The GPML Toolbox version 4.2

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Abstract

The GPML toolbox is an Octave 3.2.x and Matlab 7.x implementation of inference and prediction in Gaussian process (GP) models. It implements algorithms discussed in Rasmussen & Williams: Gaussian Processes for Machine Learning, the MIT press, 2006 and Nickisch & Rasmussen: Approximations for Binary Gaussian Process Classification, JMLR, 2008.

The strength of the function lies in its flexibility, simplicity and extensibility. The function is flexible as firstly it allows specification of the properties of the GP through definition of mean function and covariance functions. Secondly, it allows specification of different inference procedures, such as e.g. exact inference and Expectation Propagation (EP). Thirdly it allows specification of likelihood functions e.g. Gaussian or Laplace (for regression) and e.g. cumulative Logistic (for classification). Simplicity is achieved through a single function and compact code. Extensibility is ensured by modular design allowing for easy addition of extension for the already fairly extensive libraries for inference methods, mean functions, covariance functions and likelihood functions.

This document is a technical manual for a developer containing many details. If you are not yet familiar with the GPML toolbox, the *user documentation* and examples therein are a better way to get started.

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1 Gaussian Process Training and Prediction

The gpml toolbox contains a single user function gp described in section 2. In addition there are a number of supporting structures and functions which the user needs to know about, as well as an internal convention for representing the posterior distribution, which may not be of direct interest to the casual user.

Inference Methods: An inference method is a function which computes the (approximate) posterior, the (approximate) negative log marginal likelihood and its partial derivatives w.r.t.. the hyperparameters, given a model specification (i.e., GP mean and covariance functions and a likelihood function) and a data set. Inference methods are discussed in section 3. New inference methods require a function providing the desired inference functionality and possibly extra functionality in the likelihood functions applicable.

Hyperparameters: The hyperparameters is a struct controlling the properties of the model, i.e., the GP mean and covariance function and the likelihood function. The hyperparameters is a struct with the three fields mean, cov and lik, each of which is a vector. The number of elements in each field must agree with number of hyperparameters in the specification of the three functions they control (below). If a field is either empty or non-existent it represents zero hyperparameters. When working with FITC approximate inference, the inducing inputs xu can also be treated as hyperparameters for some common stationary covariances.

Hyperparameter Prior Distributions: When optimising the marginal likelihood w.r.t. hyperparameters, it is sometimes useful to softly constrain the hyperparameters by means of prior knowledge. A prior is a probability distribution over individual or a group of hyperparameters, section 7.

Likelihood Functions: The likelihood function specifies the form of the likelihood of the GP model and computes terms needed for prediction and inference. For inference, the required properties of the likelihood depend on the inference method, including properties necessary for hyperparameter learning, section 4.

Mean Functions: The mean function is a cell array specifying the GP mean. It computes the mean and its derivatives w.r.t.. the part of the hyperparameters pertaining to the mean. The cell array allows flexible specification and composition of mean functions, discussed in section 5. The default is the zero function.

Covariance Functions: The covariance function is a cell array specifying the GP covariance function. It computes the covariance and its derivatives w.r.t.. the part of the hyperparameters pertaining to the covariance function. The cell array allows flexible specification and composition of covariance functions, discussed in section 6.

Inference methods, see section 3, compute (among other things) an approximation to the posterior distribution of the latent variables f_i associated with the training cases, $i=1,\ldots,n$. This approximate posterior is assumed to be Gaussian, and is communicated via a struct post with the fields post.alpha, post.sW and post.L. Often, starting from the Gaussian prior $p(f) = \mathcal{N}(f|m,K)$ the approximate posterior admits the form

$$q(f|\mathcal{D}) = \mathcal{N}\big(f|\mu = m + K\alpha, \ V = (K^{-1} + W)^{-1}\big), \ \text{where} \ \ W \ \text{diagonal with} \ \ W_{\mathfrak{i}\mathfrak{i}} = s_{\mathfrak{i}}^2. \tag{1}$$

In such cases, the entire posterior can be computed from the two vectors post.alpha and post.sW; the inference method may optionally also return $L = \text{chol}(\text{diag}(s) \times \text{diag}(s) + I)$.

If on the other hand the posterior doesn't admit the above form, then post. L returns the matrix $L = -(K + W^{-1})^{-1}$ (and post.sW is unused). In addition, a sparse representation of the posterior may be used, in which case the non-zero elements of the post.alpha vector indicate the active entries.

2 The gp Function

The gp function is typically the only function the user would directly call.

It offers facilities for training the hyperparameters of a GP model as well as predictions at unseen inputs as detailed in the following help.

```
\langle gp \ function \ help \ 4b \rangle \equiv
4b
                                                                  (4a)
     1 % Gaussian Process inference and prediction. The gp function provides a
     2 % flexible framework for Bayesian inference and prediction with Gaussian
     3 % processes for scalar targets, i.e. both regression and binary
     4 % classification. The prior is Gaussian process, defined through specification
     5~\% of its mean and covariance function. The likelihood function is also
     6 % specified. Both the prior and the likelihood may have hyperparameters
     7 % associated with them.
     8 %
     9 % Two modes are possible: training or prediction: if no test cases are
     10 % supplied, then the negative log marginal likelihood and its partial
     11 % derivatives w.r.t. the hyperparameters is computed; this mode is used to fit
     12 % the hyperparameters. If test cases are given, then the test set predictive
    13 % probabilities are returned. Usage:
    14 %
    15 %
          training: [nlZ dnlZ
                                         ] = gp(hyp, inf, mean, cov, lik, x, y);
     16 % prediction: [ymu ys2 fmu fs2 ] = gp(hyp, inf, mean, cov, lik, x, y, xs);
     17 %
                 or: [ymu ys2 fmu fs2 lp] = gp(hyp, inf, mean, cov, lik, x, y, xs, ys);
    18 %
     19 % where:
    20 %
     21 %
                    struct of column vectors of mean/cov/lik hyperparameters
     22 %
                    function specifying the inference method
          inf
    23 %
          mean prior mean function
    24 % cov
                    prior covariance function
    25 %
                    likelihood function
          lik
    26 %
          X
                  n by D matrix of training inputs
           xs ns by D matrix of test inputs
ys column vector of
    27 %
                   column vector of length n of training targets
          У
    28 %
    29 %
                    column vector of length nn of test targets
    30 %
    31 %
           nlZ returned value of the negative log marginal likelihood
    32 %
           dnlZ
                    struct of column vectors of partial derivatives of the negative
    33 %
                        log marginal likelihood w.r.t. mean/cov/lik hyperparameters
     34 %
                    column vector (of length ns) of predictive output means
          ymu
    35 %
                    column vector (of length ns) of predictive output variances
           vs2
    36 %
                    column vector (of length ns) of predictive latent means
           fmu
     37 %
           fs2
                    column vector (of length ns) of predictive latent variances
    38 %
                    column vector (of length ns) of log predictive probabilities
    39 %
```

```
40 %
                      struct representation of the (approximate) posterior
            post
     41 %
                      3rd output in training mode or 6th output in prediction mode
     42 %
                      can be reused in prediction mode gp(.., cov, lik, x, post, xs,..)
     43 %
     44 % See also infMethods.m, meanFunctions.m, covFunctions.m, likFunctions.m.
     45 %
     46 (gpml copyright 5a)
     \langle gpml\ copyright\ 5a \rangle \equiv
                                             (4b 9 10 19 22 24a 36 38 39 44 46 49)
5a
      1 % Copyright (c) by Carl Edward Rasmussen and Hannes Nickisch, 2018-06-15.
                                                  File automatically generated using noweb.
```

Depending on the number of input parameters, gp knows whether it is operated in training or in prediction mode. The highlevel structure of the code is as follows. After some initialisations, we perform inference and decide whether test set predictions are needed or only the result of the inference is demanded.

```
5b ⟨initializations 5b⟩≡ (4a)

1 ⟨minimalist usage 5c⟩

2 ⟨multivariate output by recursion 5d⟩

3 ⟨process input arguments 6a⟩

4 ⟨check hyperparameters 6b⟩
```

If the number of input arguments is incorrect, we echo a minimalist usage and return.

If there is more than a single output dimension in y, we call multiple instances of gp with shared hyperparameters and settings each computing the corresponding result with scalar output.

```
\langle multivariate \ output \ by \ recursion \ 5d \rangle \equiv
5d
                          % deal with (independent) multivariate output y by recursing
     1 if size(y,2)>1
         d = size(y,2); varargout = cell(nargout,1); out = cell(nargout,1); % allocate
     3
         for i=1:d
     4
           in = {hyp, inf, mean, cov, lik, x, y(:,i)};
     5
           if nargin>7, in = {in{:}, xs}; end
     6
           if nargin>8, in = {in{:}, ys(:,i)}; end
     7
           8
                         [out{:}] = gp(in{:});
                                                          % .. number i in the output
     9
             if nargin==7, no = 2;
    10
               varargout{1} = varargout{1} + out{1};
                                                                            % sum nlZ
                                                                            % sum dnlZ
    11
                 varargout{2} = vec2any(hyp,any2vec(varargout{2})+any2vec(out{2}));
    12
    13
               end
                                                      \% concatenate ymu ys2 fmu fs2 lp
    14
    15
               for j=1:min(nargout,no), varargout{j} = [varargout{j},out{j}]; end
    16
             end
    17
             if nargout>no
                                                                    % concatenate post
               if i==2, varargout{no+1} = {varargout{no+1}},out{no+1}};
    18
               else varargout{no+1} = {varargout{no+1}{:},out{no+1}}; end
    19
    20
             end
    21
           end
    22
         end, return
                                                         % return to end the recursion
    23 end
```

Next, we set some useful default values for empty arguments, and convert inf and lik to function handles and mean and cov to cell arrays if necessary. Initialize variables.

```
\langle process\ input\ arguments\ 6a \rangle \equiv
6a
     1 if isempty(mean), mean = {@meanZero}; end
                                                                      % set default mean
     2 if ischar(mean) || isa(mean, 'function_handle'), mean = {mean}; end % make cell
     3 if isempty(cov), error('Covariance function cannot be empty'); end % no default
     4 if ischar(cov) || isa(cov, 'function_handle'), cov = {cov}; end
     5 cstr = cov{1}; if isa(cstr,'function_handle'), cstr = func2str(cstr); end
     6 if (strcmp(cstr,'covFITC') || strcmp(cstr,'apxSparse')) && isfield(hyp,'xu')
     7 \quad cov{3} = hyp.xu;
                                                                              %use hyp.xu
     8 \text{ end}
    9 if isempty(inf), inf = {@infGaussLik}; end
10 if ischar(inf), inf = str2func(inf); end
                                                        % set default inference method % convert into function handle
    11 if ischar(inf) || isa(inf, 'function_handle'), inf = {inf}; end
    12 istr = inf{1}; if isa(istr, 'function_handle'), istr = func2str(istr); end
    13 if strcmp(istr, 'infPrior')
    istr = inf{2}; if isa(istr,'function_handle'), istr = func2str(istr); end
    15 end
    16 if isempty(lik), lik = {@likGauss}; end
                                                                       % set default lik
    18 lstr = lik{1}; if isa(lstr,'function_handle'), lstr = func2str(lstr); end
    19
    20 D = size(x,2);
    21 if strncmp(cstr,'covGrid',7) || strcmp(cstr,'apxGrid') % only some inf* possible
    22 D = 0; xg = cov{3}; p = numel(xg); for i=1:p, D = D+size(xg{i},2); end % dims
    23 end
```

Check that the sizes of the hyperparameters supplied in hyp match the sizes expected. The three parts hyp.mean, hyp.cov and hyp.lik are checked separately, and define empty entries if they don't exist.

Inference is performed by calling the desired inference method inf. In training mode, we accept a failure of the inference method (and issue a warning), since during hyperparameter learning, hyperparameters causing a numerical failure may be attempted, but the minimize function may gracefully recover from this. During prediction, failure of the inference method is an error.

```
9
         end
10
       end
     end
11
12
     if nargin>7 % compute marginal likelihood and its derivatives only if needed
13
       if isstruct(y)
14
                              % reuse a previously computed posterior approximation
         post = y;
15
       else
16
         post = feval(inf{:}, hyp, mean, cov, lik, x, y);
17
18
     else
19
       if nargout <= 1</pre>
20
         [post nlZ] = feval(inf{:}, hyp, mean, cov, lik, x, y); dnlZ = {};
21
22
         [post nlZ dnlZ] = feval(inf{:}, hyp, mean, cov, lik, x, y);
23
24
     end
25 catch
26
    msgstr = lasterr;
    if nargin>7, error('Inference method failed [%s]', msgstr); else
27
28
       warning('Inference method failed [%s] .. attempting to continue', msgstr)
29
       varargout = {NaN, vec2any(hyp,zeros(numel(any2vec(hyp)),1))}; return % go on
30
     end
31 end
```

We copy the already computed negative log marginal likelihood to the first output argument, and if desired report its partial derivatives w.r.t. the hyperparameters if running in inference mode.

Predictions are computed in a loop over small batches to avoid memory problems for very large test sets.

7

```
\langle compute\ test\ predictions\ 7 \rangle \equiv
                                                               (4a)
1 alpha = post.alpha; L = post.L; sW = post.sW;
2 if issparse(alpha)
                                        % handle things for sparse representations
   nz = alpha \sim 0;
                                                        % determine nonzero indices
   if issparse(L), L = full(L(nz,nz)); end
                                                   % convert L and sW if necessary
   if issparse(sW), sW = full(sW(nz)); end
6 else nz = true(size(alpha,1),1); end
                                                        % non-sparse representation
7 if isempty(L)
                                        % in case L is not provided, we compute it
   K = feval(cov\{:\}, hyp.cov, x(nz,:));
9 L = chol(eye(sum(nz))+sW*sW'.*K);
10 end
11 %verify whether L contains valid Cholesky decomposition or something different
12 Lchol = isnumeric(L) && all(all(tril(L,-1)==0)&diag(L)'>0&isreal(diag(L))');
13 \text{ ns} = \text{size}(xs, 1);
                                                            % number of data points
14 if strncmp(cstr, 'apxGrid',7), xs = apxGrid('idx2dat',cov{3},xs); end % expand
                                            % number of data points per mini batch
15 nperbatch = 1000;
                                   % number of already processed test data points
16 \text{ nact} = 0;
17 ymu = zeros(ns,1); ys2 = ymu; fmu = ymu; fs2 = ymu; lp = ymu; % allocate mem
18 while nact<ns
                                % process minibatches of test cases to save memory
19
    id = (nact+1):min(nact+nperbatch,ns);
                                                           % data points to process
20
    (make predictions 8)
21
    nact = id(end);
                            % set counter to index of last processed data point
22 end
23 if nargin<9
   varargout = {ymu, ys2, fmu, fs2, [], post};
                                                        % assign output arguments
    varargout = {ymu, ys2, fmu, fs2, lp, post};
27 end
```

In every iteration of the above loop, we compute the predictions for all test points of the batch.

```
\langle make\ predictions\ 8\rangle \equiv
8
    1 kss = feval(cov{:}, hyp.cov, xs(id,:), 'diag');
                                                             % self-variance
    3 Ks = feval(cov{:}, hyp.cov, x, xs(id,:)); Ks = Ks(nz,:); % res indep. of x
    4 else
    5 Ks = feval(cov{:}, hyp.cov, x(nz,:), xs(id,:)); % avoid computation
    7 ms = feval(mean{:}, hyp.mean, xs(id,:));
    8 N = size(alpha,2); % number of alphas (usually 1; more in case of sampling)
    9 Fmu = repmat(ms,1,N) + Ks'*full(alpha(nz,:)); % conditional mean fs|f
   10 fmu(id) = sum(Fmu, 2)/N;
                                                            % predictive means
   11 if Lchol
                 % L contains chol decomp => use Cholesky parameters (alpha,sW,L)
   12 V = L'\(repmat(sW,1,length(id)).*Ks);
   13 fs2(id) = kss - sum(V.*V,1);
                                                        % predictive variances
                        % L is not triangular => use alternative parametrisation
   15 if isnumeric(L), LKs = L*Ks; else LKs = L(Ks); end % matrix or callback
   16 fs2(id) = kss + sum(Ks.*LKs,1);
                                                        % predictive variances
   17 end
   18 fs2(id) = max(fs2(id),0); % remove numerical noise i.e. negative variances
   19 Fs2 = repmat(fs2(id),1,N); % we have multiple values in case of sampling
   20 if nargin<9
   21 [Lp, Ymu, Ys2] = feval(lik{:},hyp.lik,[], Fmu(:),Fs2(:));
   22 else
   23     Ys = repmat(ys(id),1,N);
   24 [Lp, Ymu, Ys2] = feval(lik{:},hyp.lik,Ys(:),Fmu(:),Fs2(:));
   25 end
   26 lp(id) = sum(reshape(Lp, [],N),2)/N;
                                            % log probability; sample averaging
   27 ymu(id) = sum(reshape(Ymu,[],N),2)/N;
                                                  % predictive mean ys|y and ...
   28 ys2(id) = sum(reshape(Ys2,[],N),2)/N;
                                                                % .. variance
```

3 Inference Methods

Inference methods are responsible for computing the (approximate) posterior post, the (approximate) negative log marginal likelihood n1Z and its partial derivatives dn1Z w.r.t. the hyperparameters hyp. The arguments to the function are hyperparameters hyp, mean function mean, covariance function cov, likelihood function lik and training data x and y. Several inference methods are implemented and described this section.

```
\langle infMethods.m 9 \rangle \equiv
 1 % Inference methods: Compute the (approximate) posterior for a Gaussian process.
 2 % Methods currently implemented include:
 3 %
4 %
                        Exact inference (only possible with Gaussian likelihood)
      infGaussLik
 5 %
      infLaplace
                       Laplace's Approximation
 6 %
      infEP
                       Expectation Propagation
 7 %
      infVB
                        Variational Bayes Approximation
 8 %
      infKL
                        Kullback-Leibler optimal Approximation
9 %
10 %
      infMCMC
                Markov Chain Monte Carlo and Annealed Importance Sampling
11 %
                   We offer two samplers.
12 %
                     - hmc: Hybrid Monte Carlo
13 %
                     - ess: Elliptical Slice Sampling
14 %
                   No derivatives w.r.t. to hyperparameters are provided.
15 %
16 %
     infL00
                  Leave-One-Out predictive probability and Least-Squares Approxim.
17 %
      infPrior
                  Perform inference with hyperparameter prior.
18 %
19 % The interface to the approximation methods is the following:
20 %
21 %
      function [post nlZ dnlZ] = inf..(hyp, mean, cov, lik, x, y)
22 %
23 % where:
24 %
25 %
      hyp
               is a struct of hyperparameters
26 %
      mean
               is the name of the mean function
                                                      (see meanFunctions.m)
27 %
               is the name of the covariance function (see covFunctions.m)
     cov
28 %
               is the name of the likelihood function (see likFunctions.m)
      lik
29 %
               is a n by D matrix of training inputs
      X
30 %
               is a (column) vector (of size n) of targets
31 %
             is the returned value of the negative log marginal likelihood
32 %
      nlZ
33 %
      dnlZ
               is a (column) vector of partial derivatives of the negative
34 %
                   log marginal likelihood w.r.t. each hyperparameter
35 %
               struct representation of the (approximate) posterior containing
     post
      alpha is a (sparse or full column vector) containing inv(K)*(mu-m),
36 %
37 %
                   where K is the prior covariance matrix, m the prior mean,
38 %
                   and mu the approx posterior mean
39 %
               is a (sparse or full column) vector containing diagonal of sqrt(W)
       sW
40 %
                   the approximate posterior covariance matrix is inv(inv(K)+W)
41 %
               is a (sparse or full) triangular matrix, L = chol(sW*K*sW+eye(n)),
        L
42 %
                   or a full matrix, L = -inv(K+inv(W)),
43 %
                   or a function L(A) of a matrix A such that -(K+inv(W))*L(A) = A
45 % Usually, the approximate posterior to be returned admits the form
46 % N(mu=m+K*alpha, V=inv(inv(K)+W)), where alpha is a vector and W is diagonal.
```

48 % For more information on the individual approximation methods and their

```
49 % implementations, see the separate inf??.m files. See also gp.m. 50 % 51 \ \langle gpml \ copyright \ 5a \rangle
```

Not all inference methods are compatible with all likelihood functions, e.g., exact inference is only possible with Gaussian likelihood. In order to perform inference, each method needs various properties of the likelihood functions, section 4.

3.1 Exact Inference with Gaussian likelihood

10

For Gaussian likelihoods, GP inference reduces to computing mean and covariance of a multivariate Gaussian which can be done exactly by simple matrix algebra. The program inf/infExact.m does exactly this. If it is called with a likelihood function other than the Gaussian, it issues an error. The Gaussian posterior $q(f|\mathcal{D}) = \mathcal{N}(f|\mu, V)$ is exact.

```
\langle inf/infGaussLik.m \ 10 \rangle \equiv
1 function [post nlZ dnlZ] = infGaussLik(hyp, mean, cov, lik, x, y, opt)
3 % Exact inference for a GP with Gaussian likelihood.
5 % Compute a parametrization of the posterior, the negative log marginal
6 % likelihood and its derivatives w.r.t. the hyperparameters. The function takes
7 % a specified covariance function (see covFunctions.m) and likelihood function
8 \% (see likFunctions.m), and is designed to be used with gp.m.
9 %
10 \(\langle gpml \) copyright 5a\(\rangle \)
11 %
12 % See also INFMETHODS.M, APX.M.
13
14 if nargin<7, opt = []; end
                                                           % make sure parameter exists
15 if iscell(lik), likstr = lik{1}; else likstr = lik; end
16 if ~ischar(likstr), likstr = func2str(likstr); end
17 if ~strcmp(likstr,'likGauss')
                                                % NOTE: no explicit call to likGauss
    error('Exact inference only possible with Gaussian likelihood');
19 end
20
21 [n, D] = size(x);
22 [m,dm] = feval(mean{:}, hyp.mean, x);
                                                     % evaluate mean vector and deriv
23 sn2 = \exp(2*hyp.lik); W = ones(n,1)/sn2;
                                                           % noise variance of likGauss
24 \text{ K} = \text{apx}(\text{hyp,cov,x,opt});
                                                     % set up covariance approximation
25 [ldB2,solveKiW,dW,dhyp,post.L] = K.fun(W); % obtain functionality depending on W
26
27 alpha = solveKiW(y-m);
28 post.alpha = K.P(alpha);
                                                     \% return the posterior parameters
29 \text{ post.sW} = \text{sqrt(W)};
                                                      % sqrt of noise precision vector
30 if nargout>1
                                                 % do we want the marginal likelihood?
31 nlZ = (y-m)'*alpha/2 + ldB2 + n*log(2*pi*sn2)/2; % -log marginal likelihood
32
    if nargout>2
                                                             % do we want derivatives?
33
       dnlZ = dhyp(alpha); dnlZ.mean = -dm(alpha);
34
       dnlZ.lik = -sn2*(alpha'*alpha) - 2*sum(dW)/sn2 + n;
35
     end
36 \text{ end}
```

3.2 Laplace's Approximation

For differentiable likelihoods, Laplace's approximation, approximates the posterior by a Gaussian centered at its mode and matching its curvature inf/infLaplace.m.

