

Imaging with multi-source experiments

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Outline

The problem

Standard solution techniques

Reformulation and some new insight

Numerical experiments

Extensions to moving sources/receivers

The problem

We consider an inverse problem solved by the optimization of

$$\begin{aligned} \min_m \quad & \frac{1}{2} \sum_j \|P^\top u_j - d_j\|^2 + R(m) \\ \text{s.t} \quad & A(m)u_j = Q_j \quad j = 1, \dots, N_s \end{aligned}$$

- $A(m)$ - a discretization of a parameter dependent differential operator
- Q_j - source
- P - observation matrix
- u_j - field
- d_j - data vector
- $R(m)$ - regularization

The problem

$$\begin{aligned} \min_m \quad & \frac{1}{2} \sum_j \|P^\top u_j - d_j\|^2 + R(m) \\ \text{s.t} \quad & A(m)u_j = Q_j \quad j = 1, \dots, N_s \end{aligned}$$

- The number of sources is **LARGE**
- The discretized PDE $A(m)$ is large and ill-conditioned
- Special structure - all sources share the same receivers

The problem

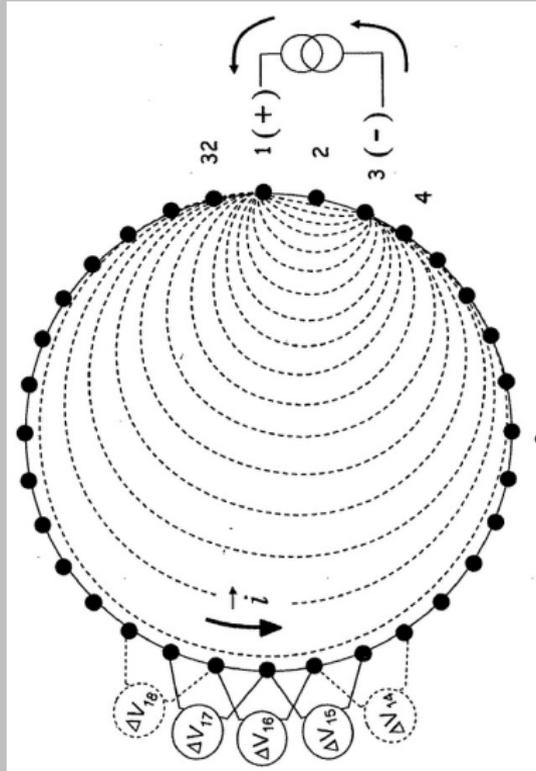
Medical examples include

- Electrical Impedance Tomography (EIT)
- Magnetic Induction Tomography (MIT)
- Microwave Imaging
- 3D Ultrasound

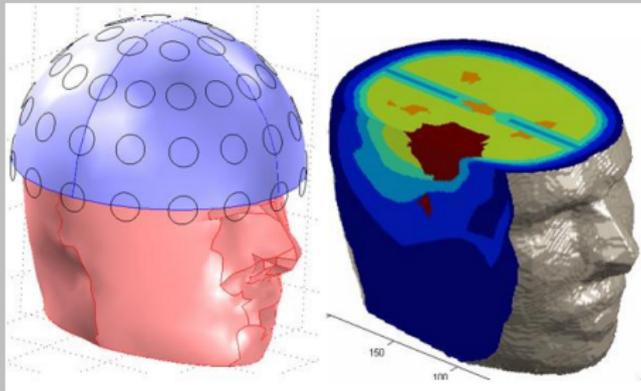
Geophysical examples include

- DC resistivity
- Electromagnetics
- Seismic imaging

Electrical Impedance Tomography



Magnetic Induction Tomography



Solution technique

$$\begin{aligned} \min_m \quad & \frac{1}{2} \sum_j \|P^\top u_j - d_j\|^2 + R(m) \\ \text{s.t.} \quad & A(m)u_j = Q_j \quad j = 1, \dots, N_s \end{aligned}$$

Impossible to store all fields, use unconstrained approach [H. Oldenburg, Ascher 2000]

$$\min_m \mathcal{J}(m) = \frac{1}{2} \sum_{j=1}^{N_s} \|P^\top A(m)^{-1} Q_j - d_j\|^2 + R(m)$$

Solution technique

$$\min_m \mathcal{J}(m) = \frac{1}{2} \sum_{j=1}^{N_s} \|P^\top A(m)^{-1} Q_j - d_j\|^2 + R(m)$$

