \text{MSE} \\ & \overline{\alpha}_2^t = \overline{\gamma}_2^t \mathrm{E}\left\{\left[S^2/\tau_w + \overline{\gamma}_2^t\right]^{-1}\right\} & \text{divergence} \\ & \overline{\gamma}_1^{t+1} = \overline{\gamma}_2^t \frac{1 - \overline{\alpha}_2^t}{\overline{\alpha}_2^t}, \quad \tau_1^{t+1} = \frac{1}{(1 - \overline{\alpha}_2^t)^2} \left[\mathcal{E}_2^t - \left(\overline{\alpha}_2^t\right)^2 \tau_2^t\right] \end{aligned}$$

More complicated expressions for \mathcal{E}_2^t and $\overline{\alpha}_2^t$ apply when $\widehat{\tau}_w \neq \tau_w$.

¹⁴Rangan, Schniter, Fletcher-arXiv:1610.03082

Connections to the replica prediction

- The replica method from statistical physics is often used to characterize the average behavior of large disordered systems.
- Although not fully rigorous, replica predictions are usually correct.
- For SLR under large right-rotationally invariant A and matched priors, The MMSE $\mathcal{E}_1(\overline{\gamma}_1)$ should satisfy the fixed-point equation¹⁵

$$\overline{\gamma}_1 = R_{\boldsymbol{A}^{\mathsf{T}}\boldsymbol{A}/\tau_w}(-\mathcal{E}_1(\overline{\gamma}_1)),$$

where $\mathbf{R}_{\mathbf{C}}(\cdot)$ denotes the R-transform of matrix \mathbf{C} and $\mathcal{E}_1(\overline{\gamma}_1) \triangleq \mathrm{E}\left\{ \left[g_{mmse} \left(X_o + \mathcal{N}(0, 1/\overline{\gamma}_1); \overline{\gamma}_1 \right) - X_o \right]^2 \right\}.$

- It can be shown that VAMP's matched SE obeys the above equation.
- Thus, if the replica prediction is correct, then VAMP's estimates will be MMSE whenever the replica fixed-point equation has a unique solution.

¹⁵Tulino,Caire,Verdu,Shamai–TIT'13

Experiment with Matched Priors I



$$\begin{split} N &= 1024 \\ M/N &= 0.5 \end{split}$$

 $A = U \operatorname{Diag}(s) V^{\mathsf{T}}$ U, V drawn uniform $s_n/s_{n-1} = \phi \ \forall n$ $\phi \text{ determines } \kappa(A)$

 $X_o \sim \text{Bernoulli-Gaussian}$ $\Pr\{X_0 \neq 0\} = 0.1$

 $\mathsf{SNR} = 40\mathsf{dB}$

Note robustness w.r.t. condition number of A.

Experiment with Matched Priors II



$$\begin{split} N &= 1024 \\ M/N &= 0.5 \end{split}$$

 $A = U \operatorname{Diag}(s) V^{\mathsf{T}}$ U, V drawn uniform $s_n/s_{n-1} = \phi \ \forall n$ $\phi \text{ determines } \kappa(A)$

 $X_o \sim \text{Bernoulli-Gaussian}$ $\Pr\{X_0 \neq 0\} = 0.1$

 $\mathsf{SNR} = 40\mathsf{dB}$

Note convergence speed relative to (damped) EM-AMP.

Non-parametric (model-free) regression

So far we considered recovering x_o from

$$\boldsymbol{y} = \boldsymbol{A}\boldsymbol{x}_o + \boldsymbol{w}, \quad \boldsymbol{x}_o \sim p(\boldsymbol{x}), \quad \boldsymbol{w} \sim \mathcal{N}(\boldsymbol{0}, \tau_w \boldsymbol{I}),$$

when $p(\boldsymbol{x})$ and τ_w are known.

Can we learn τ_w ? Yes, through an EM procedure.¹⁶ Can we learn $p(\boldsymbol{x})$? Yes if $p(\boldsymbol{x}) = \prod_j p(x_j)$.

• Why is $p(x_j)$ learnable with VAMP?

- Recall that $m{r}_1^t = m{x}_o + \mathcal{N}(m{0}, au_1^tm{I}).$
- Thus \boldsymbol{r}_1^t contains i.i.d. samples of $p(x_j) * \mathcal{N}(x_j; 0, \tau_1^t)$.
- Should be able to deconvolve $p(x_j)$ from the empirical distribution of r_1^t .
- A practical method: Model $p(x_j) = \text{GMM}(x_j; \theta_x)$. Learn parameters θ_x using EM.

¹⁶Fletcher,Schniter-arXiv:1602.08207

EM-VAMP

- Recall $\begin{cases} \text{prior } p(\boldsymbol{x}; \boldsymbol{\theta}_x) \\ \text{likelihood } \ell(\boldsymbol{x}; \tau_w) \end{cases} \rightarrow \text{Learn parameters } \boldsymbol{\theta} \triangleq (\boldsymbol{\theta}_x, \tau_w). \end{cases}$
- EM: iterate $Q(\boldsymbol{\theta}; \widehat{\boldsymbol{\theta}}^k) = \int p(\boldsymbol{x} | \boldsymbol{y}; \widehat{\boldsymbol{\theta}}^k) \ln p(\boldsymbol{x}, \boldsymbol{y}; \boldsymbol{\theta}) d\boldsymbol{x}$ "expectation" $\widehat{\boldsymbol{\theta}}^{k+1} = \arg \max_{\boldsymbol{\theta}} Q(\boldsymbol{\theta}; \widehat{\boldsymbol{\theta}}^k)$ "maximization" which uses the posterior $p(\boldsymbol{x} | \boldsymbol{y}; \widehat{\boldsymbol{\theta}}^k)$ in the E step.