More concretely, the mean of the posterior $q(f|\mathcal{D}) = \mathcal{N}\big(f|\mu,V\big)$ is – defining $\ell_i(f_i) = \ln p(y_i|f_i)$ and $\ell(f) = \sum_{i=1}^n \ell_i(f_i)$ – given by

$$\mu = \arg\min_{\mathbf{f}} \phi(\mathbf{f}), \text{ where } \phi(\mathbf{f}) = \frac{1}{2} (\mathbf{f} - \mathbf{m})^{\top} \mathbf{K}^{-1} (\mathbf{f} - \mathbf{m}) - \ell(\mathbf{f}) \stackrel{c}{=} -\ln[p(\mathbf{f})p(\mathbf{y}|\mathbf{f})], \tag{2}$$

which we abbreviate by $\mu \leftarrow \mathcal{L}(\ell)$. The curvature $\frac{\partial^2 \varphi}{\partial f_1^{\dagger}} = K^{-1} + W$ with $W_{ii} = -\frac{\partial^2}{\partial f_i^2} \ln p(y_i|f_i)$ serves as precision for the Gaussian posterior approximation $V = (K^{-1} + W)^{-1}$ and the marginal likelihood $Z = \int p(f)p(y|f)df$ is approximated by $Z \approx Z_{LA} = \int \tilde{\varphi}(f)df$ where we use the 2nd order Taylor expansion at the mode μ given by $\tilde{\varphi}(f) = \varphi(\mu) + \frac{1}{2}(f - \mu)^{\top}V^{-1}(f - \mu) \approx \varphi(f)$.

Laplace's approximation needs derivatives up to third order for the mode fitting procedure (Newton method)

$$d_k = \frac{\partial^k}{\partial f^k} \log p(y|f), \quad k = 0, 1, 2, 3$$

and

$$d_k \; = \; \frac{\partial}{\partial \rho_i} \frac{\partial^k}{\partial f^k} \log p(y|f), \quad k = 0, 1, 2$$

evaluated at the latent location f and observed value y. The likelihood calls (see section 4)

and

return exactly these values.

3.3 Expectation Propagation

The basic idea of Expectation Propagation (EP) as implemented in inf/infEP.m. is to replace the non-Gaussian likelihood terms $p(y_i|f_i)$ by Gaussian functions $t(f_i; \nu_i, \tau_i) = \exp(\nu_i f_i - \frac{1}{2}\tau_i f_i^2)$ and to adjust the natural parameters ν_i, τ_i such that the following identity holds:

$$\frac{1}{Z_{t,i}}\int f^kq_{-i}(f)\cdot t(f;\nu_i,\tau_i)df = \frac{1}{Z_{t,i}}\int f^kq_{-i}(f)\cdot p(y_i|f)df,\quad k=1,2$$

with the so-called cavity distributions $q_{-i}(f) = \mathcal{N}(f|m,K) \prod_{j \neq i} t(f_j; \nu_j, \tau_j) \propto \mathcal{N}(f|\mu,V)/t(f_i; \nu_i, \tau_i)$ equal to the posterior divided by the ith Gaussian approximation function and the two normalisers $Z_{t,i} = \int q_{-i}(f) \cdot t(f_i; \nu_i, \tau_i) df$ and $Z_{p,i} = \int q_{-i}(f) \cdot p(y_i|f_i) df$. The moment matching corresponds to minimising the following local KL-divergence

$$\nu_i, \tau_i = arg \min_{\nu,\tau} KL[q_{-i}(f)p(y_i|f_i)/Z_{p,i} \|q_{-i}(f)t(f_i;\nu,\tau)/Z_{t,i}].$$

In order to apply the moment matching steps in a numerically safe way, EP requires the deriviatives of the expectations w.r.t. the Gaussian mean parameter μ

$$d_k = \frac{\partial^k}{\partial \mu^k} \log \int p(y|f) \mathcal{N}(f|\mu, \sigma^2) df, \quad k = 0, 1, 2$$

and the ith likelihood hyperparameter ρ_i

$$d = \frac{\partial}{\partial \rho_i} \log \int p(y|f) \mathcal{N}(f|\mu, \sigma^2) df$$

which can be obtained by the likelihood calls (see section 4)

• [d0, d1, d2] = lik(hyp, y, mu, s2, 'infEP')

and

• d = lik(hyp, y, mu, s2, 'infEP', i).

3.4 Kullback Leibler Divergence Minimisation

Another well known approach to approximate inference implemented inf/infKL.m in attempts to directly find the closest Gaussian $q(f|\mathcal{D}) = \mathcal{N}(f|\mu, V)$ to the exact posterior $p(f|\mathcal{D})$ w.r.t. to some proximity measure or equivalently to maximise a lower bound $Z(\mu, V)$ to the marginal likelihood Z as described in Nickisch & Rasmussen *Approximations for Binary Gaussian Process Classification*, JMLR, 2008. In particular, one minimises $KL(\mathcal{N}(f|\mu, V)||p(f|\mathcal{D}))$ which amounts to minimising $-\ln Z(\mu, V)$ as defined by:

$$\begin{split} -\ln \mathsf{Z} &= -\ln \int p(f)p(y|f)df = -\ln \int q(f|\mathcal{D}) \frac{p(f)}{q(f|\mathcal{D})} p(y|f)df \\ &\overset{Jensen}{\leqslant} \int q(f|\mathcal{D}) \ln \frac{q(f|\mathcal{D})}{p(f)} df - \int q(f|\mathcal{D}) \ln p(y|f)df =: -\ln \mathsf{Z}(\mu, V) \\ &= \mathsf{KL}\left(\mathcal{N}(f|\mu, V)||\mathcal{N}(f|m, K)\right) - \sum_{i=1}^n \int \mathcal{N}(f_i|\mu_i, \nu_{i\,i}) \ln p(y_i|f_i) df_i, \ \nu_{i\,i} = [V]_{i\,i} \\ &= \frac{1}{2} \left(\operatorname{tr}(VK^{-1} - I) - \ln |VK^{-1}| \right) + \frac{1}{2} (\mu - m)^\top K^{-1} (\mu - m) - \sum_{i=1}^n \ell^{KL}(\mu_i, \nu_{i\,i}) \right) \end{split}$$

where $\ell_v^{KL}(\mu_i) = \int \mathcal{N}(f_i|\mu_i,\nu_{ii})\ell_i(f_i)df_i$ is the convolution of the log likelihood ℓ_i with the Gaussian \mathcal{N} and $\mathbf{v} = dg(\mathbf{V})$. Equivalently, one can view ℓ^{KL} as a smoothed version of ℓ with univariate smoothing kernel \mathcal{N} .

From Challis & Barber Concave Gaussian Variational Approximations for Inference in Large Scale Bayesian Linear Models, AISTATS, 2011 we know that the mapping $(\mu, L) \mapsto -\ln Z(\mu, L^{\top}L)$ is jointly convex whenever the likelihoods $f_i \mapsto \mathbb{P}(y_i|f_i)$ are log concave. In particular, this implies that every $(\mu_i, s_i) \mapsto -\ell^{KL}(\mu_i, s_i^2)$ is jointly convex.

We use an optimisation algorithm similar to EP (section 3.3) where we minimise the local KL-divergence the other way round μ_i , $s_i = \arg\min_{\mu,s} \text{KL}[\mathcal{N}(f|\mu,s^2)||q_{-i}(f)p(y_i|f_i)/Z_{p,i}]$. This view was brought forward by Tom Minka Convex Divergence measures and message passing, MSR-TR, 2005. The KL minimisation constitutes a jointly convex 2d optimisation problem solved by klmin using a scaled Newton approach which is included as a sub function in inf/infKL.m. The smoothed likelihood $\ell^{\text{KL}}(\mu_i,\nu_{ii})$ is implemented as a meta likelihood in likKL; it uses Gaussian-Hermite quadrature to compute the required integrals. Note that – as opposed to EP – Gaussian-Hermite quadrature is appropriate since we integrate against the $\ln \mathbb{P}(y_i|f_i)$ (which can be well approximated by a polynomial) instead of $\mathbb{P}(y_i|f_i)$ itself. The algorithm is – again unlike EP – provably convergent for log-concave likelihoods (e.g. likGauss, likLaplace, likSech2, likLogistic, likPoisson) since it can be regarded as coordinate descent with guaranteed decrease in the objective in every step. Due to the complex update computations, infKL can be quite slow although it has the same $\mathfrak{O}(\mathfrak{n}^3)$ asymptotic complexity as EP and Laplace.

3.5 Variational Bayes

One can drive the bounding even further by means of local quadratic lower bounds to the log likelihood $\ell(f) = \ln p(y|f)$. Suppose that we use a super-Gaussian likelihood p(y|f) i.e. likelihoods that can be lower bounded by Gaussians of any width w (e.g. likLaplace, likT, likLogistic, likSech2). Formally, that means that there are $b, z \in \mathbb{R}$ such that

$$\rho(f) = \ln p(y|f - z) - bf$$

is symmetric and $\sqrt{f} \mapsto \rho(f)$ is a convex function for all $f \geqslant 0$. As a result, we obtain the following exact representation of the likelihood

$$\ell(f) = \ln p(y|f) = \max_{w>0} \left((b+wz)f - \frac{wf^2}{2} - \frac{1}{2}h(\gamma) \right),$$

which can be derived by convex duality and assuming the likelihoods to be super-Gaussian. Details can be found in papers by Palmer et al. *Variational EM Algorithms for Non-Gaussian Latent Variable Models*, NIPS, 2006 and Nickisch & Seeger *Convex Variational Bayesian Inference for Large Scale Generalized Linear Models*, ICML, 2009.

The bottom line is that we can treat the variational bounding as a sequence of Laplace approximations with the "variational Bayes" log likelihood

$$\ell^{VB}(f_{\mathfrak{i}}) = \ell(g_{\mathfrak{i}}) + b_{\mathfrak{i}}(f_{\mathfrak{i}} - g_{\mathfrak{i}}), \ g = sgn(f - z) \odot \sqrt{(f - z)^2 + v} + z$$

instead of the usual likelihood $\ell(f_i) = \ln p(y_i|f_i)$ i.e. we solve $\mu \leftarrow \mathcal{L}(\ell_v^{VB})$ instead of $\mu \leftarrow \mathcal{L}(\ell)$. See section 3.2. In the code of inf/infVB.m, the likelihood is implemented in the function likVB.

At the end, the optimal value of **W** can be obtained analytically via $w_i = |b_i - \ell'(g_i)|/|g_i - z_i|$.

For the minimisation in inf/infVB.m, we use a provably convergent double loop algorithm, where in the inner loop a nonlinear least squares problem (convex for log-concave likelihoods) is solved using inf/infLaplace.m such that $\mu \leftarrow \mathcal{L}(\ell_v^{VB})$ and in the outer loop, we compute $\mathbf{v} \leftarrow \mathrm{dg}((\mathbf{K}^{-1} + \mathbf{W})^{-1})$. The only requirement to the likelihood function is that it returns the values z and b required by the bound which are delivered by the call (see section 4)

The negative marginal likelihood upper bound $-\ln Z_{VB}$ is obtained by integrating the prior times the exact representation of the likelihood

$$p(y|f) = \max_{\gamma > 0} q(y|f, \gamma), \ q(y|f, \gamma) = \mathcal{N}(f|\nu, \gamma) \exp\left(-\frac{h(\gamma)}{2} - \frac{\nu^2}{2\gamma}\right) \sqrt{2\pi\gamma}, \ \gamma = \frac{1}{w}, \ \nu = b\gamma + z$$

w.r.t. the latent variables f yielding

$$\begin{split} -\ln Z_{\mathrm{VB}} &= -\ln \int \mathcal{N}(\mathbf{f}|\mathbf{m},\mathbf{K}) \prod_{i=1}^{n} q_{i}(y_{i}|f_{i},\gamma_{i}) d\mathbf{f} \\ &= -\ln \mathcal{N}(\mathbf{m}|\mathbf{v},\mathbf{K}+\boldsymbol{\Gamma}) + \frac{1}{2} \left(h(\boldsymbol{\gamma}) - \mathbf{w}^{\top} \boldsymbol{\nu}^{2} - \mathbf{1}^{\top} \ln 2\pi \boldsymbol{\gamma}\right). \end{split}$$

3.6 Compatibility Between Inference Methods and Covariance Approximations

Another kind of approximation is needed to render an inference method scalable. We have two approximation schemes which in fact approximate the covariance to make it amenable to large number of training data points. The following table shows the compatibility between some inference methods and two major groups of covariance approximations we will discuss in the next two sections.

Approximation \ Inference	Exact	Laplace	VB	EP	KL, MCMC, LOO	Implementation	
Approximation (interence	infGaussLik	infLaplace	infVB	infEP	inf{KL,MCMC,LOO}	implementation	
variational free energy (VFE)	✓	✓	✓			apxSparse, opt.s=0.0	
fully independent training conditionals (FITC)	✓	✓	✓	✓		apxSparse, opt.s=1.0	
hybrid between VFE and FITC, $s_0 \in [0, 1]$	✓	✓	✓			apxSparse, opt.s=s0	
Kronecker/Toeplitz/BTTB grid (KISS-GP)	✓	✓	✓			apxGrid	
State space representation	✓	✓	✓	(√, ADF)		apxState	

3.7 Sparse Covariance Approximations

One of the main problems with GP models is the high computational load for inference computations. In a setting with n training points x, exact inference with Gaussian likelihood requires $O(n^3)$ effort; approximations like Laplace or EP consist of a sequence of $O(n^3)$ operations.

There is a line of research with the goal to alleviate this burden by using approximate covariance functions \tilde{k} instead of k. A review is given by Candela and Rasmussen¹. One basic idea in those approximations is to work with a set of m inducing inputs \mathbf{u} with a reduced computational load of $O(nm^2)$. In the following, we will provide a rough idea of the FITC approximation used in the toolbox. Let K denote the $n \times n$ covariance matrix between the training points \mathbf{x} , $K_{\mathbf{u}}$ the $m \times n$ covariance matrix between the n training points and the m inducing points, and $K_{\mathbf{u}\mathbf{u}}$ the $m \times m$ covariance matrix between the m inducing points. The FITC approximation to the covariance is given by

$$K \approx \tilde{K} = Q + G, \ G = \text{diag}(g), \ g = \text{diag}(K - Q), \ Q = K_u^\top Q_{uu}^{-1} K_u, \ Q_{uu} = K_{uu} + \sigma_{n_u}^2 I,$$

where σ_{n_u} is the noise from the inducing inputs. Note that \tilde{K} and K have the same diagonal elements diag(\tilde{K}) = diag(K); all off-diagonal elements are the same as for Q. Internally, the necessary covariance evaluations are performed by a meta covariance function cov/apxSparse.m. The toolbox offers FITC versions for regression with Gaussian likelihood inf/infGaussLik.m, as well as for Laplace's approximation inf/infLaplace.m.

The user can decide whether to treat the inducing inputs **u** as fixed or as hyperparameters. The latter allows to adjust the inducing inputs **u** w.r.t. the marginal likelihood. As detailed in the documentation of inf/apx.m, **u** is treated as fixed if it is passed as the 2nd parameter of apxSparse(cov,xu,..). If the hyperparameter structure hyp contains a field hyp.xu in inference method calls such as infGaussLik(hyp,..) or inference/prediction calls like gp(hyp,@infGaussLik,..) the inducing inputs **u** are treated as hyperparameters and can be optimised. See doc/demoSparse.m for an illustration.

3.8 Grid-Based Covariance Approximations

Another way to bring down computational costs is to take advantage of grid structure \mathbf{x} . For example, in geostatistics or image processing, the training data $\mathbf{x} \in \mathbb{R}^{n \times D}$ could be a complete 2d lattice of size $n_1 \times n_2$ as given by the axes $\mathbf{g}_1 \in \mathbb{R}^{n_1}$, $\mathbf{g}_2 \in \mathbb{R}^{n_2}$ so that $\mathbf{n} = \mathbf{N} = \mathbf{n}_1 \cdot \mathbf{n}_2$, $\mathbf{D} = 2$ and $\mathbf{x} = [\text{vec}(\mathbf{g}_1 \mathbf{1}^\top), \text{vec}(\mathbf{1}\mathbf{g}_2^\top)]$. In general, a p-dimensional grid $\mathbf{U} \in \mathbb{R}^{N \times D}$ is specified by a set of axis matrices $\{\mathbf{g}_i \in \mathbb{R}^{n_i \times D_i}\}_{i=1..p}$ so that $\mathbf{N} = \prod_{i=1}^p \mathbf{n}_i$ and $\mathbf{D} = \sum_{i=1}^p \mathbf{D}_i$ where the axes do not need to be 1d nor do their components need to be sorted. As a consequence, \mathbf{U} represents a Cartesian product of its axes $\mathbf{U} = \mathbf{g}_1 \times \mathbf{g}_2 \times ... \times \mathbf{g}_p$. The cov/apxGrid.m covariance function represents a Kronecker product covariance matrix

$$K_{U,U} = K_p \otimes .. \otimes K_2 \otimes K_1$$

whose factorisation structure is given by the grid \mathbf{x}_g . The gain in computationial efficiency is due to the fact that matrix-vector product, determinant, inverse and eigenvalue computations decompose so that many operations with an overall cost of $\mathcal{O}(N^3)$ now only cost $\mathcal{O}(\sum_{i=1}^p n_i^3)$.

¹A Unifying View of Sparse Approximate Gaussian Process Regression, JMLR, 2005

For off-grid data points, we can still take advantage of the computational properties of a grid-based covariance matrix $K_{U,U}$ via the structured kernel interpolation (SKI) framework aka KISS-GP by Wilson and Nickisch² with extensions³. Here, the $n \times n$ covariance K is obtained from the $N \times N$ grid covariance $K_{U,U}$ by interpolation $K \approx W_X K_{U,U} W_X^{\top}$, where $K_{U,U}$ is a covariance matrix formed by evaluating the user-specified kernel over a set of latent inducing inputs U, with locations that have been chosen to create algebraic structure in $K_{U,U}$ that we can exploit for efficiency. Here, the interpolation matrix $W_X \in \mathbb{R}^{n \times N}$ is extremely sparse; i.e., for local cubic interpolation W_X contains only 4^D nonzeros per row, where D is the diata dimension. In addition W_X is row-normalised $1_n = W_X 1_N$. The structure in $K_{U,U}$ alongside the sparsity of W_X , allows for very fast MVMs with the SKI approximate covariance matrix K over the inputs x enabling fast inference and prediction.

Internally, we use a meta covariance function cov/apxGrid.m to represent the Kronecker covariance matrix and a Gaussian regression inference method inf/infGaussLik.m. We also support incomplete grids where n < N. A good starting point is Yunus Saatçi's PhD thesis⁴. For incomplete grids, we use the interpolation-based extensions by Wilson et al.⁵ where conjugate gradients and a determinant approximations are used. See doc/demoGrid1d.m and doc/demoGrid2d.m for an illustration. We also offer non-Gaussian likelihoods as described by Seth Flaxman⁶ so that inf/infLaplace.m can be used.

3.9 State Space Representation of GPs

GP models with covariance functions with a Markovian structure can be transformed into equivalent discrete state space models where inference can be done in linear time O(n). Exact models can be derived for sum, product, linear, noise, constant, Matérn (half-integer), Ornstein–Uhlenbeck, and Wiener covariance functions. Other common covariance functions can be approximated by their Markovian counterparts, including squared exponential, rational quadratic, and periodic covariance functions.

A state space model describes the evolution of a dynamical system at different time instances t_i , i = 1, 2, ... by

$$\mathbf{f_i} \sim \mathbb{P}(\mathbf{f_i}|\mathbf{f_{i-1}}), \quad \mathbf{y_i} \sim \mathbb{P}(\mathbf{y_i}|\mathbf{f_i}),$$

where $f_i := f(t_i) \in \mathbb{R}^d$ and $f_0 \sim \mathbb{P}(f_0)$ with f_i being the latent (hidden/unobserved) variable and y_i being the observed variable. In continuous time, a simple dynamical system able to represent many covariance functions is given by the following linear time-invariant stochastic differential equation:

$$\dot{f}(t) = F f(t) + L w(t), \quad y_i = H f(t_i) + \varepsilon_i,$$

where $\mathbf{w}(t)$ is an s-dimensional white noise process, the measurement noise $\varepsilon_i \sim \mathcal{N}(0, \sigma_n^2)$ is Gaussian, and $\mathbf{F} \in \mathbb{R}^{d \times d}$, $\mathbf{L} \in \mathbb{R}^{d \times s}$, $\mathbf{H} \in \mathbb{R}^{1 \times d}$ are the feedback, noise effect, and measurement matrices, respectively. The initial state is distributed according to $\mathbf{f}_0 \sim \mathcal{N}(\mathbf{0}, \mathbf{P}_0)$.

The latent GP is recovered by f(t) = Hf(t) and $w(t) \in \mathbb{R}^s$ is a multivariate white noise process with spectral density matrix $Q_c \in \mathbb{R}^{s \times s}$. For discrete values, this translates into

$$f_i \sim \mathcal{N}(A_{i-1}f_{i-1}, Q_{i-1}), \quad y_i \sim \mathbb{P}(y_i|H\,Lf_i),$$

²Kernel Interpolation for Scalable Structured Gaussian Processes, ICML, 2015

³Thoughts on Massively Scalable Gaussian Processes, TR, 2015.

