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Ample Manual

Usage

ample can be called from Matlab using the following command.

[a,c] = ample(A,y,priorHandle,<options>);

- y is the set of observations.
- A is the system used to obtain y from the unknown signal. This system must be in the ample-system format, which will be described below.
- priorHandle is a function handle which takes two inputs, the hidden variables $\{R, \Sigma\}$.
- <options> is a Matlab-style list of varargin options in the pairs of
 'opt_name',opt_value.

The demos directory contains a number of useful examples for different signal priors and ample option modes. The best place to start is with demos/ample_gb_demo.m.

Ample-System Format

In order to find the factorized probabilities, ample needs to know the system which generated the observations, y. This information is given by the variable A in the description in the previous section. ample will operate on two different methods of specifying A:

Explicit

In this mode, A is assumed to be a linear system which is defined by an MxN matrix where y is of dimensionality M and the observed signals are of dimensionality x. I.e. a set of observations of a given signal is given by y = Ax.

Functional

This is a more general mode for defining the observational model of A. Here, both linear and non-linear observations can be constructed. Instead of an explicit matrix defining a linear system, in functional mode, the variable A is given as a structure with four fields, each of which is a function handle of a single vector input to produce each of the following four operations...

- A.forward: Computes the forward (signal->observation) mapping.
- A.adjoint: Computes the adjoint (observation->kernel-space) mapping.
- A.squared_forward: The squared operation is particular to the AMP algorithm. For a linear system, the squared forward mapping is equivalent to

$$(A \circ A) x$$
,

where \circ is the element-wise, or Hadamaard, product.

• A.squared_adjoint: The squared adjoint is similar to the above, but for the adjoint operation. For a linear system, this would be equivalent to

$$(A \circ A)^T y$$
.

Here is an example of a random linear system using this format.

```
% Signal dimensionality
N
     = 128;
                                   % Number of observations
Μ
     = 64;
     = randn(N,1);
                                   % Random signal
Phi = randn(M,N) ./ sqrt(N);
                                   % Random system
PhiT = Phi';
                                   % Defining the adjoint
Phi2 = Phi.*Phi;
                                   % Defining the squared forward
Phi2T= Phi2';
                                   % Defining the squared adjoint
A.forward
                  = 0(x_) Phi *x_;
                  = @(y_) PhiT *y_;
A.adjoint
A.squared_forward = @(x_) Phi2 *x_;
A.squared_adjoint = @(y_) Phi2T*y_;
y = A.forward(x);
z = A.adjoint(y);
```

Prior Modules

One of the nice aspects of the AMP framework is its applicability to a wide range of possible signal priors. With the right approach, even non-iid (i.e. structured) signal priors can even be used to great effect with AMP. The ample package provides this same versatility to the user by modularizing prior-specific calculations. The prior modules follow a standard format.

Module Format

The prior modules must follow the same general format.

```
function [a,c,learned_params] = prior_module(r,s,params)
    % Must return new learned prior parameters if requested
    learn_prior = 0;
    if nargout > 2
        learn_prior = 1;
    end

% 1. Calculate factorized means of variational distribution, `a`.
    % 2. Calculate factorized means of variational distribution, `c`.
    if learn_prior
        % 3. If new learned parameters are requested, update accordingly
    end
```

Here, the values **r** and **s** are the AMP hidden variables for the means and variances associated with the variational Gaussian. The structure of the **params** variable is entirely up to the module writer, **ample** will support both vector-valued and cell-valued **params**.

Important Note: If params is specified as a cell variable, then params{1} must contain the vector of learned variables!