The gradient

$$\sum_{j=1}^{N_s} -G(m, u_j)^\top A(m)^{-\top} P (P^\top A(m)^{-1} Q_j - d_j) + \nabla R(m)$$

where $G(m, u_j) = \nabla_m (A(m)u_j)$

Solution technique

Computing the misfit and gradient

- Set $\text{misfit} = 0$ $\nabla \text{misfit} = 0$
- For $j = 1, \dots, N_s$
 - Solve $A(m)u_j = Q_j$
 - $r_j = P^\top u_j - d_j$
 - $\text{misfit} \leftarrow \text{misfit} + r_j^\top r_j$
 - Solve $A^\top \lambda = P r_j$
 - $\nabla \text{misfit} \leftarrow \nabla \text{misfit} - G^\top \lambda$

Computation of misfit and its derivative require $2N_s$ solutions of the forward/adjoint problem.

For large scale problems difficult if not impossible

Solution technique

- Set $\text{misfit} = 0$ $\nabla \text{misfit} = 0$
- For $j = 1, \dots, N_s$
 - Solve $\overbrace{A(m)u_j = Q_j}^{\text{expensive!}}$
 - $r_j = P^\top u_j - d_j$
 - $\text{misfit} \leftarrow \text{misfit} + r_j^\top r_j$
 - Solve $\overbrace{A^\top \lambda = Pr_j}^{\text{expensive}}$
 - $\nabla \text{misfit} \leftarrow \nabla \text{misfit} - G^\top \lambda$

Solution technique

Current methods to deal with multiple rhs

- Factor the system if possible [Pratt, 2000, H. & Oldenburg, 2006]
- Almost factor the system (ILU, domain decomposition with large domains) [Ascher & van den Doel, 2009]
- Recycle right hand sides [Kilmer & de Sturler 2006]

Issues - complexity, storage

Solution technique

For the computation of a Gauss-Newton step similar calculations are needed.

Typically, avoid Gauss-Newton and use L-BFGS, nonlinear CG and steepest descent (storage).

Converges can be slow

A different point of view

The difficulty: computing the misfit.

Can we do this cheaper?

A different point of view

The difficulty: computing the misfit.

Can we do this cheaper?

Recall that

$$\begin{aligned}\text{misfit} &= \frac{1}{2} \sum_j \|P^\top A(m)^{-1} Q_j - d_j\|^2 \\ &= \frac{1}{2} \|P^\top A(m)^{-1} Q - D\|_F^2 = \\ &= \frac{1}{2} \text{trace} \left((P^\top A(m)^{-1} Q - D)^\top (P^\top A(m)^{-1} Q - D) \right) \\ &= \frac{1}{2} \mathbf{E}_w \| (P^\top A(m)^{-1} Q - D) w \|^2\end{aligned}$$

where w is a random variable with

$$\mathbf{E}(w) = 0 \quad \text{Cov}(w) = I$$

A different point of view

The original (deterministic) optimization problem is therefore equivalent to the (stochastic) optimization problem

$$\begin{aligned}\hat{m} &= \arg \min_m \frac{1}{2} \sum_{j=1}^{N_s} \|P^\top A(m)^{-1} Q_j - d_j\|^2 + R(m) \\ &= \arg \min_m \frac{1}{2} \mathbf{E}_w \|P^\top A(m)^{-1} Qw - Dw\|^2 + R(m)\end{aligned}$$

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So What?

A different point of view

$$\hat{m} = \arg \min_m \frac{1}{2} \mathbf{E}_w \|P^\top A(m)^{-1} Qw - Dw\|^2 + R(m)$$

- This is a stochastic optimization problem [Shapiro 09] and has been treated extensively in the literature
- We can capitalize on the structure of the problem to obtain cheap algorithms

Main point - Given a realization w_i a **Single** PDE solve is required to evaluate $\text{misfit}(m; w_i)$

A different point of view

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Two methods for stochastic optimization [Shapiro 05]

- SAA - Sample Average Approximation
Discretize the Expectation THEN optimize
- SA - Stochastic Approximation
Optimize AND Discretize