⁴Scalable Inference for Structured Gaussian Process Models, University of Cambridge, 2011

⁵Fast Kernel Learning for Multidimensional Pattern Extrapolation, NIPS, 2014

⁶Fast Kronecker inference in Gaussian processes with non-Gaussian likelihoods, ICML, 2015

Algorithm 1 Kalman (forward) filtering.

```
Input: \{t_i\}, y
                                                                                                                                                  # training inputs and targets
                                                                                                                                                                     # state space model
               \{A_i\}, \{Q_i\}, H, P_0
               W,b
                                                                                                                           # likelihood eff. precision and location
for i = 1 to n do
     if i == 1 then
         m_i \leftarrow 0; P_i \leftarrow P_0
                                                                                                                                                                                                   # init
          \mathbf{m_i} \leftarrow \mathbf{A_i} \mathbf{m_{i-1}}; \ \mathbf{P_i} \leftarrow \mathbf{A_i} \mathbf{P_{i-1}} \mathbf{A_i}^\top + \mathbf{Q_i}
                                                                                                                                                                                           # predict
     end if
     if has label yi then
         \mu_f \leftarrow Hm_i; \ u \leftarrow P_iH^\top; \ \sigma_f^2 \leftarrow Hu
                                                                                                                                                                                             # latent
          z_i \leftarrow W_{ii}\sigma_f^2 + 1
         \mathbf{k}_{\mathfrak{i}} \leftarrow \mathbf{W}_{\mathfrak{i}\mathfrak{i}}\mathbf{u}/z_{\mathfrak{i}}; \ \mathbf{P}_{\mathfrak{i}} \leftarrow \mathbf{P}_{\mathfrak{i}} - \mathbf{k}_{\mathfrak{i}}\mathbf{u}^{\top}
                                                                                                                                                                                       # variance
          c_i \leftarrow W_{ii}\mu_f - b_i; m_i \leftarrow m_i - uc_i/z_i
                                                                                                                                                                                               # mean
     end if
end for
\log \det(\mathbf{I} + \mathbf{W}^{\frac{1}{2}} \mathbf{K} \mathbf{W}^{\frac{1}{2}}) \leftarrow \sum_{i} \log z_{i}
```

Algorithm 2 Rauch-Tung-Striebel (backward) smoothing.

```
\begin{array}{ll} \text{Input: } \{m_i\}, \{P_i\} & \text{\# Kalman filter output} \\ \{A_i\}, \{Q_i\} & \text{\# state space model} \\ \text{for } i = n \text{ down to 2 do} \\ m \leftarrow A_i m_{i-1}; \ P \leftarrow A_i P_{i-1} A_i^\top + Q_i & \text{\# predict} \\ G_i \leftarrow P_{i-1} A_i^\top P^{-1}; \ \Delta m_{i-1} \leftarrow G_i (m_i - m) \\ P_{i-1} \leftarrow P_{i-1} + G_i (P_i - P) G_i^\top & \text{\# variance} \\ m_{i-1} \leftarrow m_{i-1} + \Delta m_{i-1} & \text{\# mean} \\ \text{end for} & & \text{\# mean} \\ \end{array}
```

with $f_0 \sim \mathcal{N}(0, P_0)$. The discrete-time matrices are

$$\begin{split} \mathbf{A}_{i} &= \mathbf{A}[\Delta t_{i}] = e^{\Delta t_{i}F}, \\ \mathbf{Q}_{i} &= \int_{0}^{\Delta t_{i}} e^{(\Delta t_{k} - \tau)F} \mathbf{L} \, \mathbf{Q}_{c} \, \mathbf{L}^{\top} e^{(\Delta t_{i} - \tau)F^{\top}} d\tau, \end{split}$$

where $\Delta t_i = t_{i+1} - t_i \geqslant 0$.

For stationary covariances k(t,t')=k(t-t'), the stationary state is distributed by $f_{\infty}\sim \mathcal{N}(0,P_{\infty})$ and the stationary covariance can be found by solving the Lyapunov equation

$$\dot{P}_{\infty} = F P_{\infty} + P_{\infty} F^{\top} + L Q_{c} L^{\top} = 0,$$

which leads to the identity $Q_i = P_{\infty} - A_i P_{\infty} A_i^{\top}$.

In practice, the evaluation of the n discrete-time transition matrices $A_i = e^{\Delta t_i F}$ and the noise covariance matrices Q_i (in the stationary case) for different values of Δt_i is a computational challenge. Since the matrix exponential $\psi: s \mapsto e^{sX}$ is smooth, its evaluation can be accurately approximated by convolution interpolation.

From the discrete set of matrices, all the necessary computations can be done using Kalman filtering and Kalman smoothing as detailed in Algorithms 1 and 2.

Internally, we use a meta covariance function cov/apxState.m to represent the state space represen-

tation. A good starting point is Arno Solin's PhD thesis⁷. See doc/demoState.m for an illustration. We also offer non-Gaussian likelihoods 8 so that inf/infLaplace.m and inf/infVB.m can be used. EP is not fully functional; we offer single-sweep EP aka assumed density filtering (ADF).

⁷Stochastic Differential Equation Methods for Spatio-Temporal Gaussian Process Regression, Aalto University, 2016

⁸State Space Gaussian Processes with Non-Gaussian Likelihood, ICML, 2018

4 Likelihood Functions

A likelihood function $p_{\rho}(y|f)$ (with hyperparameters ρ) is a conditional density $\int p_{\rho}(y|f) dy = 1$ defined for scalar latent function values f and outputs y. In the GPML toolbox, we use iid. likelihoods $p_{\rho}(y|f) = \prod_{i=1}^{n} p_{\rho}(y_i|f_i)$. The approximate inference engine does not explicitly distinguish between classification and regression likelihoods: it is fully generic in the likelihood allowing to use a single code in the inference step.

Likelihood functionality is needed both during inference and while predicting.

4.1 Prediction

A prediction at \mathbf{x}_* conditioned on the data $\mathcal{D}=(X,y)$ (as implemented in gp.m) consists of the predictive mean μ_{y_*} and variance $\sigma_{y_*}^2$ which are computed from the the latent marginal moments μ_{f_*} , $\sigma_{f_*}^2$ i.e. the Gaussian marginal approximation $\mathcal{N}(f_*|\mu_{f_*},\sigma_{f_*}^2)$ via

$$p(y_*|\mathcal{D}, \mathbf{x}_*) = \int p(y_*|f_*)p(f_*|\mathcal{D}, \mathbf{x}_*)df_* \approx \int p(y_*|f_*)\mathcal{N}(f_*|\mu_{f_*}, \sigma_{f_*}^2)df_*.$$
 (3)

The moments are given by $\mu_{y_*}=\int y_*p(y_*|\mathcal{D},\mathbf{x}_*)dy_*$ and $\sigma_{y_*}^2=\int (y_*-\mu_{y_*})^2p(y_*|\mathcal{D},\mathbf{x}_*)dy_*$. The likelihood call

• [lp,ymu,ys2] = lik(hyp, [], fmu, fs2)

does exactly this. Evaluation of the logarithm of $p_{y_*} = p(y_*|\mathcal{D}, \mathbf{x}_*)$ for values y_* can be done via

• [lp,ymu,ys2] = lik(hyp, y, fmu, fs2)

where 1p contains the number $\ln p_{u_*}$.

Using the moments of the likelihood $\mu(f_*) = \int y_* p(y_*|f_*) dy_*$ and $\sigma^2(f_*) = \int (y_* - \mu(f_*))^2 p(y_*|f_*) dy_*$ we obtain for the predictive moments the following (exact) expressions

$$\begin{array}{lcl} \mu_{y_*} & = & \int \mu(f_*) p(f_* | \mathcal{D}, \mathbf{x}_*) df_*, \text{ and} \\ \\ \sigma_{y_*}^2 & = & \int \left[\sigma^2(f_*) + (\mu(f_*) - \mu_{y_*})^2 \right] p(f_* | \mathcal{D}, \mathbf{x}_*) df_*. \end{array}$$

1. The binary case is simple since $y_* \in \{-1, +1\}$ and $1 = p_{y_*} + p_{-y_*}$. Using $\pi_* = p_{+1}$, we find

$$\begin{array}{lll} p_{y_*} &=& \begin{cases} \pi_* & y_* = +1 \\ 1 - \pi_* & y_* = -1 \end{cases} \\ \mu_{y_*} &=& \sum_{y_* = \pm 1} y_* p(y_* | \mathcal{D}, \mathbf{x}_*) = 2 \cdot \pi_* - 1 \in [-1, 1], \quad \text{and} \\ \sigma_{y_*}^2 &=& \sum_{y_* = \pm 1} (y_* - \mu_{y_*})^2 p(y_* | \mathcal{D}, \mathbf{x}_*) = 4 \cdot \pi_* (1 - \pi_*) \in [0, 1]. \end{array}$$

2. The continuous case for homoscedastic likelihoods depending on $r_* = y_* - f_*$ only and having noise variance $\sigma^2(f_*) = \sigma_n^2$ is also simple since the identity $p(y_*|f_*) = p(y_* - f_*|0)$ allows to substitute $y_* \leftarrow y_* + f_*$ yielding $\mu(f_*) = f_* + \int y_* p(y_*|0) dy_*$ and assuming $\int y_* p(y_*|0) dy_* = 0$ we arrive at

$$\mu_{y_*} = \mu_{f_*}, \text{ and }$$
 $\sigma^2_{u_*} = \sigma^2_{f_*} + \sigma^2_{n}.$

3. The generalised linear model (GLM) case is also feasible. Evaluation of the predictive distribution is done by quadrature

$$p_{y_*} = \int p(y_*|f_*)p(f_*|\mathcal{D}, \mathbf{x}_*)df_* \approx \int p(y_*|f_*)\mathcal{N}(f_*|\mu_{f_*}, \sigma_{f_*}^2)df_*.$$

For GLMs the mean is given by $\mu(f_*) = g(f_*)$ and the variance is usually given by a simple function of the mean $\sigma^2(f_*) = \nu(g(f_*))$, hence we use Gaussian-Hermite quadrature with $\mathcal{N}(f_*|\mu_{f_*},\sigma_{f_*}^2) \approx p(f_*|\mathcal{D},\mathbf{x}_*)$ to compute

$$\begin{array}{lcl} \mu_{y_*} & = & \int g(f_*) p(f_* | \mathbb{D}, \mathbf{x}_*) df_*, \text{ and} \\ \\ \sigma_{y_*}^2 & = & \int \left[\nu(g(f_*)) + (g(f_*) - \mu_{y_*})^2 \right] p(f_* | \mathbb{D}, \mathbf{x}_*) df_* \neq \nu(\mu_{y_*}). \end{array}$$

4. Finally the warped Gaussian likelihood predictive distribution with strictly monotonically increasing warping function g is given by the expression

$$p(y_*|\mathcal{D}, \mathbf{x}_*) = g'(y_*) \mathcal{N}\left(g(y_*)|\mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2\right)$$

so that the predictive moments can be computed by Gaussian-Hermite quadrature.

In the following, we will detail how and which likelihood functions are implemented in the GPML toolbox. Further, we will mention dependencies between likelihoods and inference methods and provide some analytical expressions in addition to some likelihood implementations.

4.2 Interface

19

The likelihoods are in fact the most challenging object in our implementation. Different inference algorithms require different aspects of the likelihood to be computed, therefore the interface is rather involved as detailed below.

 $\langle likFunctions.m \ 19 \rangle \equiv$ 1 % likelihood functions are provided to be used by the gp.m function: 2 % 3 % likErf (Error function, classification, probit regression) 4 % likLogistic classification, logit regression) 5 % (Uniform likelihood, classification) likUni 6 % 7 % likGauss (Gaussian, regression) 8 % likGaussWarp (Warped Gaussian, regression) 9 % (Gumbel likelihood for extremal values) likGumbel 10 % likLaplace (Laplacian or double exponential, regression) 11 % likSech2 (Sech-square, regression) 12 % likT (Student's t, regression) 13 % 14 % likPoisson (Poisson regression, count data) 15 % (Negativ binomial regression, count data) likNegBinom likGamma 16 % (Nonnegative regression, positive data) 17 % likExp (Nonnegative regression, positive data) 18 % likInvGauss (Nonnegative regression, positive data) 19 % (Beta regression, interval data) likBeta 20 % 21 % (Mixture of individual likelihood functions) likMix 22 %

```
23 % The likelihood functions have three possible modes, the mode being selected
24 % as follows (where "lik" stands for any likelihood function in "lik/lik*.m".):
25 %
                                               [REPORT NUMBER OF HYPERPARAMETERS]
26 % 1) With one or no input arguments:
28 %
       s = lik OR s = lik(hyp)
29 %
30 % The likelihood function returns a string telling how many hyperparameters it
31 % expects, using the convention that "D" is the dimension of the input space.
32 % For example, calling "likLogistic" returns the string '0'.
33 %
34 %
35~\% 2) With three or four input arguments:
                                                                  [PREDICTION MODE]
37 %
       lp = lik(hyp, y, mu) OR [lp, ymu, ys2] = lik(hyp, y, mu, s2)
38 %
39 % This allows to evaluate the predictive distribution. Let p(y_*|f_*) be the
40 % likelihood of a test point and N(f_*|mu,s2) an approximation to the posterior
41 % marginal p(f_*|x_*,x,y) as returned by an inference method. The predictive
42 % distribution p(y_*|x_*,x,y) is approximated by.
43 %
      q(y_*) = \inf N(f_*|mu,s2) p(y_*|f_*) df_*
44 %
45 %
     lp = log(q(y)) for a particular value of y, if s2 is [] or 0, this
46 %
                        corresponds to log(p(y|mu))
47 %
     ymu and ys2
                       the mean and variance of the predictive marginal q(y)
48 %
                       note that these two numbers do not depend on a particular
49 %
                       value of y
50 % All vectors have the same size.
51 %
52 %
53 % 3) With five or six input arguments, the fifth being a string [INFERENCE MODE]
55 % [varargout] = lik(hyp, y, mu, s2, inf) OR
56 % [varargout] = lik(hyp, y, mu, s2, inf, i)
58 % There are three cases for inf, namely a) infLaplace, b) infEP and c) infVB.
59 % The last input i, refers to derivatives w.r.t. the ith hyperparameter.
60 %
61 % a1) [lp,dlp,d2lp,d3lp] = lik(hyp, y, f, [], 'infLaplace')
62 % lp, dlp, d2lp and d3lp correspond to derivatives of the log likelihood
63 % log(p(y|f)) w.r.t. to the latent location f.
64 \% lp = log(p(y|f))
65 \% dlp = d log(p(y|f)) / df
66 \% d21p = d^2 log(p(y|f)) / df^2
67 \% d31p = d^3 log(p(y|f)) / df^3
69 % a2) [lp_dhyp,dlp_dhyp,d2lp_dhyp] = lik(hyp, y, f, [], 'infLaplace', i)
70 % returns derivatives w.r.t. to the ith hyperparameter
     lp_dhyp = d log(p(y|f)) / (
72 % dlp_dhyp = d^2 log(p(y|f)) / (df dhyp_i)
73 \% d2lp_dhyp = d^3 log(p(y|f)) / (df^2 dhyp_i)
74 %
75 %
76 % b1) [1Z,d1Z,d21Z] = lik(hyp, y, mu, s2, 'infEP')
77 % let Z = \inf p(y|f) N(f|mu,s2) df then
78 \% 1Z =
              log(Z)
79 \% d1Z = d log(Z) / dmu
80 \% d21Z = d^2 \log(Z) / dmu^2
```

```
82 % b2) [dlZhyp] = lik(hyp, y, mu, s2, 'infEP', i)
 83 % returns derivatives w.r.t. to the ith hyperparameter
 84 \% dlZhyp = d log(Z) / dhyp_i
 85 %
 86 %
 87 % c1) [b,z] = lik(hyp, y, [], ga, 'infVB')
 88 % ga is the variance of a Gaussian lower bound to the likelihood p(y|f).
       p(y|f) \ge exp(b*(f+z) - (f+z).^2/(2*ga) - h(ga)/2) \ge N(f|b*ga-z,ga)
 90 % The function returns the linear part b and z.
 91 %
 92 % Cumulative likelihoods are designed for binary classification. Therefore, they
 93 % only look at the sign of the targets y; zero values are treated as +1.
 95 % Some examples for valid likelihood functions:
 96 %
         lik = @likLogistic;
 97 %
          lik = {'likMix',{'likUni',@likErf}}
 98 %
          lik = {@likPoisson,'logistic'};
99 %
100 % See the help for the individual likelihood for the computations specific to
101 % each likelihood function.
102 %
103 (gpml copyright 5a)
```

4.3 Implemented Likelihood Functions

The following table enumerates all (currently) implemented likelihood functions that can be found at lik/lik<NAME>.m and their respective set of hyperparameters ρ .

lik <name></name>	regression $y_i \in \mathbb{R}$	$p_{\rho}(y_i f_i) =$	$\rho =$	
Gauss	Gaussian	$\mathcal{N}(y_i f_i,\sigma^2) = \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{(y_i-f_i)^2}{2\sigma^2}\right)$	{ln σ}	
GaussWarp	Warped Gaussian	$\mathcal{N}(g_{\theta},(y_i) f_i,\sigma^2)g_{\theta}'(y_i)$	$\{\theta_1,,\theta_{n_g},\ln\sigma\}$	
Gumbel	Gumbel	$\frac{\pi}{\sigma\sqrt{6}}\exp(-z_{i}-e^{-z_{i}}), z_{i}=\gamma+\frac{s\cdot\pi(y_{i}-f_{i})}{\sigma\sqrt{6}}, s =1$	{ln σ}	
Sech2	Sech-squared	$\frac{\pi}{\frac{\sigma\sqrt{6}}{\sqrt{6}}} \exp\left(-z_{i} - e^{-z_{i}}\right), z_{i} = \gamma + \frac{s \cdot \pi(y_{i} - f_{i})}{\sigma\sqrt{6}}, s = 1$ $\frac{\tau}{2\cosh^{2}(\tau(y_{i} - f_{i}))}, \tau = \frac{\pi}{2\sigma\sqrt{3}}$	{ln σ}	
Laplace	Laplacian	$\left(\frac{1}{2b}\exp\left(-\frac{ y_i-f_i }{b}\right),\ b=\frac{\sigma}{\sqrt{2}}\right)$	$\{ln \ \sigma\}$	
Т	Student's t	$\frac{\Gamma(\frac{\nu+1}{2})}{\Gamma(\frac{\nu}{2})} \frac{1}{\sqrt{\nu\pi}\sigma} \left(1 + \frac{(y_i - f_i)^2}{\nu\sigma^2}\right)^{-\frac{\nu+1}{2}}$	$\{\ln(\nu-1), \ln\sigma\}$	
lik <name></name>	classification $y_i \in \{\pm 1\}$	$p_{\rho}(y_i f_i) =$	$\rho =$	
Erf	Error function	$\int_{-\infty}^{y_i f_i} \mathcal{N}(t) dt$	Ø	
Logistic	Logistic function	$\frac{1}{1+\exp(-y_i f_i)}$	Ø	
Uni	Label noise	$\frac{1}{2}$	Ø	
lik <name></name>	count data $y_i \in \mathbb{N}$	$p_{\rho}(y_i f_i) = $ where $\mu = g(f_i)$	$\rho =$	
Poisson	Poisson	$\mu^{y_i}e^{-\mu}/y_i!$	Ø	
NegBinom	Negative Binomial	${y_i+r-1 \choose y_i} r^r \mu^{y_i}/(r+\mu)^{r+y_i}$	$\{\ln r\}$	
lik <name></name>	nonnegative data $y_i \in \mathbb{R}_+ \setminus \{0\}$	$p_{\rho}(y_i f_i) = $ where $\mu = g(f_i)$	ρ =	
Weibull	Weibull, $\gamma_1 = \Gamma(1 + 1/\kappa)$	$\kappa \gamma_1 / \mu (y_i \gamma_1 / \mu)^{\kappa - 1} \exp(-(y_i \gamma_1 / \mu)^{\kappa})$	{ln κ}	
Gamma	Gamma	$\frac{\alpha^{\alpha}y_{i}^{\alpha-1}}{\Gamma(\alpha)}\mu^{-\alpha}\exp\left(-\frac{y_{i}\alpha}{\mu}\right)$	{ln α}	
Exp	Exponential	$\mu^{-1} \exp\left(-\frac{y_i}{\mu}\right)$	Ø	
InvGauss	Inverse Gaussian	$\sqrt{\frac{\lambda}{2\pi y_1^3}} \exp\left(-\frac{\lambda (y_1-\mu)^2}{2\mu^2 y_1}\right)$	$\{\ln\lambda\}$	
lik <name></name>	interval data $y_i \in [0, 1]$	$p_{\rho}(y_i f_i) = $ where $\mu = g(f_i)$	ρ =	
Beta	Beta	$\begin{array}{ll} p_{\rho}(y_i f_i) = & \text{where } \mu = g(f_i) \\ \frac{\Gamma(\varphi)}{\Gamma(\mu\varphi)\Gamma((1-\mu)\varphi)} y_i^{\mu\varphi-1} (1-y_i)^{(1-\mu)\varphi-1} \end{array}$	{ln φ}	
Composite l	ikelihood functions $[p_1(y_i f_i), p_1$	$(y_i f_i),]\mapsto p_{\rho}(y_i f_i)$		
Mix	Mixture	$\sum_{j} \alpha_{j} p_{j}(y_{i} f_{i})$	$\{\ln \alpha_1, \ln \alpha_2,\}$	

4.4 Usage of Implemented Likelihood Functions

Some code examples taken from doc/usageLik.m illustrate how to use simple and composite likelihood functions to specify a GP model.

Syntactically, a likelihood function 1f is defined by

```
lk := 'func' | @func // simple
lf := {lk} | {param, lk} | {lk, {lk, .., lk}} // composite
```

i.e., it is either a string containing the name of a likelihood function, a pointer to a likelihood function or one of the former in combination with a cell array of likelihood functions and an additional list of parameters.

```
\langle doc/usageLik.m 22 \rangle \equiv
22
     1 % demonstrate usage of likelihood functions
     3 % See also likFunctions.m.
     4 %
     5 (gpml copyright 5a)
     6 clear all, close all
     7 n = 5; f = randn(n,1);
                               % create random latent function values
     9 % set up simple classification likelihood functions
     10 \text{ yc} = \text{sign}(f);
     11 lc0 = {'likErf'};
                              hypc0 = [];  % no hyperparameters are needed
     12 lc1 = {@likLogistic}; hypc1 = []; % also function handles are OK
     13 lc2 = {'likUni'};
                              hypc2 = [];
     14 lc3 = {'likMix', {'likUni', @likErf}}; hypc3 = log([1;2]); %mixture
     1.5
     16 % set up simple regression likelihood functions
     17 \text{ yr} = f + randn(n,1)/20;
     18 \text{ sn} = 0.1;
                                                  % noise standard deviation
     19 lr0 = {'likGauss'}; hypr0 = log(sn);
     20 lr1 = {'likLaplace'}; hypr1 = log(sn);
     21 lr2 = {'likSech2'}; hypr2 = log(sn);
     22 \text{ nu} = 4;
                                              % number of degrees of freedom
     23 lr3 = {'likT'}; hypr3 = [log(nu-1); log(sn)];
     24 lr4 = {'likMix',{lr0,lr1}}; hypr4 = [log([1,2]);hypr0;hypr1];
     26 a = 1; % set up warped Gaussian with g(y) = y + a*sign(y).*y.^2
     27 1r5 = {'likGaussWarp',['poly2']}; hypr5 = log([a;sn]);
     28 lr6 = {'likGumbel', '+'}; hypr6 = log(sn);
     29
     30 % set up Poisson/negative binomial regression
     31 \text{ yp} = \text{fix}(abs(f)) + 1;
     32 lp0 = {@likPoisson,'logistic'};
                                                hypp0 = [];
     33 lp1 = {@likPoisson,{'logistic2',0.1}}; hypp1 = [];
     34 lp2 = {@likPoisson,'exp'};
                                                hypp2 = [];
     35 ln1 = {@likNegBinom, 'logistic2'};
                                                hypn1 = [];
     37 % set up other GLM likelihoods for positive or interval regression
    38 lg1 = {@likGamma, 'logistic'}; al = 2; hyp.lik = log(al);
    39 lg2 = {@likInvGauss,'exp'};
                                                lam = 1.1; hyp.lik = log(lam);
                                       phi = 2.1; hyp.lik = log(phi);
     40 lg3 = {@likBeta,'expexp'};
    41 lg4 = {@likBeta, 'logit'};
                                                phi = 4.7; hyp.lik = log(phi);
    42 lg5 = {@likWeibull,{'logistic2',0.01}}; ka = 0.5; hyp.lik = log(ka);
```

44 % 0) specify the likelihood function

```
45 lik = 1c0; hyp = hypc0; y = yc;
46 % lik = lr4; hyp = hypr4; y = yr;
47 \% lik = lp1; hyp = hypp1; y = yp;
49 % 1) query the number of parameters
50 feval(lik{:})
52 % 2) evaluate the likelihood function on f
53 exp(feval(lik{:},hyp,y,f))
55 % 3a) evaluate derivatives of the likelihood
56 [lp,dlp,d2lp,d3lp] = feval(lik{:}, hyp, y, f, [], 'infLaplace');
58 % 3b) compute Gaussian integrals w.r.t. likelihood
59 \text{ mu} = f; s2 = rand(n,1);
60 [1Z,d1Z,d21Z] = feval(lik{:}, hyp, y, mu, s2, 'infEP');
61
62 % 3c) obtain lower bound on likelihood
63 \text{ ga} = \text{rand}(n,1);
64 [b,z] = feval(lik{:}, hyp, y, [], ga, 'infVB');
```

4.5 Compatibility Between Likelihoods and Inference Methods

The following table lists all possible combinations of likelihood function and inference methods.