The module writer must be familiar with the calculation of the means and variances of the variational form of their desired prior distribution,

$$a_i = \langle x_i \rangle = \frac{1}{Z} \int dx \ x \ P_0(x; \theta) \exp \left\{ \frac{(x - R_i)^2}{2\Sigma_i^2} \right\},$$

$$c_i = \left\langle x_i^2 \right\rangle - a_i^2 = \frac{1}{Z} \int \mathrm{d}x \ x^2 \ P_0(x;\theta) \exp\left\{ \frac{(x - R_i)^2}{2\Sigma_i^2} \right\},$$

where P_0 is the desired prior distribution and θ are the parameters associated with that distribution. This will generally also include the calculation of the

proper partition function. Depending on the prior, this calculation may be extremly non-trivial. If numerical integration is required to solve for these variables, ample will still support this, but at the cost computation time.

The determination of the prior parameter update may be equally non-trivial, in general.

Included Modules

For ease of use, some common prior modules are already included in the ample package. These serve as good examples for the expected format of the prior modules. ample currently provides the following signal priors:

• Binary

- priors/prior_binary.m
- Signal coefficients are discrete values are in $\{0,1\}$ and have value 1 with probability ρ .
- params format:
 - * params(1) $-\rho$

• Gauss-Bernoulli

- priors/prior_gb.m
- Signal coefficients are real values given by

$$Prob[x_i = x] \propto (1 - \rho) \delta(x) + \rho \mathcal{N}(\mu, \sigma^2).$$

This distribution has been shown to be very good for modelling "sparse" signals, such as those desired for compressed sensing.

- params format:
 - * params(1) $-\mu$
 - * params(2) σ^2
 - * params(3) $-\rho$

• L1-Sparse

- priors/prior_l1sparse.m
- Signal coefficients are assumed to be sparse, or at least compressible.
 That is, there are assumed to be very few non-zero coefficients. Using the L1-Sparse module isn't necessarily defining a "prior," per-se, but it does solve the equivalent LASSO problem when using ample for inverse-problems such as compressed sensing,

$$a = \arg\min_{x} ||x||_1 \text{ s.t. } ||y - Ax||_2^2 < \epsilon$$

- params format:

- * params(1) Minimum coefficient value
- * params(2) Maximum coefficient value

• Q-ary

- priors/prior_qary.m
- Signal coefficients are discrete values in $\{\tau_0,\tau_1,\ldots,\tau_{Q-1}\}$ with probabilities defined according to the PMF

$$Prob[x = \tau_q] = \rho_q.$$

- params format:
 - * params{1} PMF vector, $[\rho_0, \rho_1, \dots, \rho_{Q-1}]$
 - * params{2} Alphabet vector, $[\tau_0, \tau_1, \dots, \tau_{Q-1}]$

Table of Options & Default Values

Option Name	Values	Default Value
convergence_tolerance	Positive Small Real	1e-10
convergence_type	{'residual','iteration'}	'iteration'
damp	Real in $[0,1)$	0.0
debug	Boolean	false
delta	Positive Real	1.0
image_mode	Boolean	false
init_a	Vector of reals	zeros(N,1)
init_c	Vector of positive reals	ones(N,1)
learn_delta	Boolean	true
learning_mode	{'track','em'}	'track'
learn_prior_params	Boolean	false
max_em_iterations	Positive Integer	20
max_iterations	Positive Integer	250
mean_approximation	Boolean	false
pause_mode	Boolean	false
prior_damp	Real in $[0,1)$	0.0
prior_params	Function Handle	See 'Prior-Handle Format
report_history	Boolean	true
true_solution	Vector of Reals	[]

Option Name	Values	Default Value
verbose_mode	{0,1}	1

Option Descriptions

Required

• prior_params:

General

- convergence_tolerance:
- convergence_type:
- max_iterations:
- verbose_mode:
- delta:

Debug

- true_solution:
- debug:
- pause_mode:
- image_mode:
- report_history:

Fine-Tuning

- init_a: Initial value for the factorized means. The dimensionality of init_a should be equal to that of the original signal.
- init_c: Initial value for the factorized variances. The dimensionality of init_c should be equal to that of the original signal.
- damp:
- prior_damp:

Learning Control

- learn_delta:
- learn_prior_params:
- learning_mode:
- max_em_iterations: