

Surrogate modeling-based optimization for the integration of static and dynamic data into a reservoir description

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Abstract

This paper presents a solution methodology for the inverse problem of estimating the distributions of permeability and porosity in heterogeneous and multiphase petroleum reservoirs by matching the static and dynamic data available. The solution methodology includes, the construction of a “fast surrogate” of an objective function whose evaluation involves the execution of a time-consuming mathematical model (i.e., reservoir numerical simulator) based on neural networks, DACE (design and analysis of computer experiment) modeling, and adaptive sampling. Using adaptive sampling, promising areas are searched considering the information provided by the surrogate model and the expected value of the errors. The proposed methodology provides a global optimization method, hence avoiding the potential problem of convergence to a local minimum in the objective function exhibited by the commonly Gauss–Newton methods. Furthermore, it exhibits an affordable computational cost, is amenable to parallel processing, and is expected to outperform other general-purpose global optimization methods such as, simulated annealing, and genetic algorithms. The methodology is evaluated using two case studies of increasing complexity (from 6 to 23 independent parameters). From the results, it is concluded that the methodology can be used effectively and efficiently for reservoir characterization purposes. In addition, the optimization approach holds promise to be useful in the optimization of objective functions involving the execution of computationally expensive reservoir numerical simulators, such as those found, not only in reservoir characterization, but also in other areas of petroleum engineering (e.g., EOR optimization). © 2002 Elsevier Science B.V. All rights reserved.

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1. Introduction

The identification of the permeability and porosity parameters that best match the data (static and dynamic) available for a given reservoir is critical for devising an optimal strategy for the development of oil and gas fields. The static data makes reference to those

originated from geology, electrical logs, core analysis, fluid properties, seismic and geostatistics; while the dynamic data is represented by field measurements such as, production history, bottom hole pressures from permanent gauges, water-cut, and gas/oil ratio.

Estimating permeability and porosity parameters from available data is difficult because of the following reasons: (i) in general, the number of parameters to be estimated are very high, since data is scarce, and the reservoirs are heterogeneous (permeability and porosity have spatial variability), (ii) the available

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data may have very different scope and nature, and (iii) the numerical simulation of the reservoir, necessary to assess how well given permeability and porosity parameters match the available data are computationally expensive.

This paper presents a solution methodology, called NEGO (neural network-based efficient global optimization), for the problem of estimating the distributions of permeability and porosity in heterogeneous and multiphase petroleum reservoirs by matching the static and dynamic data available. The solution methodology includes the construction of a “fast surrogate” of an objective function whose evaluation involves the execution of a time-consuming mathematical model (i.e., reservoir numerical simulator) based on neural networks, DACE (Sacks et al., 1989) modeling, and adaptive sampling. Using adaptive sampling, promising areas are searched considering the information provided by the surrogate model and the expected value of the errors.

The DACE surrogate model is initially constructed using sample data generated from the execution of mathematical models with parameters given by a Latin hypercube experimental design, and a neural network, and provides error estimates at any point. Additional points are obtained balancing the exploitation of the information provided by the surrogate model (where the surface is minimized) with the need to improve the surface (where error estimates are high). The proposed methodology provides a global optimization method, hence avoiding the potential problem of convergence to a local minimum in the objective function exhibited by the commonly used Gauss–Newton methods (Tan, 1995; Landa and Horne, 1997) and computational cost involved in numerically estimating derivatives, and in the step by step movement along given trajectories. Furthermore, it exhibits an affordable computational cost, is amenable to parallel processing, and is expected to outperform other general purpose global optimization methods such as, simulated annealing, and genetic algorithms (Huang and Kelkar, 1994; Datta Gupta et al., 1992).

2. Problem definition

The problem under consideration is an optimization problem (inverse parameter estimation) with

typically a high number of parameters and computationally expensive objective function evaluations. Formally, it can be written as:

$$\text{find } \mathbf{x} \in X \subseteq R^p$$

such that

$f(\mathbf{x})$ is minimized

where f is an objective function of \mathbf{x} , the permeability and porosity parameter vector, and X is the set constraint.

