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Sequential dimension reduction
for learning features of expensive black-box functions *

Malek Ben Salem †, François Bachoc‡, Olivier Roustant §, Fabrice Gamboa †, and Lionel Tomaso ¶

Abstract. Learning a feature of an expensive black-box function (optimum, contour line,...) is a difficult
task when the dimension increases. A classical approach is two-stage. First, sensitivity analysis
is performed to reduce the dimension of the input variables. Second, the feature is estimated by
considering only the selected influential variables. This approach can be computationally expensive
and may lack flexibility since dimension reduction is done once and for all. In this paper, we
propose a so called Split-and-Doubt algorithm that performs sequentially both dimension reduction
and feature oriented sampling. The ‘split’ step identifies influential variables. This selection relies
on new theoretical results on Gaussian process regression. We prove that large correlation lengths
of covariance functions correspond to inactive variables. Then, in the ‘doubt’ step, a doubt function
is used to update the subset of influential variables. Numerical tests show the efficiency of the
Split-and-Doubt algorithm.

Key words. Variable selection, Surrogate modeling, Design of experiments, Bayesian optimization

AMS subject classifications. 62L05, 62K99, 62F15

1. Introduction. In design problems, the goal may be the estimation of a feature of an
expensive black-box function (optimum, probability of failure, level set, ...). Several methods
have been proposed to achieve this goal. Nevertheless, they generally suffer from the curse of
dimensionality. Thus, their usage is limited to functions depending on a moderate number of
variables. Meanwhile, most of real life problems are complex and may involve a large number
of variables.

Let us focus first on high-dimensional optimization problems. In this context, we look for
a good approximation of a global minimum of an expensive-to-evaluate black-box function
\( f : \Omega = [0, 1]^D \to \mathbb{R} \) using a limited number of evaluations of \( f \). That is, we aim at approximating
\( x^\star \in \Omega \) such that:

\[
(1.1) \quad x^\star \in \arg\min_{x \in \Omega} f(x).
\]

Bayesian optimization (BO) techniques have been successfully used in various problems
[19, 18, 17, 27, 31]. These methods give interesting results when the number of evaluations

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of the function $f$ is relatively low [15]. They are generally limited to problems of moderate dimension, typically up to about 10 [32]. Here, we are particularly interested in the case where the dimension $D$ is large and the number of influential variables $d$, also called effective dimension, is much smaller: $d << D$. In this case, there are different approaches to tackle the dimensionality problem.

A direct approach consists in first performing global sensitivity analysis. Then, the most influential variables are selected and used in the parametric study. Chen et al. [11] stated that “Variables selection and optimization have both been extensively studied separately from each other”. Most of these methods are two-stage: First, the influential variables are selected and then optimization is performed on these influential variables. These strategies are generally computationally expensive. Furthermore, the set of selected variables does not take into account the new data. However, this new information may modify the results of the sensitivity analysis study. For an overview of global sensitivity analysis methods, one may refer to [16].

Some Bayesian optimization techniques are designed to handle the dimensionality problem. For instance, the method called Random EMbedding Bayesian Optimization (REMBO) selects randomly the subspace of influential variables [32, 9]. The main strengths of REMBO are that the selected variables are linear combinations of the input variables and that it works for huge values of $D$. However, the effective dimension $d$ must be specified.

In this paper, we propose a versatile sequential dimension reduction method called Split-and-Doubt. The design is sequentially generated in order to achieve jointly two goals. The first goal is the estimation of the optimum (in the optimization case). The second one is the learning of the influential variables. In the “split” step, the algorithm selects the set of influential variables based on the values of the correlation lengths of Automatic Relevance Determination (ARD) covariances. We show theoretical results that support the intuition that large correlation lengths correspond to inactive variables. The “doubt” step questions the “split” step and helps correcting the estimation of the correlation lengths.

The paper is organized as follows. Section 2 presents the background and the notations. Section 3 introduces the Split-and-Doubt. The algorithm is based on theoretical results stated in Section 4. Finally, Section 5 illustrates the performance of the algorithm on various test functions. For readability, proofs are postponed to Section 6. Concluding remarks are given in Section 7.

2. General notations and background.

2.1. Gaussian Process Regression (GPR). Kriging or Gaussian process regression (GPR) models predict the outputs of a function $f: \Omega = [0, 1]^D \rightarrow \mathbb{R}$, based on a set of $n$ observations [30, 23]. It is a widely used surrogate modeling technique. Its popularity is mainly due to its statistical nature and properties. Indeed, it is a Bayesian inference technique that provides an estimate of the prediction error distribution. This uncertainty is an efficient tool to construct strategies for various problems such as prediction refinement, optimization or inversion.

The GPR framework uses a centered real-valued Gaussian Process (GP) $Y$ over $\Omega$ as a prior distribution for $f$. The predictions are given by the conditional distribution of $Y$ given the observations $y = (y_1, \ldots, y_n)^\top$ where $y_i = f(x^{(i)})$ for $1 \leq i \leq n$. We denote by $k_\theta : \Omega \times \Omega \rightarrow \mathbb{R}$ the covariance function (or kernel) of $Y$: $k_\theta(x, x') = \text{Cov}[Y(x), Y(x')]$ ($((x, x') \in \Omega^2$), by $X = (x^{(1)}, \ldots, x^{(n)})^\top \in \Omega^n$ the matrix of observation locations and
by $Z = (X \ y)$ the matrix of observation locations and values where $x^{(i)} = (x_1^{(i)}, \ldots, x_D^{(i)})$ for $1 \leq i \leq n$. $\theta$ is a parameter that will be discussed later. Without loss of generality, we consider the simple kriging framework. The a posteriori conditional mean $m_{\theta,Z}$ and the a posteriori conditional variance $\hat{\sigma}^2_{\theta,Z}$ are given by:

\begin{equation}
(2.1) \quad m_{\theta,Z}(x) = k_\theta(x, X) \top K^{-1}_\theta y
\end{equation}

\begin{equation}
(2.2) \quad \hat{\sigma}^2_{\theta,Z}(x) = k_\theta(x, x) - k_\theta(x, X) \top K^{-1}_\theta k_\theta(x, X).
\end{equation}

Here, $k_\theta(x, X)$ is the vector $(k_\theta(x, x^{(1)}), \ldots, k_\theta(x, x^{(n)}))^\top$ and $K_\theta = K_\theta(X, X)$ is the invertible matrix with entries $k_\theta(x^{(i)}, x^{(j)})$, for $1 \leq i, j \leq n$.

Several methods are useful to select the covariance function. A common approach consists in assuming that the covariance function belongs to a parametric family. In this paper, we consider the Automatic Relevance Determination (ARD) kernels defined in (2.3). A review of classical covariance functions is given in [1].

\begin{equation}
(2.3) \quad k_\theta(x, y) = \sigma^2 \prod_{p=1}^D k\left(\frac{d(x_p, y_p)}{\theta_p}\right), \text{ for } x, y \in \Omega.
\end{equation}

Here, $d(\cdot)$ is a distance on $\Omega \times \Omega$ and $k : \mathbb{R} \rightarrow \mathbb{R}$ is a fixed stationary covariance function. Without loss of generality, we suppose that the hyper-parameter $\sigma$ is fixed while $\theta_1, \ldots, \theta_D$ have to be estimated. The ARD kernels include most popular kernels such as the exponential kernel, the Matérn 5/2 kernel and the squared exponential kernel.

