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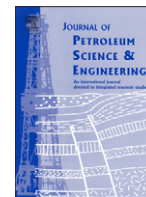
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## A geostatistical perspective for the surrogate-based integration of variable fidelity models

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### ABSTRACT

When constructing surrogate models, time/cost constraints make the designer frequently face the dilemma of whether to use a small sample of data obtained from, for example, high fidelity/computationally expensive computer simulations, or, a large one but with low fidelity values. More generally, variable fidelity samples can be the result of: i) different physical/mathematical representations (e.g., inviscid/Euler versus viscous/Navier–Stokes calculations), ii) alternative resolution models (e.g., fine/coarse grids), or, iii) experiments. Ideally, surrogate models should allow: a) the integration of variable fidelity samples, and, b) provide estimation and appraisal (error) information consistent with the amount and fidelity level of the available data. While there have been significant progress in this area through deterministic modeling and optimization approaches (e.g., correction surfaces, and space mapping), a spatial-stochastic perspective such as those provided by the branch of spatial statistics known as geostatistics offers distinctive advantages when satisfying the above referenced requirements (a and b). This paper discusses the effectiveness and requirements of geostatistical methods such as classic and collocated cokriging for the integration of variable fidelity models. The discussion is illustrated using well-known analytical functions and, alternative resolution models, in the surrogate-based modeling of a field scale alkali–surfactant–polymer (ASP) enhanced oil recovery (EOR) process.

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### 1. Introduction

Surrogate modeling is increasingly popular and has been successfully used in the analysis and optimization of computationally expensive models in, for example, the aerospace (Giunta et al., 1997; Balabanov et al., 1998; Li and Padula, 2004; Queipo et al., 2005), automotive (Craig et al., 2002; Kurtaran et al., 2002), and oil industries (Queipo et al., 2002a, b). Recent review papers on the subject are those of Li and Padula (2004), Wang and Shan (2007), Queipo et al. (2005). When constructing surrogate models, time/cost constraints make the designer frequently face the dilemma of whether to use a small sample of data obtained from, for example, high fidelity/computationally expensive computer simulations, or, a large one but with low fidelity values. More generally, variable fidelity samples can be the result of: i) different physical/mathematical representations (e.g., inviscid/Euler versus viscous/Navier–Stokes calculations), ii) alternative resolution models (e.g., fine/coarse grids), or, iii) experiments. Ideally, surrogate models should allow: a) the integration of variable fidelity samples, and, b) provide estimation and appraisal (error) information consistent with the amount and fidelity level of the available data.

There have been significant progress in this area through deterministic modeling and optimization approaches such as correction surfaces (Chang et al., 1993; Toropov et al., 1999; Vitali et al., 2002; Alexandrov and Lewis, 2002) and space mapping (Bakr et al., 2002; Bandler et al., 2004); however, these alternatives do not provide appraisal (error) information. From a spatial-stochastic perspective, geostatistical methods offer distinctive advantages for addressing issues a) and b), most notably: they are sound (founded on solid statistical principles), flexible (e.g., can be extended to higher dimensions), and efficient (e.g., can be implemented through matrix operations). The required covariance models (structure and parameters) of the variable fidelity data can be identified either through the so called DACE (Sacks et al., 1989b) or, variogram approaches.

While there is a variety of geostatistical methods (Isaaks and Srivastava, 1989; Goovaerts, 1997; Chiles and Delfiner, 1999) potentially useful for addressing the problem of interest such as (increasingly complex): kriging with external drift (KED), collocated cokriging with Markov models (CCM), collocated cokriging (CC), classic cokriging (CLC), in this work, the more general CC and CLC methods are selected for investigation. Related works include those of Kennedy and O'Hagan (2000) which was restricted to a CCM method, and those reported in Koh-Sung and Tapabrata (2004) and Chung and Alonso (2002a,b) where cokriging approaches use gradient (secondary) information in surrogate-based estimations.

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This paper discusses the effectiveness and requirements of geostatistical methods such as collocated (two versions) and classic cokriging for the integration of two level (low and high) fidelity models. The discussion is illustrated using a well-known analytical function and, two distinct resolution models, in the surrogate-based modeling of a field scale alkali–surfactant–polymer (ASP) enhanced oil recovery (EOR) process. ASP flooding is the most promising EOR solution for one of the greatest challenges facing the oil industry worldwide: after conventional water flooding the residual oil (drops trapped by capillary forces) in reservoirs around the world is likely to be around 70% of the original oil in place.

## 2. Problem definition

Given  $Z = \{z_1, z_2, \dots, z_n\}$ , and,  $V = \{v_1, v_2, \dots, v_m\}$ , representing high and low fidelity model input/output pairs, respectively, with  $m \gg n$ , build a surrogate of the high fidelity model that provides estimation and appraisal (error) information consistent with the amount and fidelity level of the available data. It is assumed that: i) the high and low fidelity models are correlated, and, ii) each model output is a scalar. Fig. 1 shows a particular case of the problem of interest where high (circle) and low (diamond) fidelity model samples are available, and an estimation of the high fidelity model output and corresponding appraisal (error) information is sought at a specific location (marked with an “X”).

## 3. Geostatistical methods

Kriging and related methods represent, so called, BLUE-best (minimum error variance) linear unbiased estimators (Isaaks and Srivastava, 1989; Goovaerts, 1997; Chiles and Delfiner, 1999). From the variety of geostatistical methods potentially useful for addressing the problem of interest such as (increasingly complex): kriging with external drift (KED), collocated cokriging with Markov models (CCM), collocated cokriging (CC), classic cokriging (CLC), in this work, the more general CC and CLC methods are selected for investigation.

### 3.1. Kriging and cokriging

Table 1 summarizes the optimization problem associated with ordinary kriging, and classic and collocated cokriging, and the corresponding closed-form solutions. In general, the cokriging methods consider only a small sample associated with a high fidelity model, and, an abundant sample associated with a low fidelity model one.

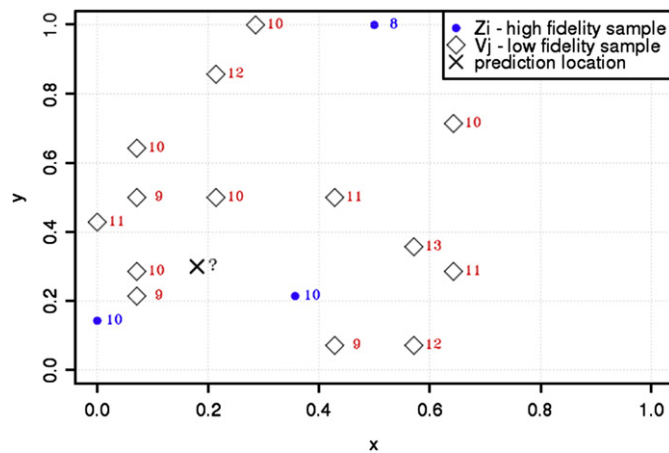


Fig. 1. Distribution of low (diamonds) and high (circles) fidelity model sample for a 2D illustration of the problem of interest.

Table 1

Optimization problem associated with kriging, and cokriging methods, and corresponding closed-form solutions.