The objective function is a measure of the discrepancy between the data (static and dynamic) available and the response of the mathematical models using the current set of parameters. Eq. (1) shows a commonly used form of the objective function (weighted least square version):

$$f(\mathbf{x}) = (d_{\text{obs}} - d_{\text{calc}})' \mathbf{W} (d_{\text{obs}} - d_{\text{calc}}) \quad (1)$$

where \mathbf{W} is a weighting matrix, d_{obs} makes reference to static and dynamic data available (normalized), and d_{cal} denotes the corresponding data obtained using a mathematical model.

Hence, the problem of interest is one of finding the vector of parameters \mathbf{x} that minimizes the difference between the available data (d_{obs}), and the values calculated (d_{cal}) substituting \mathbf{x} in the appropriate mathematical model. Note that a reservoir numerical simulator is necessary for calculating the response associated with dynamic data (e.g., production history). As mentioned before, the reservoir numerical simulator, in general, is computationally expensive and the number of elements in \mathbf{x} is usually high, since the data is scarce and the reservoirs are heterogeneous. These two issues place restrictions on the solution approach, given that the number of objective function evaluations are limited to a relatively low value considering the time restrictions typically present in the oil industry.

3. Solution methodology

The proposed solution approach called NEGO, neural network-based efficient global optimization, is an improved version of the EGO algorithm (Jones et al., 1998) for the optimization of computationally expensive black-box functions.

The proposed solution methodology involves the following five steps.

(1) Construct a sample of the parameter space using the Latin hypercube method. The Latin hypercube sampling procedure has been shown to be very effective for selecting input variables for the analysis of the output of a computer code (McKay et al., 1979).

(2) Conduct mathematical simulations using the sample from the previous step and record the response values associated with static and dynamic data available, and the objective function values.

(3) Construct a parsimonious neural network using the data from the previous step. The purpose of this neural network is to capture the general trends observed in the data; no rigorous performance criterion is placed on the neural network. The neural network model used in this study is the so-called multilayer (three layers) perceptron. The number of neurons in the input and output layers are given by the number of input (porosity and permeability parameters) and output variables (objective function value). The number of neurons in the hidden layer is settled by limiting the number of parameters of the model (weights) to be a fraction of the total number of data points available to the learning process. This process makes reference to the identification of a set of weights that, for a given architecture, minimize the sum of the square of the model errors; the errors represent the difference between the neural network model output value and the objective function value. The learning algorithm used is a gradient-based optimization procedure called

Backpropagation (Rumelhart and McClelland, 1986). For an introduction to neural network modeling, see, for example, Bishop (1995).

(4) The input variables of the neural network are the permeability and porosity parameters and the output variable is the corresponding objective function value.

(5) Construct a DACE model for the residuals, that is, the difference between the observed objective function values, and the neural network responses using the sample data. These models provide not only estimates of the residuals values but also of the respective errors. The surrogate model for the evaluation of the objective function is the sum of the neural network and DACE models. Details of this step will be given later in this section.

(6) Additional points are obtained balancing the exploitation of the information provided by the surrogate model (where the surface is minimized) with the need to improve the surface (where error estimates are high), until a stopping criteria has been met. This balance is achieved by sampling where a figure of merit is maximized. Details of the figure of merit will be given later in this section.

3.1. DACE models

These models owe their name, design and analysis of computer experiments, to the title of an article that popularized the approach (Sacks et al., 1989). This approach differs from Kriging (in the geostatistics

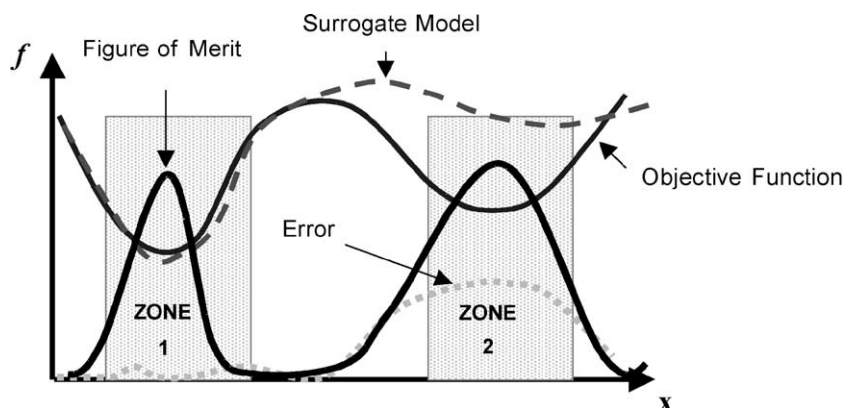


Fig. 1. Illustration of the purpose of the figure of merit.