Stochastic Optimization

SAA - Sample Average Approximation

Approximate the expectation using Monte-Carlo

$$\min_m \sum_j \frac{1}{2} \|P^\top A(m)^{-1} Q w_j - D w_j\|^2 + R(m)$$

Stochastic Optimization

SAA - Sample Average Approximation

Approximate the expectation using Monte-Carlo

$$\min_m \sum_j \frac{1}{2} \|P^\top A(m)^{-1} Q w_j - D w_j\|^2 + R(m)$$

SA - Stochastic approximation

for $j = 1, \dots$

$$\begin{aligned} \hat{s} &= \arg(\text{aprox}) \min \frac{1}{2} \|P^\top A(m_j + s)^{-1} Q w_j - D w_j\|^2 + R(s) \\ s_{j+1} &= \text{average}(s_{1:j}, \hat{s}) \end{aligned}$$

end

SAA - Sample Average Approximation

Approximate the expectation using Monte-Carlo

$$\min_m \mathcal{J}(m; w) = \sum_j \frac{1}{2} \|P^\top A(m)^{-1} Q w_j - D w_j\|^2 + R(m)$$

- How to pick w ?
- How many w 's?

Any distribution with $\mathbf{E}(w) = 0$ and $\mathbf{Cov}(w) = I$ has

$$\mathbf{E}(w^\top H w) = \text{trace}(H)$$

Choose the distribution such that [Hutchinson 93]

$$\text{Var}(w^\top H w) \rightarrow \min$$

$$w = \text{rand}(\pm 1)$$

SAA - Sample Average Approximation

Approximate the expectation using Monte-Carlo

$$\min_m \mathcal{J}(m; w) = \sum_j \frac{1}{2} \|P^\top A(m)^{-1} Q w_j - D w_j\|^2 + R(m)$$

- How to pick w ?
- How many w 's?

The number of w 's depends on the variability of the unbiased estimator

$$\mathbf{E}(\mathcal{J}(m; w)) \approx \frac{1}{N} \sum_{j=1}^N \mathcal{J}(m; w_j)$$

and the accuracy we would like to obtain.

SAA - Example

Generate $A(m)$ by discretizing the PDE

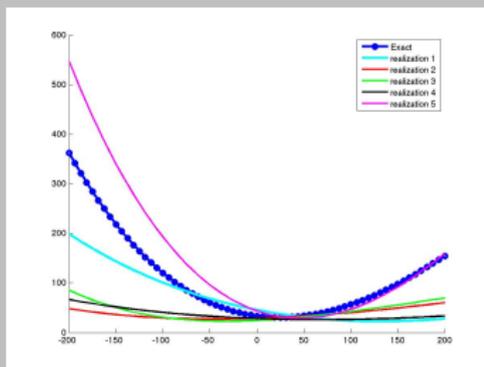
$$\nabla \cdot \exp(m) \nabla u$$

Assume 1089 sources (right hand sides) and 1089 receivers
Look at

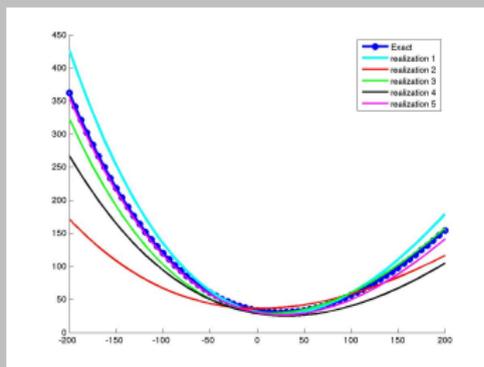
$$f(\alpha) = \frac{1}{2N} \sum_j \|P^\top A(m + \alpha s)^{-1} Q - D\|_F^2$$

SAA - Sample Average Approximation

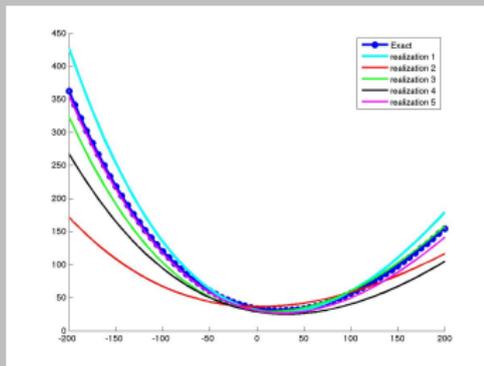
1 vector



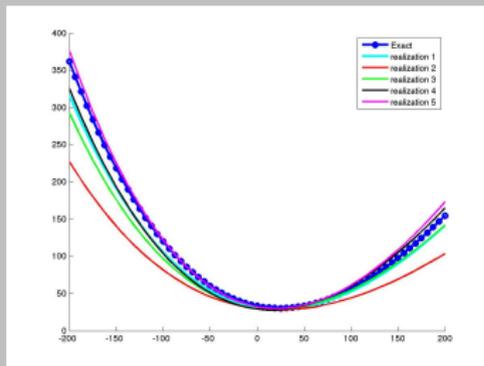
10 vectors



25 vectors



50 vectors



SAA - Sample Average Approximation

Controlling the quality of the approximation can be done by repeating the minimization with different samples