Likelihood \ Inference	Gaussian Likelihood	Laplace	VB	EP	KL	MCMC	LOO	Type, Output Domain	Alternative Names
	approx. cov.	possible, T	able3.6	no approx. cov. possible, Table3.6					
Gaussian	✓	✓	√	✓	√	✓	✓	regression, R	
Warped Gaussian		✓	✓	✓	√	✓	✓	regression, R	
Gumbel		✓			√	✓	✓	regression, R	
Sech-squared		✓	✓	✓	√	✓	✓	regression, R	logistic distribution
Laplacian		✓	✓	✓	√	✓	✓	regression, R	double exponential
Student's t		✓	✓		√	✓	✓	regression, R	
Mixture		✓		✓	√	✓	✓		mixing meta likelihood
Error function		✓		✓	√	✓	✓	classification, {±1}	probit regression
Logistic function		✓	✓	✓	√	✓	✓	classification, {±1}	logit regression
Uniform		√	✓	✓	√	✓	✓	classification, {±1}	label noise
Weibull		✓				✓	✓	positive data, $\mathbb{R}_+ \setminus \{0\}$	nonnegative regression
Gamma		✓				✓	✓	positive data, $\mathbb{R}_+ \setminus \{0\}$	nonnegative regression
Exp		✓				✓	✓	positive data, $\mathbb{R}_+ \setminus \{0\}$	nonnegative regression
Inverse Gaussian		✓				✓	✓	positive data, $\mathbb{R}_+ \setminus \{0\}$	nonnegative regression
Poisson		√		(√)*	√	√	√	count data, N	Poisson regression
Negative Binomial		✓				✓	✓	count data, N	negative binomial regression
Beta		✓				√	√	interval data, [0, 1]	beta regression

$(\checkmark)^*$ EP might not converge in some cases since quadrature is used.

Exact inference is only tractable for Gaussian likelihoods. Expectation propagation together with Student's t likelihood is inherently unstable due to non-log-concavity. Laplace's approximation for Laplace likelihoods is not sensible because at the mode the curvature and the gradient is undefined due to the non-differentiable peak of the Laplace distribution. Special care has been taken for the non-convex optimisation problem imposed by the combination Student's t likelihood and Laplace's approximation.

4.6 Gaussian Likelihood

The Gaussian likelihood is the simplest likelihood because the posterior distribution is not only Gaussian but can be computed analytically. In principle, the Gaussian likelihood would only be

operated in conjunction with the exact inference method but we chose to provide compatibility with all other inference algorithms as well because it enables code testing and allows to switch between different regression likelihoods very easily.

```
24a
       \langle lik/likGauss.m 24a \rangle \equiv
       1 function [varargout] = likGauss(hyp, y, mu, s2, inf, i)
       3 % likGauss - Gaussian likelihood function for regression. The expression for the
       4 % likelihood is
       5 \% likGauss(t) = exp(-(t-y)^2/2*sn^2) / sqrt(2*pi*sn^2),
       6 \% where y is the mean and sn is the standard deviation.
       7 %
       8 % The hyperparameters are:
      9 %
      10 \% \text{ hyp} = [\log(\text{sn})]
      12 % Several modes are provided, for computing likelihoods, derivatives and moments
      13 % respectively, see likFunctions.m for the details. In general, care is taken
      14 % to avoid numerical issues when the arguments are extreme.
      16 \(\langle gpml \copyright 5a\rangle
      17 %
      18 % See also LIKFUNCTIONS.M.
      20 if nargin<3, varargout = {'1'}; return; end % report number of hyperparameters
      22 sn2 = exp(2*hyp);
      23
      24 if nargin<5
                                                      % prediction mode if inf is not present
         (Prediction with Gaussian likelihood 24b)
      25
      26 else
      27 switch inf
      28 case 'infLaplace'
      29
           (Laplace's method with Gaussian likelihood 25a)
      30 case 'infEP'
      31
            (EP inference with Gaussian likelihood 25b)
      32
         case 'infVB'
      33
             (Variational Bayes inference with Gaussian likelihood 25c)
      34
           end
      35 \text{ end}
24b
      ⟨Prediction with Gaussian likelihood 24b⟩≡
                                                                       (24a)
       1 if isempty(y), y = zeros(size(mu)); end
       2 s2zero = 1; if nargin>3&&numel(s2)>0&&norm(s2)>eps, s2zero = 0; end % s2==0 ?
       3 if s2zero
                                                                           % log probability
          lp = -(y-mu).^2./sn2/2-log(2*pi*sn2)/2; s2 = 0;
           lp = likGauss(hyp, y, mu, s2, 'infEP');
                                                                                 % prediction
       7 end
       8 \text{ ymu} = \{\}; \text{ ys2} = \{\};
       9 if nargout>1
      10 ymu = mu;
                                                                             % first y moment
      if nargout>2
      12
           ys2 = s2 + sn2;
                                                                            % second y moment
      13 end
      14 end
      15 varargout = {lp,ymu,ys2};
```

The Gaussian likelihood function has a single hyperparameter ρ , the log of the noise standard deviation σ_n .

4.6.1 Exact Inference

Exact inference doesn't require any specific likelihood related code; all computations are done directly by the inference method, section 3.1.

4.6.2 Laplace's Approximation

```
\langle Laplace's method with Gaussian likelihood 25a \rangle \equiv
25a
                                                                 (24a)
      1 if nargin<6
                                                                % no derivative mode
          if isempty(y), y=0; end
          ymmu = y-mu; dlp = {}; d2lp = {}; d3lp = {};
          lp = -ymmu.^2/(2*sn2) - log(2*pi*sn2)/2;
      5
         if nargout>1
      6
          dlp = ymmu/sn2;
                                                 % dlp, derivative of log likelihood
      7
          if nargout>2
                                            % d2lp, 2nd derivative of log likelihood
      8
            d2lp = -ones(size(ymmu))/sn2;
      9
             if nargout>3
                                            % d3lp, 3rd derivative of log likelihood
               d3lp = zeros(size(ymmu));
     10
     11
             end
     12
            end
     13
         end
     varargout = {lp,dlp,d2lp,d3lp};
     15 else
                                                                   % derivative mode
          lp_dhyp = (y-mu).^2/sn2 - 1;  % derivative of log likelihood w.r.t. hypers
     16
     17
          dlp_dhyp = 2*(mu-y)/sn2;
                                                                % first derivative,
     d2lp_dhyp = 2*ones(size(mu))/sn2; % and also of the second mu derivative
     19
          varargout = {lp_dhyp,dlp_dhyp,d2lp_dhyp};
     20 end
```

4.6.3 Expectation Propagation

```
25b
      \langle EP \text{ inference with Gaussian likelihood 25b} \rangle \equiv
                                                                (24a)
      1 if nargin<6
                                                               % no derivative mode
        1Z = -(y-mu).^2./(sn2+s2)/2 - log(2*pi*(sn2+s2))/2; % log part function
         d1Z = {}; d21Z = {};
      4
        if nargout>1
                                             % 1st derivative w.r.t. mean
      5
           dlZ = (y-mu)./(sn2+s2);
            if nargout>2
      6
      7
                                                % 2nd derivative w.r.t. mean
             d21Z = -1./(sn2+s2);
      8
            end
      9
        end
     10  varargout = {1Z,d1Z,d21Z};
                                                                  % derivative mode
     11 else
          d1Zhyp = ((y-mu).^2./(sn2+s2)-1)./(1+s2./sn2); % deriv. w.r.t. hyp.lik
     12
     varargout = {dlZhyp};
     14 end
```

4.6.4 Variational Bayes

25c $\langle Variational \ Bayes \ inference \ with \ Gaussian \ likelihood \ 25c \rangle \equiv (24a)$

```
1 % variational lower site bound

2 % t(s) = \exp(-(y-s)^2/2sn^2)/sqrt(2*pi*sn^2)

3 % the bound has the form: (b+z/ga)*f - f.^2/(2*ga) - h(ga)/2

4 n = numel(s2); b = zeros(n,1); y = y.*ones(n,1); z = y;

5 varargout = \{b,z\};
```

4.7 Warped Gaussian Likelihood

Starting from the likelihood p(y|f) we are sometimes facing the situation where the data $y \in \mathcal{Y} \subseteq \mathbb{R}$ is not distributed according to p(y|f) but some nonlinear transformation of the data g(y) = z so that $z \sim p(z|f)$. Here, the warping function $g: \mathcal{Y} \to \mathbb{R}$ needs to be strictly monotonically increasing i.e. g'(y) > 0. Formally, we start from the fact that p(z|f) integrates to one and use the derivative dz = g'(y)dy to substitute

$$\int p(z|f)dz = 1 = \int p_g(y|f)dy, \ p_g(y|f) = p(g(y)|f)g'(y)$$

where we have defined the log warped likelihood $\ln p_g(y|f) = \ln p(g(y)|f) + \ln g'(y)$. The interesting bit is that approximate inference methods such as infExact, infLaplace, infEP, infVB, infKL remain fully feasible; only prediction and derivatives become more involved. The usual GP inference is recovered by using the identity warping function $g: y \mapsto y$. The construction works in princple for any likelihood but our implementation in likGaussWarp is limited to the Gaussian likelihood.

Hyperparameter derivatives

Hyperparameter derivatives for infLaplace are obtained as follows

$$\begin{split} \frac{\partial}{\partial \theta} \ln \frac{\partial^k}{\partial f^k} p_g(y|f) &= \frac{\partial}{\partial \theta} \ln \frac{\partial^k}{\partial f} p(g(y)|f) + \frac{\partial}{\partial \theta} \frac{\partial^k}{\partial f^k} \ln g'(y), \ k = 0, 1, 2 \\ &= -\frac{\partial^{k+1}}{\partial f^{k+1}} \ln p(g(y)|f) \frac{\partial}{\partial \theta} g(y) + \frac{\partial}{\partial \theta} \frac{\partial^k}{\partial f^k} \ln g'(y). \end{split}$$

Similarly for infEP the derivatives are given by

$$\begin{split} \frac{\partial}{\partial \theta} \ln \int p_g(y|f) \mathcal{N}(f|\mu,\sigma^2) df &= \frac{\partial}{\partial \theta} \ln \int p(g(y)|f) \mathcal{N}(f|\mu,\sigma^2) df + \frac{\partial}{\partial \theta} \ln g'(y) \\ &= -\frac{\partial}{\partial \mu} \ln \int p(g(y)|f) \mathcal{N}(f|\mu,\sigma^2) df \frac{\partial}{\partial \theta} g(y) + \frac{\partial}{\partial \theta} \ln g'(y). \end{split}$$

This trick above works for any homoscedastic likelihood where $p(y|f) = p(y + \beta|f + \beta)$ such as likGauss, likLaplace, likSech2 and likT.

Predictive moments

As detailed in 4, the predictive distribution is - for Gaussian likelihood - given by

$$\begin{split} p(z_*|\mathcal{D}, \mathbf{x}_*) &= \int p(z_*|f_*) p(f_*|\mathcal{D}, \mathbf{x}_*) df_* = \int \mathcal{N}(z_*|f_*, \sigma_n^2) \mathcal{N}(f_*|\mu_{f_*}, \sigma_{f_*}^2) df_* \\ &= \mathcal{N}(z_*|\mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2), \text{ where } z_* = g(y_*) \\ p(y_*|\mathcal{D}, \mathbf{x}_*) &= g'(y_*) \mathcal{N}(g(y_*)|\mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2). \end{split}$$

Hence, the predictive moments are obtained by the 1d integrals

$$\begin{split} \mu_{y_*} &= \int y_* g'(y_*) \mathcal{N}(g(y_*) | \mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2) dy_* \\ &= \int g^{-1}(z_*) \mathcal{N}(z_* | \mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2) dz_*, \text{ and} \\ \sigma_{y_*}^2 &= \int (y_* - \mu_{y_*})^2 g'(y_*) \mathcal{N}(g(y_*) | \mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2) dy_* \\ &= \int (g^{-1}(z_*) - \mu_{y_*})^2 \mathcal{N}(z_* | \mu_{f_*}, \sigma_n^2 + \sigma_{f_*}^2) dz_*. \end{split}$$

4.8 Gumbel Likelihood

Distributions of extrema are well captured by the Gumbel distribution

$$p(y) = \frac{1}{\beta} \exp(-z - e^{-z}), z = s \frac{y - \eta}{\beta}, s \in \{\pm 1\}$$

with mean $\mu=\eta+\beta\gamma$ and variance $\sigma^2=\pi^2\beta^2/6$ where $\gamma=0.57721566490153$ denotes Euler–Mascheroni's constant. Skewness is approximately given by 1.1395s where s is a sign switching between left and right skewness and kurtosis is 12/5. The final expression for the Gumbel likelihood is

$$p(y|f) = \frac{\pi}{\sigma\sqrt{6}} \exp\left(-z - e^{-z}\right), \ z = \gamma + s \frac{\pi}{\sigma\sqrt{6}} (y - f), \ s \in \{\pm 1\}.$$

4.9 Laplace Likelihood

Laplace's Approximation

The following derivatives are needed:

$$\ln p(y|f) = -\ln(2b) - \frac{|f-y|}{b}$$

$$\frac{\partial \ln p}{\partial f} = \frac{\operatorname{sign}(f-y)}{b}$$

$$\frac{\partial^2 \ln p}{(\partial f)^2} = \frac{\partial^3 \ln p}{(\partial f)^3} = \frac{\partial^3 \ln p}{(\partial \ln \sigma_n)(\partial f)^2} = 0$$

$$\frac{\partial \ln p}{\partial \ln \sigma_n} = \frac{|f-y|}{b} - 1$$

Expectation Propagation

Expectation propagation requires integration against a Gaussian measure for moment matching,

We need to evaluate $\ln Z = \ln \int \mathcal{L}(y|f,\sigma_n^2) \mathcal{N}(f|\mu,\sigma^2) df$ as well as the derivatives $\frac{\partial \ln Z}{\partial \mu}$ and $\frac{\partial^2 \ln Z}{\partial \mu^2}$ where $\mathcal{N}(f|\mu,\sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(f-\mu)^2}{2\sigma^2}\right)$, $\mathcal{L}(y|f,\sigma_n^2) = \frac{1}{2b} \exp\left(-\frac{|y-f|}{b}\right)$, and $b = \frac{\sigma_n}{\sqrt{2}}$. As a first

step, we reduce the number of parameters by means of the substitution $\tilde{f}=\frac{f-y}{\sigma_{n}}$ yielding

$$\begin{split} Z &= \int \mathcal{L}(y|f,\sigma_n^2) \mathcal{N}(f|\mu,\sigma^2) df \\ &= \frac{1}{\sqrt{2\pi}\sigma} \frac{\sqrt{2}}{2\sigma_n} \int \exp\left(-\frac{(f-\mu)^2}{2\sigma^2}\right) \exp\left(-\sqrt{2}\frac{|f-y|}{\sigma_n}\right) df \\ &= \frac{\sqrt{2}}{2\sigma\sqrt{2\pi}} \int \exp\left(-\frac{(\sigma_n\tilde{f}+y-\mu)^2}{2\sigma^2}\right) \exp\left(-\sqrt{2}|\tilde{f}|\right) d\tilde{f} \\ &= \frac{\sigma_n}{\sigma\sigma_n\sqrt{2\pi}} \int \exp\left(-\frac{\sigma_n^2\left(\tilde{f}-\frac{\mu-y}{\sigma_n}\right)^2}{2\sigma^2}\right) \mathcal{L}(\tilde{f}|0,1) d\tilde{f} \\ &= \frac{1}{\sigma_n} \int \mathcal{L}(f|0,1) \mathcal{N}(f|\tilde{\mu},\tilde{\sigma}^2) df \\ \ln Z &= \ln \tilde{Z} - \ln \sigma_n = \ln \int \mathcal{L}(f|0,1) \mathcal{N}(f|\tilde{\mu},\tilde{\sigma}^2) df - \ln \sigma_n \end{split}$$

with $\tilde{\mu} = \frac{\mu - y}{\sigma_n}$ and $\tilde{\sigma} = \frac{\sigma}{\sigma_n}$. Thus, we concentrate on the simpler quantity $\ln \tilde{Z}$.

$$\begin{split} \ln \mathsf{Z} &= \ln \int \exp \left(-\frac{(f-\tilde{\mu})^2}{2\tilde{\sigma}^2} - \sqrt{2} |f| \right) df - \ln \tilde{\sigma} \sqrt{2\pi} - \ln \sqrt{2} \sigma_n \\ &= \ln \left[\int_{-\infty}^0 \exp \left(-\frac{(f-\tilde{\mu})^2}{2\tilde{\sigma}^2} + \sqrt{2} f \right) df + \int_0^\infty \exp \left(-\frac{(f-\tilde{\mu})^2}{2\tilde{\sigma}^2} - \sqrt{2} f \right) df \right] + C \\ &= \ln \left[\int_{-\infty}^0 \exp \left(-\frac{f^2 - 2(\tilde{\mu} + \tilde{\sigma}^2 \sqrt{2})f + \tilde{\mu}^2}{2\tilde{\sigma}^2} \right) df + \int_0^\infty \exp \left(-\frac{f^2 - 2(\tilde{\mu} - \tilde{\sigma}^2 \sqrt{2})f + \tilde{\mu}^2}{2\tilde{\sigma}^2} \right) df \right] + C \\ &= \ln \left[\exp \left(\frac{m_-^2}{2\tilde{\sigma}^2} \right) \int_{-\infty}^0 \exp \left(-\frac{(f-m_-)^2}{2\tilde{\sigma}^2} \right) df + \exp \left(\frac{m_+^2}{2\tilde{\sigma}^2} \right) \int_0^\infty \exp \left(-\frac{(f-m_+)^2}{2\tilde{\sigma}^2} \right) df \right] - \frac{\tilde{\mu}^2}{2\tilde{\sigma}^2} + C \\ &= \ln \left[\exp \left(\frac{m_-^2}{2\tilde{\sigma}^2} \right) \int_{-\infty}^0 \mathcal{N}(f|m_-, \tilde{\sigma}^2) df + \exp \left(\frac{m_+^2}{2\tilde{\sigma}^2} \right) \left(1 - \int_{-\infty}^0 \mathcal{N}(f|m_+, \tilde{\sigma}^2) df \right) \right] - \frac{\tilde{\mu}^2}{2\tilde{\sigma}^2} - \ln \sqrt{2} \sigma_n \\ &= \ln \left[\exp \left(\frac{m_-^2}{2\tilde{\sigma}^2} \right) \Phi \left(\frac{m_-}{\tilde{\sigma}} \right) - \exp \left(\frac{m_+^2}{2\tilde{\sigma}^2} \right) \Phi \left(\frac{m_+}{\tilde{\sigma}} \right) + \exp \left(\frac{m_+^2}{2\tilde{\sigma}^2} \right) \right] - \frac{\tilde{\mu}^2}{2\tilde{\sigma}^2} - \ln \sqrt{2} \sigma_n \end{split}$$