With VAMP's posterior approx, EM is an alternating approach to

$$\min_{b_1,b_2,\boldsymbol{\theta}} \max_{q} D(b_1 \| p(\boldsymbol{\theta}_{\boldsymbol{x}})) + D(b_2 \| \ell(\tau_w)) + H(q)$$
s.t.
$$\begin{cases} \mathrm{E}\{\boldsymbol{x}|b_1\} = \mathrm{E}\{\boldsymbol{x}|b_2\} = \mathrm{E}\{\boldsymbol{x}|q\} \\ \mathrm{Tr}[\mathrm{Cov}\{\boldsymbol{x}|b_1\}] = \mathrm{Tr}[\mathrm{Cov}\{\boldsymbol{x}|b_2\}] = \mathrm{Tr}[\mathrm{Cov}\{\boldsymbol{x}|q\}] \end{cases}$$

Can make faster by putting θ optimization in the inner loop.

Experiment with Learned Parameters I

Learning both τ_w and θ_x :



EM-VAMP achieves oracle performance at all condition numbers.

Experiment with Learned Parameters II

Learning both τ_w and θ_x :



M/N = 0.5

 $\boldsymbol{A} = \boldsymbol{U} \operatorname{Diag}(\boldsymbol{s}) \boldsymbol{V}^{\mathsf{T}}$ U, V drawn uniform $s_n/s_{n-1} = \phi \ \forall n$ ϕ determines $\kappa(A)$

 $X_{0} \sim \mathsf{Bernoulli-Gaussian}$ $\Pr\{X_0 \neq 0\} = 0.1$

SNR = 40 dB

EM-VAMP nearly as fast as VAMP and much faster than EM-AMP.

Noiseless Image Recovery with BM3D

		PSNR	time
10%	L1-AMP	17.7 dB	0.5 s
	L1-VAMP	17.6 dB	0.5 s
	BM3D-AMP	25.2 dB	10.1 s
	BM3D-VAMP	25.2 dB	10.4 s
20%	L1-AMP	20.2 dB	1.0 s
	L1-VAMP	20.2 dB	0.9 s
	BM3D-AMP	30.0 dB	8.8 s
	BM3D-VAMP	30.0 dB	8.5 s
30%	L1-AMP	22.4 dB	1.6 s
	L1-VAMP	22.4 dB	1.4 s
	BM3D-AMP	32.5 dB	8.6 s
	BM3D-VAMP	32.5 dB	8.2 s
40%	L1-AMP	24.6 dB	2.3 s
	L1-VAMP	24.8 dB	1.8 s
	BM3D-AMP	35.1 dB	9.1 s
	BM3D-VAMP	35.2 dB	8.5 s
50%	L1-AMP	27.0 dB	3.1 s
	L1-VAMP	27.2 dB	2.3 s
	BM3D-AMP	37.4 dB	9.8 s
	BM3D-VAMP	37.7 dB	8.8 s

Avg results for recovering 128×128 lena, barbara, boat, fingerprint, house, and peppers from $y = Ax_o$ with i.i.d. Gaussian A at various sampling ratios.

All algorithms use 20 iterations and learn the noise variance τ_w .

VAMP slightly outperforms AMP in accuracy and runtime.

Noiseless Image Recovery with BM3D (cont.)



Generalized linear models

- Until now we have considered SLR, $y = Ax_o + w$.
- VAMP can also support the generalized linear model (GLM) $\boldsymbol{u} \sim p(\boldsymbol{u}|\boldsymbol{z})$ with hidden $\boldsymbol{z} = \boldsymbol{A}\boldsymbol{x}_{o}$

which supports, e.g.,

- $y_i = z_i + w_i$: additive, possibly non-Gaussian noise
- $y_i = \operatorname{sgn}(z_i + w_i)$: binary classification / one-bit sensing
- $y_i = |z_i + w_i|$: phase retrieval in noise
- Poisson y_i : photon-limited imaging

Trick:
$$z = Ax \quad \Leftrightarrow \quad \underbrace{0}_{\widetilde{z}} = \underbrace{[A-I]}_{\widetilde{A}} \underbrace{\begin{bmatrix} x \\ z \end{bmatrix}}_{\widetilde{x}}$$

One-bit compressed sensing / Probit regression

Learning both τ_w and θ_x :



VAMP and EM-VAMP robust to ill-conditioned A.

Phil Schniter (Ohio State)

Vector Approximate Message Passing

iTWIST — Aug'16 32 / 35

One-bit compressed sensing / Probit regression

Learning both τ_w and θ_x :



EM-VAMP mildly slower than VAMP but much faster than damped AMP.

Conclusions

AMP exhibits some remarkable properties

- low cost-per-iteration and relatively few iterations to convergence,
- intermediate estimates of form $m{r}^t = m{x}_o + \mathcal{N}(m{0}, au_r^tm{I})$,
- rigorous state evolution,
- easy tuning of prior & likelihood,
- compatibility with plug-in denoisers like BM3D,

but those properties are guaranteed only under large i.i.d. Gaussian A.

Vector AMP has the same properties, but for a much larger class of A.

Ongoing work: analysis of EM procedure, bilinear extensions, connections with deep learning, various applications...

Thanks for listening!