The hyper-parameters of these parametric families can be estimated by maximum Likelihood (ML) or cross validation (CV). Both methods have interesting asymptotic properties [2, 4, 5]. Nevertheless, when the number of observations is relatively low, the estimation can be misleading. These methods are also computationally demanding when the number of observations is large.

On one hand, estimating the correlation lengths by the maximum likelihood estimator gives the estimator $\hat{\theta}_{MLE}^* \in \arg \max_{\theta} l_Z(\theta)$ where the likelihood $l_Z(\theta)$ is given in (2.4).

\begin{equation}
(2.4) \quad l_Z(\theta) = \frac{1}{(2\pi)^n/2} \frac{1}{\sqrt{\det(k_\theta(X, X))}} \exp\left(-y^\top k_\theta(X, X)^{-1} y\right).
\end{equation}

On the other hand, the idea behind Cross-validation (CV) is to estimate the prediction errors by splitting the observations once or several times. One part is used as a test set while the remaining parts are used to construct the model. The Leave-One-Out Cross-Validation (LOO-CV) consists in dividing the $n$ points into $n$ subsets of one point each. Then, each subset plays the role of test set while the remaining points are used together as the training set. Using Dubrule’s formula [13], the LOO-CV estimator is given in (2.5).
\[
\hat{\theta}_{CV}^* \in \arg \min_{\theta} \frac{1}{n} y^\top K_{\theta}^{-1} \text{diag}(K_{\theta}^{-1})^{-1} K_{\theta}^{-1} y.
\]

For more insight on these estimators, one can refer to [3].

2.2. Derivative based global sensitivity measures: DGSM. Sobol and Kucherenko [28, 29] proposed the so-called Derivative-based Global Sensitivity Measures (DGSM) to estimate the influence of an input variable on a function \( f : \Omega = [0, 1]^D \rightarrow \mathbb{R} \). For each variable \( x_i \), the index \( \vartheta_i \) is the global energy of the corresponding partial derivatives.

\[
\vartheta_i(f) = \int_{\Omega} \left( \frac{\partial f(x)}{\partial x_i} \right)^2 dx, \quad i = 1, \ldots, D.
\]

DGSM provides a quantification of the influence of a single input on \( f \). Indeed, assuming that \( f \) is of class \( C^1 \), then \( x_i \) is not influential iff \( \frac{\partial f}{\partial x_i}(x) = 0, \forall x \in \Omega \) iff \( \vartheta_i = 0 \). DGSM has recently shown its efficiency for the identification of non-influential inputs [24]. We further define the normalized DGSM \( \tilde{\vartheta}_i \) in (2.7). \( \tilde{\vartheta}_i \) measures the influence of \( x_i \) with regard to the total energy.

\[
\tilde{\vartheta}_i(f) = \frac{\vartheta_i(f)}{\sum_{p=1}^{D} \vartheta_p(f)}, \quad i = 1, \ldots, D.
\]


3.1. Definitions.

Variable splitting. Let us consider the framework of a GPR using a stationary ARD kernel. Intuitively, large correlation length values correspond to inactive variables in the function. We prove this intuition in Proposition 4.1. The influential variables are selected in our algorithm according to the estimated values of their corresponding correlation lengths. We show also that the ML (and CV) estimator is able to assign asymptotically large correlation length values to the inactive variables (Propositions 4.3 and 4.4).

Let \( \hat{\theta}^* = (\hat{\theta}_1^*, \ldots, \hat{\theta}_D^*) \) be the ML estimation of the correlation lengths:

\[
\hat{\theta}^* \in \arg \max_{\theta} l_{\mathcal{Z}}(\theta).
\]

The influential variables are then selected according to the estimated values of their corresponding correlation lengths. We split the indices into a set of influential variables \( I_M \) and a set of minor variables \( I_m \) as follows:

- \( I_M = \{ i : \hat{\theta}_i^* < T \} \)
- \( I_m = \{ i : \hat{\theta}_i^* \geq T \} \)

where \( T \in \mathbb{R} \) is a suitable threshold. Let \( d_M \) (resp. \( d_m \)) be the size of \( I_M \) (resp. \( I_m \)). We further call \( \Omega_m := [0, 1]^{d_m} \) the minor subspace, that is the space of minor variables and
Ω_M := [0, 1]^d_M the major subspace, that is the subspace of major variables. We will use the set notation: for a set I of \{1, \ldots, D\}, x_I will denote the vector extracted from x with coordinates x_i, i \in I. Hence, x_{IM} (resp. x_{Im}) denotes the sub-vector of x whose coordinates are in the major (resp. minor) subspace. For simplicity, we will also write x = (x_{IM}, x_{Im}), without specifying the re-ordering used to obtain x by gathering x_{IM} and x_{Im}.

**Doubt.** The doubt of a correlation length \( \theta \) measures the influence on \( m_{\theta,Z} \) of the variables from the minor subspace \( \Omega_m \). It is a decreasing function of the correlation lengths. We will use it to question the variable splitting.

**Definition 3.1 (Doubt).** Let \( \delta \) be the following function associated with a variable splitting \((I_m, I_M)\). For all vector \( \theta = (\theta_1, \ldots, \theta_D) \in \mathbb{R}^D \):

\[
\delta(\theta) = \sum_{i \in I_m} \max(\theta_i^{-1} - T^{-1}, 0).
\]

**Contrast.** Given two different correlation lengths \( \theta^{(1)} \) and \( \theta^{(2)} \) and a location \( x \), the contrast measures the discrepancy between the predictions using \( \theta^{(1)} \) and \( \theta^{(2)} \) at \( x \). It will be used to build a sequential design in the minor subspace.

**Definition 3.2 (Prediction contrast).** For a point \( x \) and two correlation lengths \( \theta^{(1)} \) and \( \theta^{(2)} \), the prediction contrast \( PC(x, \theta^{(1)}, \theta^{(2)}) \) is

\[
PC(x, \theta^{(1)}, \theta^{(2)}) = |m_{\theta^{(1)},Z}(x) - m_{\theta^{(2)},Z}(x)|.
\]

### 3.2. The algorithm.

The **Split-and-Doubt** algorithm performs a new variable selection at each iteration. It samples a point in two steps: a goal-oriented sampling in the major subspace and a sampling of the minor variables to question the variable selection done at the previous step. The **Split-and-Doubt** algorithm for optimization using the expected improvement (EI) criterion [18] is described in Algorithm 3.1.

Here, the algorithm is applied for optimization (Step 6). We used the Expected Improvement criterion (3.1).

\[
EI_{Z}(x) = \mathbb{E} \left[ \max(\min_i y_i - Y(x), 0) | Z \right].
\]

(3.1)

It is possible to use any other classical optimization criterion to sample \( x^*_M \). We can use other criteria for other purposes such as contour estimation [21, 22, 8], probability of failure estimation [6] or surrogate model refinement [10].

The settings of the algorithm are mainly the kernel \( k \), the limit \( \ell \) and the threshold \( T \). An other hidden setting is the search space for the ML estimator. We use a Matérn 5/2 kernel and we set \( \ell = \text{erf}(\frac{1}{\sqrt{2}}) \) and an adaptive threshold \( T = 20 \min_{i \in \{1, \ldots, D\}} (\hat{\theta}_i) \).