Here, $\Phi(z) = \int_{-\infty}^{z} \mathcal{N}(f|0,1) df$ denotes the cumulative Gaussian distribution. Finally, we have

$$\begin{split} \ln \mathsf{Z} &= & \ln \left[\exp \left(- \sqrt{2} \tilde{\mu} \right) \Phi \left(\frac{m_-}{\tilde{\sigma}} \right) + \exp \left(\sqrt{2} \tilde{\mu} \right) \Phi \left(- \frac{m_+}{\tilde{\sigma}} \right) \right] + \tilde{\sigma}^2 - \ln \sqrt{2} \sigma_n \\ &= & \ln \left[\exp \left(\underbrace{\ln \Phi(-z_+) + \sqrt{2} \tilde{\mu}}_{\alpha_+} \right) + \exp \left(\underbrace{\ln \Phi(z_-) - \sqrt{2} \tilde{\mu}}_{\alpha_-} \right) \right] + \tilde{\sigma}^2 - \ln \sqrt{2} \sigma_n \\ &= & \ln (e^{\alpha_+} + e^{\alpha_-}) + \tilde{\sigma}^2 - \ln \sqrt{2} \sigma_n \end{split}$$

where
$$z_{+} = \frac{\tilde{\mu}}{\tilde{\sigma}} + \tilde{\sigma}\sqrt{2} = \frac{\mu - y}{\sigma} + \frac{\sigma}{\sigma_{n}}\sqrt{2}$$
, $z_{-} = \frac{\tilde{\mu}}{\tilde{\sigma}} - \tilde{\sigma}\sqrt{2} = \frac{\mu - y}{\sigma} - \frac{\sigma}{\sigma_{n}}\sqrt{2}$ and $\tilde{\mu} = \frac{\mu - y}{\sigma_{n}}$, $\tilde{\sigma} = \frac{\sigma}{\sigma_{n}}$.
Now, using $\frac{d}{d\theta} \ln \Phi(z) = \frac{1}{\Phi(z)} \frac{d}{d\theta} \Phi(z) = \frac{\mathcal{N}(z)}{\Phi(z)} \frac{dz}{d\theta}$ we tackle first derivative

$$\begin{array}{ll} \frac{\partial \ln Z}{\partial \mu} &=& \frac{e^{\alpha_{+}} \frac{\partial \alpha_{+}}{\partial \mu} + e^{\alpha_{-}} \frac{\partial \alpha_{-}}{\partial \mu}}{e^{\alpha_{+}} + e^{\alpha_{-}}} \\ \frac{\partial \alpha_{+}}{\partial \mu} &=& \frac{\partial}{\partial \mu} \ln \Phi(-z_{+}) + \frac{\sqrt{2}}{\sigma_{n}} \\ &=& -\frac{\mathcal{N}(-z_{+})}{\sigma \Phi(-z_{+})} + \frac{\sqrt{2}}{\sigma_{n}} = -\frac{q_{+}}{\sigma} + \frac{\sqrt{2}}{\sigma_{n}} \\ \frac{\partial \alpha_{-}}{\partial \mu} &=& \frac{\partial}{\partial \mu} \ln \Phi(z_{-}) - \frac{\sqrt{2}}{\sigma_{n}} \\ &=& \frac{\mathcal{N}(z_{-})}{\sigma \Phi(z_{-})} - \frac{\sqrt{2}}{\sigma_{n}} = \frac{q_{-}}{\sigma} - \frac{\sqrt{2}}{\sigma_{n}} \\ \frac{\partial \alpha_{\pm}}{\partial \mu} &=& \mp \frac{q_{\pm}}{\sigma} \pm \frac{\sqrt{2}}{\sigma_{n}}. \end{array}$$

as well as the second derivative

$$\begin{split} \frac{\partial^2 \ln Z}{\partial \mu^2} &= \frac{\frac{\partial}{\partial \mu} \left(e^{\alpha_+} \frac{\partial \alpha_+}{\partial \mu}\right) + \frac{\partial}{\partial \mu} \left(e^{\alpha_-} \frac{\partial \alpha_-}{\partial \mu}\right)}{e^{\alpha_+} + e^{\alpha_-}} - \left(\frac{\partial \ln Z}{\partial \mu}\right)^2 \\ \frac{\partial}{\partial \mu} \left(e^{\alpha_\pm} \frac{\partial \alpha_\pm}{\partial \mu}\right) &= e^{\alpha_\pm} \left[\left(\frac{\partial \alpha_\pm}{\partial \mu}\right)^2 + \frac{\partial^2 \alpha_\pm}{\partial \mu^2}\right] \\ \frac{\partial^2 \alpha_+}{\partial \mu^2} &= -\frac{1}{\sigma} \frac{\frac{\partial}{\partial \mu} \mathcal{N}(-z_+) \Phi(-z_+) - \frac{\partial}{\partial \mu} \Phi(-z_+) \mathcal{N}(-z_+)}{\Phi^2(-z_+)} \\ &= -\frac{1}{\sigma} \frac{\mathcal{N}(-z_+) \Phi(-z_+) \frac{\partial - z_+^2/2}{\partial \mu} - \mathcal{N}^2(-z_+) \frac{\partial - z_+}{\partial \mu}}{\Phi^2(-z_+)} \\ &= \frac{\mathcal{N}(-z_+)}{\sigma^2} \cdot \frac{\Phi(-z_+)z_+ - \mathcal{N}(-z_+)}{\Phi^2(-z_+)} = -\frac{q_+^2 - q_+ z_+}{\sigma^2} \\ \frac{\partial^2 \alpha_-}{\partial \mu^2} &= \frac{1}{\sigma} \frac{\frac{\partial}{\partial \mu} \mathcal{N}(z_-) \Phi(z_-) - \frac{\partial}{\partial \mu} \Phi(z_-) \mathcal{N}(z_-)}{\Phi^2(z_-)} \\ &= \frac{1}{\sigma} \frac{\mathcal{N}(z_-) \Phi(z_-) \frac{\partial - z_-^2/2}{\partial \mu} - \mathcal{N}^2(z_-) \frac{\partial z_-}{\partial \mu}}{\Phi^2(z_-)} \\ &= \frac{\mathcal{N}(z_-)}{\sigma^2} \cdot \frac{-\Phi(z_-)z_- - \mathcal{N}(z_-)}{\Phi^2(z_-)} = -\frac{q_-^2 + q_- z_-}{\sigma^2} \\ \frac{\partial^2 \alpha_\pm}{\partial \mu^2} &= -\frac{q_\pm^2 \mp q_\pm z_\pm}{\sigma^2} \end{split}$$

which can be simplified to

$$\frac{\partial^2 \ln Z}{\partial \mu^2} = \frac{e^{\alpha_+}b_+ + e^{\alpha_-}b_-}{e^{\alpha_+} + e^{\alpha_-}} - \left(\frac{\partial \ln Z}{\partial \mu}\right)^2$$

using

$$\begin{split} b_{\pm} &= \left(\frac{\partial \alpha_{\pm}}{\partial \mu}\right)^2 + \frac{\partial^2 \alpha_{\pm}}{\partial \mu^2} &= \left(\mp \frac{q_{\pm}}{\sigma} \pm \frac{\sqrt{2}}{\sigma_n}\right)^2 - \frac{q_{\pm}^2 \mp q_{\pm} z_{\pm}}{\sigma^2} \\ &= \left(\frac{q_{\pm}}{\sigma} - \frac{\sqrt{2}}{\sigma_n}\right)^2 - \frac{q_{\pm}^2}{\sigma^2} \pm \frac{q_{\pm} z_{\pm}}{\sigma^2} \\ &= \frac{2}{\sigma_n^2} - \left(\frac{\sqrt{8}}{\sigma \sigma_n} \mp \frac{z_{\pm}}{\sigma^2}\right) q_{\pm}. \end{split}$$

We also need

$$\frac{\partial \ln Z}{\partial \ln \sigma_n} = \frac{e^{\alpha_+} \frac{\partial \alpha_+}{\partial \ln \sigma_n} + e^{\alpha_-} \frac{\partial \alpha_-}{\partial \ln \sigma_n}}{e^{\alpha_+} + e^{\alpha_-}} - \frac{2\sigma^2}{\sigma_n^2} - 1.$$

Variational Bayes

We need $h(\gamma)$ and its derivatives as well as $\beta(\gamma)$:

$$h(\gamma) = \frac{2}{\sigma_n^2} \gamma + \ln(2\sigma_n^2) + y^2 \gamma^{-1}$$

$$h'(\gamma) = \frac{2}{\sigma_n^2} - y^2 \gamma^{-2}$$

$$h''(\gamma) = 2y^2 \gamma^{-3}$$

$$\beta(\gamma) = y\gamma^{-1}$$

4.10 Student's t Likelihood

The likelihood has two hyperparameters (both represented in the log domain to ensure positivity): the degrees of freedom ν and the scale σ_n with mean y (for $\nu > 1$) and variance $\frac{\nu}{\nu-2}\sigma_n^2$ (for $\nu > 2$).

$$p(y|f) = Z \cdot \left(1 + \frac{(f - y)^2}{v\sigma_n^2}\right)^{-\frac{v + 1}{2}}, \quad Z = \frac{\Gamma\left(\frac{v + 1}{2}\right)}{\Gamma\left(\frac{v}{2}\right)\sqrt{v\pi\sigma_n^2}}$$

Laplace's Approximation

For the mode fitting procedure, we need derivatives up to third order; the hyperparameter derivatives at the mode require some mixed derivatives. All in all, using r = y - f, we have

$$\begin{split} &\ln p(y|f) &= &\ln \Gamma \left(\frac{\nu+1}{2} \right) - \ln \Gamma \left(\frac{\nu}{2} \right) - \frac{1}{2} \ln \nu \pi \sigma_n^2 - \frac{\nu+1}{2} \ln \left(1 + \frac{r^2}{\nu \sigma_n^2} \right) \\ &\frac{\partial \ln p}{\partial f} &= (\nu+1) \frac{r}{r^2 + \nu \sigma_n^2} \\ &\frac{\partial^2 \ln p}{(\partial f)^2} &= (\nu+1) \frac{r^2 - \nu \sigma_n^2}{(r^2 + \nu \sigma_n^2)^2} \\ &\frac{\partial^3 \ln p}{(\partial f)^3} &= 2(\nu+1) \frac{r^3 - 3r \nu \sigma_n^2}{(r^2 + \nu \sigma_n^2)^3} \\ &\frac{\partial \ln p}{\partial \ln \nu} &= \frac{\partial Z}{\partial \ln \nu} - \frac{\nu}{2} \ln \left(1 + \frac{r^2}{\nu \sigma_n^2} \right) + \frac{\nu+1}{2} \cdot \frac{r^2}{r^2 + \nu \sigma_n^2} \\ &\frac{\partial Z}{\partial \ln \nu} &= \frac{\nu}{2} \frac{d \ln \Gamma \left(\frac{\nu+1}{2} \right)}{d \ln \nu} - \frac{\nu}{2} \frac{d \ln \Gamma \left(\frac{\nu}{2} \right)}{d \ln \nu} - \frac{1}{2} \\ &\frac{\partial^3 \ln p}{(\partial \ln \nu)(\partial f)^2} &= \nu \frac{r^2 (r^2 - 3(\nu+1)\sigma_n^2) + \nu \sigma_n^2}{(r^2 + \nu \sigma_n^2)^3} \\ &\frac{\partial \ln p}{\partial \ln \sigma_n} &= (\nu+1) \frac{r^2}{r^2 + \nu \sigma_n^2} - 1 \\ &\frac{\partial^3 \ln p}{(\partial \ln \sigma_n)(\partial f)^2} &= 2\nu \sigma_n^2 (\nu+1) \frac{\nu \sigma_n^2 - 3r^2}{(r^2 + \nu \sigma_n^2)^3} \end{split}$$

4.11 Cumulative Logistic Likelihood

The likelihood has one hyperparameter (represented in the log domain), namely the standard deviation σ_n

$$p(y|f) = Z \cdot cosh^{-2}\left(\tau(f-y)\right), \ \tau = \frac{\pi}{2\sigma_n\sqrt{3}}, \ Z = \frac{\pi}{4\sigma_n\sqrt{3}}$$

Laplace's Approximation

The following derivatives are needed where $\phi(x) \equiv \ln(\cosh(x))$

$$\begin{array}{rcl} & \ln p(y|f) & = & \ln(\pi) - \ln(4\sigma_n\sqrt{3}) - 2\varphi\left(\tau(f-y)\right) \\ & \frac{\partial \ln p}{\partial f} & = & 2\tau\varphi'\left(\tau(f-y)\right) \\ & \frac{\partial^2 \ln p}{(\partial f)^2} & = & -2\tau^2\varphi''\left(\tau(f-y)\right) \\ & \frac{\partial^3 \ln p}{(\partial f)^3} & = & 2\tau^3\varphi'''\left(\tau(f-y)\right) \\ & \frac{\partial^3 \ln p}{(\partial \ln \sigma_n)(\partial f)^2} & = & 2\tau^2\left(2\varphi''\left(\tau(f-y)\right) + \tau(f-y)\varphi'''\left(\tau(f-y)\right)\right) \\ & \frac{\partial \ln p}{\partial \ln \sigma_n} & = & 2\tau(f-y)\varphi'\left(\tau(f-y)\right) - 1 \end{array}$$

4.12 GLM Likelihoods: Poisson, Negative Binomial, Weibull, Gamma, Exponential, Inverse Gaussian and Beta

Data y from a space other than \mathbb{R} e.g. \mathbb{N} , \mathbb{R}_+ or [0,1] can be modeled using generalised linear model likelihoods p(y|f) where the expected value $\mathbb{E}[y] = \mu$ is related to the underlying Gaussian process f by means of an inverse link function $\mu = g(f)$. Typically, the likelihoods are from an exponential family, hence the variance $\mathbb{V}[y] = \nu(\mu)$, is a simple function of the mean μ as well as higher order moments such as skewness $\mathbb{S}[y] = s(\mu)$ and kurtosis $\mathbb{K}[y] = k(\mu)$.

Here, we directly specify the inverse link function $\mu = g(f)$ defining the mapping from the GP f to the mean intensity μ . For numerical reasons, we work with the log of the inverse link function $h(f) = \ln g(f)$ and use its derivatives h', h'' and h''' for subsequent computations. In the table below, we have summarised the GLM likelihood expressions, the moments, the range of their variables and the applicable inverse link functions.

Likelihood	$\rho =$	$\nu(\mu) =$	$s(\mu) =$	k(μ) =	p(y f) =	y ∈	$\mu \in$	Inverse Links
Poisson	Ø	μ	$1/\sqrt{\mu}$	1/μ	$\mu^{y} \exp(-\mu)/y!$	N	\mathbb{R}_{+}	exp, logistic*
Neg. Binomial	{ln r}	$\mu(\mu/r+1)$		$\frac{6}{r} + \frac{r}{\mu(\mu + r)}$	$\tbinom{y+r-1}{y}r^r\mu^y/(r+\mu)^{r+y}$	N	\mathbb{R}_{+}	exp, logistic*
Weibull	{ln κ}	$\mu^2(\gamma_2/\gamma_1^2-1)$	$\frac{\gamma_3 - 3\gamma_1\gamma_2 + 2\gamma_1^3}{(\gamma_2 - \gamma_1^2)^{3/2}}$	$\frac{\gamma_4 - 4\gamma_1\gamma_3 + 12\gamma_1^2\gamma_2 - 3\gamma_2^2 - 6\gamma_1^4}{(\gamma_2 - \gamma_1^2)^2}$	$\kappa \gamma_1/\mu \left(y\gamma_1/\mu\right)^{\kappa-1} exp\left(-(y\gamma_1/\mu)^\kappa\right)$	$\mathbb{R}_{+}\backslash\{0\}$	$\mathbb{R}_+\backslash\{0\}$	exp, logistic*
Gamma	$\{ln \ \alpha\}$	μ^2/α	$2/\sqrt{\alpha}$	6/α	$\frac{\alpha^{\alpha}y^{\alpha-1}}{\Gamma(\alpha)}\mu^{-\alpha}\exp\left(-\frac{y\alpha}{\mu}\right)$	$\mathbb{R}_{+}\backslash\{0\}$	$\mathbb{R}_+\backslash\{0\}$	exp, logistic*
Exponential	Ø	μ^2	2	6	$\mu^{-1} \exp\left(-\frac{y}{\mu}\right)$	$\mathbb{R}_{+}\backslash\{0\}$	$\mathbb{R}_+\backslash\{0\}$	exp, logistic*
Inv. Gauss	$\{\ln\lambda\}$	μ^3/λ	$3\sqrt{\mu/\lambda}$	$15\mu/\lambda$	$\sqrt{\frac{\lambda}{2\pi y^3}} \exp\left(-\frac{\lambda(y-\mu)^2}{2\mu^2 y}\right)$	$\mathbb{R}_{+}\backslash\{0\}$	$\mathbb{R}_{+}\backslash\{0\}$	exp, logistic*
Beta	{ln φ}	$\mu(1-\mu)/(1+\varphi)$	$\frac{(2-4\mu)(1+\phi)}{\sqrt{\nu(\mu)}(2+\phi)}$	$6\frac{(\phi+1)^2-\nu(\mu)(5\phi+6)}{\nu(\mu)(\phi+2)(\phi+3)}$	$\frac{\Gamma(\varphi)}{\Gamma(\mu\varphi)\Gamma((1-\mu)\varphi)}y^{\mu\varphi-1}(1-y)^{(1-\mu)\varphi-1}$	[0, 1]	[0, 1]	expexp, logit

4.12.1 Inverse Link Functions

Possible inverse link functions and their properties (\cup convex, \cap concave, \uparrow monotone) are summarised below:

util/glm_invlink_*	$g(f) = \mu =$	$g:\mathbb{R} o$	g is	$h(f) = \ln \mu =$	h is
exp	e ^f	\mathbb{R}_{+}	∪,↑	f	∪,∩,↑
logistic	$\ell(f) = \ln(1 + e^f)$	\mathbb{R}_{+}	∪,↑	$ln(ln(1+e^f))$	∩,↑
logistic2	$\ell(f + \alpha f \ell(f))$	\mathbb{R}_{+}	↑	$ln(\ell(f + \alpha f \ell(f)))$	∩,↑
expexp	$\exp(-e^{-f})$	[0, 1]	1	$-e^{-f}$	∪,↑
logit	$1/(1+e^{-f})$	[0, 1]	↑	$-\ln(1+e^{-f})$	∪,↑

Please see doc/usageLik.m for how to specify the pair likelihood function and link function in GPML.

Exponential inverse link: exp

For $g(f) = e^f$ things are simple since h(f) = f, h'(f) = 1 and h''(f) = h'''(f) = 0.

Logistic inverse link: logistic

For $g(f) = \ln(1 + e^f)$ the derivatives of h(f) are given by

$$\begin{split} h(f) &= \ln(\ln(1+e^f)) \\ h'(f) &= \frac{1}{\ln(1+e^f)}s(-f), \, s(f) = \frac{1}{1+e^f}, \, s'(f) = \frac{-e^f}{(1+e^f)^2} = -s(-f)s(f) \\ h''(f) &= \frac{1}{\ln(1+e^f)}\frac{e^{-f}}{(1+e^{-f})^2} - \frac{1}{\ln^2(1+e^f)}\frac{e^f}{1+e^f}\frac{1}{1+e^{-f}} \\ &= h'(f)\left[s(f) - h'(f)\right] \\ h'''(f) &= h''(f)\left[s(f) - h'(f)\right] + h'(f)\left[\frac{-e^f}{(1+e^f)^2} - h''(f)\right] \\ &= h''(f)\left[s(f) - 2h'(f)\right] - h'(f)s(f)s(-f). \end{split}$$

Note that $g(f) = e^{h(f)} = \ln(1 + e^f)$ is convex and $h(f) = \ln(\ln(1 + e^f))$ with

$$h''(f) = \frac{1}{\ln(1+e^f)} \left(1 - \frac{e^f}{\ln(1+e^f)} \right) \frac{1}{1+e^f} \frac{1}{1+e^{-f}} \leqslant 0$$

is concave since $e^f \ge \ln(1 + e^f)$ for all $f \in \mathbb{R}$.

Twice logistic inverse link: logistic2

Note that $h(f) = \ln(\ell(f + \alpha f \ell(f)))$ is – according to Seeger et al.⁹ – concave.

Double negative exponential inverse link: expexp

For $g(f) = \exp(-e^{-f})$ the derivatives of h(f) are given by

$$h(f) = -e^{-f}$$
 $h'(f) = -h(f)$
 $h'''(f) = h(f)$
 $h''''(f) = -h(f)$

Logit regression inverse link: logit

For $g(f) = 1/(1+e^{-f})$ the derivatives of h(f) can be computed using the logistic inverse link function $h_{\ell}(f)$ since $h(f) = f - \exp(h_{\ell}(f))$

$$\begin{array}{lll} h(f) & = & f - e^{h_{\ell}(f)} \\ h'(f) & = & 1 - e^{h_{\ell}(f)} h'_{\ell}(f) \\ h''(f) & = & -e^{h_{\ell}(f)} [h'_{\ell}(f)^2 + h''_{\ell}(f)] = e^{h_{\ell}(f)} s_{\ell}(-f) s_{\ell}^2(f) \\ h'''(f) & = & -e^{h_{\ell}(f)} [h'_{\ell}(f)^3 + 3h''_{\ell}(f)h'_{\ell}(f) + h'''_{\ell}(f)] \end{array}$$

⁹Bayesian Intermittent Demand Forecasting, NIPS, 2016

4.12.2 Poisson Likelihood

Count data $y \in \mathbb{N}^n$ can be modeled in the GP framework using the Poisson distribution $p(y) = \mu^y e^{-\mu}/y!$ with mean/variance $\mathbb{E}[y] = \mathbb{V}[y] = \mu$, skewness $\mathbb{S}[y] = 1/\sqrt{\mu}$ and kurtosis $\mathbb{K}[y] = 1/\mu$ leading to the likelihood

$$\begin{array}{rcl} p(y|f) & = & \mu^y \exp(-\mu)/y!, \; \mu = g(f) \\ \Leftrightarrow \ln p(y|f) & = & y \cdot \ln g(f) - g(f) - \ln \Gamma(y+1). \end{array}$$

For Laplace's method to work, we need the first three derivatives of the log likelihood $\ln p(y|f)$, where $h(f) = \ln q(f)$

Note that if $\ln \mu = h(f)$ is concave and $\mu = g(f)$ is convex then the Poisson likelihood p(y|f) is log-concave in f which is the case for both exp and logistic.

4.12.3 Weibull Likelihood

Nonnegative data $y \in \mathbb{R}_+$ such as time-to-failure can be modeled in the GP framework using the Weibull distribution $p(y) = \kappa/\lambda(y/\lambda)^{\kappa-1}e^{-(y/\lambda)^{\kappa}}$ with shape parameter $\kappa > 0$, scale parameter $\lambda > 0$, mean $\mathbb{E}[y] = \lambda \gamma_1 = \mu$ where $\gamma_j = \Gamma(1+j/\kappa)$, variance $\mathbb{V}[y] = \lambda^2 \gamma_2 - \mu^2 = \mu^2(\gamma_2/\gamma_1^2 - 1)$, skewness $\mathbb{S}[y] = (\gamma_3 - 3\gamma_1\gamma_2 + 2\gamma_1^3)/(\gamma_2 - \gamma_1^2)^{3/2}$ and kurtosis $\mathbb{K}[y] = (\gamma_4 - 4\gamma_1\gamma_3 + 12\gamma_1^2\gamma_2 - 3\gamma_2^2 - 6\gamma_1^4)/(\gamma_2 - \gamma_1^2)^2$. Using the substitution $\mu = \lambda \gamma_1 \Leftrightarrow 1/\lambda = \gamma_1/\mu$, we obtain

$$\begin{split} p(y|f) &= \gamma_1 \frac{\kappa}{\mu} \left(\gamma_1 \frac{y}{\mu} \right)^{\kappa-1} exp \left(- \left(\gamma_1 \frac{y}{\mu} \right)^{\kappa} \right), \; \mu = g(f) > 0 \\ \Leftrightarrow \ln p(y|f) &= \ln \left(\gamma_1 \frac{\kappa}{\mu} \right) + (\kappa - 1) \ln \left(\gamma_1 \frac{y}{\mu} \right) - \left(\gamma_1 \frac{y}{\mu} \right)^{\kappa}. \end{split}$$

Note that the Weibull likelihood p(y|f) is log-concave in f neither for the exp nor for the logistic inverse link.

4.12.4 Gamma Likelihood

Nonnegative data $y \in \mathbb{R}_+$ can be modeled in the GP framework using the Gamma distribution $p(y) = \theta^{-\alpha}/\Gamma(\alpha)y^{\alpha-1}e^{-y/\theta}$ with shape parameter $\alpha > 0$, scale parameter $\theta > 0$, mean $\mathbb{E}[y] = \alpha\theta = \mu$, variance $\mathbb{V}[y] = \alpha\theta^2 = \mu^2/\alpha$, skewness $\mathbb{S}[y] = 2/\sqrt{\alpha}$ and kurtosis $\mathbb{K}[y] = 6/\alpha$. Using the substitution $\mu = \alpha\theta \Leftrightarrow \alpha/\mu = 1/\theta$, we obtain

$$\begin{split} p(y|f) &=& \frac{\alpha^{\alpha}y^{\alpha-1}}{\Gamma(\alpha)}\mu^{-\alpha}\exp\left(-\frac{y\alpha}{\mu}\right),\; \mu=g(f)>0\\ \Leftrightarrow& \ln p(y|f) &=& -\alpha\left(\ln \mu+\frac{y}{\mu}\right)-\ln Z_{\alpha}(y),\; \ln Z_{\alpha}(y)=\ln \Gamma(\alpha)-\alpha\ln \alpha+(1-\alpha)\ln y. \end{split}$$

Note that if $\ln \mu = h(f)$ was convex and $\mu = g(f)$ was concave then the Gamma likelihood p(y|f) would be log-concave in f which is not the case for both exp and logistic.