Table 1
Test function (without noise) results for the NEGO algorithm

Test function	Dimensions	NIO	OFE	RE (%)	$ \mathbf{x} - \mathbf{x}_{\text{opt}} $
Branin-Hoo	2	20	42	0.35	0.0161
Hock-Shittkowski 5	2	20	25	0.08	0.0371
HGO 468:1	2	20	22	0.22	0.0023
HGO 468:2	2	20	22	0.021	0.0173
Six-Hump Camel	2	29	59	1.69	0.0476
Hartman 3	3	30	36	0.11	0.0110
Hartman 6	6	72	85	4.08	1.1151

literature; see, for example, Matheron, 1963; Cressie, 1991) in that, the latter, is limited to two or three dimensions, and does not have a fixed correlation structure.

These models suggest to estimate deterministic functions as shown in Eq. (2).

$$y(\mathbf{x}_j) = \mu + \varepsilon(\mathbf{x}_j) \quad (2)$$

where y is the function to be modeled, μ is the mean of the population, and ε is the error with zero expected value, and with a correlation structure given by Eq. (3).

$$\text{cov}(\varepsilon(\mathbf{x}_i), \varepsilon(\mathbf{x}_j)) = \sigma^2 \exp \left(- \sum_{h=1}^p \theta_h \left(\mathbf{x}_i^h - \mathbf{x}_j^h \right)^2 \right) \quad (3)$$

where p denotes the number of dimensions in the vector \mathbf{x} , σ identifies the standard deviation of the population, and θ_h is a correlation parameter, which is a measure of the degree of correlation among the data along the h direction.

Specifically, given a set of n input/output pairs (x, y) , the parameters μ , σ , and θ are estimated such that the likelihood function is maximized (Sacks et

al., 1989). Having estimated these values, the function estimate for new points is given by Eq. (4).

$$\bar{y}(\mathbf{x}) = \bar{\mu} + \mathbf{r}' \mathbf{R}^{-1} (\mathbf{y} - L \bar{\mu}) \quad (4)$$

where the line above the letters denote *estimates*, \mathbf{r}' identifies the correlation vector between the new point and the points used to construct the model, \mathbf{R} is the correlation matrix among the n sample points, and L denotes an n -vector of ones.

The mean square error of the estimate is given by Eq. (5).

$$s^2(\mathbf{x}^*) = \sigma^2 \left[1 - \mathbf{r}' \mathbf{R}^{-1} \mathbf{r} + \frac{(1 - L' \mathbf{R}^{-1} \mathbf{r})}{L' \mathbf{R}^{-1} L} \right] \quad (5)$$

The model is validated through a cross-validation procedure, that essentially makes sure that the estimates using all but the point being tested and the actual response values are within a specified number of standard deviations. The original EGO algorithm may not cross-validate properly if there are trends in the data, in contrast to NEGO, which is expected to subtract any significant trends in the data.

The benefits of modeling deterministic functions using this probabilistic approach are: (i) represents a best linear unbiased estimator, (ii) interpolates the data, and (iii) provides error estimates.

3.2. Figure of merit

With reference to Fig. 1, there are two zones where it is desirable to add additional points. The zone (left) where the objective function is minimized and the zone (right) where there is a significant error in the prediction. Hence, the figure of merit for adding sample points should be high in either of these

Table 2
Test function (with noise) results for the NEGO algorithm ($f_{\eta} = f(1 + \eta * \text{randn}())$)

Test function (f)	η	Dimensions	NIO	OFE	$ \mathbf{x} - \mathbf{x}_{\text{opt}} _{\eta}$	$ \mathbf{x} - \mathbf{x}_{\text{opt}} $
Branin-Hoo	0.001	2	20	35	0.2459	0.0161
Branin-Hoo	0.010	2	20	39	0.2242	0.0161
Branin-Hoo	0.100	2	20	50	0.0616	0.0161
Hartman 3	0.001	3	30	42	0.1195	0.0110
Hartman 3	0.010	3	30	36	0.0822	0.0110
Hartman 3	0.100	3	30	60	0.0791	0.0110