Method	Optimization formulation
Ordinary kriging	<p>Find <math>\alpha</math> in:</p> $\hat{z}_0 = \alpha^T Z$ <p>such that:</p> $\text{Var}(z_0 - \hat{z}_0) = \sigma_z^2 + \alpha^T \text{Cov}(Z)\alpha - 2\alpha^T \text{Cov}(Z, z_0)$ <p>is minimized subject to the restriction:</p> $\sum \alpha = 1$ <p>to ensure unbiasedness.</p> <p>Closed-form solution is:</p> $\hat{z}_0 = w^T IC (I - h L L^T IC) Z + h L^T IC Z$ <p>with</p> <p><math>L</math> = vector of ones with the same length of <math>Z</math></p> <p><math>IC = \text{Cov}(Z)^{-1}</math></p> <p><math>h = (L^T IC L)^{-1}</math>.</p>
Classic cokriging	<p>Find <math>\alpha</math> and <math>\beta</math> in:</p> $\hat{z}_0 = \alpha^T Z + \beta^T V$ <p>such that:</p> $\text{Var}(z_0 - \hat{z}_0) = \sigma_z^2 + \alpha^T \text{Cov}(Z)\alpha + \beta^T \text{Cov}(V)\beta + 2\alpha^T \text{Cov}(Z, V)\beta - 2\alpha^T \text{Cov}(Z, z_0) - 2\beta^T \text{Cov}(V, z_0)$ <p>is minimized subject to the restrictions: <math>\sum \alpha = 1</math> and <math>\sum \beta = 0</math> to ensure unbiasedness.</p> <p>Closed-form solution is:</p> $\hat{z}_0 = \begin{bmatrix} 0 \\ 0 \\ 1 \\ 0 \end{bmatrix}^T \text{COVAMP}^{-1} \begin{bmatrix} Z \\ V \\ 0 \\ 0 \end{bmatrix} + \begin{bmatrix} \text{Cov}(Z, z_0) \\ \text{Cov}(V, z_0) \\ 0 \\ 0 \end{bmatrix}^T \text{COVAMP}^{-1} \begin{bmatrix} Z \\ V \\ 0 \\ 0 \end{bmatrix}$ <p>with</p> $\text{COVAMP} = \begin{bmatrix} \text{Cov}(Z) & \text{Cov}(Z, V) & L_n & 0 \\ \text{Cov}(V, Z) & \text{Cov}(V) & 0 & L_m \\ L_n^T & 0 & 0 & 0 \\ 0 & L_m^T & 0 & 0 \end{bmatrix}$ <p><math>L_n</math> and <math>L_m</math> are vectors of ones with the same length of <math>Z</math> and <math>V</math> respectively.</p>
Collocated cokriging	<p>Find <math>\alpha</math> and <math>\beta</math> in:</p> $\hat{z}_0 = \alpha^T Z + \beta v_0$ <p>such that:</p> $\text{Var}(z_0 - \hat{z}_0) = \sigma_z^2 + \beta^2 \sigma_v^2 + \alpha^T \text{Cov}(Z)\alpha + 2\alpha^T \text{Cov}(Z, v_0)\beta - 2\alpha^T \text{Cov}(Z, z_0) - 2\beta \text{Cov}(v_0, z_0)$ <p>is minimized subject to the restrictions: <math>(\sum \alpha) + \beta = 1</math> to ensure unbiasedness.</p> <p>Closed-form solution is:</p> $\hat{z}_0 = \begin{bmatrix} 0 \\ 0 \\ 1 \\ 0 \end{bmatrix}^T \text{COVAMP}^{-1} \begin{bmatrix} Z \\ v_0 \\ 0 \\ 0 \end{bmatrix} + \begin{bmatrix} \text{Cov}(Z, z_0) \\ \text{Cov}(v_0, z_0) \\ 0 \\ 0 \end{bmatrix}^T \text{COVAMP}^{-1} \begin{bmatrix} Z \\ v_0 \\ 0 \\ 0 \end{bmatrix}$ <p>with</p> $\text{COVAMP} = \begin{bmatrix} \text{Cov}(Z) & \text{Cov}(Z, v_0) & L_n \\ \text{Cov}(v_0, Z) & \sigma_v^2 & 1 \\ L_n^T & 1 & 0 \end{bmatrix}$ <p><math>L_n</math> = vectors of ones with the same length of <math>Z</math>.</p>

Note that: i) the geostatistical method estimates are linear combinations of the available data, and the main goal of the optimization problem is to find the weights to ensure the best unbiased estimation, and ii) the optimization problems and closed-form solutions assume that a covariance function (Cov) has been identified (this issue is addressed later in this section).

Kriging is a spatial prediction method from geostatistics developed by Matheron (1963) and named after the pioneering work of D.G. Krige (a South African mining engineer); Sacks et al. (1989a,b), and Jones et al. (1998) made it well-known in the context of the modeling and optimization of deterministic functions, respectively. In this context, the prediction is modeled as the sum of two effects, a mean (constant or a linear trend), and fluctuations around the mean, and it represents a so called BLUE-best (minimizes prediction error variance) linear unbiased (expected value of the error equal to zero) estimator. Depending on the mean assumption, there are several kriging formulations, namely: simple, ordinary, universal and kriging with external drift; the assumptions: a known constant, a constant (unknown), a linear polynomial on the input variables, and a linear polynomial on an external (secondary) variable, respectively. The fluctuations around the mean are represented by a zero mean, second order, stationary process.

Note that the geostatistical methods are: i) founded on solid statistical principles (*sound*), ii) can be extended to higher dimensions, provide both estimation and error information, and allow for non-collocated low and high fidelity data (*flexible*), and iii) represent linear estimators which make them easy to implement and update through matrix operations (*efficient*). In addition, the geostatistical methods exhibit the following features: interpolates the high fidelity data, estimated values not constrained by maximum and minimum data values, and the so called *declustering*, *screening*, and *smoothing* effects. The declustering effect refers to the fact that a cluster of closely located samples will have collectively the weight of a single sample located near the centroid of the cluster. On the other hand, the screening effect reduces the influence of a sample (original) by the addition of one or more samples at intermediate locations between the original sample and the prediction location. Furthermore, the considered geostatistical methods behave as low-pass filters away from the available data and their effect grows with the prediction error variance.

Fig. 2 illustrates the ordinary kriging weights of sample values for estimating the high fidelity model output value at a prediction location (marked with an "X"). Note that even though samples denoted as a, b and c, are equally distant from the prediction location, the kriging weights for the closely located samples b and c are reduced because of the so called declustering effect; the figure also illustrates

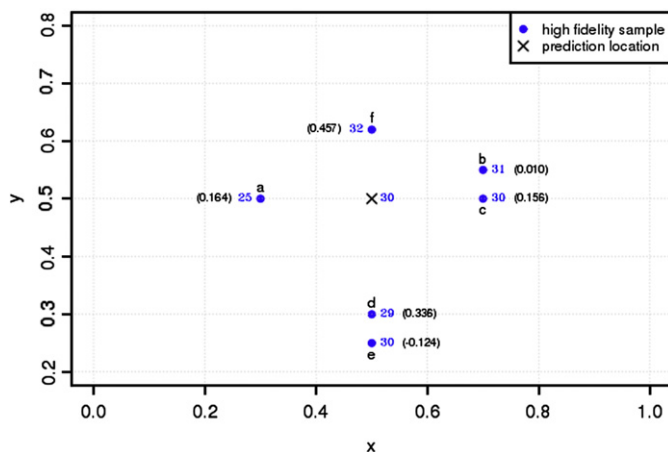


Fig. 2. Ordinary kriging weights (in parenthesis) of sample values for estimating the high fidelity model output at a prediction site ("X"). The circle mark represents sample locations, and the number aside the mark denotes high fidelity model output values.

the screening effect in samples d and e, having the latter a negative weight. Notice how the sum of all the weights is one hence satisfying the ordinary kriging restriction to ensure unbiasedness.

Additionally, these methods give appraisal (error) information with the estimation, as it is shown in Fig. 3. Fig. 3(a) shows the estimation and prediction error standard deviation obtained at a prediction location using the KRI method. Fig. 3(b), on the other hand, depicts the effect of the addition of low fidelity samples when using the CLC method; in this scenario, the prediction error standard deviation is considerably reduced at the prediction location.