### 3.3. Remarks on the steps of the Split-and-Doubt algorithm.

**Remark on the doubt.** When the observations do not carry enough information, it is hard to estimate accurately the correlation lengths. The use of such values can lead to unsatisfactory results [14, 7]. In our algorithm, the estimated correlation lengths are used to select the major variables. If this estimation is done once and for all, poor estimation can lead to considering
**Algorithm 3.1** Split-and-Doubt-EGO (f)

1. **Algorithm parameters:** \( \ell, k, T \).
2. **Start:** Inputs: \( Z = (X, y) \).
3. **while** Stop conditions are not satisfied **do**
   4. Estimate the correlation lengths: \( \hat{\theta}^* \in \arg \max_{\theta \in \mathbb{R}^+} l_Z(\theta) \) Eq (2.4).
   5. Split the variables: Define the major set \( I_M = \{ i; \hat{\theta}^*_i < T \} \) and the minor set \( I_m = \{ i; \hat{\theta}^*_i \geq T \} \), \( d_m = |I_m| \).
   6. Design in the major subspace: Compute \( x_M^* \) according to the objective function in the major subspace (by EI for instance): We compute a new GPR considering only the major variables to compute the EI. Let \( Z_M = (X_{I_M}, y) \)
   \[ x_M^* \in \arg \max_{x_M \in \Omega_M} \text{EI}_{Z_M}(x_M). \]
   7. Doubt the variable splitting: Compute a challenger \( \theta' \) for correlation lengths.
   \[ \theta' \in \arg \max_{\theta \in \mathbb{R}^+} \delta(\theta) \text{ subject to } 2 \left| \frac{\ln l_Z(\theta)}{\ln l_Z(\hat{\theta}^*)} \right| < \chi^2(\ell, d_m). \]
   8. Design in the minor subspace: Compute \( x_m^* \) by maximum contrast with the challenger \( \theta' \)
   \[ x_m^* \in \arg \max_{x_m \in \Omega_m} \text{PC} \left( x = (x_M^*, x_m), \hat{\theta}^*, \theta' \right). \]
   9. Update: Evaluate the new point output \( y_{n+1} = f(x^{(n+1)}) \) with \( x_m^{(n+1)} = x_{m}^* \) and \( x_M^{(n+1)} = x_{M}^* \) and add the new point to the design:
   \[ X^T \leftarrow (X^T \ x^{(n+1)}), \quad y^T \leftarrow (y^T, y_{n+1}). \]
10. **end while**
11. **return** Outputs: \( Z = (X, y) \).

A major variable inactive. So, it is important to always question the estimation. Therefore, we look for a "challenger kernel" at each iteration. Specifically, we are looking for correlation lengths that maximize the doubt and that are accepted by a likelihood ratio test. Indeed, this is why we limit the search space by a likelihood ratio deviation from the estimated correlation lengths \( \hat{\theta}^* \): \( \Theta_l = \{ \theta; 2 \left| \frac{\ln l_Z(\theta)}{\ln l_Z(\hat{\theta}^*)} \right| < l \} \). Notice that we used \( l = \chi^2(\ell, d_m). \) Following [14, 11], the likelihood ratio test is compared to the \( \chi^2 \) distribution to decide whether the correlation lengths are allowable or not.

**Remark on the contrast.** Sampling the coordinates in the non-influential variable subspace \( \{ x_M^* \} \times \Omega_m = \{ (x_M^*, x_m), x_m \in \Omega_m \} \) aims at revealing the contrast between the maximum
likelihood correlation lengths $\hat{\theta}^\star$ and a challenging correlation length parameter is $\theta'$. The main idea is to sample the point that helps either correcting the first estimation or reducing the allowable doubt space $\Theta$ in order to strengthen the belief in the estimated kernel.

We could have used an alternative direct approach. It consists in maximizing the likelihood ratio between two estimations of the correlation lengths in the future iterations.

**Definition 3.3 (likelihood contrast).** For a point $x$ and two correlation lengths $\theta(1)$ and $\theta(2)$, the likelihood contrast $LC$ is:

$$LC(x, \theta(1), \theta(2)) = \mathbb{E}[\left|\ln\left(\frac{L(\theta(1), Z \cup (x, \hat{Y}(x)))}{L(\theta(2), Z \cup (x, \hat{Y}(x)))}\right)\right|]$$

where $\hat{Y}(x) \sim \mathcal{N}(m_{\theta_2,B}(x), (\hat{\sigma}_{\theta_2,B}(x))^2)$.

However, this approach is computationally more expensive. Therefore, we prefer to use the prediction contrast (Definition 3.2).

### 3.4. Example: Illustration of the contrast effect.**

We illustrate here how the Doubt/Contrast strategy can help correcting an inaccurate variable splitting. To do so, let us consider the following example. Let $f(x_1, x_2) = \cos(2\pi x_2)$. We assume that we have at hands four design points $x^{(1)} = (0, \frac{2}{3})$, $x^{(2)} = (\frac{1}{3}, 0)$, $x^{(3)} = (\frac{2}{3}, 1)$ and $x^{(4)} = (1, \frac{1}{3})$ and their corresponding responses $y_1 = y_4 = f(x^{(1)}) = f(x^{(4)}) = -0.5$ and $y_2 = y_3 = f(x^{(2)}) = f(x^{(3)}) = 1$. Here, the search space for the correlation lengths is $[0.5, 10]^2$.

**Misleading estimation.** The log-likelihood of the correlation lengths in the search space $[0.5, 10]^2$ for the Matérn 5/2 kernel is displayed in Figure 1. Notice that the likelihood is maximized for different values of $\theta$ and that $\hat{\theta}^\star = (0.5, 10)$ is among these values:

$$(0.5, 10) \in \arg\max_{\theta \in \mathbb{R}_+^D} l_Z(\theta).$$

**Figure 1.** Left: $f(x_1, x_2) = \cos(2\pi x_2)$, the color code indicates the values of $f$ and solid black circles indicate the design points. Middle: log-likelihood of the correlation lengths, solid black triangle: $\hat{\theta}^\star$. Right: The predictions given by the GPR using $k_{\hat{\theta}}$. 
We also display in Figure 1 the function \( f \), the design points, and the predictions using \( k_{\theta^*} \). This example shows that a limited number of observations may lead to inaccurate correlation lengths and consequently inaccurate predictions.

**Doubt/Contrast strategy.** Adding more points will arguably improve the quality of the correlation lengths estimation. Here, we want to bold that the improvement due to the Doubt/Contrast strategy is not only related to the fact that more points are sampled. To do so, we set \( T = 10 \). So, \( I_M = \{1\} \) and \( I_m = \{2\} \). Notice that the challenger correlation lengths is \( \theta' = (0.5, 0.5) \). It gives the maximum doubt \( \delta(\theta') = \frac{1}{0.5} - \frac{1}{10} = 1.9 \). In this example, we are sampling the fifth point \( x^{(5)} \). The value sampled by the EI in the major space is \( x^*_{M} = 0.64 \).

We display in Figure 2 the prediction contrast \( PC((x^*_M, x_2), \hat{\theta^*}, \theta') \) as a function of \( x_2 \).

**Figure 2.** The prediction contrast in function of \( x_2 \).

Let us now consider two cases: a) we add the point sampled by the maximum contrast \( (x^*_M, x^*_m) \) and b) we add the point with the minimum contrast \( (x^*_M, 1) \). For both cases, we display the updated likelihood function in Figure 3.