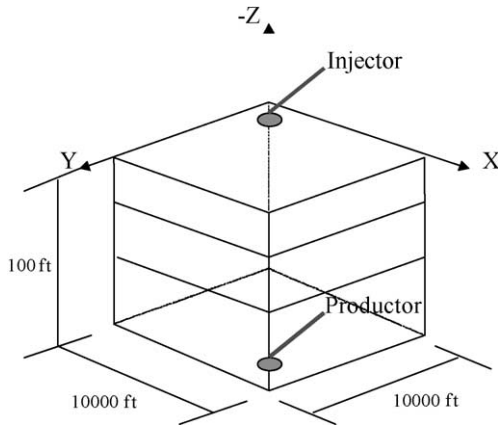


Fig. 2. Reservoir illustration (Case Study No. 1).

situations. Specifically, the figure of merit (Jones et al., 1998) used in this work, is given by Eq. (6).

$$\text{fom}(\mathbf{x}) = (f_{\min} - \hat{f})\Phi\left(\frac{f_{\min} - \hat{f}}{s}\right) + s\phi\left(\frac{f_{\min} - \hat{f}}{s}\right) \quad (6)$$

where Φ and ϕ are the cumulative and density normal distribution functions, respectively; and f_{\min} denotes the minimum current objective function value. Eq. (6) establishes the desired balance of sampling where the response surface (the predictor)

Table 3

Permeability and porosity parameters (Case Study No. 1)

Variable	Description	Range	Units
K_1	Permeability in the x and y directions	450–650	md
K_2	Permeability in the x and y directions	35–55	md
K_3	Permeability in the x and y directions	160–240	md
K_4	Permeability in the z direction	52–82	md
K_5	Permeability in the z direction.	22–32	md
	Layers		
P	Porosity. Layers 1, 2	0.25–0.37	–

is minimized (left term) and in zones where error estimates are high (right term). Note that the figure of merit makes reference to the objective function so it includes the sum of the output of both the neural network and the residual models.

This surface response approach for global optimization is expected to outperform competing methods, in terms of necessary computationally expensive objective function evaluations, to meet a stopping criteria. It can identify promising areas without the need of moving step by step along a given trajectory. In addition, by providing estimates of the errors at unsampled points, it is possible to establish a reasonable stopping criterion. Furthermore, provides a fast surrogate model that could be used to visualize the relationship between the sought parameters and the

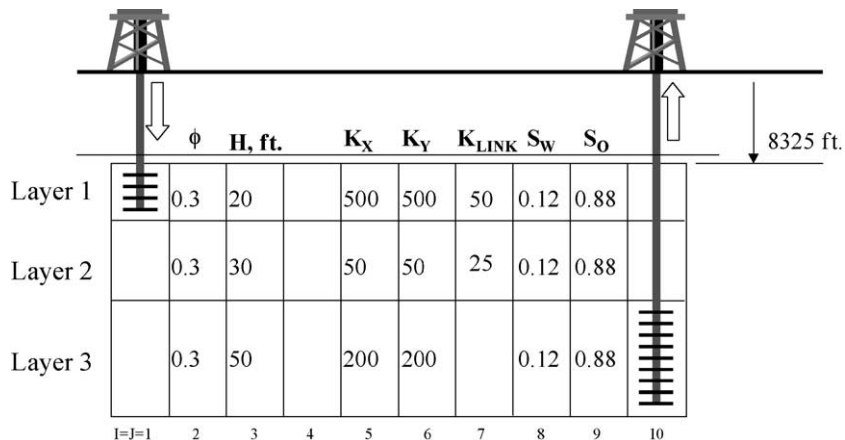


Fig. 3. Schematic representation of the reservoir considered in (Case Study No. 1).

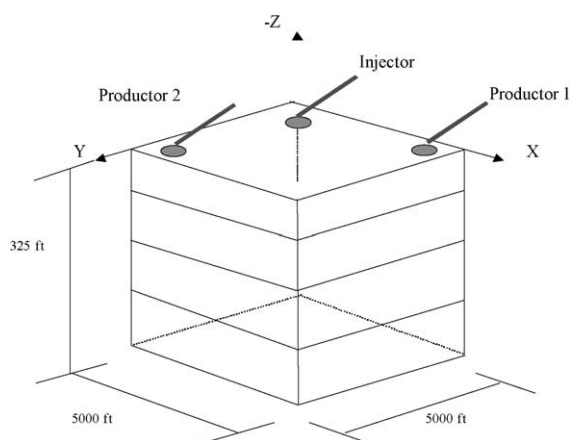


Fig. 4. Reservoir illustration (Case Study No. 2).

objective function values and to identify the relative significance of each of the parameters.