The CC and CLC methods differ in whether they place restrictions on the relative position of the variable fidelity data, and on the covariance functions required. In the CC method values of the low fidelity model output have to be available at the prediction sites (exhaustive). The CLC method does not have that requirement. On the other hand, the CLC method has to fully specify a covariance function for the high and low fidelity model outputs, and for the cross-covariance between the high and low fidelity models. In contrast, the CC method does not require a covariance function for the low fidelity model. When using the CC method two different strategies are used to estimate the values of the low fidelity model at prediction locations (if necessary), namely; ordinary kriging (CC-Kri) and the  $k$ -nn (with  $k=1$ ) classifier algorithm (CC-1nn).

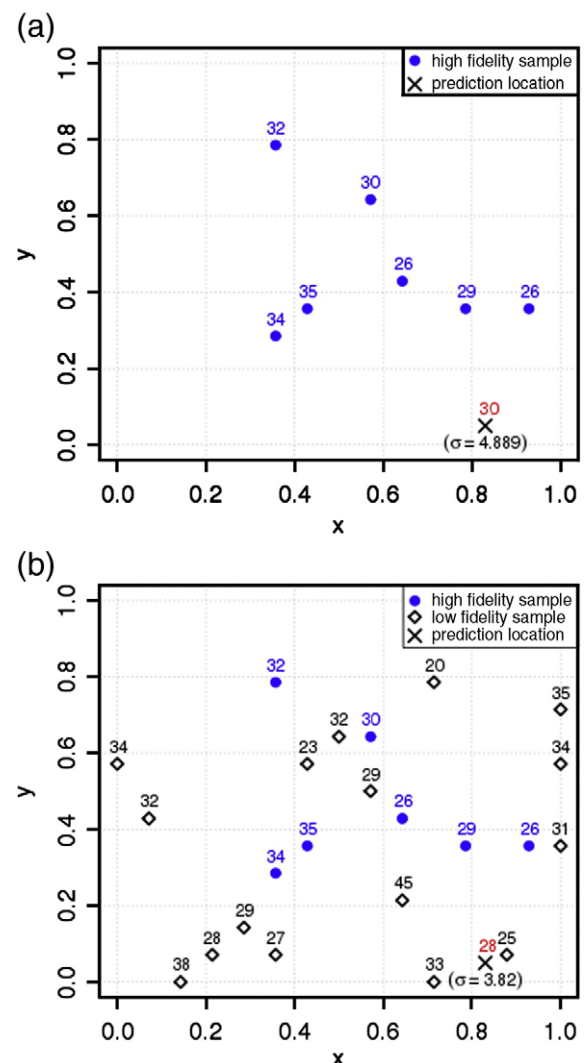


Fig. 3. Estimation and prediction error standard deviation obtained at a prediction location using (a) KRI and (b) CLC methods.



### 3.2. Covariance function identification

The covariance function is identified assuming the spatial-stochastic process associated with the kriging estimates is stationary, hence,  $\text{Cov}(Z_j, Z_k)$  is the same for any two points  $x_j, x_k$ , with  $x_j - x_k = h$ , a constant vector. More formally, in this context, the covariance between two random variables  $Z_j$  and  $Z_k$  is defined by the following expression ( $R^p \rightarrow R$ ):

$$\text{Cov}(Z_j, Z_k) = \text{Cov}(h) \quad \text{where } h = x_j - x_k$$

When  $\text{Cov}(h)$  only depends on the modulus  $|h|$  and not on the direction of vector  $h$ , the spatial-stochastic process is considered isotropic and  $\text{Cov}(h)$  is a scalar function ( $R \rightarrow R$ ). In order to identify the covariance function (structure and parameters) from only a sample of model input/output pairs some additional assumptions are required; in particular, an assumption regarding the principal directions of continuity. This fact lead to the two most well-known methods for covariance function estimation: the so called DACE (Sacks et al., 1989b) and variogram modeling (from geostatistics) approaches (see, for example, Isaaks and Srivastava, 1989; Goovaerts, 1997; Chiles and Delfiner, 1999).

The DACE approach assumes that the principal directions of continuity coincide with the Cartesian axes, then, for every variable component  $x_j$  from  $x$  there is a parameter  $\theta_j$  that adjust the continuity (smoothness) along the  $j$  Cartesian direction, and the correlation function is considered the product of a correlation function for each of the variable component in vector  $x$ , consequently, the covariance function can be specified as:

$$\text{Cov}(h) = \text{Cov}(0) \prod_{j=1}^p R(\theta_j, h_j, y)$$

where  $\text{Cov}(0)$  is the process variance, the  $R$  function makes reference to the correlation structure (specified by the user) that may depend on some parameters denoted here as  $y$ . The parameters  $\theta$  and those in  $y$  can be specified using, for example, maximum likelihood estimates.

In contrast to the DACE approach, the variogram modeling approach looks to explore the principal directions of continuity before identifying covariance parameters. For this purpose, for vector  $h$ , a so called variogram is defined as:

$$\gamma(h) = \frac{1}{2} E(Z_j - Z_k)^2 \quad \text{where } h = x_j - x_k$$

where  $E$  denotes expected value. The basic parameters that describe a variogram are the *sill* and *range*. The sill is the (special) asymptotic variogram value associated to large values of  $h$  (lag) and the range is the lag distance ( $h$ ) for the sill (spherical model) or 95% of the sill value (Gaussian and exponential models). Fig. 4 is an example of an experimental variogram (circles) with three different theoretical model structures; the sill and range are also shown.

Considering the stationarity assumption, the covariance and variogram are related by the following expression  $\text{Cov}(0) = \text{Cov}(h) + \gamma(h)$ ; so if you identify the variogram, the covariance function is also established. In order to identify  $\text{Cov}(h)$  using this approach another postulate needs to be made regarding the covariance anisotropy: the iso-contours of  $\text{Cov}(h)$ , that is,  $\text{Cov}(h) = c$ , are similar (with an elliptical shape) regardless of the constant  $c$ , and, as in principal component analysis, the orthogonal principal directions and axis sizes need to be established (e.g., using so called directional variograms). Having accomplished this task, a linear transformation takes the ellipsoids into spheres and the covariance problem is now isotropic and all the samples can be used to estimate the  $\text{Cov}(h)$ . The covariance model structures used are parsimonious and rely at most on only two or three parameters. Once the structure is established, the parameters (a lower number than

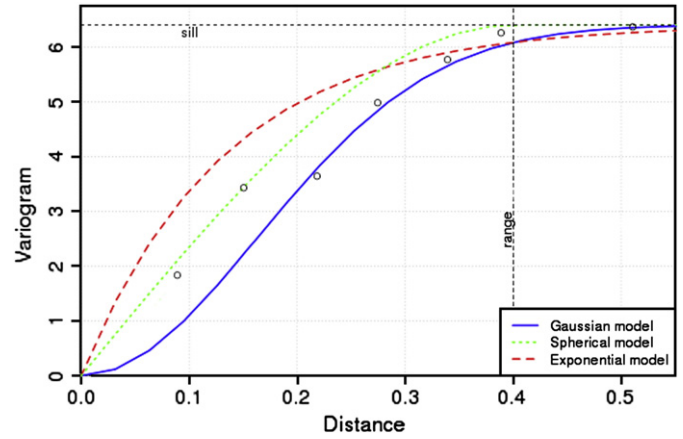


Fig. 4. Example of an experimental variogram and three theoretical variograms. The sill ( $\gamma = 6.4$ ) and range ( $h = 0.4$ ) are represented with dotted lines.

the  $p$  parameters  $\theta_j$  in the DACE approach) can be more robustly identified using weighted mean squares, maximum likelihood, etc.

### 4. Solution methodology

The proposed approach to build a surrogate of a high fidelity model from high and low fidelity data, as specified in Section 2, includes the following steps:

- (1) Identify the covariance models (structure and associated parameters) required for the selected geostatistical method. In this work, for such purpose, the variogram approach is favored. The CC method requires the identification of a covariance model associated with the high fidelity data, and, a cross-covariance model for the high and low fidelity data. On the other hand, the CLC method requires not only the above referenced covariance models but also a covariance model for the low fidelity data. Once the theoretical variogram model is identified (Table 2), the covariance model is obtained through the following expression:  $\text{Cov}(h) = \sigma^2 - \gamma(h)$ , where  $\sigma^2$  represents the sill,  $\gamma(h)$  the theoretical variogram model, and,  $h$  the distance between the points whose covariance is sought.
- (2) If necessary, extend the low fidelity data to the prediction sites (CC method). The cited extension can be done using kriging, or using the 1-nn classifier. In the case of kriging, a covariance function for the low fidelity model needs to be identified. Both approaches are used in the context of this work.
- (3) Build a surrogate of the high fidelity model using the covariance structures/parameters identified in step (1) as specified by the CC and CLC methods. For a given set of prediction sites, the cokriging models require the solution of constrained optimization problems established in Table 1.

Table 2  
Description of common theoretical variogram models.

Model	$\gamma(h)/\sum^2$	$h$
Nugget	0	$h = 0$
	1	$h > 0$
Spherical	$\frac{3h}{2a} - \frac{1}{2} \left(\frac{h}{a}\right)^3$	$0 \leq h \leq a$
	1	$h > a$
Exponential	$1 - \exp\left(-\frac{h}{a}\right)$	$h \geq 0$
Gaussian	$1 - \exp\left(-\left(\frac{h}{a}\right)^2\right)$	$h \geq 0$

Given a sample of high fidelity values, the relative performance of the CC and CLC methods will be established by measuring the number of additional low fidelity values required by each of the methods to achieve different error reduction percentages with respect to the error obtained using only the initial high fidelity sample. The number of additional high fidelity values that would be required to achieve similar error reductions will also be available.

The methodology was implemented using a well-known geostatistical software package named Gstat ([www.gstat.org](http://www.gstat.org)). Available since 1997, Gstat is an open source (GPL) computer code for multivariable geostatistical modeling, prediction and simulation. As of 2003, the Gstat functionally is also available as an S extension, either as an R package or S-Plus library. Details of Gstat can be found in Pebesma (2004).

## 5. Case studies

Analytical and industrial case studies are used to illustrate the proposed approach and to assess the relative performance of the cokriging methods under consideration. The analytical case study corresponds to a well-known optimization test function (Jin and Chen, 2000) denoted as F1, and the industrial case study refers to a modeling problem in the area of enhanced oil recovery.

### 5.1. Analytical test function F1

The function of interest is represented in Eq. (1). The domains of interest for the input variables  $x_1$  and  $x_2$  are given by the intervals (0.9), and, (0.6) respectively; the function range is 88.8.

$$F_1 = [30 + x_1 \sin(x_1)] [4 + \exp(-x_2)^2] \quad (1)$$

The equation above will represent the so called high fidelity model. The low fidelity model corresponds to a filtered version of the function denoted by Eq. (1) obtained using Wavelets Daubechies #4 (Daubechies, 1988). This filtering process is expected to preserve the basic features of the original function so that the high and low fidelity models are correlated. Fig. 5 illustrates both the high and low fidelity models.

The test set includes 4096 input/output pairs from a grid of  $64 \times 64$ .

### 5.2. Alkali-surfactant-polymer (ASP) modeling problem

The ASP enhanced oil recovery modeling problem addressed here is to build a surrogate model of a computationally expensive numerical simulator, that will take as input: surfactant (S) and polymer (P) concentrations, ASP slug size (expressed in the form of the injection time), and as output the cumulative oil production (NP) in bbls. The ranges of the input variables surfactant and polymer concentrations are given by  $0 \leq S \leq 0.005$  vol.fract. and  $0 \leq P \leq 0.10$  wt. %. The injection time is 194 days and the cumulative oil production is calculated at 800 days. As illustrated in Fig. 6, the ASP flooding pilot has an inverted five-spot pattern and a total of 5 vertical wells, 4 producers and 1 injector. The high and low fidelity models (Fig. 7) are associated with two different numerical grid resolutions, namely,  $18 \times 18 \times 3$  (range: 41,360 bbls), and,  $9 \times 9 \times 3$  (range: 27,850 bbls) in the x, y, and z directions, respectively. The cited ranges were calculated on a  $15 \times 15$  grid for the input variables (S and P). The numerical grid associated with the low fidelity model was obtained using an appropriate upscaling process.

The reservoir is at a depth of 4150 ft., has an average initial pressure of 1740 psi, and the porosity is assumed to be constant throughout the reservoir and equal to 0.3. The crude oil viscosity is 40 cp, the initial brine salinity is 0.0583 meq/ml and the initial brine divalent cation concentration is 0.0025 meq/ml. A summary of the

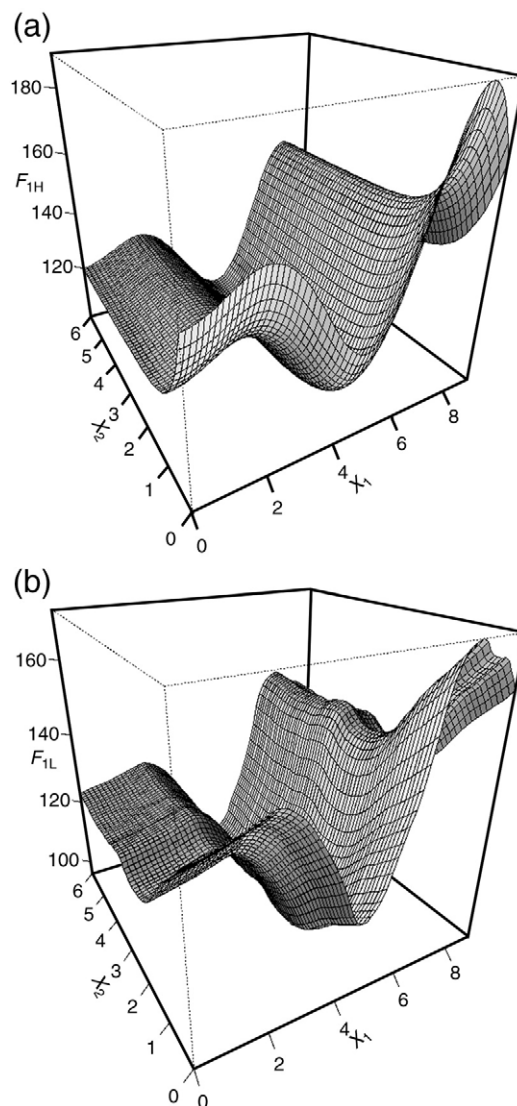


Fig. 5. High (a) and low (b) fidelity models — F1 case study.

reservoir and fluid properties is presented in Table 3. The injection scheme and other reference configuration details can be found in the sample data files of the UTCHEM program (UTCHEM, 2000).

The UTCHEM program is a three-dimensional, multiphase, multi-component reservoir simulator of chemical flooding processes

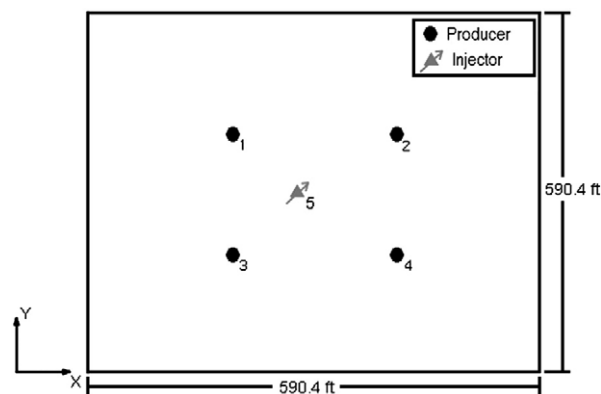


Fig. 6. Inverted five-spot in the ASP flooding pilot case study.

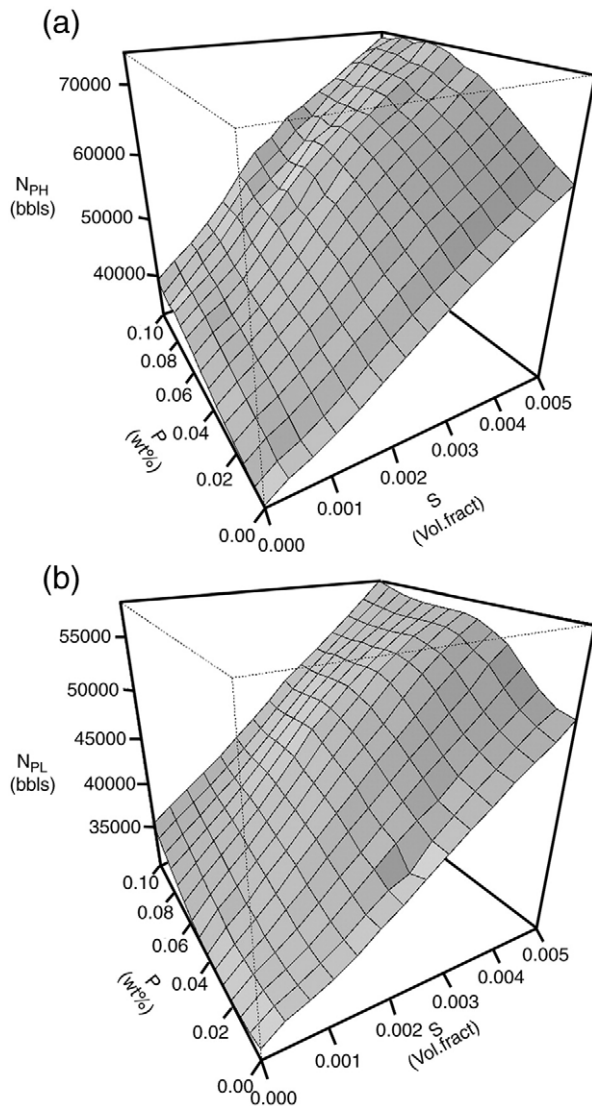


Fig. 7. High (a) and low (b) fidelity models — ASP case study.

developed at the University of Texas at Austin (Pope and Nelson, 1978; Engelsen et al., 1987; Lake et al., 1990). The basic governing differential equations consist of: a mass conservation equation for each component, an overall mass conservation equation that determines the pressure (the pressure equation), an energy balance, and Darcy's Law generalized for multiphase flow. The resulting flow equations are solved using a block-centered finite-difference scheme. The solution method is implicit in pressure and explicit in concentration, similar to the well-known IMPES method used in blackoil reservoir simulators. A Jacobi conjugate gradient method is used to

solve the system of finite-difference equations resulted from the discretization of the pressure equation.

Three flowing phases and eleven components are considered in the numerical simulations. The phases are water, oil and microemulsion, while the components are water, oil, surfactant, polymer, chloride anions, divalent cations ( $\text{Ca}^{2+}$ ,  $\text{Mg}^{2+}$ ), carbonate, sodium, hydrogen ion, and oil acid. The ASP interactions are modeled using the reactions: in situ generated surfactant, precipitation and dissolution of minerals, cation exchange with clay and micelle, and chemical adsorption. Note the detailed chemical reaction modeling, and the heterogeneous and multiphase petroleum reservoir under consideration.

The test set includes 225 input/output pairs from a grid of  $15 \times 15$ .

## 6. Results and discussion

Figs. 8 and 9 show the theoretical variograms associated with the high, low, and cross-covariance functions adjusted to the experimental variograms in the F1 and ASP modeling case studies, respectively. The experimental variograms were constructed using an extended sample to make sure the correlation models were properly identified; specifically, for the F1 (ASP modeling) case study 90 (60) and 80 (50) high and low fidelity sample sizes were used. In all instances a single model type and range were used to assure the positive definiteness of the covariance matrices and hence a proper solution of the cokriging optimization problems. The model type that provided the best fit was the spherical model; the sill and range values for all the theoretical variograms are shown in Table 4. As expected, the sills for the low and high fidelity samples are similar, and their arithmetic mean is higher than the sill for the cross-covariance model.

Figs. 10 and 11 show for the F1, and ASP modeling case studies, respectively, the relative performance of the CC and CLC methods by establishing the number of additional low fidelity values (for a given initial sample of high fidelity values) required by each of the methods to achieve a particular RMSE. The number of additional high fidelity values required to achieve similar RMSE is also shown. Additionally, Table 5 shows the number of additional high or low fidelity values necessary to reach a particular percentage reduction of the RMSE on a test data. It also shows the number of additional low fidelity values equivalent to a single high fidelity value in terms of percentage reduction of RMSE.

With reference to Figs. 10 and 11, and Table 5, all the cokriging models (CLC, CC) allow improving performance by integrating low fidelity samples to an existing high fidelity one. For example, a RMSE of six (6) in the F1 case study can be achieved by the integration of approximately thirty (30) low fidelity samples using CLC or CC-Kri, or approximately fifteen (15) additional samples if only high fidelity samples are used. Similarly, in the case of the ASP modeling case study a RMSE of five hundred and fifty (550) bbls can be achieved by the integration of approximately twenty two (22) low fidelity samples using CLC or CC-Kri, or approximately seventeen (17) additional samples if only high fidelity samples are used.

In terms of percentage reduction of RMSE (Table 5) observe that, in the F1 case study, when using CLC an eight percent (8%) reduction of the RMSE can be achieved by either eight (8) additional high fidelity values, or, eighteen (18) low fidelity ones. Hence, in this context, a high fidelity value is worth 2.25 times a low fidelity one, and, in general, the worthiness of high fidelity values decreases with increasing values of percentage reduction of the RMSE. The range of the worthiness of high fidelity values versus low fidelity ones was between 1.25 and 3.70. Estimates of the worthiness of high fidelity values can be useful when combining a small sample of data obtained from, for example, high fidelity/computationally expensive computer simulations, and, a larger one but with low fidelity values.

Additionally, note that for the case of low fidelity samples, after a certain sample size (problem dependent), no error reduction is observed (asymptotic behavior), which indicates that no significant

Table 3  
Reservoir and fluid properties. ASP modeling case study.

Property	Value	Unit
Reservoir depth	4150 (1265)	ft (m)
Oil viscosity	40	cp
Porosity	0.3	fraction
Average initial pressure	1740	psi
Well ratio	0.49 (15)	ft (m)
Skin factor	0	adim
Water salinity	$C_{Na}$ 0.0583	meq/ml
	$C_{Ca}$ 0.0025	meq/ml



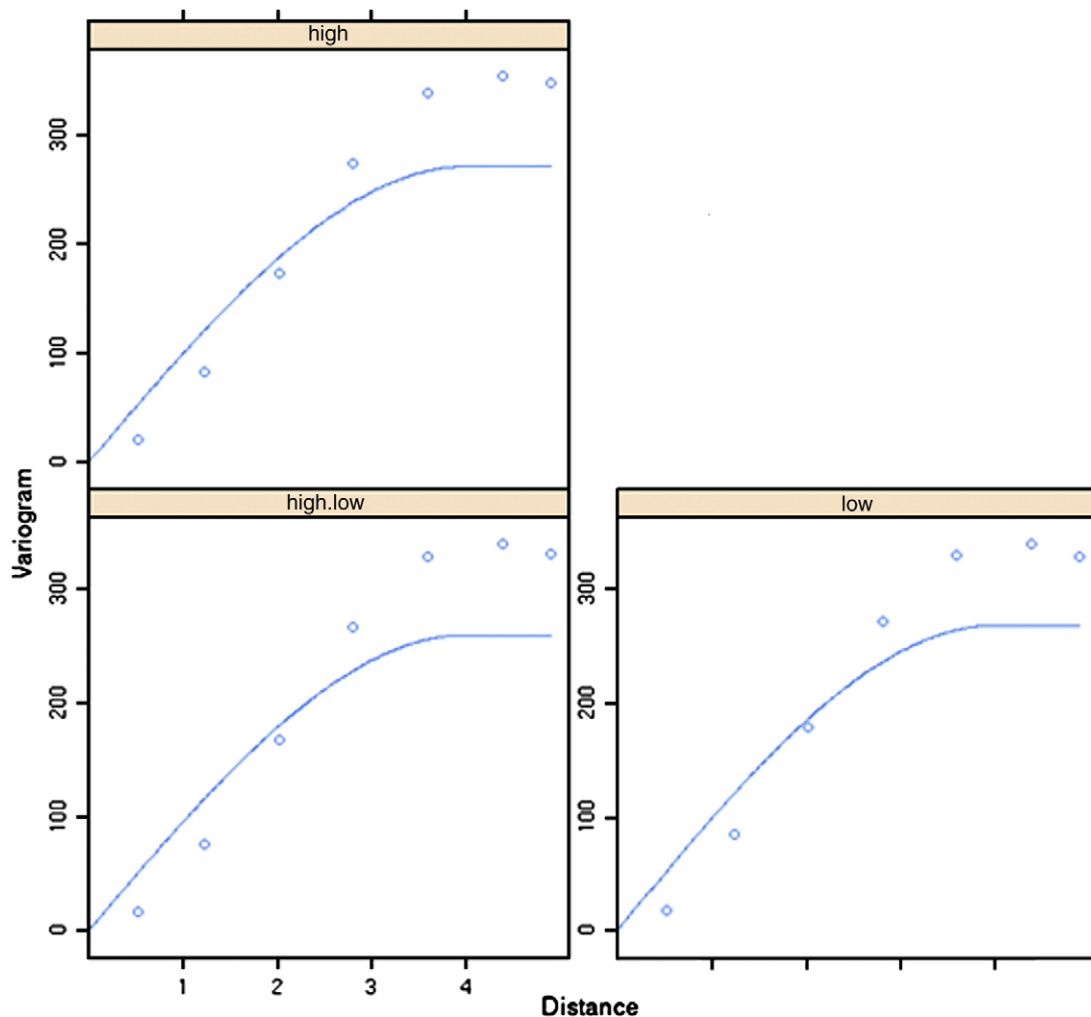


Fig. 8. Theoretical variogram adjusted to the experimental variogram – F1 case study.

modeling information is added. In general, though, this is not the case when incorporating high fidelity samples. It is also observed that the CC-Kri and CLC methods exhibited similar performance, outperforming in all instances, the CC-1nn method. For example, in the context of the ASP case study, a RMSE of five hundred and fifty (550) bbls can be achieved by the integration of approximately twenty two (22) low fidelity samples using CLC or CC-Kri, or thirty five (35) using CLC-1nn. The differences in the performance of CC-Kri vs. CC-1nn (more than 50% additional low fidelity samples required) can be explained by the latter method giving potentially inaccurate estimates when extending the low fidelity values to prediction sites.

The fact that CLC and CC-Kri exhibit similar performance is noteworthy since the former method requires the additional effort of estimating the covariance model associated with the low fidelity values; however, the performance of the latter can deteriorate if the low fidelity sample is too small to reasonably estimate the low fidelity values at prediction sites. Another consideration when selecting between the CLC and CC methods, is that if the sample size of the low fidelity variable is relatively high, the corresponding optimization problem can be significantly harder to solve in the case of the CLC method, although this latter issue can be overcome by using a restricted set of the low fidelity sample close to the prediction site.

For all case studies and cokriging approaches uncertainty estimates were consistent with amount and fidelity of the available data. For example, Figs. 12 and 13 display standard deviation estimates, and uncertainty reductions throughout the input space as the result of the addition of low fidelity samples for the F1 and ASP case studies,

respectively. In both cases, CLC is used and two scenarios are considered: only high fidelity simulations (10) are used (a), and both high (10) and low (10) fidelity simulations. The final picture (c) within the figures shows the uncertainty reduction as the result of the addition of low fidelity samples.

With references to Figs. 12(a) and 13(a) note that the standard deviation estimates goes from zero (white color) at the available data to the highest values of the scale in regions away from the available data. Also observe that the addition of low fidelity values (green circles) as shown in Figs. 12(b) and 13(b) translates into lower uncertainty values (whitening of previously blue regions). Figs. 12(c) and 13(c) depict that, as expected, the largest uncertainty reductions associated with the addition of low fidelity samples corresponds to regions where high fidelity values were not available.

## 7. Conclusions

Ideally, surrogate models should allow: a) the integration of variable fidelity samples, and, b) provide estimation and appraisal (error) information consistent with the amount and fidelity level of the available data. The branch of spatial statistics known as geostatistics offers considerable advantages when satisfying the above referenced requirements (a and b). This paper discussed the effectiveness and requirements of geostatistical methods such as classic cokriging and two variants of collocated cokriging for the integration of two level fidelity models; these methods can be shown to give unbiased and optimal estimates among linear models. Two



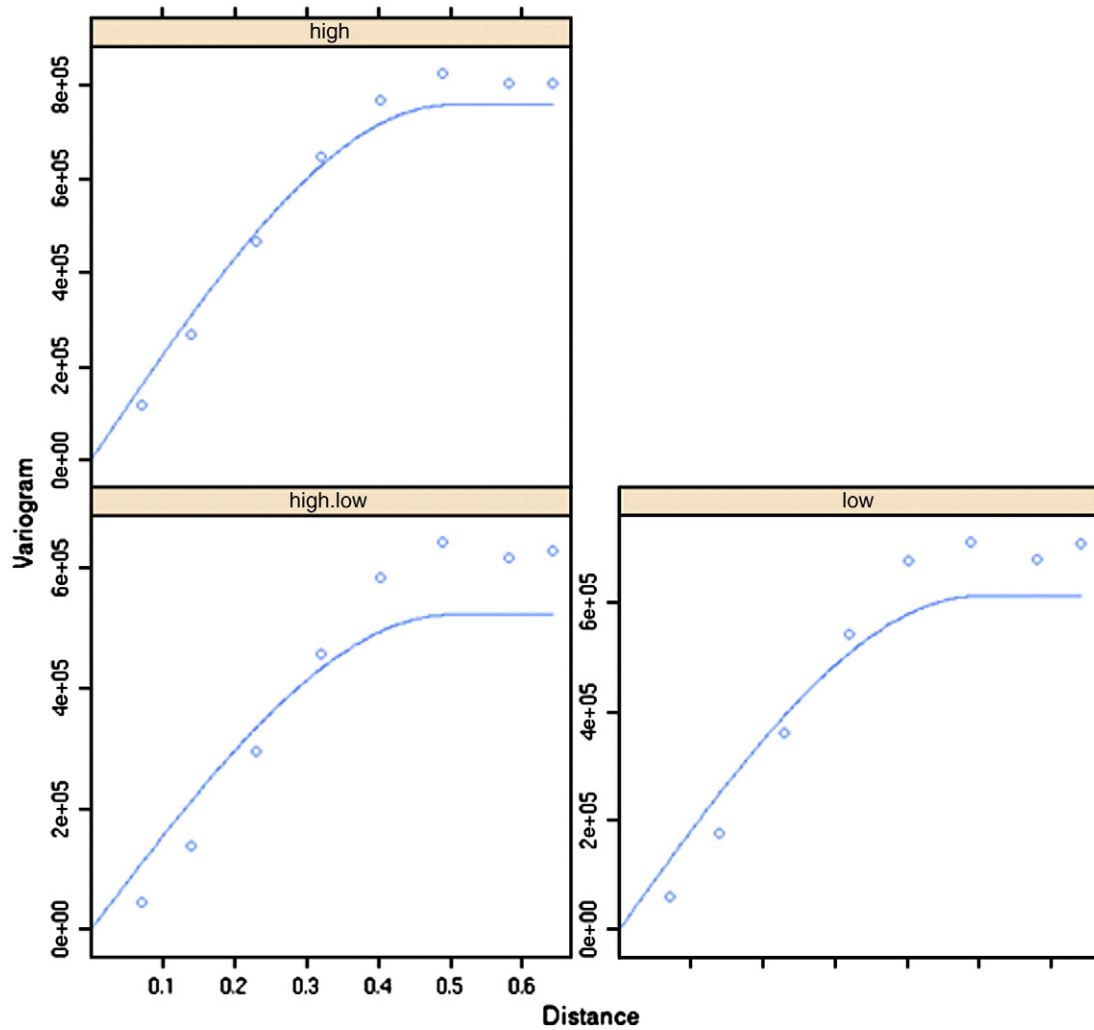


Fig. 9. Theoretical variogram adjusted to the experimental variogram — ASP case study.

case studies are considered: a well-known analytical function and, distinct resolution models, in the surrogate-based modeling of a field scale alkali–surfactant–polymer (ASP) enhanced oil recovery (EOR) process.

All cokriging methods considered allowed improving performance by integrating low fidelity samples to an existing high fidelity one. Classic cokriging offered the best overall performance but it has the drawback that it requires the specification of three covariance functions and the optimization formulation associated with predicting values may be hard to solve for large low fidelity sample sizes. As an alternative, the collocated cokriging (CC) approach does not need to specify the correlation model associated with the low fidelity values, and the optimization problem associated with the estimations are easier to solve. However, the CC approach require estimates of the

low fidelity variable at prediction sites; properly extending the low fidelity value to prediction sites is sensitive to the method used and sample size. For the latter task, the kriging method showed to be

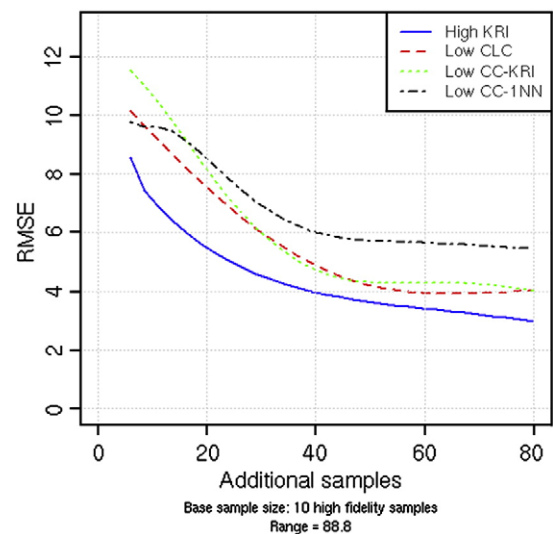


Fig. 10. Number of additional low fidelity samples required for a given RMSE on a test data set for the different cokriging models. The number of additional high fidelity samples to achieve a given RMSE is also shown (high KRI) — F1 case study.

Table 4

Sill and range values for the theoretical variograms identified in the F1 and ASP modeling case studies.

	Models	Sill	Range
F1	High	271.4	4
	Low	268.3	4
	Crossed	259.8	4
ASP	High	756,977.1	0.5
	Low	612,088.3	0.5
	Crossed	521,801.9	0.5

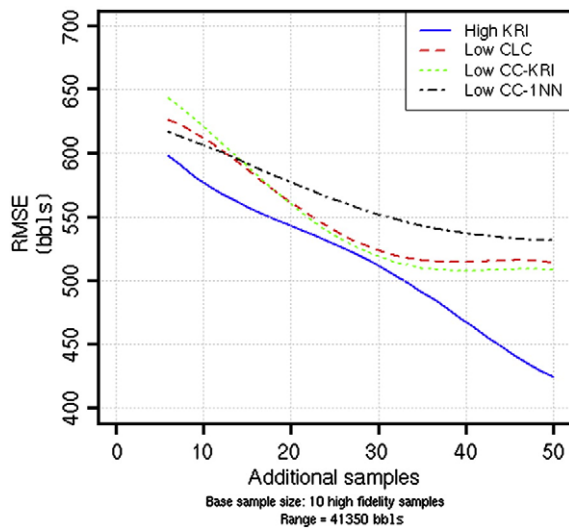


Fig. 11. Number of additional low fidelity samples required for a given RMSE on a test data set for the different cokriging models. The number of additional high fidelity samples to achieve a given RMSE is also shown (high KRI) — ASP case study.

much more effective than 1-nn for all samples sizes considered in this work. The classic and collocated cokriging (with kriging for extending the low fidelity values to prediction sites) showed similar performance with the latter offering predictions at a significantly lower computational cost.

The cokriging approaches uncertainty estimates were consistent with amount and fidelity of the available data with the greatest uncertainty reductions associated with the addition of low fidelity samples to regions where high fidelity values were not available.

The effectiveness of the approach in the context of modeling problems in higher dimensions, the accuracy of the error variance predictions, and practical specifications of the covariance models are the subject of current research efforts.

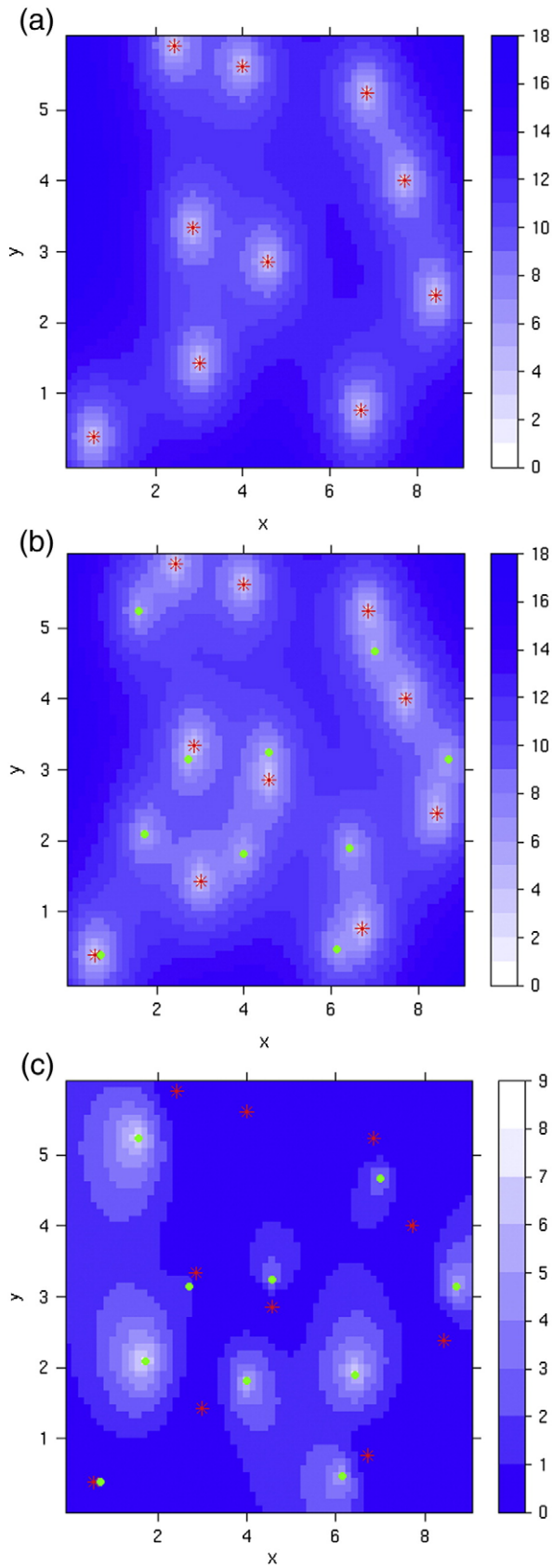
#### Nomenclature

KRI	Ordinary kriging
CLC	Classic cokriging
CC	Collocated cokriging
Cov	Covariance function
Var	Variance function
R	Correlation function
$\gamma$	Variogram function

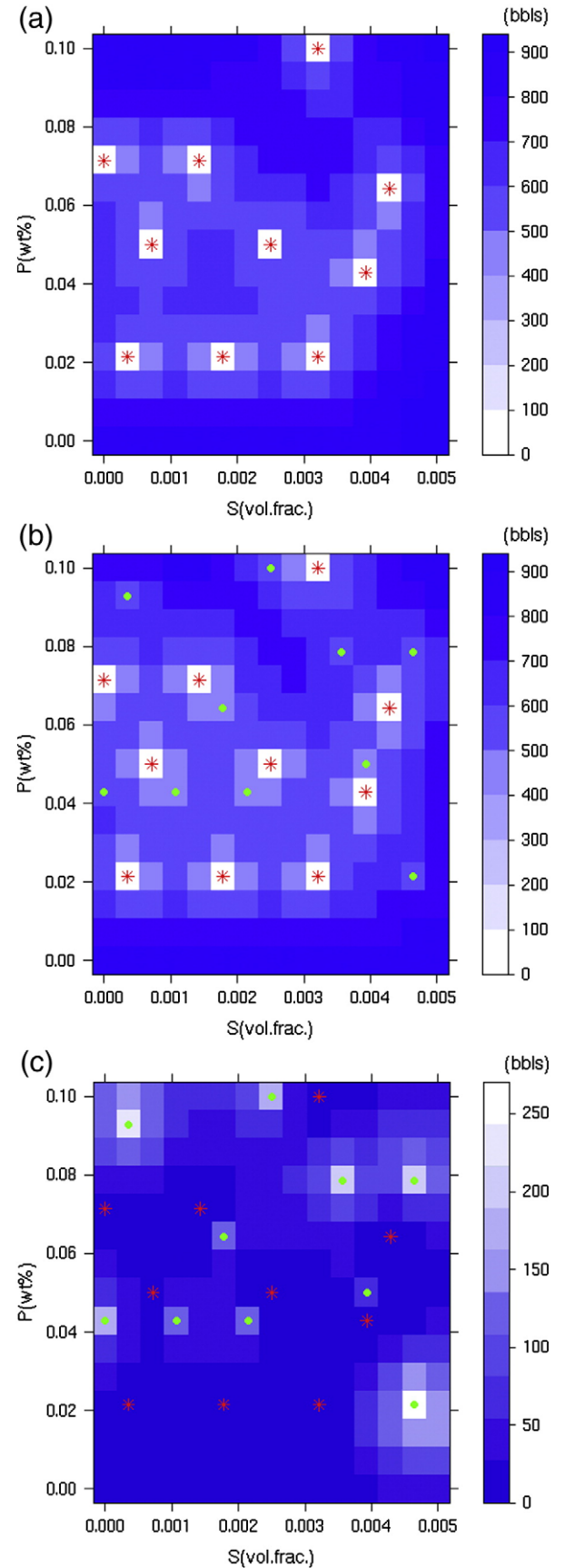
Table 5

The number of additional high or low fidelity values necessary to reach a particular percentage reduction of the RMSE on a test set. It also shows the number of additional low fidelity values equivalent to a single high fidelity value in terms of percentage reduction of RMSE. The initial high fidelity sample size was equal to ten (10).

Model:			CLC		CC-KRI		CC-1NN	
Case study	% of error reduction	High fidelity samples	Low fidelity samples	Worthiness	Low fidelity samples	Worthiness	Low fidelity samples	Worthiness
F1	2	7	15	2.14	20	2.86	21	3.00
	4	7	16	2.29	20	2.86	22	3.14
	6	8	17	2.13	21	2.63	23	2.88
	8	8	18	2.25	21	2.63	24	3.00
	10	8	19	2.38	22	2.75	25	3.13
	12	9	20	2.22	23	2.56	26	2.89
	14	9	21	2.33	24	2.67	27	3.00
	16	10	22	2.20	25	2.50	28	2.80
	18	10	24	2.40	25	2.50	30	3.00
	20	11	25	2.27	26	2.36	31	2.82
	22	11	26	2.36	27	2.45	32	2.91
	24	12	27	2.25	28	2.33	34	2.83
	26	13	28	2.15	29	2.23	36	2.77
	28	15	30	2.00	30	2.00	38	2.53
	30	16	31	1.94	31	1.94	40	2.50
	32	17	32	1.88	31	1.82	44	2.59
	34	18	33	1.83	32	1.78	60	3.33
	36	20	35	1.75	34	1.70	74	3.70
	38	22	36	1.64	35	1.59	–	–
	40	24	38	1.58	37	1.54	–	–
	42	25	39	1.56	38	1.52	–	–
ASP	44	27	41	1.52	40	1.48	–	–
	46	29	43	1.48	41	1.41	–	–
	48	31	45	1.45	44	1.42	–	–
	50	33	49	1.48	49	1.48	–	–
	52	37	52	1.41	76	2.05	–	–
	54	41	60	1.46	–	–	–	–
	56	45	–	–	–	–	–	–
	1	7	14	2.00	15	2.14	15	2.14
	2	9	17	1.89	17	1.89	19	2.11
	3	10	18	1.80	18	1.80	21	2.10
	4	11	19	1.73	19	1.73	24	2.18
	5	12	20	1.67	20	1.67	26	2.17
	6	14	21	1.50	21	1.50	28	2.00
	7	16	23	1.44	22	1.38	31	1.94
	8	18	24	1.33	24	1.33	34	1.89
	9	20	26	1.30	25	1.25	38	1.90
	10	22	28	1.27	27	1.23	43	1.95
	11	24	30	1.25	29	1.21	50	2.08
	12	26	33	1.27	31	1.19	–	–
	13	28	38	1.36	33	1.18	–	–
	14	29	47	1.62	36	1.24	–	–
	15	31	–	–	45	1.45	–	–
	16	33	–	–	–	–	–	–



**Fig. 12.** Standard deviation estimates (a, and b) and uncertainty reductions (c) throughout the input space as the result of the addition of low fidelity samples. The red stars represent the location of high fidelity samples, while the green circles depict the location of low fidelity samples (F1 case study).



**Fig. 13.** Standard deviation estimates (a, and b) and uncertainty reductions (c) throughout the input space as the result of the addition of low fidelity samples. The red stars represent the location of high fidelity samples, while the green circles depict the location of low fidelity samples (ASP case study).

F1	Analytical case study
ASP	Alkali–surfactant–polymer
EOR	Enhanced oil recovery
NP	Cumulative oil production
RMSE	Root mean square error
$\alpha, \beta$	Weights of the high and low fidelity samples
$Z, V$	High and low fidelity model samples
$\sigma_Z^2, \sigma_V^2$	High and low fidelity model sample variances
$\hat{z}_0$	High fidelity model estimate
$v_0$	Low fidelity model output at prediction sites

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