4.12.5 Exponential Likelihood

Nonnegative data $y \in \mathbb{R}_+$ can be modeled in the GP framework using the Exponential distribution $p(y) = \theta^{-1}e^{-y/\theta}$ with scale parameter $\theta > 0$, mean $\mathbb{E}[y] = \theta = \mu$, variance $\mathbb{V}[y] = \mu^2$, skewness $\mathbb{S}[y] = 2$ and kurtosis $\mathbb{K}[y] = 6$. We obtain

$$\begin{split} p(y|f) &= & \mu^{-1} \exp\left(-\frac{y}{\mu}\right), \; \mu = g(f) > 0 \\ \Leftrightarrow & \ln p(y|f) &= & -\ln \mu - \frac{y}{\mu}. \end{split}$$

Note that for exp (but not for logistic) the likelihood is log-concave. The exponential distribution corresponds to the Gamma distribution with $\alpha = 1$ and the Weibull distribution with $\kappa = 1$.

4.12.6 Inverse Gaussian Likelihood

Nonnegative data $y \in \mathbb{R}^n_+$ can be modeled in the GP framework using the Inverse Gaussian distribution $p(y) = \sqrt{\lambda/(2\pi y^3)} \exp(-\lambda(y-\mu)^2/(2\mu^2 y))$ with shape parameter $\lambda > 0$, mean parameter $\mu > 0$, mean $\mathbb{E}[y] = \mu$, variance $\mathbb{V}[y] = \mu^3/\lambda$, skewness $\mathbb{S}[y] = 3\sqrt{\mu/\lambda}$ and kurtosis $\mathbb{K}[y] = 15\mu/\lambda$. We obtain

$$\begin{split} p(y|f) &= \sqrt{\frac{\lambda}{2\pi y^3}} \exp\left(-\frac{\lambda (y-\mu)^2}{2\mu^2 y}\right), \; \mu = g(f) > 0 \\ \Leftrightarrow \ln p(y|f) &= -\frac{\lambda (y-\mu)^2}{2\mu^2 y} - \ln Z_\alpha(y), \; \ln Z_\alpha(y) = -\frac{1}{2} (\ln \lambda - \ln 2\pi y^3). \end{split}$$

The inverse Gaussian likelihood is in general not log-concace in f for both exp and logistic.

4.12.7 Beta Likelihood

Interval data $y \in [0,1]^n$ can be modeled in the GP framework using the Beta distribution $p(y) = y^{\alpha-1}(1-y)^{\beta-1}/B(\alpha,\beta)$ with shape parameters $\alpha,\beta>0$, mean $\mathbb{E}[y]=\alpha/(\alpha+\beta)$ and variance $\mathbb{V}[y]=\alpha\beta/[(\alpha+\beta)^2(\alpha+\beta+1)]$ and $1/B(\alpha,\beta)=\Gamma(\alpha+\beta)/[\Gamma(\alpha)\Gamma(\beta)]$. Reparametrising using the mean parameter $\mu=\mathbb{E}[y]=\alpha/(\alpha+\beta)$, the shape parameter $\varphi=\alpha+\beta$, the variance $\mathbb{V}[y]=\mu(1-\mu)/(1+\varphi)$ and hence

$$\begin{split} p(y|f) &=& \frac{\Gamma(\varphi)}{\Gamma(\mu\varphi)\Gamma((1-\mu)\varphi)} y^{\mu\varphi-1} (1-y)^{(1-\mu)\varphi-1}, \; \mu=g(f)>0 \\ \Leftrightarrow & \ln p(y|f) &=& \ln \Gamma(\varphi) - \ln \Gamma(\mu\varphi) - \ln \Gamma((1-\mu)\varphi) + (\mu\varphi-1) \ln y + ((1-\mu)\varphi-1) \ln (1-y). \end{split}$$

The Beta likelihood is in general not log-concace in f for both exp and logistic.

5 Mean Functions

A mean function $\mathfrak{m}_{\Phi}: \mathfrak{X} \to \mathbb{R}$ (with hyperparameters Φ) of a GP f is a scalar function defined over the whole domain \mathfrak{X} that computes the expected value $\mathfrak{m}(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ of f for the input \mathbf{x} .

5.1 Interface

In the GPML toolbox, a mean function $m: \mathcal{X} \to \mathbb{R}$ needs to implement evaluation $m = m_{\varphi}(X)$ and first derivatives $m_i = \frac{\partial}{\partial \varphi_i} m$ with respect to the components i of the parameter $\varphi \in \Phi$ as detailed below.

```
\langle meanFunctions.m \ 36 \rangle \equiv
36
      1 % Mean functions to be use by Gaussian process functions. There are two
      2 % different kinds of mean functions: simple and composite:
      3 %
      4 % Simple mean functions:
      5 %
      6 % meanZero - zero mean function
      7 % meanOne
                             - one mean function
      8 % meanConst
                            - constant mean function
     9 % meanLinear - linear mean function
10 % meanPoly - polynomial mean function
     11 % meanDiscrete - precomputed mean for discrete data
     12 % meanGP - predictive mean of another GP
     13 % meanGPexact - predictive mean of a regression GP
     14\ \% meanNN - nearest neighbor mean function 15\ \% meanWSPC - weighted sum of projected cosines
     16 %
     17 % Composite mean functions (see explanation at the bottom):
     18 %
     19 % meanScale - scaled version of a mean function
20 % meanSum - sum of mean functions
21 % meanProd - product of mean functions
22 % meanPow - power of a mean function
     23 % meanMask - mask some dimensions of the data
24 % meanPref - difference mean for preference learning
25 % meanWarp - warped mean function
     26 %
     27 % Naming convention: all mean functions are named "mean/mean*.m".
     28 %
     29 %
     30 % 1) With no or only a single input argument:
     31 %
     32 %
              s = meanNAME or s = meanNAME(hyp)
     33 %
     34 % The mean function returns a string s telling how many hyperparameters hyp it
     35 % expects, using the convention that "D" is the dimension of the input space.
     36 % For example, calling "meanLinear" returns the string 'D'.
     37 %
     38 % 2) With two input arguments and one output argument:
     39 %
     40 %
           m = meanNAME(hyp, x)
     41 %
     42 % The function computes and returns the mean vector m with components
     43 % m(i) = m(x(i,:)) where hyp are the hyperparameters and x is an n by D matrix
```

44 % of data, where D is the dimension of the input space. The returned mean

```
45 % vector m is of size n by 1.
46 %
47 % 3) With two input arguments and two output arguments:
48 %
49 % [m,dm] = meanNAME(hyp, x)
50 %
51 % The function computes and returns the mean vector m as in 2) above.
52 % In addition to that, the (linear) directional derivative function dm is
53 % returned. The call dhyp = dm(q) for a direction vector q of size n by 1
54 % returns a vector of directional derivatives dhyp = d (q'*m(x)) / d hyp of
55 % the same size as the hyperparameter vector hyp. The components of dhyp are
56 % defined as follows: dhyp(i) = q'*(d m(x) / d hyp(i)).
57 %
58 % See also doc/usageMean.m.
59 %
60 \( \lambda \text{gpml copyright 5a} \rangle \)
```

5.2 Implemented Mean Functions

We offer simple and composite mean functions producing new mean functions $m(\mathbf{x})$ from existing mean functions $\mu_j(\mathbf{x})$. All code files are named according to the pattern mean/mean<NAME>.m for simple identification. This modular specification allows to define affine mean functions $m(\mathbf{x}) = c + \mathbf{a}^{\top}\mathbf{x}$ or polynomial mean functions $m(\mathbf{x}) = (c + \mathbf{a}^{\top}\mathbf{x})^2$. All currently available mean functions are summarised in the following table.

Simple mean functions $m(x)$				
<name></name>	Mmeaning	$m(\mathbf{x}) =$	ф	
Zero	mean vanishes always	0	Ø	
One	mean equals 1	1	Ø	
Const	mean equals a constant	С	$c\in\mathbb{R}$	
Linear	mean linearly depends on $\mathbf{x} \in \mathcal{X} \subseteq \mathbb{R}^D$	$\mathbf{a}^{T}\mathbf{x}$	$\mathbf{a} \in \mathbb{R}^{\mathrm{D}}$	
Poly	mean polynomially depends on $\mathbf{x} \in \mathcal{X} \subseteq \mathbb{R}^D$	$\sum_{\mathbf{d}} \mathbf{a}_{\mathbf{d}}^{T} \mathbf{x}^{\mathbf{d}}$	$\mathbf{a} \in \mathbb{R}^{D \times d}$	
Discrete	precomputed mean for discrete data $\mathbf{x} \in \mathcal{X} \subseteq \mathbb{N}$	$\mu_{\mathbf{x}}$	$\mu \in \mathbb{R}^s$	
GP	predictive mean of another GP	$\int \mathbf{y} \cdot \mathbf{p}(\mathbf{y} \mathcal{D}, \mathbf{x}) d\mathbf{y}$	Ø	
GPexact	predictive mean of a regression GP	$\int \mathbf{y} \cdot \mathbf{p}(\mathbf{y} \mathcal{D}, \mathbf{x}) d\mathbf{y}$	ρ, ψ, σ_n	
NN	nearest neighbor for a set $(\mathbf{z}_j, \mathbf{m}_j) \in \mathcal{X} \times \mathbb{R}$	m_i , $i = arg min_j d(x, z_j)$	Ø	
WSPC	weighted sum of d projected cosines $\mathbf{x} \in \mathcal{X} \subseteq \mathbb{R}^D$	$\mathbf{a}^{\top}\cos(\mathbf{W}\mathbf{x}+\mathbf{b})$	$\mathbf{W} \in \mathbb{R}^{d \times D}, \mathbf{a}, \mathbf{b} \in \mathbb{R}^d$	
Composite mean functions $[\mu_1(\mathbf{x}), \mu_2(\mathbf{x}),] \mapsto \mathfrak{m}(\mathbf{x})$				
<name></name>	meaning	$m(\mathbf{x}) =$	ф	
Scale	scale a mean	$\alpha\mu(\mathbf{x})$	$lpha \in \mathbb{R}$	
Sum	add up mean functions	$\sum_{j} \mu_{j}(\mathbf{x})$	Ø	
Prod	multiply mean functions	$\prod_{j} \mu_{j}(\mathbf{x})$	Ø	
Pow	raise a mean to a power	$\mu(\mathbf{x})^{d}$	Ø	
Mask	act on components $I \subseteq [1, 2,, D]$ of $\mathbf{x} \in \mathcal{X} \subseteq \mathbb{R}^D$ only	$\mu(\mathbf{x}_{\mathrm{I}})$	Ø	
Pref	preference learning mean $\mathbf{x} = [\mathbf{x}_1; \mathbf{x}_2], \ \mathbf{x}_i \subseteq \mathbb{R}^{D/2}$	$\mu(\mathbf{x}_1) - \mu(\mathbf{x}_2)$	Ø	
Warp	warped mean	$g[\mu(\mathbf{x})]$	Ø	

5.3 Usage of Implemented Mean Functions

Some code examples taken from doc/usageMean.m illustrate how to use simple and composite mean functions to specify a GP model.

Syntactically, a mean function mf is defined by

```
mn := 'func' | @func // simple
```

 $mf := \{mn\} \mid \{mn, \{param, mf\}\} \mid \{mn, \{mf, ..., mf\}\} // composite$ i.e., it is either a string containing the name of a mean function, a pointer to a mean function or one of the former in combination with a cell array of mean functions and an additional list of parameters.

```
\langle doc/usageMean.m \ 38 \rangle \equiv
38
      1 % demonstrate usage of mean functions
      3 % See also meanFunctions.m.
      4 %
      5 \(\langle gpml \copyright 5a\rangle
      6 clear all, close all
      7 \text{ n} = 5; D = 2; x = \text{randn}(n,D); % create a random data set
      9 % set up simple mean functions
     12 mc = {@meanConst}; hypc = 2; % also function handles are possible
     13 ml = {@meanLinear}; hypl = [2;3];
                                                         % m(x) = 2*x1 + 3*x2
     14 mp = {@meanPoly,2}; hypp = [1;1;2;3]; % m(x) = x1+x2+2*x1^2+3*x2^2
     15 mn = {@meanNN,[1,0; 0,1],[0.9,0.5]}; hypn = []; % nearest neighbor
     16 s = 12; hypd = randn(s,1); % discrete mean with 12 hypers
     17 md = {'meanDiscrete',s};
                                                          % GP predictive mean
     18 \text{ hyp.cov} = [0;0]; \text{ hypg} = [];
     19 xt = randn(2*n,D); yt = sign(xt(:,1)-xt(:,2)); % training data
     20 mg = {@meanGP,hyp,@infEP,@meanZero,@covSEiso,@likErf,xt,yt};
     21 hype = [0;0; log(0.1)]; % regression GP predictive mean
     22 xt = randn(2*n,D); yt = xt(:,1).*xt(:,2);
                                                              % training data
     23 me = {@meanGPexact,@meanZero,@covSEiso,xt,yt};
     24
     25 % set up composite mean functions
     26 msc = {'meanScale',{m1}}; hypsc = [3; hyp1]; % scale by 3
     27 msu = {'meanSum', {m0,mc,ml}}; hypsu = [hyp0; hypc; hyp1]; % sum
     28 mpr = {@meanProd,{mc,ml}}; hyppr = [hypc; hypl]; % product
29 mpo = {'meanPow',3,msu}; hyppo = hypsu; % third power
     30 mask = [false,true]; % mask excluding all but the 2nd component
     31 mma = {'meanMask', mask, ml}; hypma = hypl(mask);
32 mpf = {@meanPref, ml}; hyppf = 2; % linear pref with slope
     33 mwp = {@meanWarp,ml,@sin,@cos};hypwp = 2; % sin of linear
     35~\% 0) specify mean function
     36 \% \text{ mean} = \text{md}; \text{ hyp} = \text{hypd}; \text{ x} = \text{randi}([1,s],n,1);
     37 \% \text{ mean} = mn; \text{ hyp} = \text{hypn};
     38 % mean = mg; hyp = hypg;
     39 mean = me; hyp = hype;
     40 \% \text{ mean} = m0; \text{ hyp} = \text{hyp0};
     41 % mean = msu; hyp = hypsu;
     42 % mean = mpr; hyp = hyppr;
     43 % mean = mpo; hyp = hyppo;
     44 % mean = mpf; hyp = hyppf;
     46 % 1) query the number of parameters
     47 feval(mean{:})
     49 \% 2) evaluate the function on x
     50 feval(mean{:},hyp,x)
     51
     52 % 3) compute the derivatives w.r.t. to hyperparameter i
     53 i = 2; feval(mean{:},hyp,x,i)
```

6 Covariance Functions

A covariance function $k_{\psi}: \mathfrak{X} \times \mathfrak{X} \to \mathbb{R}$ (with hyperparameters ψ) of a GP f is a scalar function defined over the whole domain \mathcal{X}^2 that computes the covariance $k(\mathbf{x}, \mathbf{z}) = \mathbb{V}[f(\mathbf{x}), f(\mathbf{z})] = \mathbb{E}[(f(\mathbf{x}) - \mathbf{z})]$ m(x)(f(z) - m(z)) of f between the inputs x and z.

6.1 Interface

43 %

Again, the interface is simple since only evaluation of the full covariance matrix $K = k_{\psi}(X)$ and its derivatives $K_i = \frac{\partial}{\partial \psi} K$ as well as cross terms $k_* = k_{\psi}(X, x_*)$ and $k_{**} = k_{\psi}(x_*, x_*)$ for prediction are required.

```
\langle covFunctions.m \ 39 \rangle \equiv
39
           1 % Covariance functions to be use by Gaussian process functions. There are two
           2 % different kinds of covariance functions: simple and composite:
           3 %
           4 % 1) Elementary and standalone covariance functions:
           5 % covZero - zero covariance function
           6 % covEye
                                               - unit covariance function
           7 % covOne - unit constant covariance function
           8 % covDiscrete - precomputed covariance for discrete data
           9 %
         10 % 2) Composite covariance functions:
         11 % covScale - scaled version of a covariance function
         12 % covSum
                                               - sums of covariance functions
        12 % covSum - sums of covariance functions
13 % covProd - products of covariance functions
14 % covMask - mask some dimensions of the data
15 % covPref - difference covariance for preference learning
16 % covPER - make stationary covariance periodic
17 % covADD - additive covariance function
18 % covWarp - warp input to a covariance function
         19 %
         20 % 3) Mahalanobis distance based covariances and their modes
         21 % covMaha - generic "mother" covariance
        21 % covMaha - generic "mother" covariance
22 % * eye - unit length scale
23 % * iso - isotropic length scale
24 % * ard - automatic relevance determination
25 % * prot - (low-rank) projection in input space
26 % * fact - factor analysis covariance
27 % * vlen - spatially varying length scale
28 % covGE - Gamma exponential covariance
29 % covMatern - Matern covariance function with nu=1/2, 3/2 or 5/2
30 % covPP - piecewise polynomial covariance function (compact support)
31 % covRQ - rational quadratic covariance function
32 % covSE - squared exponential covariance function
33 %
         33 %
         34 % 4) Dot product based covariances and their modes
         35~\% covDot - generic "mother" covariance
        30 % * eye - unit length scale
37 % * iso - isotropic length scale
38 % * ard - automatic relevance determination
39 % * pro - (low-rank) projection in input space
40 % * fac - factor analysis covariance
41 % covLIN - linear covariance function
42 % covPoly - polynomial covariance function
43 %
```

```
44 % 5) Time series covariance functions on the positive real line
 45 % covFBM - fractional Brownian motion covariance
46 % covULL - underdamped linear Langevin process covariance
 47 % covW
                                - i-times integrated Wiener process covariance
 48 % covOU
                                - i-times integrated Ornstein-Uhlenbeck process covariance
 49 %
 50 % 6) Standalone covariances
 51 % covNNone - neural network covariance function
 52 % covLINone - linear covariance function with bias
 53 % covPeriodic - smooth periodic covariance function (1d)
 54 % covPeriodicNoDC - as above but with zero DC component and properly scaled
 55 \% covCos - sine periodic covariance function (1d) with unit period
 56 %
         covGabor - Gabor covariance function
 57 %
 58 % 7) Shortcut covariances assembled from library
 59~\% covConst - covariance for constant functions - independent covariance function (i.e. white noise)
 61 % covPERiso - make isotropic stationary covariance periodic
62 % covPERard - make ARD stationary covariance periodic
 63 % covMaterniso - Matern covariance function with nu=1/2, 3/2 or 5/2
 64 \% covMaternard - Matern covariance function with nu=1/2, 3/2 or 5/2 with ARD
 65 % covPPiso - piecewise polynomial covariance function (compact support)
- piecewise polynomial covariance function
- piecewise polynomial covariance function
- piecewise polynomial covariance function
- projection squared exponential covariance function
- piecewise polynomial covariance function
- projection squared exponential covariance function
 76 % covGaborard - Gabor covariance function with ARD
 77 % covGaborsio - isotropic Gabor covariance function
 78 % covSM
                              - spectral mixture covariance function
 79 %
 80 % 8) Special purpose (approximation) covariance functions
 81 % apxSparse - sparse approximation: to be used for large scale inference
 82 %
                                     problems with inducing points aka FITC
 83 %
         apxGrid
                                 - grid interpolation: to be used for large scale inference
 84 %
                                     problems with Kronecker/Toeplitz/BTTB covariance matrix
 85 %
 86 % The covariance functions are written according to a special convention where
 87 % the exact behaviour depends on the number of input and output arguments
 88 % passed to the function. If you want to add new covariance functions, you
 89 % should follow this convention if you want them to work with the function gp.
 90 % There are four different ways of calling the covariance functions:
 92 % 1) With no (or one) input argument(s):
 93 %
 94 %
             s = cov
 95 %
 96 % The covariance function returns a string s telling how many hyperparameters it
 97 % expects, using the convention that "D" is the dimension of the input space.
 98 % For example, calling "covRQard" returns the string '(D+2)'.
100 % 2) With two input arguments and one output argument:
101 %
```

```
102 %
        K = cov(hyp, x) equivalent to K = cov(hyp, x, [])
103 %
104 % The function computes and returns the covariance matrix where hyp are
105 % the hyperparameters and x is an n by D matrix of cases, where
106 % D is the dimension of the input space. The returned covariance matrix is of
107 % size n by n.
108 %
109 % 3) With three input arguments and one output argument:
110 %
111 %
      Kz = cov(hyp, x, z)
112 %
        kx = cov(hyp, x, 'diag')
113 %
114 % The function computes test set covariances; kx is a vector of self covariances
115 % for the test cases in x (of length n) and Kz is an (n by nz) matrix of cross
116 % covariances between training cases x and test cases z.
117 %
118 % 4) With two output arguments:
119 %
120 %
         [K,dK] = cov(hyp, x) equivalent to [K,dK] = cov(hyp, x, [])
121 %
         [K,dK] = cov(hyp, x, z)
122 %
         [K,dK] = cov(hyp, x, 'diag')
123 %
124 % The function computes and returns the covariances K as in 3) above.
125 % In addition to that, the (linear) directional derivative function dK is
126 % returned. The two possible calls dhyp = dK(Q) and [dhyp,dx] = dK(Q) for a
127 % direction Q of the same size as K are possible. The return arguments dhyp
128 % and dx are the directional derivatives dhyp = d trace(Q'*K) / d hyp and
129 \% dx = d trace(Q'*K) / d x are of the same size as the hyperparameter
130 % vector hyp and the input data x, respectively. The components of dhyp and
131 % dx are defined as follows: dhyp(i) = trace(Q**(dK/dhyp(i)))
132 % and dx(i,j) = trace(Q'*(dK/dx(i,j))).
133 %
134 % Covariance functions can be specified in two ways: either as a string
135 % containing the name of the covariance function or using a cell array. For
136 % example:
137 %
138 %
      cov = 'covRQard';
139 % cov = {'covRQard'};
140 %
      cov = {@covRQard};
141 %
142 % are supported. Only the second and third form using the cell array can be used
143 % for specifying composite covariance functions, made up of several
144 % contributions. For example:
145 %
146 %
           cov = {'covScale', {'covRQiso'}};
147 %
            cov = {'covSum', {'covRQiso','covSEard','covNoise'}};
148 %
            cov = {'covProd',{'covRQiso','covSEard','covNoise'}};
149 %
            cov = {'covMask',{mask,'covSEiso'}}
150 % q=1; cov = {'covPPiso',q};
151 %
      d=3; cov = {'covPoly',d};
            cov = {'covADD',{[1,2],'covSEiso'}};
152 %
153 %
            cov = {@apxSparse, {@covSEiso}, u}; where u are the inducing inputs
154 %
155 % specifies a covariance function which is the sum of three contributions. To
156 % find out how many hyperparameters this covariance function requires, we do:
157 %
      feval(cov{:})
158 %
159 %
```

```
160 % which returns the string '3+(D+1)+1' (i.e. the 'covRQiso' contribution uses 161 % 3 parameters, the 'covSEard' uses D+1 and 'covNoise' a single parameter). 162 % 163 % See also doc/usageCov.m. 164 % 165 \langle gpml\ copyright\ 5a \rangle
```

6.2 Implemented Covariance Functions

Similarly to the mean functions, we provide a whole algebra of covariance functions $k: \mathcal{X} \times \mathcal{X} \to \mathbb{R}$ with the same generic name pattern cov/cov<NAME>.m as before.

Besides a long list of simple covariance functions, we also offer a variety of composite covariance functions as shown in the following table.

Const		σf^2	ln σf
1) Elementary a	and standalone covariance functions k(x,x')		
<name></name>	meaning	$k(\mathbf{x}, \mathbf{z}) =$	ψ
Zero	covariance vanishes always	0	Ø
Eye	unit additive measurement noise	$\delta(\mathbf{x} - \mathbf{z})$	Ø
One	unit constant	1	Ø
Discrete	precomputed covariance for discrete data $\mathbf{x} \in \mathfrak{X} \subseteq \mathbb{N}$	$k_{xx'}$ where $K = L^{\top}L$ is Cholesky decomposition, $L \in \mathbb{R}^{s \times s}$	L
2) Composite c	ovariance functions $[\kappa_1(\mathbf{x}, \mathbf{z}), \kappa_2(\mathbf{x}, \mathbf{z}),] \mapsto k(\mathbf{x}, \mathbf{z})$		
<name></name>	meaning	$k(\mathbf{x}, \mathbf{z}) =$	ψ
Scale	modulate covariance function by a scalar	$\sigma_f^2 \kappa(\mathbf{x}, \mathbf{z})$	ln σ _f
Scale	modulate covariance function depending on input	$\sigma_f(\mathbf{x})\kappa(\mathbf{x},\mathbf{z})\sigma_f(\mathbf{z})$	Фℓ
Sum	add up covariance functions	$\sum_{i} \kappa_{i}(\mathbf{x}, \mathbf{z})$	Ø
Prod	multiply covariance functions	$\prod_{i} \kappa_{i}(\mathbf{x}, \mathbf{z})$	Ø
Mask	act on components $I \subseteq [1, 2,, D]$ of $\mathbf{x} \in \mathcal{X} \subseteq \mathbb{R}^D$ only	$\kappa(\mathbf{x}_{\mathrm{I}}, \mathbf{z}_{\mathrm{I}})$	Ø
Pref	preference learning covariance $\mathbf{x} = [\mathbf{x}_1; \mathbf{x}_2], \ \mathbf{x}_i \subseteq \mathbb{R}^{D/2}$	$\kappa(\mathbf{x}_1, \mathbf{z}_1) + \kappa(\mathbf{x}_2, \mathbf{z}_2) - \kappa(\mathbf{x}_1, \mathbf{z}_2) - \kappa(\mathbf{x}_2, \mathbf{z}_1)$	Ø
PER	turn covariance into a periodic, $\mathfrak{X} \subseteq \mathbb{R}^D$	$\kappa(u(x),u(z)) \text{ , } u(x) = \left[\begin{array}{c} \sin 2\pi x_p \\ \cos 2\pi x_p \end{array}\right], x_p = \left\{\begin{array}{c} x & \text{eye} \\ x/p & \text{iso} \\ x/p & \text{ard} \end{array}\right.$	ψ_{u}
ADD	additive, $\mathfrak{X} \subseteq \mathbb{R}^{D}$, index degree set $\mathfrak{D} = \{1,, D\}$	$\begin{array}{c} \sum_{\mathbf{d} \in \mathcal{D}} \sigma_{f_{\mathbf{d}}}^{2} \sum_{ I = \mathbf{d}} \prod_{i \in I} \kappa(x_{i}, z_{i}; \psi_{i}) \\ \kappa(p(\mathbf{x}), p(\mathbf{z})), \text{ where } p : \mathbb{R}^{D} \rightarrow \mathbb{R}^{D_{p}} \end{array}$	$\{\psi_1,, \psi_D, \ln \sigma_{f_1},, \ln \sigma_{f_{ D }}\}$
Warp	warp inputs to a covariance function	$\kappa(p(x), p(z)), \text{ where } p : \mathbb{R}^D \to \mathbb{R}^{D_p}$	Ø
3) Covariance f	unctions based on Mahalanobis distances $k(\mathbf{x}, \mathbf{z}) = k(\mathbf{r}), \ \mathbf{r}^2 = (\mathbf{x} - \mathbf{z})^T$	$\mathbf{P}^{-1}(\mathbf{x} - \mathbf{z})$, where $\mathbf{x}, \mathbf{z} \in \mathfrak{X} \subseteq \mathbb{R}^{D}$	
Maha	shared generic "mother" covariance function	k(r) =	0
GE	gamma exponential gamma越接近于2,光顺性越好	$\exp(- \mathbf{r} ^{\gamma}), \gamma \in (0,2]$	$\ln(\gamma/(2-\gamma))$
Matern	Matérn, $f_1(t) = 1$, $f_3(t) = 1 + t$, $f_5(t) = f_3(t) + \frac{t^2}{3}$	$f_d(\sqrt{dr}) \exp(-\sqrt{dr})$	0
PP	compact support, piecewise polynomial $f_{\nu}(r)$	$\max(0, 1-r)^{j+\nu} \cdot f_{\nu}(r)$	Ø
RQ	rational quadratic	$\left(1+\frac{1}{2}r^2/\alpha\right)^{-\alpha}$	ln α
SE	squared exponential	$\exp(-r^2/2)$	Ø
Allowable mode	parameter values for Mahalanibis distance covariances		
'eye'	unit lengthscale	P = I	Ø
'iso'	isotropic lengthscale	$\mathbf{P} = \ell^2 \mathbf{I}, \ \ell \in \mathbb{R}_+$	In ℓ
'ard'	automatic relevance determination	$P = \Lambda^2$, $\Lambda = \text{diag}(\lambda)$, $\lambda \in \mathbb{R}^D_+$	lnΛ
'proj'	(low-rank) projection in input space	$P^{-1} = L^{\top}L, L \in \mathbb{R}^{d \times D}$	L
'fact'	factor analysis	$\mathbf{P}^{-1} = \mathbf{L}^{T} \mathbf{L} + \mathtt{diag}(\mathbf{f}), \mathbf{f} \in \mathbb{R}^{D}, \mathbf{L} \in \mathbb{R}^{d imes D}$	{L, ln f}
'vlen'	spatially varying lengthscale,	$k_{\ell}(\mathbf{x}, \mathbf{z}) = \left(\frac{\ell(\mathbf{x})\ell(\mathbf{z})}{l^2(\mathbf{x}, \mathbf{z})}\right)^{D/2} k\left(\frac{r^2}{l^2(\mathbf{x}, \mathbf{z})}\right), \ l^2(\mathbf{x}, \mathbf{z}) = \frac{\ell^2(\mathbf{x}) + \ell^2(\mathbf{z})}{2}$	Фε
4) Covariance f	unctions based on Euclidean dot products $k(\mathbf{x}, \mathbf{z}) = k(s)$, $s = \mathbf{x}^{\top} \mathbf{P}^{-1} \mathbf{z}$,	where $\mathbf{x}, \mathbf{z} \in \mathfrak{X} \subseteq \mathbb{R}^D$	
Dot	shared generic "mother" covariance function	k(s) =	Ø
LIN	linear covariance function	s	Ø
Poly	polynomial covariance	$\sigma_f^2(s+c)^d$	ln c
Allowable mode	parameter values for Euclidean dot product covariances		
'eye'	unit lengthscale	P = I	Ø
'iso'	isotropic lengthscale	$\mathbf{P} = \ell^2 \mathbf{I}, \ \ell \in \mathbb{R}_+$	ln ℓ
'ard'	automatic relevance determination	$\begin{aligned} \mathbf{P} &= \mathbf{\Lambda}^2, \ \ \boldsymbol{\Lambda} = \operatorname{diag}(\boldsymbol{\lambda}), \ \ \boldsymbol{\lambda} \in \mathbb{R}_+^D \\ \mathbf{P}^{-1} &= \boldsymbol{L}^\top \boldsymbol{L}, \boldsymbol{L} \in \mathbb{R}^{d \times D} \end{aligned}$	ln Λ
'proj'	(low-rank) projection in input space	$\mathbf{P}^{-1} = \mathbf{L}^{T} \mathbf{L}, \mathbf{L} \in \mathbb{R}^{\mathbf{d} \times \mathbf{D}}$	L
'fact'	factor analysis	$\mathbf{P}^{-1} = \mathbf{L}^{ op} \mathbf{L} + \mathtt{diag}(\mathbf{f}), \mathbf{f} \in \mathbb{R}^{\mathrm{D}}, \mathbf{L} \in \mathbb{R}^{\mathbf{d} imes \mathrm{D}}$	{L, ln f}
5) Time series c	ovariance functions $k(\mathbf{x}, \mathbf{z}) = k(\mathbf{x}, \mathbf{z})$, where $\mathbf{x}, \mathbf{z} \in \mathcal{X} = \mathbb{R}_+$		
<name></name>	meaning	k(x, z) =	ψ
FBM	fractional Brownian motion covariance with Hurst index h	$\frac{1}{2}\sigma_{\rm f}^2\left[{\bf x} ^{2{\bf h}}- {\bf x}-{\bf z} ^{2{\bf h}}+ {\bf z} ^{2{\bf h}}\right]$	$\{\ln \sigma_f, -\ln(1/h-1)\}$
ULL	underdamped linear Langevin process covariance	$k(t) = a^2 e^{-\mu t} \left(\frac{\sin(\omega t)}{\omega} + \frac{\cos(\omega t)}{\mu} \right), t = x - z $	$\{\ln \mu, \ln \omega, \ln \alpha\}$
W	i-times integrated Wiener process covariance, $i \in \{-1,, 3\}$	$\kappa_0(\mathbf{x}, \mathbf{z}) = \sigma_{\mathbf{f}}^2 \min(\mathbf{x}, \mathbf{z}), \ \kappa_{\mathbf{i}}(\mathbf{x}, \mathbf{z}) = \int_0^{\mathbf{x}} \int_0^{\mathbf{z}} \kappa_{\mathbf{i}}(\tilde{\mathbf{x}}, \tilde{\mathbf{z}}) d\tilde{\mathbf{x}} d\tilde{\mathbf{z}}$	ln σ _f
OU	i-times integrated Ornstein-Uhlenbeck process covariance, $i \in \{0, 1\}$	$\kappa_0(x,z) = \sigma_f^2 \exp(-\frac{ x-z }{\theta}) + (\sigma_{f_0}^2 - \sigma_f^2) \exp(-\frac{ x+z }{\theta})$	$\{\ln \ell, \ln \sigma_f, \ln \sigma_{fo}\}$
	ovariance functions	1 occurs -1 f 1 + (-10 ot) orb(f)	, 1, • 101
<name></name>	meaning meaning	$k(\mathbf{x}, \mathbf{z}) =$	ψ
	neural net, $X \subseteq \mathbb{R}^D$		
NNone	, –	$\sigma_{\rm f}^2 \sin^{-1} \left(\mathbf{x}^\top \mathbf{z} / \sqrt{(\ell^2 + \mathbf{x}^\top \mathbf{x})(\ell^2 + \mathbf{z}^\top \mathbf{z})} \right)$	$\{\ln \ell, \ln \sigma_f\}$
LINone	linear with bias, $\mathfrak{X} \subseteq \mathbb{R}^{D}$	$(\mathbf{x}^{T}\mathbf{z}+1)/t^2$	ln t
Periodic	periodic, $\mathfrak{X} \subseteq \mathbb{R}$	$ \sigma_f^2 \exp\left(-\frac{2}{\ell^2} \sin^2[\pi(x-z)/p]\right) $ $ 2\kappa(x-z) - \frac{1}{2} \int_0^{\pi} \kappa(t) dt $	$\{\ln \ell, \ln p, \ln \sigma_f\}$
PeriodicNoDC	periodic, rescaled, DC component removed, $\mathfrak{X} \subseteq \mathbb{R}$	$\sigma_f^2 \frac{\kappa(x-z) - \frac{1}{\pi} \int_0^\pi \kappa(t) dt}{\kappa(0) - \frac{1}{\pi} \int_0^\pi \kappa(t) dt}, \kappa(x-z) = \exp\left(-\frac{2}{\ell^2} \sin^2[\pi(x-z)/p]\right)$	$\{\ln \ell, \ln p, \ln \sigma_f\}$
Cos	periodic cosine, $\mathfrak{X} \subseteq \mathbb{R}$	$\sigma_f^2 \cos(\pi(x-z)/p)$	$\{\ln p, \ln \sigma_f\}$
Gabor	Gabor function, $\mathfrak{X} \subseteq \mathbb{R}^D$, $\boldsymbol{\lambda}, p \in \mathbb{R}^D_+$	$\exp\left(-\frac{1}{2}t^{\top}\Lambda^{-2}t\right)\cos\left(2\pi t_{p}^{\top}1\right), t_{p} = \begin{cases} t & \text{eye} \\ t/p & \text{iso}, t = x - z \end{cases}$	$\{\ln\lambda,\psi_p\}$

$$\begin{split} & gamma=1.5;cge=\{'covGE','eye',[]\};\\ & hypge=log(gamma/(2-gamma)); cov=cge;hyp=hypge;\\ & \raiset{ff} T\\ & cgu=\{'covSEisoU'\}; \ \% \ k(x,z)=exp(-(x-z))*inv(P)*(x-z)/2); P=ell^2*I;\\ & ell=1/sqrt(2);hypgu=log(ell);cov=cgu;hyp=hypgu; \end{split}$$

1D

7) Shortcut	7) Shortcut covariance functions assembled from library $k(\mathbf{x}, \mathbf{z})$				
<name></name>	Meaning	$k(\mathbf{x}, \mathbf{z}) =$	ψ		
Const	covariance equals a constant	$\sigma_{\rm f}^2$	ln σ _f		
Noise	additive measurement noise	$\sigma_f^2 \delta(\mathbf{x} - \mathbf{z})$	ln σ _f		
PERiso	turn covariance into a periodic, $\mathfrak{X} \subseteq \mathbb{R}^D$	$\kappa(\mathbf{u}(\mathbf{x}), \mathbf{u}(\mathbf{z}))$, $\mathbf{u}(\mathbf{x}) = [\sin \mathbf{x}_p, \cos \mathbf{x}_p]$, $\mathbf{x}_p = 2\pi \mathbf{x}/p$	ln p		
PERard	turn covariance into a periodic, $\mathfrak{X} \subseteq \mathbb{R}^D$	$\kappa(\mathbf{u}(\mathbf{x}), \mathbf{u}(\mathbf{z}))$, $\mathbf{u}(\mathbf{x}) = [\sin \mathbf{x}_p, \cos \mathbf{x}_p]$, $\mathbf{x}_p = 2\pi \mathbf{x}/p$	$\{\ln p_1,, \ln p_D\}$		
Materniso	Matérn, $\mathfrak{X}\subseteq\mathbb{R}^D,$ $f_1(t)=1,$ $f_3(t)=1+t,$ $f_5(t)=f_3(t)+\frac{t^2}{3}$	$\sigma_f^2 f_d(r_d) \exp(-r_d), \ r_d = \sqrt{\tfrac{d}{\ell^2} (\mathbf{x} - \mathbf{z})^\top (\mathbf{x} - \mathbf{z})}$	$\{\ln\ell,\ln\sigma_f\}$		
Maternard	Matérn, $\mathfrak{X} \subseteq \mathbb{R}^{D}$, $f_{1}(t) = 1$, $f_{3}(t) = 1 + t$, $f_{5}(t) = f_{3}(t) + \frac{t^{2}}{3}$	$\sigma_f^2 f_d(r_d) \exp(-r_d), r_d = \sqrt{d(\mathbf{x} - \mathbf{z})^{\top} \mathbf{\Lambda}^{-2} (\mathbf{x} - \mathbf{z})}$	$\{\ln \lambda_1,,\ln \lambda_D,\ln \sigma_f\}$		
PPiso	compact support, piecewise polynomial $f_{\nu}(r)$	$\sigma_f^2 \max(0, 1-r)^{j+\nu} \cdot f_{\nu}(r), r = \frac{\ \mathbf{x}-\mathbf{z}\ }{\ell}, j = \left \frac{\mathbf{D}}{2}\right + \nu + 1$	$\{\ln \ell, \ln \sigma_f\}$		
PPard	compact support, piecewise polynomial $f_{\nu}(r)$	$\sigma_f^2 \max(0, 1-r)^{j+\nu} \cdot f_{\nu}(r), r^2 = (\mathbf{x} - \mathbf{z})^{\top} \mathbf{\Lambda}^{-2} (\mathbf{x} - \mathbf{z})$	$\{\ln \lambda_1,, \ln \lambda_D, \ln \sigma_f\}$		
RQiso	rational quadratic, $\mathfrak{X} \subseteq \mathbb{R}^D$	$\sigma_f^2 \left(1 + \frac{1}{2\alpha\ell^2} (\mathbf{x} - \mathbf{z})^\top (\mathbf{x} - \mathbf{z})\right)^{-\alpha}$	$\{\ln \ell, \ln \sigma_f, \ln \alpha\}$		
RQard	rational quadratic, $\mathfrak{X} \subseteq \mathbb{R}^D$	$\sigma_f^2 \left(1 + \frac{1}{2\alpha} (\mathbf{x} - \mathbf{z})^\top \mathbf{\Lambda}^{-2} (\mathbf{x} - \mathbf{z})\right)^{-\alpha}$	$\{\ln \lambda_1,,\ln \lambda_D,\ln \sigma_f,\ln \alpha\}$		
SEiso	diagonal squared exponential, $\mathfrak{X} \subseteq \mathbb{R}^D$	$\sigma_f^2 \exp\left(-\frac{1}{2\ell^2}(\mathbf{x}-\mathbf{z})^\top(\mathbf{x}-\mathbf{z})\right)$	$\{\ln \ell, \ln \sigma_f\}$		
SEisoU	squared exponential, $\mathfrak{X} \subseteq \mathbb{R}^{D}$	$\exp\left(-\frac{1}{2\ell^2}(\mathbf{x}-\mathbf{z})^{\top}(\mathbf{x}-\mathbf{z})\right)$	ln ℓ		
SEard	automatic relevence determination squared exponential	$\sigma_{\rm f}^2 \exp\left(-\frac{1}{2}(\mathbf{x}-\mathbf{z})^{\top} \mathbf{\Lambda}^{-2}(\mathbf{x}-\mathbf{z})\right)$	$\{\ln \lambda_1,, \ln \lambda_D, \ln \sigma_f\}$		
SEvlen	spatially varying lengthscale squared exponential	$\sigma_f^2\left(\frac{a}{b}\right)^{\frac{D}{2}}\exp\left(-\frac{\ \mathbf{x}-\mathbf{z}\ ^2}{b}\right), a = 2\ell(\mathbf{x})\ell(\mathbf{z}), b = \ell^2(\mathbf{x}) + \ell^2(\mathbf{z})$	$\{\varphi_\ell, \ln \sigma_f\}$		
SEproj	factor analysis squared exponential	$\sigma_f^2 \exp\left(-\frac{1}{2}(\mathbf{x} - \mathbf{z})^{\top} \mathbf{L}^{\top} \mathbf{L}(\mathbf{x} - \mathbf{z})\right), \mathbf{L} \in \mathbb{R}^{d \times D}$	$\{L, \ln \sigma_f\}$		
LINard	linear with diagonal weighting	$\mathbf{x}^{T} \mathbf{\Lambda}^{-2} \mathbf{z}$	$\{\ln \lambda_1,, \ln \lambda_D\}$		
LINiso	linear with isotropic weighting	$\mathbf{x}^{T}\mathbf{z}/\ell^2$	ln ℓ		
Gaborard	anisotropic Gabor function, $\mathfrak{X}\subseteq\mathbb{R}^{D}, \lambda, p\in\mathbb{R}^{D}_{+}$	$\exp\left(-\sum_{d=1}^{D}\frac{t_d^2}{2\lambda_d^2}\right)\cos\left(2\pi\sum_{d=1}^{D}t_d/p_d\right),t_d=x_d-z_d$	$\{\ln \lambda, \ln p\}$		
Gaboriso	isotropic Gabor function, $\mathfrak{X} \subseteq \mathbb{R}^D$, $\ell, p \in \mathbb{R}_+$	$\exp(-\frac{\mathbf{t}^{T}\mathbf{t}}{2\ell^2})\cos(2\pi\mathbf{t}^{T}1/p), \mathbf{t} = \mathbf{x} - \mathbf{z}$	$\{\ln \ell, \ln p\}$		
SM	spectral mixture, $\mathfrak{X} \subseteq \mathbb{R}^D$, $\mathbf{w} \in \mathbb{R}_+^Q$, $\mathbf{M}, \mathbf{V} \in \mathbb{R}_+^{D \times Q}$	$\mathbf{w}^{\top} \left(\exp(-\frac{1}{2} \mathbf{V}^{\top} \mathbf{t}^2) \odot \cos(\mathbf{M}^{\top} \mathbf{t}) \right), \mathbf{t} = 2\pi(\mathbf{x} - \mathbf{z})$	$\{\ln w, \ln M, \ln V\}$		
	spectral mixture product, $\mathfrak{X} \subseteq \mathbb{R}^{D}$, $\mathbf{W} \in \mathbb{R}_{+}^{D \times Q}$, $\mathbf{M}, \mathbf{V} \in \mathbb{R}_{+}^{D \times Q}$	$\prod_{d=1}^{D} \mathbf{w}_{d}^{\top} \left(\exp(-\frac{1}{2}\mathbf{v}_{d}t_{d}^{2}) \odot \cos(\mathbf{m}_{d}t_{d}) \right), t_{d} = 2\pi(\mathbf{x}_{d} - \mathbf{z}_{d})$	$\{\ln W, \ln M, \ln V\}$		

The spectral mixture covariance covSM was introduced by Wilson & Adams Gaussian Process Kernels for Pattern Discovery and Extrapolation, ICML, 2013.

The periodic covariance function covPER starts from a stationary covariance function that depends on the data only through a distance $\mathbf{r}^2 = (\mathbf{x} - \mathbf{x}')^{\top} \boldsymbol{\Lambda}^{-2} (\mathbf{x} - \mathbf{x}')$ such as covMatern, covPP, covRQ, covSE and turns them into a periodic covariance function by embedding the data $\mathbf{x} \in \mathbb{R}^D$ into a periodic high-dimensional space $\mathbf{x}_p = \mathbf{u}(\mathbf{x}) \in \mathbb{R}^{2D}$ by a function $\mathbf{u}(\mathbf{x}) = 2\pi \text{diag}(\mathbf{p}^{-1})\mathbf{x}$.