3.3. Implementation and validation

The following case studies were solved using an implementation of the NEGO algorithm developed by the authors (NQ, SP, and JC) in MATLAB (MATLAB, The MathWorks). The subproblems of finding near optimal values for maximizing likelihood and the figure of merit were solved using the DIRECT method (Jones et al., 1993). Note that the solution of these subproblems do not require additional computationally expensive objective function evaluations. The

reservoir numerical simulations were conducted using a commercial reservoir numerical simulator EXODUS (EXODUS, T.T. & Associates).

For validation purposes, the NEGO algorithm was tested using a set of seven standard test functions (Björkman and Holmström, 1999; Jones et al., 1993, 1998). The results (see Table 1) show that NEGO is highly competitive. In all cases, it found solutions with a relative error lower than 5%. This is remarkable considering the very limited number of objective function evaluations used. Furthermore, test function (with noise) results for the NEGO algorithm (Table 2) show that for significant levels of noise ($\eta=0.001, 0.01, 0.1$), the algorithm is still able to identify the optimum neighborhood.

4. Case study

The NEGO algorithm was evaluated using two case studies of increasing complexity (from 6 to 23 independent parameters). The case studies consider reservoirs similar to those found in well-known benchmark cases (Aziz, 1981; Quandle, 1993). In both instances, the dynamic and static data were obtained assuming the permeability and porosity parameters were known (later called “correct” values). Then, the problems were posed in inverse fashion; that is, given dynamic and static data associated with a reservoir; what are the parameter

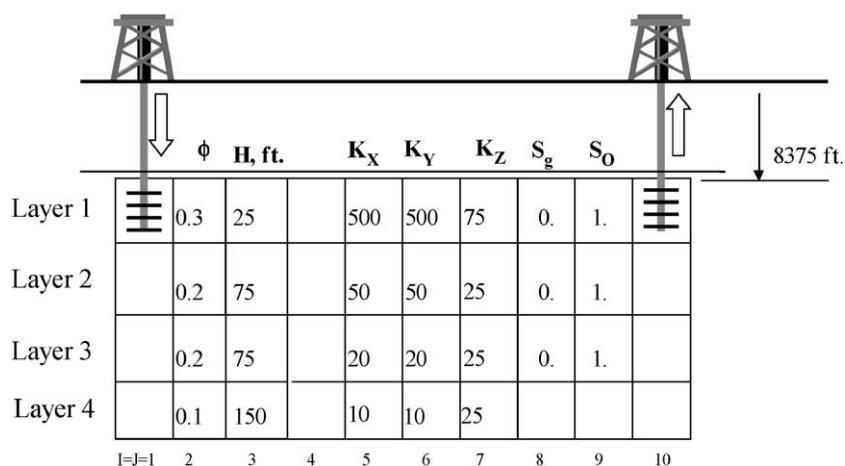


Fig. 5. Schematic representation of the reservoir considered in (Case Study No. 2).

	1	2	3	4	5	6	7	8	9	10
1	1.	1.	1.	1.	1.	1.	1.	3.	3.	3.
2	1.	1.	1.	1.	1.	1.	1.	3.	3.	3.
3	1.	1.	1.	1.	2.	2.	2.	3.	3.	3.
4	1.	1.	1.	1.	2.	2.	2.	2.	4.	4.
5	1.	1.	1.	1.	2.	2.	2.	2.	4.	4.
6	3.	3.	3.	2.	2.	2.	2.	2.	4.	4.
7	3.	3.	3.	2.	2.	2.	2.	2.	4.	4.
8	3.	3.	3.	5.	5.	5.	5.	4.	4.	4.
9	3.	3.	3.	5.	5.	5.	5.	4.	4.	4.
10	3.	3.	3.	5.	5.	5.	5.	4.	4.	4.

Fig. 6. Zones 1–5 in layer 1 (Case Study No. 2).

values that reproduce the available dynamic and static data?