Notice that:

a) For \( x_m = x^*_m \) (maximum contrast), the log-likelihood has larger value for small values of \( \theta_2 \). Thus, the same inaccurate variable splitting is prevented.

b) For \( x_m = 1 \) (small contrast value), we may still use the same misleading variable splitting.

Even if this 2-dimensional toy example is pathological, it illustrates the interests of the Doubt/Contrast strategy. This strategy can be valuable in high dimension when the number of observations is relatively small to estimate accurately the correlation lengths.

4. **Links between correlation lengths and variable importance.** In this section, we consider a deterministic function \( f : \mathbb{R}^D \to \mathbb{R} \) to be modeled as a GP path. We consider the
centered stationary GP with a covariance function \( k_\theta \) defined by

\[
k_\theta(h) = \prod_{i=1}^{D} k(h_i/\theta_i).
\]

Here \( k : \mathbb{R} \to \mathbb{R} \) is a fixed covariance function satisfying \( k(0) = 1 \) and \( \theta \in (0, \infty)^D \) is the vector of correlation lengths. As an example, \( k \) may be the function \( k(h) = e^{-h^2} \).

Intuitively, a small correlation length \( \theta_i \) for the GP should correspond to an input variable \( x_i \) that has an important impact on the function value \( f(x) \). Conversely, if the function \( f \) does not depend on \( x_i \), then the length \( \theta_i \) should ideally be infinite. This intuition is precisely the motivation for the Split-and-Doubt algorithm suggested in Section 3.

In this section, we show several theoretical results that confirm this intuition. First, we show that if the correlation length \( \theta_i \) goes to zero (respectively infinity) then the derivative-based global sensitivity measure, obtained from the GP predictor for the input \( x_i \), tends to its maximum value 1 (resp. its minimum value 0). Then, we show that an infinite correlation length \( \theta_i \) can provide an infinite likelihood or a zero LOO mean square error, for the GP model, when the function \( f \) does not depend on \( x_i \).

We use the additional following notations throughout the section. For \( D, p, q \in \mathbb{N}^* \), for a covariance function \( g \) on \( \mathbb{R}^D \), for two \( p \times D \) and \( q \times D \) matrices \( X \) and \( Z \), we denote by \( g(X, Z) \) the \( p \times q \) matrix defined by \( [g(X, Z)]_{i,j} = g(X_l, Z_j) \) where \( M_l \) is the line \( l \) of a matrix \( M \). When \( d = 1 \), \( p = 1 \) or \( q = 1 \), we identify the corresponding matrices with vectors. We further assume:

**Assumption 1 (Invertibility assumption).** For any \( p, d \in \mathbb{N} \), for any \( \theta \in (0, \infty)^D \), for any \( p \times d \) matrix \( X \) with two-by-two distinct lines, the matrix \( k_\theta(X, X) \) is invertible.

This assumption holds for instance when the spectral density of a stationary kernel is absolutely continuous. Further, for any vector \( u, u_{-i} \) is obtained from \( u \) by removing the \( i \)th
component of $u$.

4.1. Correlation lengths and derivative-based global sensitivity measures. Consider a function $f$ to be observed at the locations $x^{(1)}, \ldots, x^{(n)} \in \Omega$, with $n \in \mathbb{N}$ and for a bounded domain $\Omega \subset \mathbb{R}^D$. Let $X$ be the $n \times D$ matrix with lines given by $x^{(1)}, \ldots, x^{(n)}$, $y$ be the vector of responses $y = (f(x^{(1)}), \ldots, f(x^{(n)}))^\top$ and $Z$ the $(n + 1) \times D$ matrix $Z = (X \ y)$.

Recall that the prediction of $f$ at any line vector $x \in \Omega$, from the GP model, is given by $m_{\theta,Z}(x) = r_\theta(x)^\top K_\theta^{-1} y$, with $r_\theta(x) = k(x, X)$, $K_\theta = k_\theta(X, X)$. Then, we use the notation $\vartheta_i(\theta)$ for the DGSM index of the variable $x_i$ on the predictor function $m_{\theta,Z}(x)$:

$$\vartheta_i(\theta) = \vartheta_i(m_{\theta,Z}) = \int_\Omega \left( \frac{\partial m_{\theta,Z}(x)}{\partial x_i} \right)^2 \, dx.$$\smallskip

We also use the following notation for the normalized DGSM index of the variable $x_i$:

$$\tilde{\vartheta}_i(\theta) = \tilde{\vartheta}_i(m_{\theta,Z}) = \frac{\vartheta_i(\theta)}{\sum_{r=1}^D \vartheta_r(\theta)}.$$\smallskip

The normalized DGSM index $\tilde{\vartheta}_i(\theta)$ satisfies $0 \leq \tilde{\vartheta}_i(\theta) \leq 1$. The larger this indice is, the more important the variable $x_i$ is for $m_{\theta,Z}(x)$. In the two next propositions, we show that, under mild conditions, we have $\tilde{\vartheta}_i(\theta) \to 1$ as $\theta_i \to 0$ and $\tilde{\vartheta}_i(\theta) \to 0$ as $\theta_i \to \infty$. Hence, we give a theoretical support to the intuition that small correlation lengths correspond to important input variables.

**Proposition 4.1.** Assume that the components of $y$ are not all equal. Assume that $k$ is continuously differentiable on $\mathbb{R}$. Let $i \in \{1, \ldots, D\}$ be fixed. For $j = 1, \ldots, n$ let $v^{(j)} = x^{(j)}_i$. Assume that $v^{(1)}, \ldots, v^{(n)}$ are two by two distinct. Then, for fixed $\theta_{-i} \in (0, \infty)^{D-1}$

$$\tilde{\vartheta}_i(\theta) \xrightarrow{\theta_i \to \infty} 0.$$\smallskip

**Proposition 4.2.** Assume that the components of $y$ are not all equal. Consider the same notation as in Proposition 4.1. Assume that $k$ is continuously differentiable on $\mathbb{R}$, that $k(t) \to 0$ as $|t| \to \infty$ and that $\Omega$ is an open set. Assume also that $x^{(1)}, \ldots, x^{(n)}$ are two-by-two distinct. Let $i \in \{1, \ldots, D\}$ be fixed. Then for fixed $\theta_{-i} \in (0, \infty)^{D-1}$

$$\tilde{\vartheta}_i(\theta) \xrightarrow{\theta_i \to 0} 1.$$\smallskip

In Propositions 4.1 and 4.2, the regularity conditions on $k$ are mild, and the conditions on $x^{(1)}, \ldots, x^{(n)}$ hold in many cases, for instance when $x^{(1)}, \ldots, x^{(n)}$ are selected randomly and independently or from a latin hypercube procedure (see e.g. [26]).