For our problems we have found that a small sample size may be sufficient [Also Golub & Bai, 99, Golub & von Matt 98]

Advantage of SAA - separate the stochastic part from the optimization

Disadvantage of SAA - number of realization may be too large

SA - Stochastic Approximation

Stochastic approximation

for $j = 1, \dots$

$$\hat{s} = \arg(\text{aprox}) \min \frac{1}{2} \|P^\top A(m_j + s)^{-1} Q w_j - D w_j\|^2 + R(s)$$
$$s_{j+1} = \text{average}(s_{1:j}, \hat{s})$$

end

Questions

- How approximate?
- What methods can be used?
- Convergence?

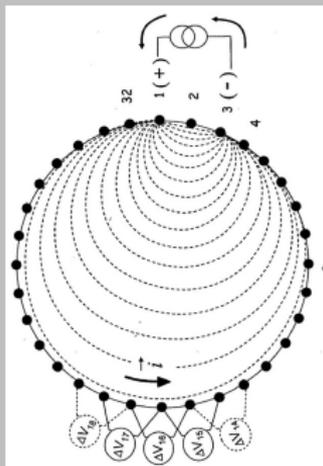
SA - Stochastic Approximation

- Proof that it works only on various flavors of steepest descent (recent work [Nemirovski and Shaoiro])
- Observed in practice - works well for L-BFGS and Gauss-Newton [Schraudolph, Yu & Gunter 10]
- Much interest in machine learning (online algorithms)

An illustrative example

Model problem - Electrical Impedance Tomography

$$\min_m \left\| P^T (G^T S(m) G)^{-1} Q - D \right\|_F^2 + \frac{\alpha}{2} \|Gm\|^2$$



An illustrative example

- Assume 1089 sources and 1089 receivers.
- Mesh $32 \times 32 \times 32$
- Number of unknowns (fields) 35,684,352
- Computation of full forward, roughly, 2 hours

Use standard, SAA and SA to solve the problem.

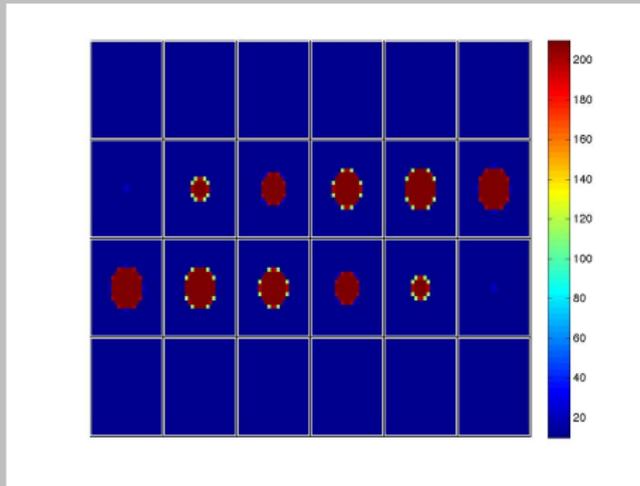
An illustrative example

Method	# iterations	# rhs/iter	Cost (pde solves)
Standard	37	1089	80,293
SSA	45	10	2757
SA	453	1	923

Computational saving of factor 100

$$\begin{aligned}\|m_{\text{SAA}} - m_{\text{Standard}}\|^2 / \|m_{\text{Standard}}\|^2 &= 2.1 \times 10^{-2} \\ \|m_{\text{SA}} - m_{\text{Standard}}\|^2 / \|m_{\text{Standard}}\|^2 &= 3.2 \times 10^{-2}\end{aligned}$$