The additive covariance function covADD starts from a one-dimensional covariance function $\kappa(x_i, x_i', \psi_i)$ acting on a single component $i \in [1, .., D]$ of \mathbf{x} . From that, we define covariance functions $\kappa_I(\mathbf{x}_I, \mathbf{x}_I) = \prod_{i \in I} \kappa(x_i, x_i', \psi_i)$ acting on vector-valued inputs \mathbf{x}_I . The sums of exponential size can efficiently be computed using the Newton-Girard formulae. Samples functions drawn from a GP with additive covariance are additive functions. The number of interacting variables |I| is a measure of how complex the additive functions are.

6.3 Usage of Implemented Covariance Functions

Some code examples taken from doc/usageCov.m illustrate how to use simple and composite covariance functions to specify a GP model.

Syntactically, a covariance function cf is defined by

```
cv := 'func' | @func // simple
cf := {cv} | {cv, {param, cf}} | {cv, {cf, ..., cf}} // composite
```

i.e., it is either a string containing the name of a covariance function, a pointer to a covariance function or one of the former in combination with a cell array of covariance functions and an additional list of parameters.

```
9 % set up simple covariance functions
10 cn = {'covNoise'}; sn = .1; hypn = log(sn); % one hyperparameter
11 cc = {@covConst}; sf = 2; hypc = log(sf); % function handles OK
12 \text{ ce} = \{@covEye\};
                               hype = [];
13 \text{ cl} = \{\text{@covLIN}\};
                                hypl = []; % linear is parameter-free
14 cla = {'covLINard'}; L = rand(D,1); hypla = log(L); % linear (ARD)
15 cli = {'covLINiso'}; 1 = rand(1); hypli = log(1); % linear iso
16 clo = {@covLINone}; ell = .9; hyplo = log(ell); % linear with bias
17 cp = \{(covPoly, 3)\}; c = 2; hypp = \log([c;sf]); % third order poly
\% Gaussian with ARD
20 cgu = {'covSEisoU'}; hypgu = log(ell); % isotropic Gauss no scale
21 cra = {'covRQard'}; al = 2; hypra = log([L;sf;al]); % ration. quad.
                             hypri = log([ell;sf;al]);  % isotropic
22 cri = {@covRQiso};
23 cma = {@covMaternard,5}; hypma = log([ell;sf]); % Matern class d=5
24 cmi = {'covMaterniso',3}; hypmi = log([ell;sf]); % Matern class d=3
25 cnn = {'covNNone'}; hypnn = log([L;sf]);
                                                     % neural network
26 cpe = {'covPeriodic'}; p = 2; hyppe = log([ell;p;sf]);  % periodic
27 cpn = {'covPeriodicNoDC'}; p = 2; hyppe = log([ell;p;sf]); % w/o DC
28 cpc = \{'covCos'\}; p = 2; hypcpc = log([p;sf]);
                                                   % cosine cov
29 cca = {'covPPard',3}; hypcc = hypgu; % compact support poly degree 3
30 cci = {'covPPiso',2}; hypcc = hypgi;% compact support poly degree 2
31 cgb = {@covGaboriso}; ell = 1; p = 1.2; hypgb=log([ell;p]); % Gabor
32 Q = 2; w = ones(Q,1)/Q; m = rand(D,Q); v = rand(D,Q);
34 cvl = {@covSEvlen, {@meanLinear}}; hypvl = [1;2;1; 0]; % var lenscal
35 s = 12; cds = {@covDiscrete,s}; % discrete covariance function
36 L = randn(s); L = chol(L'*L); L(1:(s+1):end) = log(diag(L));
37 hypds = L(triu(true(s))); xd = randi([1,s],[n,1]); xsd = [1;3;6];
38 cfa = {@covSEfact, 2}; hypfa = randn(D*2,1);
                                                  % factor analysis
39
40 % set up composite i.e. meta covariance functions
41 csc = {'covScale', {cgu}}; hypsc = [log(3); hypgu]; % scale by 9
42 csu = {'covSum', {cn,cc,cl}}; hypsu = [hypn; hypc; hypl];
43 cpr = {@covProd,{cc,cci}}; hyppr = [hypc; hypcc];
                                                            % product
44 mask = [0,1,0]; % binary mask excluding all but the 2nd component
45 cma = {'covMask', {mask, cgi{:}}}; hypma = hypgi;
46 % isotropic periodic rational quadratic
47 cpi = {'covPERiso',{@covRQiso}};
48 % periodic Matern with ARD
49 cpa = {'covPERard', {@covMaternard, 3}};
50 % additive based on SEiso using unary and pairwise interactions
51 cad = {'covADD', {[1,2], 'covSEiso'}};
52 % preference covariance with squared exponential base covariance
53 cpr = {'covPref', {'covSEiso'}}; hyppr = [0;0];
54 \text{ xp} = \text{randn}(n,2*D); \text{ xsp} = \text{randn}(3,2*D);
55
56 % 0) specify a covariance function
57 % cov = cma; hyp = hypma;
58 % cov = cci; hyp = hypcc;
59 \% \text{ cov} = \text{csm}; \text{hyp} = \text{hypsm};
60 \text{ cov} = \text{cds}; hyp = hypds; x = xd; xs = xsd;
61 % cov = cfa; hyp = hypfa;
62 \% cov = cvl; hyp = hypvl;
63 % cov = cpr; hyp = hyppr; x = xp; xs = xsp;
64
65 \% 1) query the number of parameters
66 feval(cov{:})
```

```
67
68 % 2) evaluate the function on x
69 [K,dK] = feval(cov{:},hyp,x)
70
71 % 3) evaluate the function on x and xs to get cross-terms
72 [kss,dkss] = feval(cov{:},hyp,xs,'diag')
73 [Ks, dKs] = feval(cov{:},hyp,x,xs)
```

7 Hyperpriors

A hyperprior $p(\theta)$ with $\theta = [\rho, \varphi, \psi]$ is a joint probability distribution over the likelihood hyperparameters ρ , the mean hyperparameters φ and the covariance hyperparameters ψ . We concentrate on factorial priors $p(\theta) = \prod_j p_j(\theta_j)$. Hyperpriors can be used to regularise the optimisation of the hyperparameters via the marginal likelihood $Z(\theta)$ so that $p(\theta)Z(\theta)$ is maximised instead. As we wish to perform unconstrained optimisation, we require (mainly) smooth hyperpriors with infinite support.

7.1 Interface

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In the GPML toolbox, a prior distribution $p(\theta)$ needs to implement the evaluation of the log density $\ln p(\theta)$ and its first derivative $\frac{\partial}{\partial \theta} \ln p(\theta)$. In addition, we require sampling capabilities i.e. the generation of $\theta \sim p(\theta)$.

```
\langle prior Distributions.m \ 46 \rangle \equiv
 1 % prior distributions to be used for hyperparameters of Gaussian processes
 2 % using infPrior.
 3 % There are two different kinds of prior distributions: simple and composite:
 4 %
 5 % simple prior distributions:
 6 %
 7 %
      priorGauss
                            - univariate Gaussian
 8 %
      priorLaplace
                            - univariate Laplace
9 %
      priorT
                            - univariate Student's t
10 %
      priorSmoothBox1
11 %
                            - univariate interval (linear decay in log domain)
12 %
      priorSmoothBox2
                            - univariate interval (quadr. decay in log domain)
13 %
14 %
     priorGamma
                            - univariate Gamma, IR+
15 %
      priorWeibull
                            - univariate Weibull, IR+
16 %
      priorInvGauss
                            - univariate Inverse Gaussian, IR+
17 %
      priorLogNormal
                            - univariate Log-normal, IR+
18 %
19 %
      priorClamped or
                           - fix hyperparameter to its current value by setting
20 %
      priorDelta
                           derivatives to zero, no effect on marginal likelihood
21 %
22 %
      priorGaussMulti
                            - multivariate Gauss
23 %
      priorLaplaceMulti
                            - multivariate Laplace
24 %
      priorTMulti
                            - multivariate Student's t
25 %
26 %
      priorClampedMulti or - fix hyperparameter to its current value by setting
27 %
      priorDeltaMulti
                           derivatives to zero, no effect on marginal likelihood
28 %
29 %
      priorEqualMulti or - make several hyperparameters have the same value by
30 %
      priorSameMulti same derivative, no effect on marginal likelihood
```

```
32 % composite prior distributions (see explanation at the bottom):
33 %
34 %
                            - nonnegative mixture of priors
     priorMix
35 % priorTransform
                         - prior on g(t) rather than t
36 %
37 % Naming convention: all prior distributions are named "prior/prior*.m".
38 %
39 %
40 % 1) With only a fixed input arguments:
41 %
42 %
     r = priorNAME(par1,par2,parN)
43 %
44 % The function returns a random sample from the distribution for e.g.
45 % random restarts, simulations or optimisation initialisation.
47 \% 2) With one additional input arguments:
48 %
49 %
        [lp,dlp] = priorNAME(par1,par2,parN, t)
50 %
51\ \% The function returns the log density at location t along with its first
52 % derivative.
53 %
54 % See also doc/usagePrior.m, inf/infPrior.m.
55 %
56 \(\langle gpml \copyright 5a\rangle
```

7.2 Implemented Hyperpriors

All code files are named according to the pattern prior/prior<NAME>.m for simple identification. All currently available hyperpriors are summarised in the following table.

Simple hyperpri	Simple hyperpriors $p(\theta)$					
	Univariate hyperpriors defined over the whole reals with mean μ and variace σ^2					
<name></name>	Meaning	$p(\theta) =$	τ			
Gauss	normally distributed hyperparameter $\theta \in \mathbb{R}$	$\frac{1}{\sigma\sqrt{2\pi}} \exp\left(-\frac{(\theta-\mu)^2}{2\sigma^2}\right)$ $\frac{1}{2b} \exp\left(-\frac{(\theta-\mu)}{b}\right), b = \sigma/\sqrt{2}$	$\mu \in \mathbb{R}, \sigma^2 \in \mathbb{R}_+$			
Laplace	double exponentially hyperparameter $\theta \in \mathbb{R}$		$\mu \in \mathbb{R}, \sigma^2 \in \mathbb{R}_+$			
Т	Student's t distributed hyperparameter $\theta \in \mathbb{R}$	$\frac{\Gamma(\frac{\nu+1}{2})}{\Gamma(\frac{\nu}{2})}\frac{1}{\sqrt{(\nu-2)\pi}\sigma}\left(1+\frac{(\theta-\mu)^2}{(\nu-2)\sigma^2}\right)^{-\frac{\nu+1}{2}}$	$\mu \in \mathbb{R}, \sigma^2, u \in \mathbb{R}_+$			
Univariate hype	rpriors with effective bounded support but defin	ned over the whole real line				
SmoothBox1	interval hyperparameter $\theta \in \mathbb{R}$ i.e. $\theta \stackrel{\approx}{\in} [a,b]$	$\begin{vmatrix} \frac{1 - \exp(\eta(a - b))}{b - a} \cdot \frac{1}{1 + \exp(-\eta(\theta - a))} \cdot \frac{1}{1 + \exp(\eta(\theta - b))} \\ \mu = \frac{a + b}{2}, \sigma^2 = \frac{w^2}{1 - \exp(-2a)} \frac{1 + \pi^2/g^2}{12}, g = \frac{w\eta}{2}, w = a - b \end{vmatrix}$	$a\leqslant b\in\mathbb{R},\eta\in\mathbb{R}_{+}$			
SmoothBox2	localised hyperparameter $\theta \in \mathbb{R}$ i.e. $\theta \stackrel{\approx}{\in} [\alpha,b]$	$\begin{split} &\frac{1-\exp(\eta(\alpha-b))}{b-\alpha} \cdot \frac{1}{1+\exp(-\eta(\theta-\alpha))} \cdot \frac{1}{1+\exp(\eta(\theta-b))} \\ &\mu = \frac{a+b}{2}, \ \sigma^2 = \frac{w^2}{1-\exp(-2g)} \frac{1+\pi^2/g^2}{12}, \ g = \frac{w\eta}{2}, \ w = a-b \\ &\frac{1}{(1/\eta+1)(b-\alpha)} \begin{cases} \mathcal{N}(\theta a, \sigma_{ab}^2) & t < a \\ 1 & t \in [a,b], \ \sigma_{ab} = \frac{b-\alpha}{\eta\sqrt{2\pi}} \\ \mathcal{N}(\theta a, \sigma_{ab}^2) & b < t \end{cases} \\ &\mu = \frac{a+b}{2}, \ \sigma^2 = \frac{w^2}{4} \frac{\eta^3/3+\eta^2+4\eta/\pi+2/\pi}{\eta^3+\eta^2}, \ w = a-b \end{split}$	$a\leqslant b\in\mathbb{R},\eta\in\mathbb{R}_{+}$			
Univariate hype	erpriors supported only over the positive reals					
Gamma	Gamma hyperparameter $\theta \in \mathbb{R}_+$	$\frac{1}{\Gamma(k)t^k} \exp(\frac{\theta}{t})\theta^{k-1}$	$k \in \mathbb{R}_+, t \in \mathbb{R}_+$			
Weibull	Weibull hyperparameter $\theta \in \mathbb{R}_+$	$\frac{k}{\lambda} \left(\frac{\theta}{\lambda} \right)^{k-1} \exp \left(-\left(\frac{\theta}{\lambda} \right)^k \right)$	$k \in \mathbb{R}_+, \lambda \in \mathbb{R}_+$			
InverseGauss	inverse Gaussian hyperparameter $\theta \in \mathbb{R}_+$	$\frac{1}{\frac{\lambda}{\lambda}\left(\frac{\theta}{\lambda}\right)^{k-1}} \exp\left(-\left(\frac{\theta}{\lambda}\right)^{k}\right)}{\frac{1}{2\pi\theta^{3}/\lambda}} \exp\left(-\frac{\lambda(\theta-\mu)^{2}}{2\mu^{2}\theta}\right)$	$k\in\mathbb{R}_+,\lambda\in\mathbb{R}_+$			
LogNormal	log-normal hyperparameter $\theta \in \mathbb{R}_+$	$\mathcal{N}(\theta \mu, \sigma^2) = \frac{1}{\theta \sigma \sqrt{2\pi}} \exp\left(-\frac{(\ln \theta - \mu)^2}{2\sigma^2}\right)$	$\mu \in \mathbb{R}, \sigma^2 \in \mathbb{R}_+$			
Multivariate hy	perpriors supported all over \mathbb{R}^D with mean μ as					
GaussMulti	multivariate normal distribution $\theta \in \mathbb{R}^D$	$ 2\pi\mathbf{\Sigma} ^{-\frac{1}{2}}\exp\left(-\frac{1}{2}(\mathbf{\theta}-\mathbf{\mu})^{\top}\mathbf{\Sigma}^{-1}(\mathbf{\theta}-\mathbf{\mu})\right)$	$\mu \in \mathbb{R}^{D}, \Sigma \in \mathbb{R}^{D imes D}$			
LaplaceMulti	multivariate Laplace distribution $\theta \in \mathbb{R}^D$	$ \sqrt{2}\Sigma ^{-\frac{1}{2}}\exp\left(-\sqrt{2}\left\ \mathbf{L}^{-1}(\mathbf{\theta}-\mathbf{\mu})\right\ _{1}\right), \mathbf{L}^{\top}\mathbf{L} = \mathbf{\Sigma}$	$\mu \in \mathbb{R}^{D}, \Sigma \in \mathbb{R}^{D \times D}$			
TMulti	multivariate Student's t distribution $\boldsymbol{\theta} \in \mathbb{R}^D$	$ (\mathbf{v}-2)\pi\mathbf{\Sigma} ^{-\frac{1}{2}} \frac{\Gamma(\frac{\mathbf{v}+\mathbf{D}}{2})}{\Gamma(\frac{\mathbf{v}}{2})} \left(1 + \frac{(\mathbf{\theta}-\mathbf{\mu})^{\top}\mathbf{\Sigma}^{-1}(\mathbf{\theta}-\mathbf{\mu})}{(\mathbf{v}-2)}\right)^{-\frac{\mathbf{v}+\mathbf{D}}{2}}$	$\mu \in \mathbb{R}^{D}, \Sigma \in \mathbb{R}^{D \times D}, \nu \in \mathbb{R}$			
Improper hyperpriors used to fix the value of a particular hyperparameter						
Delta Clamped	clamped hyperparameter $\theta=\theta_0\in\mathbb{R}$	$\delta(\theta - \theta_0)$	Ø			
DeltaMulti ClampedMulti	clamped hyperparameter $\boldsymbol{\theta} = \boldsymbol{\theta}_0 \in \mathbb{R}^D$	$\delta(\theta - \theta_0)$	Ø			
SameMulti EqualMulti	same hyperparameter $\theta = \theta_0 1 \in \mathbb{R}^D$	$\delta\left(\textstyle\prod_{i=1}^D(\theta_0-\theta_i)\right)$	Ø			
Composite hyperpriors $[\pi_1(\theta), \pi_2(\theta),] \mapsto p(\theta)$						
Transform	prior distribution on $g(\theta)$ instead of θ	$\pi(g(\theta))$	{g}			
Mix	mixture distribution	$\sum_{i} w_{i} \pi_{i}(\theta)$	{ w }			

The priorSmoothBox2 is a Gauss-uniform sandwich obtained by complementing a uniform distribution on $[\mathfrak{a},\mathfrak{b}]$ with two Gaussian halves at each side. The parameter \mathfrak{n} balances the probability mass between the constituents so that $\mathfrak{n}/(\mathfrak{n}+1)$ is used for the box and $1/(\mathfrak{n}+1)$ for the Gaussian sides. Its brother priorSmoothBox1 is the product of two sigmoidal functions.

The priorDelta or equivalently priorClamped can be used to exclude some hyperparameters from the optimisation. Their values are clamped to θ_0 and the derivative vanishes. There are also multivariate counterparts priorDeltaMulti and priorClampedMulti.

7.3 Usage of Implemented Hyperpriors

Some code examples taken from doc/usagePrior.m illustrate how to use univariate, multivariate and composite priors on hyperparameters. Syntactically, a hyperprior hp is defined by

i.e., it is either a string containing the name of a hyperprior function, a pointer to a hyperprior function or one of the former in combination with a cell array of hyperprior functions and an additional list of parameters. Furthermore, we have multivariate hyperprior variants and 2 (equivalent) predefined hyperpriors allowing to exclude variables from optimisation.

```
\langle doc/usagePrior.m \ 49 \rangle \equiv
49
      1 % demonstrate usage of prior distributions
      3 % See also priorDistributions.m.
      4 %
      5 (gpml copyright 5a)
      6 clear all, close all
      8 % 1) specify some priors
     9 % a) univariate priors
     10 \text{ mu} = 1.0; \text{ s2} = 0.01^2; \text{ nu} = 3;
     11 pg = {@priorGauss,mu,s2};
                                                               % Gaussian prior
     12 pl = {'priorLaplace', mu, s2};
                                                                % Laplace prior
     13 pt = {@priorT, mu, s2, nu};
                                                            % Student's t prior
     14 p1 = {@priorSmoothBox1,0,3,15}; % smooth box constraints lin decay
     15 p2 = {@priorSmoothBox2,0,2,15}; % smooth box constraints qua decay
     16 pd = {'priorDelta'}; % fix value of prior exclude from optimisation
     17 pc = {@priorClamped};
                                                          % equivalent to above
     18 \text{ lam} = 1.05; k = 2.5;
     19 pw = {@priorWeibull,lam,k};
                                                                % Weibull prior
     20
     21 % b) meta priors
     22 pmx = {@priorMix,[0.5,0.5],{pg,pl}}; % mixture of two priors
     23 g = 0exp; dg = 0exp; ig = 0log;
     24 ptr = {@priorTransform,g,dg,ig,pg}; % Gaussian in the exp domain
     26 % c) multivariate priors
     27 \text{ m} = [1;2]; V = [2,1;1,2];
     28 pG = {@priorGaussMulti,m,V};
                                                            % 2d Gaussian prior
     29 pD = {'priorDeltaMulti'};  % fix value of prior exclude from optim
     30 pC = {@priorClampedMulti};
                                                         % equivalent to above
     31
     32 % 2) evaluation
     33 % pri = pt; hp = randn(1,3);
     34 \% \text{ pri} = pmx; \text{ hp} = randn(1,3);
     35 % pri = ptr; hp = randn(1,3);
     36 \text{ pri} = pG; \quad \text{hp} = \text{randn}(2,3);
     37
     38~\% a) draw a sample from the prior
     39 feval(pri{:})
     41 % b) evaluate prior and derivative if requires
     42 [lp,dlp] = feval(pri{:},hp)
     43
     44 % 3) comprehensive example
     45 \times = (0:0.1:10); y = 2*x + randn(size(x)); % generate training data
     46 mean = {@meanSum,{@meanConst,@meanLinear}}; % specify mean function
     47 cov = {@covSEiso}; lik = {@likGauss}; % specify covariance and lik
     48 hyp.cov = [log(1);log(1.2)]; hyp.lik = log(0.9); hyp.mean = [2;3];
```

```
49 par = {mean,cov,lik,x,y}; mfun = @minimize; % input for GP function
50
51 % a) plain marginal likelihood optimisation (maximum likelihood)
52 im = @infExact;
                                                    % inference method
53 hyp_plain = feval(mfun, hyp, @gp, -10, im, par{:});
                                                            % optimise
55~\% b) regularised optimisation (maximum a posteriori) with 1d priors
56 prior.mean = {pg;pc}; % Gaussian prior for first, clamp second par
57 prior.cov = {p1;[]}; % box prior for first, nothing for second par
58 im = {@infPrior,@infExact,prior};
                                                    % inference method
59 hyp_p1 = feval(mfun, hyp, @gp, -10, im, par{:});
                                                            % optimise
61 % c) regularised optimisation (maximum a posteriori) with Nd priors
62 prior = [];
                                                 % clear the structure
63 % multivariate Student's t prior on the first and second mean hyper
64 prior.multi{1} = {@priorTMulti,[mu;mu],diag([s2,s2]),nu,...
65
                    struct('mean',[1,2])};
                                                    % use hyper struct
66 % Equivalent shortcut (same mu and s2 for all dimensions)
67 prior.multi{1} = {@priorTMulti,mu,s2,nu,struct('mean',[1,2])};
68 % multivariate Gaussian prior jointly on 1st and 3rd hyper
69 prior.multi{2} = {@priorGaussMulti,[mu;mu],diag([s2,s2]),...
70
                    [1,3]};
                                          % use unwrapped hyper vector
71 % Equivalent shortcut (same mu and s2 for all dimensions)
72 prior.multi{2} = {@priorGaussMulti,mu,s2,[1,3]};
73 im = {@infPrior,@infExact,prior};
                                                    % inference method
74 hyp_pN = feval(mfun, hyp, @gp, -10, im, par{:});
                                                            % optimise
76 [any2vec(hyp), any2vec(hyp_plain), any2vec(hyp_p1), any2vec(hyp_pN)]
```