4.2. Estimated correlation lengths and inactive variables. We first recall the likelihood function:

$$l_Z(\theta) = \frac{1}{(2\pi)^{n/2}} \frac{1}{\sqrt{\det(k_\theta(X,X))}} \exp \left( -y^\top k_\theta(X,X)^{-1} y \right).$$
Proposition 4.3 holds in particular when \( k \phi \) has a Fourier transform \( \hat{\phi} \) as in [33] and from the fact that an infinitely differentiable function with compact support \( \|\cdot\| \to \infty \). Hence, Proposition 4.3 holds. This follows from Theorem 10.12 in [33] and from the fact that an infinitely differentiable function with compact support \( \phi \) has a Fourier transform \( \hat{\phi}(w) \) satisfying \( \hat{\phi}(w)\|w\|^\gamma \to 0 \) as \( \|w\| \to \infty \) for any \( \gamma < \infty \). Hence, Proposition 4.3 holds in particular when \( k \) is the exponential covariance function with \( k(t) = e^{-|t|} \). Proposition 4.3 also holds when \( k \) is the Matérn covariance function with

\[
    k(t) = \frac{1}{\Gamma(\nu)} 2^{\nu-1} (2\sqrt{\nu}|t|)^\nu K_\nu(2\sqrt{\nu}|t|),
\]

where \( 0 < \nu < \infty \) is the smoothness parameter (see e.g. [30]). It should however be noted that the squared exponential covariance function \( k \) (defined by \( k(t) = \exp(-t^2) \) with \( t \in \mathbb{R} \)) does not satisfy the condition of Lemma 4.3. [Notice that [34] study specifically the asymptotic behavior of the maximum likelihood estimation of a variance parameter for this covariance function, when the number of observations of a smooth function goes to infinity.]

In the next proposition, we study the LOO mean square prediction error

\[
    CV_Z(\theta) = \sum_{j=1}^n (y_j - \hat{y}_{\theta,j})^2,
\]
with 
\[ \hat{y}_{\theta,j} = k_{\theta}(x^{(j)}, X_{-j})k_{\theta}(X_{-j}, X_{-j})^{-1}y_{-j}, \]
where \( X_{-j} \) and \( y_{-j} \) are obtained, respectively, by removing the line \( j \) of \( X \) and the component \( j \) of \( y \). We show that, as for the likelihood, inactive variables can be detected by this LOO criterion, since we can have \( CV_{Z}(\theta) \to 0 \) as \( \theta_{i} \to \infty \) if the function \( f \) does not depend on \( x_{i} \).

**Proposition 4.4.** Let \( k \) satisfy the same conditions as in Proposition 4.3. Let \( i \in \{1, \ldots, D\} \) be fixed.

For \( j = 1, \ldots, n \) let \( v^{(j)} = x_{-i}^{(j)} \). Assume that
i) \( x^{(1)}, \ldots, x^{(n)} \) are two-by-two distinct;
ii) \( y_{r} = y_{s} \) if \( v^{(r)} = v^{(s)} \);
iii) for all \( r \in \{1, \ldots, n\} \) there exists \( s \in \{1, \ldots, n\}, r \neq s, \) such that \( v^{(r)} = v^{(s)} \).

Then, for any fixed \( \theta_{-i} \in (0, \infty)^{D-1} \), we have
\[ CV_{Z}(\theta) \to 0. \]

In Proposition 4.4, Conditions i) and ii) are interpreted similarly as in Proposition 4.3. Condition iii), however, provides more restrictions than for the likelihood in Proposition 4.3. This condition states that for any observation point in the data set, there exists another observation point for which only the inactive input \( i \) is changed. This condition is arguably necessary to have \( CV_{Z}(\theta) \to 0 \).

**5. Numerical examples.**

**5.1. Test sets.** We illustrate the Split-and-Doubt algorithm on five benchmark optimization problems. The first four are classical synthetic functions: the two-dimensional Branin function, the general Ackley function in six dimensions, the six-dimensional Hartmann function and the general Rosenbrock function in five dimensions. The fifth one is the Borehole function [20]. It models the water-flow in a borehole. For each function, we added inactive input variables in order to embed them in a higher dimension \( D^{(i)} \). The settings are summarized in Table 1.

<table>
<thead>
<tr>
<th>( f^{(i)} )</th>
<th>( d^{(i)} )</th>
<th>( D^{(i)} )</th>
<th>( N^{(i)}_{\max} )</th>
<th>( n_{0}^{(i)} )</th>
</tr>
</thead>
<tbody>
<tr>
<td>Hartmann 6-dim</td>
<td>6</td>
<td>15</td>
<td>30</td>
<td>30</td>
</tr>
<tr>
<td>Rosenbrock</td>
<td>5</td>
<td>20</td>
<td>40</td>
<td>60</td>
</tr>
<tr>
<td>Ackley</td>
<td>6</td>
<td>20</td>
<td>45</td>
<td>40</td>
</tr>
<tr>
<td>Borehole</td>
<td>6</td>
<td>25</td>
<td>30</td>
<td>25</td>
</tr>
<tr>
<td>Branin</td>
<td>2</td>
<td>25</td>
<td>30</td>
<td>50</td>
</tr>
</tbody>
</table>

We launched the optimization process for these functions with three different optimization algorithms:

• **Split-and-Doubt** algorithm with Matérn $5/2$ covariance function.

• **Split-without-Doubt** algorithm: It uses the same variable splitting as **Split-and-Doubt** and generates the minor variables by uniform random sampling.

For each function $f^{(i)}$, we launched each optimization process for $N_{\text{max}}^{(i)}$ iterations starting with $N_{\text{seed}} = 20$ different initial Designs Of Experiments (DOE) of size $n_0^{(i)}$ generated by a maximin space-filling sampling.

### 5.2. Results.

The results are represented by boxplots in Appendix A. We also display the mean best value evolution in Figure 4.

We can see that **Split-and-Doubt** gives better results than EGO for Rosenbrock, Ackley and Borehole function. EGO does not converge for the first two functions and used more iterations for the last one. These cases illustrate the efficiency of the dimension reduction for limited budget optimization. For Branin function the convergence is relatively fast for all the three algorithms. This is due to the fact that the effective dimension is 2 and that the first design of experiments covers well these dimensions.

On one hand, sampling the minor variables at random or using the Doubt/Contrast strategy gives close results when the influential variables are easily determined. On the other hand, the efficiency of the Doubt/Contrast approach is visible on Hartmann and Ackley functions, by comparing the results of **Split-and-Doubt** and **Split-without-Doubt**. Notice that we start in general with a relatively small amount of design points. Thus, the initial estimation of the correlation lengths can be inaccurate. In these cases, the Doubt/Contrast approach is valuable to improve the estimation. To further highlight this idea, we display in Figure 5 the percentage of undetected influential variables and the miss-classification rate of all the variables for both **Split-and-Doubt** and **Split-without-Doubt** for the Rosenbrock function.

Among the 20 DOE, the **Split-and-Doubt** detects all the major variables for 19 cases starting from iteration 15 and for all the DOE starting from iteration 37. However, the **Split-without-Doubt** struggles to select properly all the influential variables even in the last iterations. Considering all the variables, the miss-classification rate decrease rapidly for the **Split-and-Doubt**. However, for one test a minor variable is considered influential until the end. This can be explained by the nature of the doubt function that aims at correcting only a miss-classification of an influential variable.

Finally, as we can see in Figure 6, **Split-and-Doubt** is faster than EGO in terms of computing time. The fact that we perform two optimization procedures in smaller spaces makes the algorithm faster than optimizing the EI in dimension $D$.

### 6. Proofs.

For the proofs of Propositions 4.1 and 4.2, we let $k'(t) = \partial k(t)/\partial t$.

**Proof of Proposition 4.1.** Without loss of generality, we consider $i = 1$ in the proof. Let $\theta_{-1} \in (0, \infty)^{D-1}$ be fixed. We have

$$\frac{\partial}{\partial x_j} m_{\theta,Z}(x) = \left( \frac{\partial r_{\theta}(x)}{\partial x_j} \right)^T K_{\theta}^{-1} y.$$ 

When $\theta_1 \rightarrow \infty$, $K_{\theta}$ converges to the $n \times n$ matrix $L_{\theta_{-1}}$ with $(L_{\theta_{-1}})_{pq} = \prod_{r=2}^{D} k([x_{r}^{(p)} - x_{r}^{(q)})/\theta_r]$, by continuity of $k$. This matrix is invertible by assumption on $k$ and $v^{(1)}, ..., v^{(n)}$. Hence
Figure 4. Comparison of 3 optimization strategies. Mean over $N_{seed}$ of the best current values as a function of the number of iterations.
(a) Miss-classification rate of major variables  (b) Miss-classification rate of all the variables

Figure 5. Solid line: mean value over 20 repetitions, colored area: 95% confidence interval. Blue: Split-and-Doubt, Red: Split-without-Doubt. x-axis, iteration number.

|\|K^{-1}\|_\Omega is bounded as \theta_1 \to \infty. We have for \( j = 1, \ldots, n \)

\[
\left( \frac{\partial r_\theta(x)}{\partial x_1} \right)_j = \frac{1}{\theta_1} k'(\frac{[x_1 - x_1^{(j)}]}{\theta_1}) \prod_{p=2}^{D} k(\frac{[x_p - x_p^{(j)}]}{\theta_p}).
\]

Observe that \( k \) is continuously differentiable and that \( \Omega \) is bounded. Hence by uniform
continuity as $\theta_1 \to \infty$

$$\sup_{x \in \Omega} \left| \frac{\partial r_\theta(x)}{\partial x_1} \right| \to 0.$$  

Hence, $\vartheta_1(\theta) \to 0$ as $\theta_1 \to \infty$. Let now for $x = (u,v)$ with $u \in \mathbb{R}$, $l_{\theta_1}(x)$ be the $n \times 1$ vector defined by $[l_{\theta_1}(x)]_j = \prod_{r=1}^{d-1} k([v_r - v_r^{(j)})/\theta_{r+1})$ (we recall that for $j = 1, ..., n$, $v^{(j)} = v^{(j)}_1$). Let $\hat{g}_{\theta}(x) = l_{\theta_2}(v)L_{\theta_1}^{-1}y$. Then for $m = 2, ..., D$, by the triangle and Cauchy-Schwarz inequalities

$$\left| \frac{\partial m_{\theta_2}(x)}{\partial x_m} - \frac{\partial \hat{g}_{\theta_1}(x)}{\partial x_m} \right| \leq \left| \frac{\partial r_\theta(x)}{\partial x_m} - \frac{\partial \hat{g}_{\theta_1}(x)}{\partial x_m} \right| \cdot \left| K_\theta^{-1}y \right| + \left| \frac{\partial l_{\theta_1}(x)}{\partial x_m} \right| \cdot \left| K_\theta^{-1}y - L_\theta^{-1}y \right|.$$  

In (6.1), the vector in the first norm has component $r \in \{1, ..., n\}$ equal to

$$(k((u - u_r)/\theta_1)) - 1 \frac{1}{\theta_2} k'(v_{m-1} - v_{m-1})/\theta_m) \prod_{p=2}^{D} k([v_{p-1} - v_{p-1})/\theta_p)$$

which goes to 0 as $\theta_1 \to \infty$, uniformly over $x \in \Omega$, by uniform continuity. The second norm in (6.1) is bounded as discussed above. The third norm in (6.1) does not depend on $\theta_1$ and is thus bounded uniformly over $x \in \Omega$ as $\theta_1 \to \infty$. The fourth norm in (6.1) goes to 0 as $\theta_1 \to \infty$ as discussed above.

Hence, uniformly over $x \in \Omega$,

$$\left| \frac{\partial m_{\theta_2}(x)}{\partial x_m} - \frac{\partial \hat{g}_{\theta_1}(x)}{\partial x_m} \right| \to 0, \quad \theta_1 \to \infty.$$  

Furthermore, the function $\hat{g}_{\theta_1}$ is continuously differentiable and non-constant on $\Omega$ because $\hat{g}_{\theta_1}(x^{(r)}) = y_r$ for $r = 1, ..., n$ and because the components of $y$ are not all equal. This implies that

$$\liminf_{\theta_1 \to \infty} \sum_{m=2}^{D} \vartheta_m(\theta) > 0,$$

which concludes the proof.

**Proof of Proposition 4.2.** As before, we consider $i = 1$ in the proof. We have for $m = 2, ..., D$ and $r = 1, ..., n$

$$\left( \frac{\partial r(x)}{\partial x_m} \right)_r = k([x_1 - x_1^{(r)}]/\theta_1) \frac{1}{\theta_2} k'(x_{m-1} - x_{m-1})/\theta_m \prod_{j=2}^{D} k([x_j - x_j^{(r)}]/\theta_j).$$

Hence, $||\partial r(x)/\partial x_m||$ is bounded as $\theta_1 \to 0^+$ uniformly in $x \in \Omega$ from the assumptions on $k$.

For $j = 1, ..., n$, let $u_j$ be the first component of $x^{(j)}$ and let $v^{(j)} = x^{(j)}_1$. As $\theta_1 \to 0^+$, the matrix $K_{\theta}$ converges to the $n \times n$ matrix $N_{\theta_1} = [1_{u_i = u_i} (L_{\theta_1})_{pi}]_{p,q=1,...,n}$ with the notation

\[1_{u_i = u_i} (L_{\theta_1})_{pi} \]
of the proof of Proposition 4.1. The matrix $N_{\theta-1}$ is invertible because its submatrices are invertible. This is so because for any $p = 1, \ldots, n$ the subset $\{v^{(q)}; q = 1, \ldots, n, u_q = u_p\}$ is composed of two-by-two distinct elements since $x^{(1)}, \ldots, x^{(n)}$ are two-by-two distinct.

Hence, $||K_{\theta}^{-1}y||$ is bounded as $\theta_1 \to 0^+$ and so $\sum_{m=2}^{D} \theta_m(\theta)$ is bounded as $\theta_1 \to 0^+$.

Let now $j \in \{1, \ldots, n\}$ for which $y_j \neq 0$. Let $\delta > 0$, not depending on $\theta_1$, be small enough so that $\prod_{r=1}^{D} [x_r^{(j)} - \delta, x_r^{(j)} + \delta] \subseteq \Omega$. Then we have

(6.2) $\sup_{s \in [-\delta, \delta] : |s| = \sqrt{\theta_1}} |m_{\theta,Z}(x^{(j)} + s)| \to 0 \quad \text{as} \quad \theta_1 \to 0^+$.

Indeed, we have

$$\left( r_{\theta}(x^{(j)} + s) \right)_p = k \left( \frac{u_p - u_j - s_1}{\theta_1} \right) \prod_{r=2}^{D} k \left( \frac{x_r^{(p)} - x_r^{(j)} - s_r}{\theta_r} \right).$$

The product above is bounded uniformly over $s \in [-\delta, \delta]^D$ by uniform continuity of $k$. Also, whether $u_p - u_j = 0$ or $u_p - u_j \neq 0$, we have

$$\sup_{|s_1| = \sqrt{\theta_1}} k \left( \frac{u_p - u_j - s_1}{\theta_1} \right) \to 0 \quad \text{as} \quad \theta_1 \to 0^+.$$

Finally, $||K_{\theta}^{-1}y||$ is bounded as $\theta_1 \to 0^+$ as discussed above. Hence (6.2) is proved. Also, let $E = \{u_j\} \times \prod_{r=2}^{D} [x_r^{(j)} - \delta, x_r^{(j)} + \delta]$. Then as $\theta_1 \to 0^+$, uniformly over $x \in E$, for $p = 1, \ldots, n$, we have

$$\left( r_{\theta}(x) \right)_p \to 1_{\{u_p = u_j\}} \prod_{r=2}^{D} k \left( \frac{x_r - (x_p)_r}{\theta_r} \right).$$

Also $K_{\theta}^{-1}y \to N_{\theta-1}y$ as discussed above. Hence, as $\theta_1 \to 0^+$, $m_{\theta,Z}(x)$ converges uniformly over $x \in E$ to a function value $\hat{g}_{\theta-1}(x)$, with $\hat{g}_{\theta-1}(x)$ continuous with respect to $x \in E$. Since $m_{\theta,Z}(x^{(j)}) = y_j$, we can choose the $\delta > 0$ (still independently of $\theta_1$) so that it also satisfies

(6.3) $\liminf_{\theta_1 \to 0^+} \inf_{x \in E} |m_{\theta,Z}(x)| \geq \frac{|y_j|}{2}.$
We have
\[
\int_{\Omega} \left( \frac{\partial m_{\theta,Z}(x)}{\partial x_1} \right)^2 dx \geq \int_{\Pi^2_{r=1}[x(i)^j - \delta x(i)^j + \delta]} \left( \frac{\partial m_{\theta,Z}(x)}{\partial x_1} \right)^2 dx
\]
\[
= \int_{\Pi^2_{r=1}[x(i)^j - \delta x(i)^j + \delta]} dx - \int_{x(i)^j} dx_1 \left( \frac{\partial m_{\theta,Z}(x)}{\partial x_1} \right)^2
\]
\[
\geq \int_{\Pi^2_{r=1}[x(i)^j - \delta x(i)^j + \delta]} dx - \int_{x(i)^j} dx_1 \left( \frac{\partial m_{\theta,Z}(x)}{\partial x_1} \right)^2
\]
\[
(\text{Jensen:}) \geq \int_{\Pi^2_{r=1}[x(i)^j - \delta x(i)^j + \delta]} dx \sqrt{\frac{1}{\theta_1}} \left( \frac{1}{\sqrt{\theta_1}} \int_{\Pi^2_{r=1}[x(i)^j - \delta x(i)^j + \delta]} dx_1 \partial m_{\theta,Z}(x) \right)^2
\]
\[
\geq (2\delta)^{D-1} \frac{1}{\sqrt{\theta_1}} \left( \inf_{x \in E} |m_{\theta,Z}(x)| - \sup_{s \in [-\delta,\delta]} |m_{\theta,Z}(x(s) + s)| \right)^2
\]
\[
\rightarrow_{\theta_1 \rightarrow 0^+} \infty,
\]
from (6.2) and (6.3). This concludes the proof.

**Proof of Proposition 4.3.** Without loss of generality, we consider \( i = 1 \) in the proof. Let us consider the \( 2 \times 2 \) submatrix of \( k_{\theta}(X, X) \) obtained by extracting the lines and columns \( a, b \), with \( a, b \) as in Condition iii) of the lemma. Then as \( \theta_1 \rightarrow \infty \) this submatrix converges to the singular matrix \( ((1, 1)^T, (1, 1)^T) \). Hence, we have, as \( \theta_1 \rightarrow \infty \), \( |k_{\theta}(X, X)| \rightarrow 0 \) (since \( k_{\theta}(X, X) \) has components bounded in absolute value by 1). Hence, it is sufficient to show that \( y^T k_{\theta}(X, X)^{-1} y \) is bounded in order to conclude the proof.

Let \( X_{\theta_1} \) be obtained from \( X \) by dividing its first column by \( \theta_1 \) and by leaving the other columns unchanged. Let \( x^{(\theta_1, j)} \) be the transpose of the line \( j \) of \( X_{\theta_1} \), for \( j = 1, ..., n \). Let \( \bar{\theta} = (1, \theta_1) \). Then, \( y^T k_{\theta}(X, X)^{-1} y = y^T k_{\bar{\theta}}(X_{\theta_1}, X_{\theta_1})^{-1} y \).

We now use tools from the theory of RKHSs and refer to, e.g., [33] for the definitions and properties of RKHSs used in the rest of the proof. Let \( \mathcal{H} \) be the RKHS of \( k_{\bar{\theta}} \). Let \( \alpha^{(\theta_1)} = k_{\bar{\theta}}(X_{\theta_1}, X_{\bar{\theta}})^{-1} y \). Then, \( f_{\bar{\theta}_1} : \mathbb{R}^D \rightarrow \mathbb{R} \) defined by \( f_{\bar{\theta}_1}(x) = \sum_{j=1}^n \alpha^{(\theta_1)}(k_{\bar{\theta}}(x - x^{(\theta_1, j)})) \) is the function of \( \mathcal{H} \) with minimal RKHS norm \( ||.||_{\mathcal{H}} \) satisfying \( f_{\bar{\theta}_1}(x^{(\theta_1, j)}) = y_j \) for \( j = 1, ..., n \).

As \( \theta_1 \rightarrow \infty \), the points \( x^{(\theta_1, 1)}, ..., x^{(\theta_1, n)} \) converge to the points \( w^{(1)}, ..., w^{(n)} \) with \( w^{(i)} = \{0, [u(i)]^T\} \). We observe that, by assumption, \( y_r = y_s \) for \( u^{(r)} = u^{(s)} \). Hence, there exists \( \epsilon > 0 \) small enough and \( p \) column vectors \( c^{(1)}, ..., c^{(p)} \) in \( \mathbb{R}^D \) with the following properties: (i) each Euclidean ball with center \( c^{(m)} \), \( m = 1, ..., p \), and radius \( 2\epsilon \) does not contain two \( u^{(r)}, u^{(s)} \) with \( y_r \neq y_s \), \( r, s \in \{1, ..., n\} \); (ii) each \( w_j \), \( j = 1, ..., n \), is contained in a ball with center \( c^{(m)} \) with \( m \in \{1, ..., p\} \) and radius \( \epsilon \); (iii) the \( p \) balls with centers \( c^{(1)}, ..., c^{(p)} \) and radii \( 2\epsilon \) are two-by-two non-intersecting. We can also assume that each ball with center \( c^{(m)} \), \( m = 1, ..., p \) and radius \( \epsilon \) contains at least one \( u^{(j(m))} \) with \( j(m) \in \{1, ..., n\} \) and we write \( z_m = u^{(j(m))} \).

Then, from Lemma 6.2, there exists an infinitely differentiable function \( g \) with compact support on \( \mathbb{R}^d \) so that for \( m = 1, ..., p \), \( g(x) = z_m \) for \( ||x - c^{(m)}|| \leq 2\epsilon \). Hence, for \( \theta_1 \) large enough, the function \( g \) satisfies \( g(x^{(\theta_1, j)}) = y_j \) for \( j = 1, ..., n \).
Hence, \( \|f_{\theta_1}\|_H \leq \|g\|_H \) for \( \theta_1 \) large enough, where \( \|g\|_H \) does not depend on \( \theta_1 \). Finally, a simple manipulation of \( \|\cdot\|_H \) (see again [33] for the definitions), provides
\[
\|f_{\theta_1}\|_H = \sum_{r,s=1}^n \alpha_r(\theta_1)\alpha_s(\theta_1)k_\theta(x_r^{(\theta_1)} - x_s^{(\theta_1)})
= y^\top k_\theta(X,X)^{-1}k_\theta(X,X)k_\theta(X,X)^{-1}y
= y^\top k_\theta(X,X)^{-1}y.
\]
This concludes the proof.

**Proof of Proposition 4.4.** Without loss of generality, we consider \( i = 1 \) in the proof. Also, up to renumbering the lines of \( X \) and components of \( y \), it is sufficient to show that, for fixed \( \theta_1 \in (0, \infty)^D \), as \( \theta_1 \to \infty \), \( \hat{y}_{\theta_1,n} \to y_n \). We use the same notation \( \hat{\theta} \), \( \mathcal{H} \) and \( x^{(\theta_1,j)} \) as in the proof of Proposition 4.3. Then, we have \( \hat{y}_{\theta_1,n} = f_{\theta_1}(x^{(\theta_1,n)}) \), where \( f_{\theta_1} \in \mathcal{H} \) is the function with minimal norm \( \|f\|_H \) satisfying \( f_{\theta_1}(x^{(\theta_1,j)}) = y_j \) for \( j = 1, \ldots, n-1 \).

Furthermore, from the proof of Proposition 4.3, there exists a function \( g \in \mathcal{H} \), not depending on \( \theta_1 \), satisfying, for \( \theta_1 \) large enough, \( g(x^{(\theta_1,j)}) = y_j \) for \( j = 1, \ldots, n \). This shows that \( \|f_{\theta_1}\|_H \) is bounded as \( \theta_1 \to \infty \). Let \( m \in \{1, \ldots, n-1\} \) be so that \( v^{(m)} = v^{(n)} \) (the existence is assumed in Condition iii)). Let also, for \( x \in \mathbb{R}^D \), \( k_{\theta,x} \in \mathcal{H} \) be the function \( k_\theta(x - \cdot) \). Then we have (see again [33]), with \( (\cdot, \cdot)_\mathcal{H} \) the inner product in \( \mathcal{H} \)
\[
|\hat{y}_n - y_n| = \left| f_{\theta_1}(x^{(\theta_1,n)}) - f_{\theta_1}(x^{(\theta_1,m)}) \right|
= \left| (f_{\theta_1}|k_{\theta,x}^{(\theta_1,n)})_\mathcal{H} - (f_{\theta_1}|k_{\theta,x}^{(\theta_1,m)})_\mathcal{H} \right|
\leq \|f_{\theta_1}\|_H \|k_{\theta,x}^{(\theta_1,n)} - k_{\theta,x}^{(\theta_1,m)}\|_H
= \|f_{\theta_1}\|_H \sqrt{k_\theta(x^{(\theta_1,n)} - x^{(\theta_1,m)}) + k_\theta(x^{(\theta_1,m)} - x^{(\theta_1,m)}) - 2k_\theta(x^{(\theta_1,n)} - x^{(\theta_1,m)})}.
\]
In the above display, the square root goes to zero as \( \theta_1 \to \infty \) because \( x^{(\theta_1,n)} - x^{(\theta_1,m)} \) goes to zero and \( k_\theta \) is continuous. This concludes the proof.

**Lemma 6.1.** For any \( 0 < \epsilon_1 < \epsilon_2 < \infty \), there exists an infinitely differentiable function \( g : \mathbb{R} \to \mathbb{R} \) satisfying \( g(u) = 1 \) for \( |u| \leq \epsilon_1 \) and \( g(u) = 0 \) for \( |u| \geq \epsilon_2 \).

**Proof.** Let \( h : \mathbb{R} \to \mathbb{R} \) be defined by \( h(t) = \exp(-1/(1 - t^2)) \mathbb{1}\{t \in [-1,1]\} \). Then \( h \) is infinitely differentiable. Hence, \( g \) can be chosen of the form
\[
g(t) = \begin{cases} A \int_{-\infty}^t h(B[u + \epsilon_1/2]) \, du & \text{if } t \leq 0 \\
A \int_t^\infty h(B[u - \epsilon_2/2]) \, du & \text{if } t \geq 0 \end{cases},
\]
with \( 2/(\epsilon_2 - \epsilon_1) < B < \infty \) and \( A = B/(\int_{-\infty}^\infty h(u)du) \). It can be checked that \( g \) is infinitely differentiable and satisfies the conditions of the lemma.

**Lemma 6.2.** Let \( d, p \in \mathbb{N} \). Let \( x^{(1)}, \ldots, x^{(p)} \) be two-by-two distinct points in \( \mathbb{R}^D \) and \( \epsilon > 0 \) be so that the \( p \) closed Euclidean balls with centers \( x^{(i)} \) and radii \( \epsilon \) are disjoint. Let \( y_1, \ldots, y_p \in \mathbb{R} \) be arbitrary. Then there exists an infinitely differentiable function \( r : \mathbb{R}^D \to \mathbb{R} \), with compact support, satisfying for \( i = 1, \ldots, p, g(u) = y_i \) when \( |u - x^{(i)}| \leq \epsilon \).
Proof. Let \( l = \min_{i \neq j} ||x^{(i)} - x^{(j)}|| \) and observe that \( \epsilon < 2l \). Let \( g \) satisfies Lemma 6.1 with \( \epsilon_1 = \epsilon^2 \) and \( \epsilon_2 = l^2 / 4 \). Then the function \( r \) defined by \( r(u) = \sum_{i=1}^{p} y_{i} g(||u - x^{(i)}||^2) \) satisfies the conditions of the lemma.

7. Conclusion. Performing Bayesian optimization in high dimension is a difficult task. In many real-life problems, some variables are not influential. Therefore, we propose the so-called Split-and-Doubt algorithm that performs sequentially both dimension reduction and feature oriented sampling. The “split” step (model reduction) is based on a property of stationary ARD kernel of Gaussian process regression. We proved that large correlation lengths correspond to inactive variables. We also showed that classical estimators such ML and CV may assign large correlation lengths to inactive variables.

The “doubt” step questions the “split” step and helps correcting the estimation of the correlation lengths. It is possible to use this strategy for different feature learning purposes such as refinement, optimization and inversion. The optimization Split-and-Doubt algorithm has been evaluated on classical benchmark functions embedded in larger dimensional spaces by adding useless input variables. The results show that Split-and-Doubt is faster than classical EGO in the whole design space and outperforms it for most of the discussed tests.

The main limitation of Split-and-Doubt is that we perform correlation length estimation in high dimension. This computation is expensive. To overcome this problem, one can use fast maximum likelihood approximation techniques [12]. Future research may investigate such methods and extend Split-and-Doubt to constrained optimization.

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Appendix A. Optimization test results. In this section, we use boxplots to display the evolution of the best value of the optimization test bench. For each iteration, we display: Left: EGO in light blue, Middle: Split-and-Doubt in dark blue, Right: Split-without-Doubt in light green.
Figure 7. Borehole: Box plots convergence.

Figure 8. Rosenbrock: Box plots convergence.

REFERENCES


Figure 9. Ackley: Box plots convergence.

Figure 10. Hartmann 6-dim: Box plots convergence.


Figure 11. Branin: Box plots convergence.