Inversion parameters

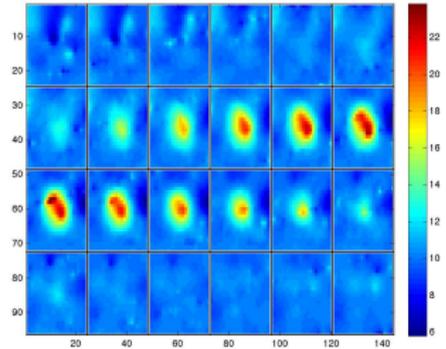
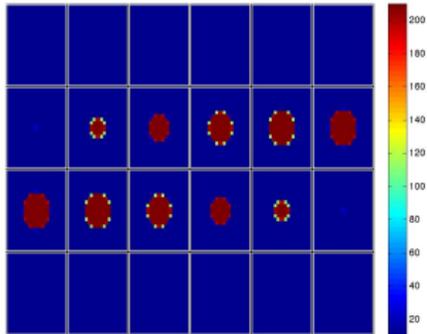


$$\alpha = 10^{-4}$$

Starting model - $m = 10^{-2} \text{S/m}$

Converges - solution does not change between iterations

Recovered solution

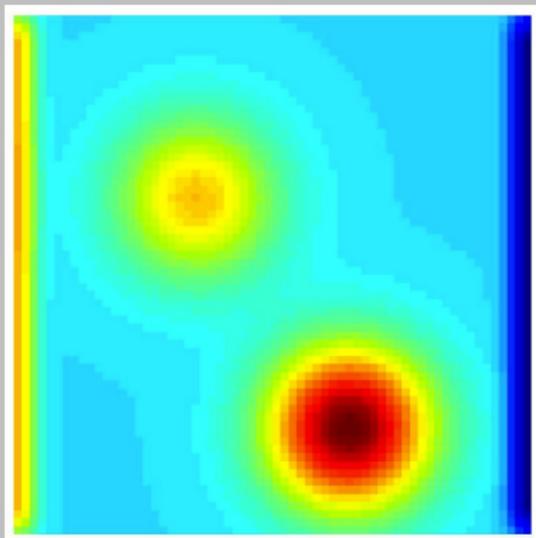


Sequential SAA

How to choose the size of the random batch?

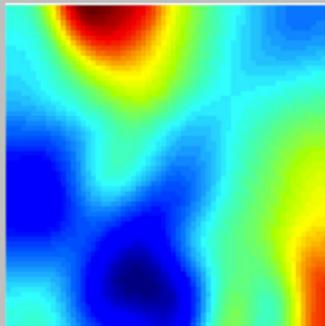
Use continuation in batch-size

Example - EIT in 2D

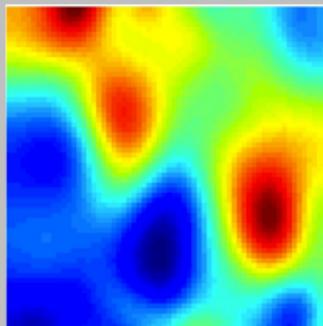


Sequential SAA

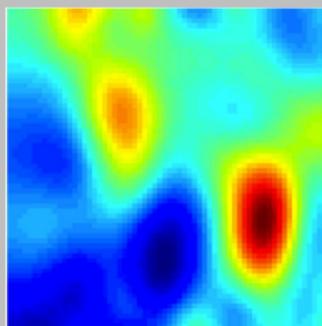
1 src



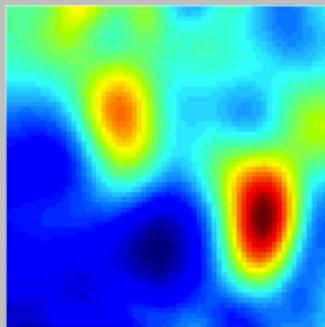
2 src



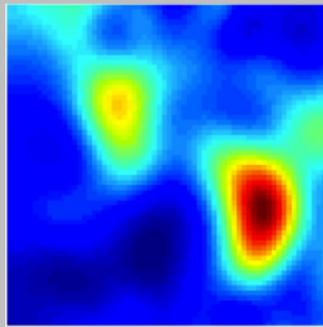
3 src



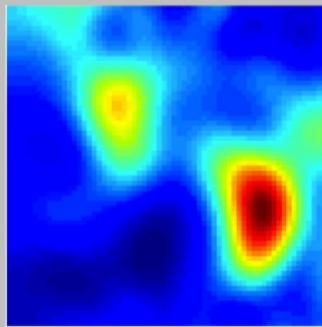
4 src



5 src



6 src



Conclusions

- Develop a new point of view for multi-source data
- Can solve the problem in a fraction of the cost of the original problem
- Key - stochastic trace estimators and stochastic optimization
- Applications in other parameter estimation problems with many sources