The first case study addresses the integration of dynamic data (cumulative oil production and gas/oil ratio), while the second case considers the integration of both dynamic (cumulative oil production and gas/oil ratio) and static data (a variogram model). In both cases, the production data is available yearly, for a

period of 10 years, and in the second case study, the covariance is calculated using 10 intervals.

4.1. Case Study No. 1

The reservoir under consideration and the coordinate system used, are illustrated in Fig. 2. It is assumed that production data (i.e., COP and GOR)

	1	2	3	4	5	6	7	8	9	10
1	6.	6.	6.	6.	6.	6.	6.	8.	8.	8.
2	6.	6.	6.	6.	6.	6.	6.	8.	8.	8.
3	6.	6.	6.	6.	7.	7.	7.	8.	8.	8.
4	6.	6.	6.	6.	7.	7.	7.	7.	9.	9.
5	6.	6.	6.	6.	7.	7.	7.	7.	9.	9.
6	8.	8.	8.	7.	7.	7.	7.	7.	9.	9.
7	8.	8.	8.	7.	7.	7.	7.	7.	9.	9.
8	8.	8.	8.	10.	10.	10.	10.	9.	9.	9.
9	8.	8.	8.	10.	10.	10.	10.	9.	9.	9.
10	8.	8.	8.	10.	10.	10.	10.	9.	9.	9.

Fig. 7. Zones 6–10 in layer 2 (Case Study No. 2).

	1	2	3	4	5	6	7	8	9	10
1	11.	11.	11.	11.	11.	11.	11.	13.	13.	13.
2	11.	11.	11.	11.	11.	11.	11.	13.	13.	13.
3	11.	11.	11.	11.	12.	12.	12.	13.	13.	13.
4	11.	11.	11.	11.	12.	12.	12.	12.	14.	14.
5	11.	11.	11.	11.	12.	12.	12.	12.	14.	14.
6	13.	13.	13.	12.	12.	12.	12.	12.	14.	14.
7	13.	13.	13.	12.	12.	12.	12.	12.	14.	14.
8	13.	13.	13.	15.	15.	15.	15.	14.	14.	14.
9	13.	13.	13.	15.	15.	15.	15.	14.	14.	14.
10	13.	13.	13.	15.	15.	15.	15.	14.	14.	14.

Fig. 8. Zones 11–15 in layer 3 (Case Study No. 2).

are available and that certain permeability and porosity parameters are unknown.

With reference to Fig. 3, the reservoir is at a depth of 8325 ft., has an initial pressure of 4800 psi, and initial oil and water saturations of 0.8, and 0.2, respectively. The numerical grid is composed of $10 \times 10 \times 3$ blocks in the x , y and z directions. The injector and producer wells are placed in the blocks denoted as (10,10,3) and (1,1,1), respectively. The porosity is assumed to be constant throughout the reservoir, the horizontal permeability is isotropic, but, as the vertical permeability, is different for each of the layers.

In this case, the unknown parameters and the restrictions on their possible values are presented in Table 3, and the objective function is given by Eq. (7).

$$f(\mathbf{x}) = \sum_{i=1}^{10} \left[\frac{\text{COP}_{\text{obs}_i} - \text{COP}_{\text{calc}_i}}{\text{COP}_{\text{obs}_i}} \right]^2 + \sum_{i=1}^{10} \left[\frac{\text{GOR}_{\text{obs}_i} - \text{GOR}_{\text{calc}_i}}{\text{GOR}_{\text{obs}_i}} \right]^2 \quad (7)$$

The neural network and DACE models were constructed using a sample of 60 points selected using a

	1	2	3	4	5	6	7	8	9	10
1	16.	16.	16.	16.	16.	16.	16.	18.	18.	18.
2	16.	16.	16.	16.	16.	16.	16.	18.	18.	18.
3	16.	16.	16.	16.	17.	17.	17.	18.	18.	18.
4	16.	16.	16.	16.	17.	17.	17.	17.	19.	19.
5	16.	16.	16.	16.	17.	17.	17.	17.	19.	19.
6	18.	18.	18.	17.	17.	17.	17.	17.	19.	19.
7	18.	18.	18.	17.	17.	17.	17.	17.	19.	19.
8	18.	18.	18.	20.	20.	20.	20.	19.	19.	19.
9	18.	18.	18.	20.	20.	20.	20.	19.	19.	19.
10	18.	18.	18.	20.	20.	20.	20.	19.	19.	19.

