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Analysis of Parameter Variability in an Integrated Wireless Power Transfer System via Partial Least Squares Regression

Mourad Larbi, Riccardo Trincherio, *Member, IEEE*, Flavio G. Canavero, *Fellow, IEEE*, Philippe Besnier, *Senior Member, IEEE*, Madhavan Swaminathan, *Fellow, IEEE*

Abstract—This paper deals with the application of the partial least squares (PLS) regression to the uncertainty quantification of an integrated wireless power transfer with 30 random variables. It considers the development of surrogate models using a limited set of training samples in order to estimate statistical quantities of the converter efficiency with a relatively low computational cost compared to the standard brute force Monte Carlo (MC) simulation. The strength, the performance and the features of the proposed modeling approach are then compared with the ones of an advanced probabilistic surrogate model combining the Least-squares Support Vector Machine (LS-SVM) and the Gaussian Process (GP) regression. The MC simulation is considered as reference.

Index Terms—Machine learning, uncertainty quantification, surrogate model, Gaussian Process regression, Least-squares Support Vector Machine regression, PLS regression, sensitivity analysis, probabilistic model, wireless power transfer (WPT).

I. INTRODUCTION

The level of complexity of electronic systems is so high nowadays that it becomes a difficult task to guarantee their reliability. Due to the fabrication process, the variability of parameters associated with the system has to be considered, since it may generate large variations of output signals and lead to a potential failure of systems. In the above scenario, uncertainty quantification (UQ) plays a key role since it allows predicting the effect of the uncertain parameters on the system performance. In advanced electronic systems, the complexity and the simulation costs are so high that the brute force Monte Carlo (MC) approach turns out to be inefficient.

Alternatively, various surrogate models have been successfully proposed. One of the most used techniques is the polynomial chaos (PC), which has shown good performance to solve different applications [1]–[3]. Despite its accuracy and its improved efficiency with respect to the plain MC simulation

in terms of simulation costs, such technique suffers the so-called *curse of dimensionality*, since the number of unknown coefficients to be estimated during the training of the PC-surrogate grows exponentially with respect to the number of uncertain parameters and the expansion order. To this aim several advanced implementations of the PC expansion such as tensor and sparse formulations have been presented to bypass the mentioned problem [4]–[7].

Recently, alternative solutions based on Machine Learning (ML) regressions have been investigated in order to mitigate the above issues. In particular, non-parametric regressions (i.e., techniques in which the regression complexity is independent from the dimensionality of the parameter space) such as support vector machine (SVM) and least-squares support vector machine (LS-SVM) regression [8]–[11], Gaussian process regression (GPR) [12]–[14] and their combinations (e.g., the LS-SVM+GPR surrogate model) [15]–[17] can be considered as valuable alternative techniques to the state-of-the-art approaches for the UQ in complex systems.

This paper presents an alternative technique for the UQ in high-dimensional problems [18]–[20], such as the partial least squares (PLS) regression [21]. The PLS regression allows to build a surrogate model of an output depending on many uncertain parameters. Similarly to the principal component analysis (PCA) [22], PLS allows to reduce the problem dimensionality using a limited number of components while mapping the relationship between input and output variables. This property of the method is very interesting as it captures the most prominent input variables of the systems. Moreover, the surrogate model can then be used to compute statistical moments and the probability density function (PDF) of the output of interest. The accuracy and the strength of the proposed modeling scheme are then investigated by comparing the results of the surrogate model built via the PLS and the ones provided by a probabilistic model based on the LS-SVM+GPR based on the results of a MC simulations. The above techniques are applied for the UQ of the efficiency of a wireless power transfer (WPT) device as a function of 30 Gaussian uncertain parameters.

This paper is structured as follows: Section II briefly presents the mathematical background of the LS-SVM+GPR regression. Section III describes the PLS regression method as well as the sensitivity analysis derived from the method. Section IV presents the numerical results and compares the performance of the PLS surrogate with the ones provided by a

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LS-SVM+GPR surrogate model by considering the WPT test-case. Section V provides additional information on the features provided by each of the considered methods. Section VI concludes the paper.

II. PROBABILISTIC MODEL COMBINING LS-SVM AND GPR

Let us consider the problem of approximating a given set of L training pairs $\{(\mathbf{x}_i, y_i)\}_{i=1}^L$ provided by a full computational model such that $y_i = \mathcal{M}(\mathbf{x}_i)$, as a function of its stochastic parameter $\mathbf{x}_i \in \mathcal{P}$ with $\mathcal{P} \subset \mathbb{R}^d$, in terms of the following GPR [13]:

$$\mathcal{M}(\mathbf{x}) \approx \mu(\mathbf{x}) + \varepsilon(\mathbf{x}) \quad (1)$$

where $\mu(\mathbf{x})$ is a deterministic function representing the mean estimated via a generic regression, also called *trend* and $\varepsilon(\mathbf{x}) \sim \mathcal{GP}(0, k(\mathbf{x}, \mathbf{x}'))$ is a Gaussian process (GP) with zero mean and covariance $k(\mathbf{x}, \mathbf{x}')$ approximating the deviation of the deterministic model from the actual function \mathcal{M} , for any configuration of the parameters \mathbf{x} . A GP can be considered as an extension of the concept of multivariate Gaussian distributions to infinite dimensionality which returns, for any arbitrary value of \mathbf{x}_* , a prediction of the mean value $\mu_{\mathbf{x}_*}$ and the variance $\sigma_{\mathbf{x}_*}^2$ of a normal distribution.

The covariance function $k(\mathbf{x}, \mathbf{x}')$ of the GP specifies the correlation between pairs of random variables under the assumption that \mathcal{M} is *smooth*, this means that points with similar predictor values are expected to have close response values. The covariance function can be defined by several kernels, however, hereafter in this paper we will consider the Matern 3/2 covariance function [13]:

$$k(\mathbf{x}, \mathbf{x}') = \sigma_f^2 \left(1 + \frac{\sqrt{3}\|\mathbf{x} - \mathbf{x}'\|}{\sigma_l} \right) \exp\left(-\frac{\sqrt{3}\|\mathbf{x} - \mathbf{x}'\|}{\sigma_l} \right) \quad (2)$$

where $\|\cdot\|$ represents the Euclidean distance and σ_f and σ_l are the hyper-parameters of the covariance estimated via an optimization procedure, e.g., by minimizing the negative log marginal likelihood of the GP.

Knowing the prior information provided by the training pairs $\{(\mathbf{x}_i, \tilde{y}_i)\}_{i=1}^L$, with $\tilde{y}_i = y_i - \mu(\mathbf{x}_i)$, for any new value of the input parameter $\mathbf{x}_* \in \mathcal{P}$ for $\mathbf{x}_* \neq \mathbf{x}_i$, the set of samples $\{\tilde{y}_1, \dots, \tilde{y}_L, \tilde{y}_*\}$ follows an $L + 1$ -dimensional joint Gaussian distribution [17], which writes:

$$\begin{bmatrix} \tilde{\mathbf{y}} \\ \tilde{y}_* \end{bmatrix} \sim \mathcal{N}\left(\begin{bmatrix} \mathbf{0} \\ 0 \end{bmatrix}, \begin{bmatrix} \mathbf{K} & \mathbf{k}_*^t \\ \mathbf{k}_* & k_{**} \end{bmatrix} \right), \quad (3)$$

where $\tilde{\mathbf{y}} = [\tilde{y}_1, \dots, \tilde{y}_L]^t$, $\mathbf{K} \in \mathbb{R}^{L \times L}$ is the correlation matrix in which each ij -element is given by $K_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$, $\mathbf{k}_* = [k(\mathbf{x}_*, \mathbf{x}_1), \dots, k(\mathbf{x}_*, \mathbf{x}_L)] \in \mathbb{R}^{1 \times L}$ and $k_{**} = k(\mathbf{x}_*, \mathbf{x}_*)$.

Therefore, for a generic configuration of the parameters $\mathbf{x}_* \in \mathcal{P}$, the probabilistic model in (1) provides as output a Gaussian distributed stochastic variable $M(\mathbf{x}_*) \sim \mathcal{N}(\mu_{\mathbf{x}_*}, \sigma_{\mathbf{x}_*}^2)$ defined by the posterior mean $\mu_{\mathbf{x}_*}$ and variance $\sigma_{\mathbf{x}_*}^2$:

$$\mu_{\mathbf{x}_*} = \mu(\mathbf{x}_*) + \mathbf{k}_* \mathbf{K}^{-1} \tilde{\mathbf{y}} \quad (4a)$$

$$\sigma_{\mathbf{x}_*}^2 = k_{**} - \mathbf{k}_* \mathbf{K}^{-1} \mathbf{k}_*^t. \quad (4b)$$

The prediction mean $\mu_{\mathbf{x}_*}$ is used instead of the deterministic model $\mu(\mathbf{x})$, whereas the variance $\sigma_{\mathbf{x}_*}^2$ gives a local error indicator about its precision. Different from the standard GP regression, hereafter in this work, we focus on the case of a GP regression with a *fixed* mean function [17].

In the proposed LS-SVM+GPR regression, we are assuming that the trend $\mu(\mathbf{x})$ of the GP in (1) is a known function previously estimated via a LS-SVM regression [8]. The LS-SVM regression is considered since it allows building accurate and compact surrogate models of the response of a generic high-dimensional nonlinear function \mathcal{M} starting from a limited number of training samples and writes:

$$\mu(\mathbf{x}) = M_{LSVM}(\mathbf{x}) = \sum_{i=1}^L \alpha_i K(\mathbf{x}_i, \mathbf{x}) + b \quad (5)$$

where $\alpha_i \in \mathbb{R}$ are scalar coefficients, $K(\cdot, \cdot) : \mathbb{R}^d \rightarrow \mathbb{R}$ is the *kernel function* and $b \in \mathbb{R}$ is the bias term.

It is important to remark that the kernel $K(\cdot, \cdot)$ in the above LS-SVM regression in the dual space is different from the covariance function $k(\cdot, \cdot)$ adopted by the GPR. In this work, the LS-SVM Lab MATLAB Toolbox, version 1.8 [31] is used to build the LS-SVM regression model.

Summarizing, in the probabilistic model given in (1), the mean function $\mu(\mathbf{x})$ is built with the LS-SVM regression, whereas the model residuals $\varepsilon(\mathbf{x})$ are approximated by a GP regression. The resulting model not only approximates the system responses for any configuration of the input parameters, but it also provides an estimation of the prediction uncertainty.

III. PARTIAL LEAST SQUARES

Partial Least Squares (PLS) [21] is a statistical method that enables to identify a linear relationship between input variables and the response variable. This technique relies on the projection of input variables onto a new space defined by new variables called principal components, which are linear combinations of the input variables. This method is well adapted when the input data set contains a small number of samples and a large number of input variables highly correlated to each other.

The principle of the PLS technique is based on the search of the best directions in the \mathbf{X} space including the input variables, in order to explain the behavior of the response \mathbf{y} . First, the input matrix \mathbf{X} of size $(N \times M)$, where N and M are the number of samples and input variables, respectively, and the response vector \mathbf{y} are centered and reduced in order to avoid a bias of the input variables having large values and strong variations. Then, the first principal component $\mathbf{t}^{(1)}$ is computed by searching the best direction $\mathbf{u}^{(1)}$ maximizing the squared covariance between $\mathbf{t}^{(1)} = \mathbf{X}\mathbf{u}^{(1)}$ and \mathbf{y} as:

$$\mathbf{u}^{(1)} = \arg \max_{\mathbf{u}^t \mathbf{u} = 1} \mathbf{u}^t \mathbf{X}^t \mathbf{y} \mathbf{y}^t \mathbf{X} \mathbf{u}. \quad (6)$$

The optimization problem (6) is maximized when $\mathbf{u}^{(1)}$ is the eigenvector of the matrix $\mathbf{X}^t \mathbf{y} \mathbf{y}^t \mathbf{X}$ associated with the eigenvalue with the largest absolute value. The vector $\mathbf{u}^{(1)}$, called *loading vector*, corresponds to the \mathbf{X} weights of the

first component. Then, the residual matrix of $\mathbf{X} = \mathbf{X}^{(0)}$ and $\mathbf{y} = \mathbf{y}^{(0)}$, which are denoted by $\mathbf{X}^{(1)}$ and $\mathbf{y}^{(1)}$, are computed as:

$$\begin{aligned}\mathbf{X}^{(1)} &= \mathbf{X}^{(0)} - \mathbf{t}^{(1)}\mathbf{b}^{(1)}, \\ \mathbf{y}^{(1)} &= \mathbf{y}^{(0)} - w_1\mathbf{t}^{(1)}\end{aligned}\quad (7)$$

with $\mathbf{b}^{(1)}$ representing a vector of size $(1 \times M)$ including the regression coefficients of the local regression of \mathbf{X} onto the first principal component $\mathbf{t}^{(1)}$, and w_1 is the regression coefficient of the local regression of \mathbf{y} onto the first principal component $\mathbf{t}^{(1)}$. The equation (7) corresponds to the local regression of \mathbf{X} and \mathbf{y} onto the first principal component.

Then, PLS computes the second principal component, which is orthogonal to $\mathbf{t}^{(1)}$, replacing \mathbf{X} by $\mathbf{X}^{(1)}$ and \mathbf{y} by $\mathbf{y}^{(1)}$ so that the problem (6) is solved again. The same technique is used to compute the rest of the components with an iterative procedure. As an illustration, Fig. 1 represents the method in a three-dimensional example by considering two principal components. The number of principal components retained is denoted by q . In the rest of the paper, the number of principal components will be selected by retaining the one minimizing the root mean square error (RMSE) given by:

$$RMSE = \frac{\sum_{i=1}^N (\hat{y}_i - y_i)^2}{N} \quad (8)$$

where N , y_i and \hat{y}_i are the sample size, the real values and the predictions of the model, respectively.

The new coordinate system, defined by the principal components, corresponds to a rotation of the initial system defined by $\mathbf{x}_1, \dots, \mathbf{x}_M$. The r -th component denoted $\mathbf{t}^{(r)}$, where $r = 1, \dots, q$, is given by:

$$\mathbf{t}^{(r)} = \mathbf{X}^{(r-1)}\mathbf{u}^{(r)} = \mathbf{X}\mathbf{u}_*^{(r)}. \quad (9)$$

The matrix $\mathbf{U}_* = [\mathbf{u}_*^{(1)}, \dots, \mathbf{u}_*^{(q)}]$ is given by:

$$\mathbf{U}_* = \mathbf{U}(\mathbf{B}^t\mathbf{U})^{-1} \quad (10)$$

with $\mathbf{U} = [\mathbf{u}^{(1)}, \dots, \mathbf{u}^{(q)}]$ and $\mathbf{B} = [\mathbf{b}^{(1)t}, \dots, \mathbf{b}^{(q)t}]$.

The vector $\mathbf{u}^{(r)}$ defines the principal direction in the initial space, maximizing the covariance of $\mathbf{X}^{(r-1)t}\mathbf{y}^{(r-1)}\mathbf{y}^{(r-1)t}\mathbf{X}^{(r-1)}$. When $q = M$, the matrix $\mathbf{U}_* = [\mathbf{u}_*^{(1)}, \dots, \mathbf{u}_*^{(M)}]$ rotates the space defined by $(\mathbf{x}_1, \dots, \mathbf{x}_M)$ to the new space defined by $(\mathbf{t}^{(1)}, \dots, \mathbf{t}^{(M)})$, oriented by the principal directions $(\mathbf{u}^{(1)}, \dots, \mathbf{u}^{(M)})$.

The variables importance projections (VIP) [23] in PLS is a variable selection method allowing to identify the importance of each input variables. It assumes that the column space of \mathbf{X} has a subspace of dimension q containing all relevant information for predicting the model response \mathbf{y} .

The VIP score for the input variable j is defined as:

$$VIP_j = \sqrt{\frac{M \cdot (\sum_{r=1}^q R^2(\mathbf{y}, \mathbf{t}^{(r)}) \cdot (\mathbf{u}^{(rj)} / \|\mathbf{u}^{(rj)}\|)^2)}{\sum_{r=1}^q R^2(\mathbf{y}, \mathbf{t}^{(r)})}} \quad (11)$$

where M is the number of input variables, $\mathbf{u}^{(rj)}$ is the weight of the j^{th} variable in component r and $R^2(\mathbf{y}, \mathbf{t}^{(r)})$ is the percentage of \mathbf{y} explained by the component r . The VIP value

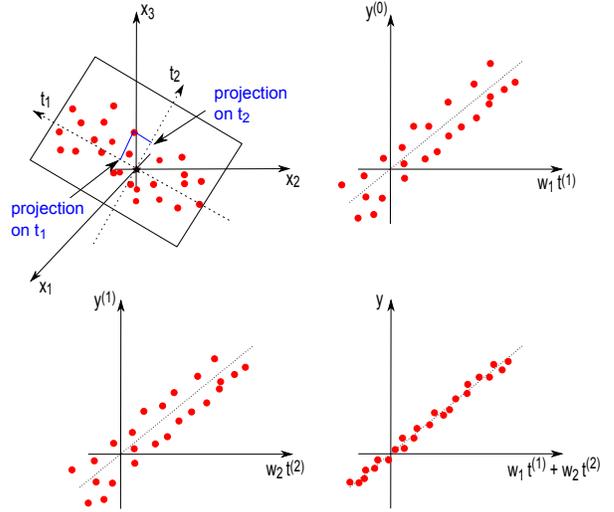


Fig. 1. Top left represents the construction of two principal components in the initial \mathbf{X} space. Top right and Bottom left illustrate the prediction of $y^{(0)}$ and $y^{(1)}$, respectively. Bottom right shows the final prediction of the output y .

is a weighted sum of squares of the PLS weights $(\mathbf{u}^{(rj)})$, which takes into account the explained variance of each PLS dimension. The input variables with a VIP score greater than one are generally considered to have a significant impact on the response \mathbf{y} [23].

IV. APPLICATION EXAMPLE: WIRELESS POWER TRANSFER

This Section investigates the feasibility and the strength of the GP+LS-SVM and the PLS regression techniques presented in Sec. II and III. Specifically, the above approaches are adopted to quantify the impact of fabrication tolerances and uncertainties on the maximal rectification efficiency, denoted E_{ff} , of the wireless power transfer (WPT) application given in Fig. 2 over [500, 1500] MHz frequency range (additional details on the considered test case are provided in [8], [29], [30]). For this purpose, the stochastic behavior of the efficiency has been analyzed by considering 30 Gaussian random variables centered at their nominal values and with a standard deviations of 5% of the nominal values, as given in Tab. I [8].

The LS-SVM+GPR and the PLS regressions are used hereafter to develop a probabilistic and deterministic surrogate model of the WPT efficiency in order to quantify the impact of uncertainties and reduce the computational cost with respect to MC simulation. A full-wave EM solver is used.

Both the LS-SVM+GPR and the PLS-based surrogate models are built by using the same set of training samples consisting of 200 realizations from Latin Hypercube sampling (LHS) [28]. As the PLS method is able to capture the behavior of an output with a small number of principal components, we construct a PLS surrogate model with a number of components varying from one to five. In order to evaluate the performance of each PLS model, a comparison with the brute force MC simulation with 10,000 realizations is carried out. Table II presents a comparison of each PLS model with respect to the

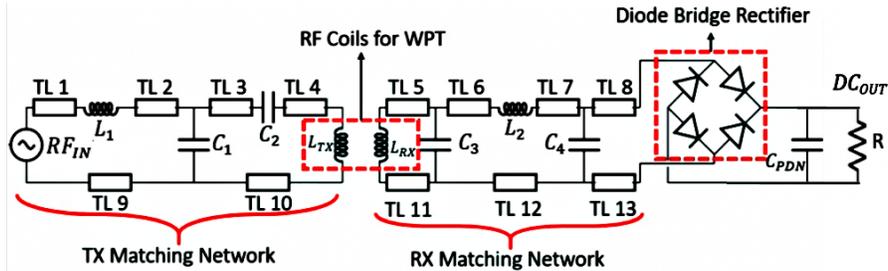


Fig. 2. Schematic of the WPT based power delivery architecture considered.

TABLE I
UNCERTAIN PARAMETERS CONSIDERED FOR THE WPT APPLICATION
EXAMPLE.

Gaussian Random Variables	Unit	μ	σ	
Feeding Gap for TX coil	$g_{f,TX}$	mm	2.46	0.1230
Feeding Gap for RX coil	$g_{f,RX}$	mm	2.3	0.1150
Width & Height of TX coil	$g_{\{x,y\},TX}$	mm	2.1	0.1050
Width & Height of RX coil	$g_{\{x,y\},RX}$	mm	3	0.1500
GND Cut-Out Ratio TX	$slot_{\{x,y\},TX}$	—	1.19	0.0595
GND Cut-Out Ratio RX	$slot_{\{x,y\},RX}$	—	1.16	0.0580
Match Capacitor TX	C_1	pF	4.09	0.2045
Resonance Capacitor TX	C_2	pF	4.59	0.2295
Resonance Capacitor RX	C_3	pF	2.16	0.1080
Match Capacitor RX	C_4	pF	0.3	0.015
Match inductor TX	L_1	nH	1.25	0.0625
Match inductor RX	L_2	nH	7.82	0.3910
Line Width TX Coil	$l_{w,TX}$	mm	1.89	0.0945
Line Width RX Coil	$l_{w,RX}$	mm	0.71	0.0355
Width TL1	w_{TL1}	mm	0.381	0.0191
Width TL2	w_{TL2}	mm	0.407	0.0204
Width TL3	w_{TL3}	mm	0.386	0.0193
Width TL4	w_{TL4}	mm	0.528	0.0264
Width TL5	w_{TL5}	mm	0.521	0.0191
Width TL6	w_{TL6}	mm	0.397	0.0198
Width TL7	w_{TL7}	mm	0.607	0.0303
Width TL8	w_{TL8}	mm	0.520	0.0260
Length TL1	l_{TL1}	mm	1.656	0.0828
Length TL2	l_{TL2}	mm	1.424	0.0712
Length TL3	l_{TL3}	mm	0.723	0.0362
Length TL4	l_{TL4}	mm	2.153	0.0828
Length TL5	l_{TL5}	mm	0.612	0.1076
Length TL6	l_{TL6}	mm	1.666	0.0833
Length TL7	l_{TL7}	mm	0.502	0.0251
Length TL8	l_{TL8}	mm	0.555	0.0278

number of components by computing the root mean square error (RMSE) over the 10,000 realizations, the mean value $\hat{\mu}$ and the standard deviation $\hat{\sigma}$ of the efficiency. We observe that the PLS model with one principal component provides a RMSE equal to 0.547. Then, the PLS model, from two to five principal components, provides a RMSE of 0.535 and 0.536. It is worth noting that the PLS surrogate model with one component achieved a quite low RMSE, i.e. 0.547, which is very close to the lowest RMSE obtained by PLS with two principal components, i.e. 0.535. In the following, this last PLS surrogate model is selected as it is the most accurate. A similar comparison has been carried out for the LS-SVM+GPR surrogate model. For this specific model the RMSE is 0.488, which is comparable with the one provided by the PLS with two components.

By comparison with several surrogate models used in [8] (see Tab. I), we notice that both the PLS surrogate model built-up using 200 LHS simulations is a bit less accurate than the LS-SVM surrogate model with a RBF kernel and more accurate than the sparse PC, which have shown their efficiency in complex problems. This highlights the capability of the PLS surrogate model to capture the main trend of a complex response based on a significant reduction of the dimensionality of the problem, i.e., from 30 initial variables to two principal components. This feature of the advocated method is very interesting and will be exploited in the following Section. On the other hand, the accuracy of the LS-SVM+GPR is the same of the one obtained by the LS-SVM surrogate model with RBF kernel.

As an illustration, Fig. 3 provides the scattering plots comparing the predictions of the LS-SVM+GPR and the PLS regression model with two principal components against the MC simulation with 10,000 realizations. We see a good agreement between the estimated efficiency by the proposed LS-SVM+GP and PLS surrogate models compared with the one obtained from the results of the MC simulation, since the realizations are quite close to the dashed line representing a perfect surrogate model. The above qualitative comment on the model accuracy is also confirmed by the correlation coefficients which take the value of 0.704 for the LS-SVM+GPR model and 0.628 for the PLS model, thus confirming the capability of the proposed techniques to deal with complex problem with a limited set of training samples (i.e., 30 variables with only 200 samples). It is worth pointing out that the low values of the WPT efficiency, where the efficiency estimated by MC simulation is below 73%, are not well predicted by the models

TABLE II

COMPARISON AMONG THE PERFORMANCES OF THE PLS, SPARSE PC, LS-SVM (RBF) AND LS-SVM (RBF)+GPR-BASED MODELS OF THE WPT EFFICIENCY BUILT WITH 200 LHS SAMPLES W.R.T THE RESULTS OF A MC SIMULATION WITH 10,000 SAMPLES. THE NUMBERS IN BOLD FONT HIGHLIGHT THE RMSE OF THE MOST ACCURATE MODELS.

Method	Number of Comp.	RMSE	$\hat{\mu}$	$\hat{\sigma}$	t_{model}^*	t_{cost}
MC	—	—	74.19	0.68	—	11 days
PLS	1	0.547	74.20	0.45	5h*	<1s
	2	0.535	74.20	0.50	5h*	<1s
	3	0.536	74.20	0.51	5h*	<1s
	4	0.536	74.20	0.51	5h*	<1s
	5	0.536	74.20	0.51	5h*	<1s
LS-SVM (RBF)+GPR	—	0.488	74.25	0.68	5h*	<1s
LS-SVM (RBF) [8]	—	0.488	74.23	0.54	5h*	<1s
Sparse PC (Order 3) [8]	—	0.565	74.25	0.52	5h*	<1s

* 200 LHS training simulations required 5 h.

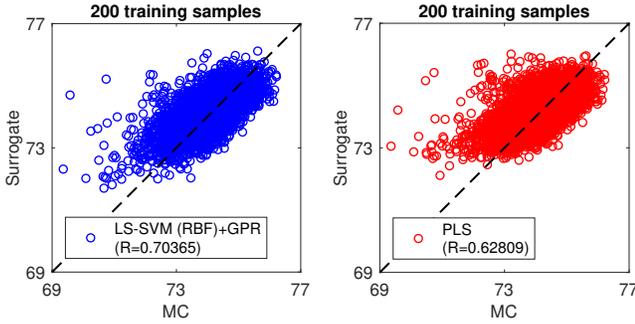


Fig. 3. Scatter plot of the WPT efficiency computed by the PLS surrogate model (red circles), the mean value of the LS-SVM (RBF)+GPR surrogate model (blue circles) vs. the actual numerical model with 10,000 MC realizations. The dashed line illustrates the perfect agreement between the models.

due to the strong non-linearity and discontinuities of the output in that region.

As the variability of the output is of interest, Fig. 4 shows the PDFs estimated by the LS-SVM+GP and PLS surrogate models and by MC simulation from 10,000 realizations. It can be observed that the LS-SVM+GP and PLS surrogate models capture well the main variability of the efficiency provided by MC simulation, although some discrepancies appear between the left tails of the probability distribution. Regarding the computational cost, 10,000 MC simulations required about 11 days while the PLS surrogate and the LS-SVM+GPR model needed less than 1 s to perform the same predictions. This comparison does not consider the computational cost required to generate the training data set (200 simulations), which took about 5 h as shown in Table II.

V. METHOD STRENGTHS

The results provided in Sec. IV show that both proposed techniques allow providing accurate surrogate models for the considered test case by using a limited set of training samples. This section aims at highlighting the advantages and differences between the two modeling techniques.

A. LS-SVM+GPR Probabilistic Model

The LS-SVM+GPR probabilistic model presented in Sec. II is extremely useful and powerful, since it provides the user

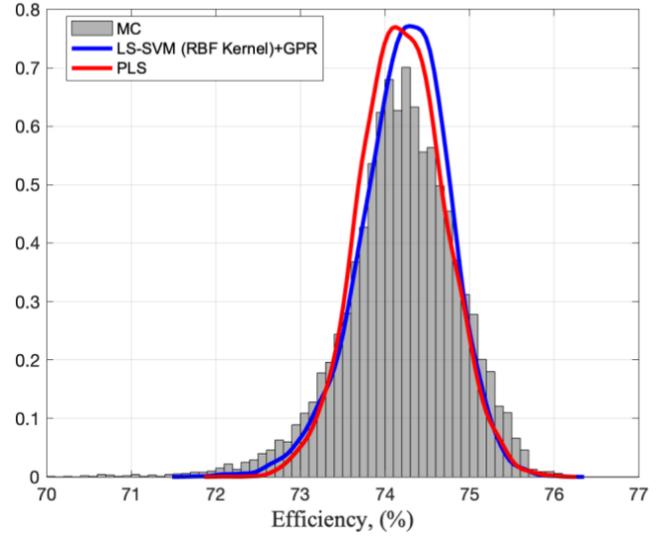


Fig. 4. PDFs of the WPT efficiency estimated by the PLS surrogate model (solid blue curve), the mean value of the LS-SVM (RBF)+GP surrogate model (solid red curve) and by the MC simulation (gray histogram) from 10,000 realizations.

with additional information on the model reliability. Indeed, for any configuration of the stochastic input variables \mathbf{x}_* , the proposed probabilistic model provides the output as a stochastic variable with a Gaussian PDF which writes:

$$\mathcal{M}(\mathbf{x}_*) \sim N(\mu_{\mathbf{x}_*}, \sigma_{\mathbf{x}_*}^2) = \mathcal{M}_{LS-SVM+GPR}(\mathbf{x}_*) \quad (12)$$

where $\mu_{\mathbf{x}_*}$ and $\sigma_{\mathbf{x}_*}^2$ are computed according to (4a) and (4b), respectively.

This means that each prediction obtained with the proposed probabilistic model is a Gaussian variable. The above description of the model output in terms of a PDF allows providing a probabilistic interpretation of the model uncertainty such as the model error, in which the prediction obtained with the model is a Gaussian variable. According to this, Fig. 5 provides a comparison between the PDF of the WPT efficiency computed from the results of a 10,000 MC simulations with the PDFs obtained by considering, for each configuration of the MC samples, 100 different realizations of the corresponding output $N(\mu_{\mathbf{x}_*}, \sigma_{\mathbf{x}_*}^2)$ in (12). The results show that the statistical information provided by the proposed probabilistic model allows providing a more conservative and reliable estimation of the actual statistical behavior of the efficiency, especially in tails of the distribution [14].

Moreover, the resulting probabilistic interpretation of the model realization allows estimating the confidence bounds on the model prediction for a given confidence interval (CI) of $100(1 - \alpha)\%$ as follows:

$$(\mu_{\mathbf{x}_*} - z_{1 - \frac{\alpha}{2}} \sigma_{\mathbf{x}_*}) \leq \mathcal{M}(\mathbf{x}_*) \leq (\mu_{\mathbf{x}_*} + z_{1 - \frac{\alpha}{2}} \sigma_{\mathbf{x}_*}), \quad (13)$$

where z denotes the inverse of the Gaussian cumulative distribution function evaluated at $1 - \frac{\alpha}{2}$.

For the considered application example, the 99% CIs predicted by the proposed LS-SVM+GPR model allow providing a conservative estimation of the actual values of the WPT

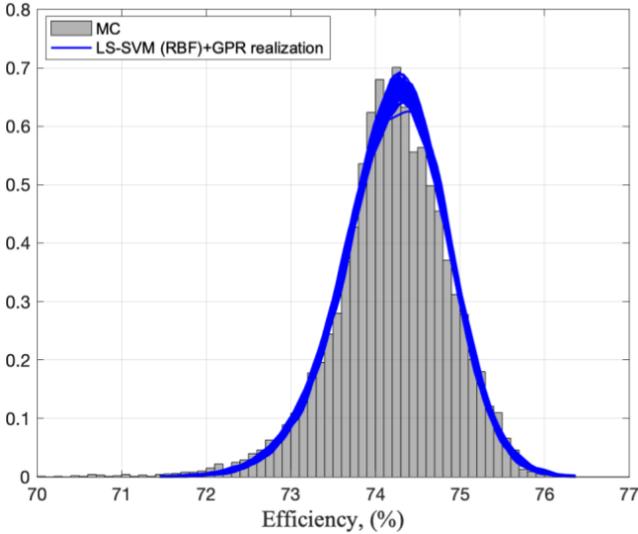


Fig. 5. Comparison between the PDFs of the WPT efficiency calculated from the results of 10,000 MC simulations of the full-computational model (gray histogram) and the PDFs obtained by considering, for each configuration of the MC samples, 100 different realizations predicted by the proposed LS-SVM+GPR probabilistic model (blue lines).

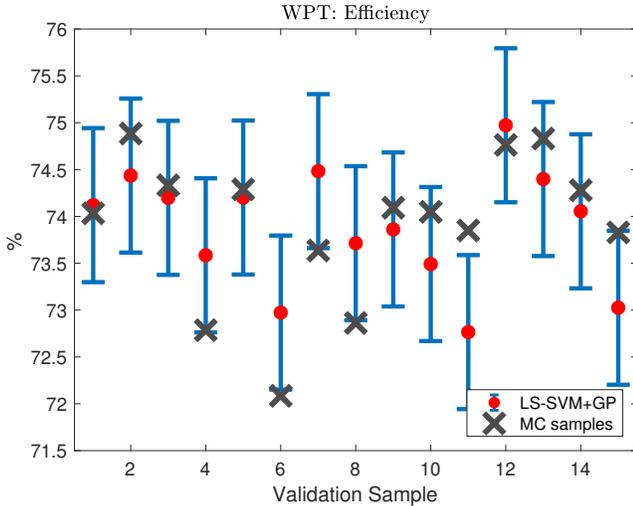


Fig. 6. Comparison between the results of 15 simulations of the full-computational model (black crosses) randomly chosen among the 10,000 MC samples with the corresponding mean value (red dots) and 99% CI (horizontal blue bars) estimated via the LS-SVM+GPR probabilistic model.

efficiency for 91,280 cases out of 10,000 validation samples (i.e., 91.3% of the overall cases). For the sake of illustration, Fig. 6 shows the CIs predicted by the model for 15 validation samples randomly selected among the 10,000 samples used in the MC simulation. According to the results, the proposed model allows to accurately predict the model uncertainty, since most of the validation samples (black crosses) fall within the CIs predicted by the proposed model (blue bars).

B. PLS & Sensitivity Analysis

Besides the estimation of the variability of the WPT efficiency, the PLS model allows to derive a sensitivity analysis of the output. Figure 7 presents a histogram of the VIP scores, which estimate the importance of each input variables in the

projection used in the PLS model. From Figure 7, we observe that, out of 30 input random variables, eight variables have a significant impact, defined by a VIP score larger than or equal to 1 (dashed line) as suggested by [23], on the variability of the efficiency E_{ff} . In particular, the most important parameters are related to the transmitter and receiver coils, i.e., $g_{\{x,y\},TX}$ and $g_{\{x,y\},RX}$, $g_{f,RX}$, $slot_{\{x,y\},RX}$, respectively, as well as some components such as L_2 , C_2 , C_3 and C_4 . The rest of the parameters, including the ones associated to the geometry of transmission lines (i.e., widths and lengths) have a negligible impact on the variability of the WPT efficiency.

The PLS surrogate also allows to analyze the correlation of such parameters on the WPT efficiency for the first two components (i.e., t_1 and t_2) via the correlation circle in Fig. 8. This correlation can be obtained by looking at the weight of each variables over the two components. More specifically, Fig. 8 shows that the projection of the output of interest (i.e. WPT efficiency E_{ff}) on the first component t_1 (which is the most important by construction) is equal to about 0.75. We also notice that the correlation with the group of input variables composed of L_2 , C_4 , $slot_{\{x,y\},RX}$, C_3 , $g_{\{x,y\},RX}$ and the first component t_1 is about 0.5. Since those variables (input and output variables) are positively correlated with the first component t_1 , and relatively close to each other, this means that a higher value of these input variables will increase the output variable (i.e., the WPT efficiency E_{ff}). The rest of the input variables are not significantly correlated with the WPT efficiency E_{ff} as their projection over the first component t_1 is quite low.

To demonstrate the validity of this information, for the input variables L_2 , C_4 , $slot_{\{x,y\},RX}$, C_3 , $g_{\{x,y\},RX}$, we increased the corresponding mean values provided in Tab. I of 2σ (i.e., $\tilde{\mu} = \mu + 2\sigma$) while choosing a standard deviation of $\tilde{\sigma} = \sigma/10$ in order to not exceed the upper bound of the input variables of the initial configuration. A new MC simulation with 10,000 samples has been run providing the new PDF (green bins) in Fig. 9. This illustration compares the PDF obtained with the optimized configuration of the system parameters (green histogram) with their initial configuration (grey histogram). The mean value and the standard deviation of the optimized and initial configurations are given by $\mu_{opt} = 74.89$ and $\sigma_{opt} = 0.74$, and $\mu_{init} = 74.19$ and $\sigma_{init} = 0.68$, respectively, confirming that controlling a small group of input parameters can significantly improve the performance of the system.

VI. CONCLUSIONS

This paper focuses on the application of the LS-SVM+GPR and the PLS technique to the development of surrogate models, with a very low computational cost compared to MC simulation, for the uncertainty quantification in high-dimensional space. The obtained models allowed to efficiently estimate the model output in a complex problem characterized by 30 uncertain input variables. The results of a 10,000 MC samples obtained with the full-computational model have been considered as reference results. The strength of the proposed technique along with their accuracy and computational cost are carefully discussed. From the results, the proposed modeling

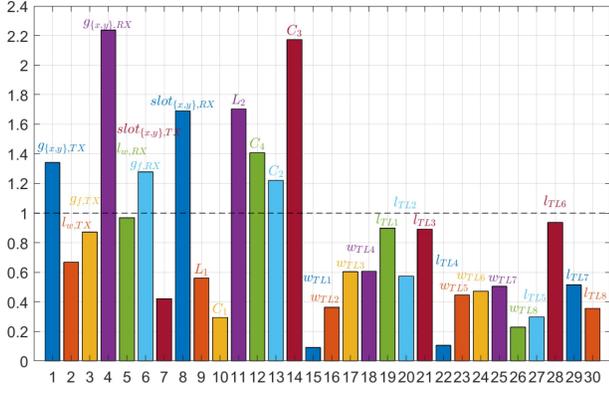


Fig. 7. Importance of the input parameters according to their impact on the behavior of the output.

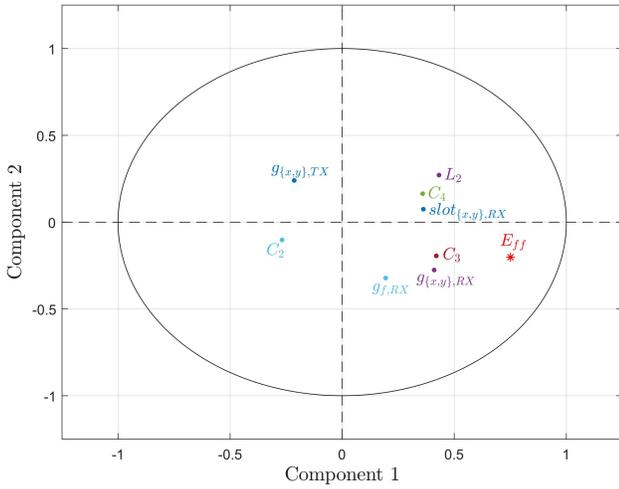


Fig. 8. Circle of correlation representing the projection of the input variables and the output variable (WPT efficiency E_{ff}) onto the plane defined by the first two PLS components. It shows a strong positive correlation between the input variables L_2 , C_4 , $slot_{\{x,y\},RX}$, C_3 , $g_{\{x,y\},RX}$ and the output variable E_{ff} since their projections are in the same direction and at a significant distance from the origin.

schemes can be considered as valuable and viable solutions for the UQ of the output of complex systems in high-dimensional parameter space.

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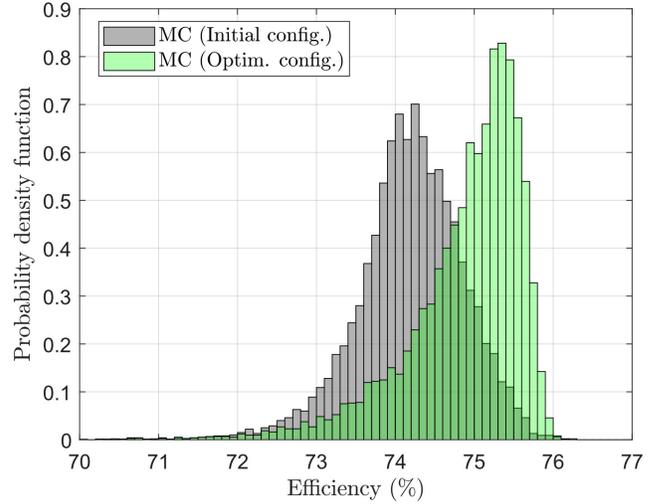


Fig. 9. PDF of the efficiency of the wireless power transfer improved (green) by increasing the nominal values of the input variables L_2 , C_4 , $slot_{\{x,y\},RX}$, C_3 , $g_{\{x,y\},RX}$ by 2σ , where σ is given Table I, according to the circle of correlation given in Fig. 8.

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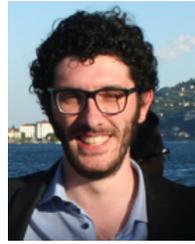
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