Fig. 9. Zones 16–20 in layer 4 (Case Study No. 2).

Table 4
Permeability and porosity parameters (Case Study No. 2)

Variable	Description	Range	Units
K_{11}	Permeability zone 1, Layer 1	432–624	md
K_{12}	Permeability zone 2, Layer 1	345–525	md
K_{13}	Permeability zone 3, Layer 1	400–600	md
K_{14}	Permeability zone 4, Layer 1	340–480	md
K_{15}	Permeability zone 5, Layer 1	376–540.5	md
K_{21}	Permeability zone 1, Layer 2	43.2–62.4	md
K_{22}	Permeability zone 2, Layer 2	31.5–49.5	md
K_{23}	Permeability zone 3, Layer 2	40–60	md
K_{24}	Permeability zone 4, Layer 2	34–48	md
K_{25}	Permeability zone 5, Layer 2	37.6–54.05	md
K_{31}	Permeability zone 1, Layer 3	16.2–23.4	md
K_{32}	Permeability zone 2, Layer 3	11.2–17.6	md
K_{33}	Permeability zone 3, Layer 3	16–24	md
K_{34}	Permeability zone 4, Layer 3	12.75–18	md
K_{35}	Permeability zone 5, Layer 3	13.6–19.55	md
K_{41}	Permeability zone 1, Layer 4	8.1–11.7	md
K_{42}	Permeability zone 2, Layer 4	4.9–7.7	md
K_{43}	Permeability zone 3, Layer 4	8–12	md
K_{44}	Permeability zone 4, Layer 4	4.25–6	md
K_{45}	Permeability zone 5, Layer 4	6.4–9.2	md
P_1	Porosity. Layer 1	0.27–0.39	–
$P_{2,3}$	Porosity. Layers 2 and 3	0.14–0.22	–
P_4	Porosity. Layer 4	0.08–0.12	–

Latin hypercube sampling procedure. Twenty additional points were added in the search of the “correct” permeability and porosity parameters.

4.2. Case study No. 2

The reservoir under consideration and the coordinate system used, are illustrated in Fig. 4. It is assumed that dynamic data (i.e., COP and CGOR) and static data (variogram model) are available and that certain permeability and porosity parameters are unknown.

With reference to Fig. 5, the reservoir is at a depth of 8375 ft, has an initial pressure of 4800 psi, and initial oil and gas saturations of 1.0, and 0.0, respectively. The numerical grid is composed of $10 \times 10 \times 4$ blocks in the x , y and z directions. The producer wells 1 and 2, are placed in the blocks denoted by (10,1,1), and (1,10,1), respectively. The injector well is placed in the block (1,1,1). The porosity varies among layers with 0.30 for layer 1, 0.20 for layers 2 and 3, and 0.10 for layer 4. The numerical grid was grouped into 20 zones, distrib-

uted among the different layers as depicted in Figs. 6–9.

Permeability is considered to be the same within a given zone and does not change with coordinate direction (isotropic).

In this case, the unknown parameters and the restrictions on their possible values are presented in Table 4, and the objective function is given by Eq. (8).

$$f(\mathbf{x}) = \sum_{i=1}^{10} \left[\frac{\text{COP}_{\text{obs}_i} - \text{COP}_{\text{calc}_i}}{\text{COP}_{\text{obs}_i}} \right]^2 + \sum_{i=1}^{10} \left[\frac{\text{CGOR}_{\text{obs}_i} - \text{CGOR}_{\text{calc}_i}}{\text{CGOR}_{\text{obs}_i}} \right]^2 + \sum_{i=1}^{10} \left[\frac{\text{COV}_{\text{obs}_i} - \text{COV}_{\text{calc}_i}}{\text{COV}_{\text{obs}_i}} \right]^2 \quad (8)$$

The neural network and DACE models were constructed using a sample of 187 points selected using a Latin hypercube sampling procedure. Eighty addi-

Table 5
Additional sampled points (Case Study No. 1)

Obs.	K_1	K_2	K_3	K_4	K_5	P	f
1	483.33	51.67	200	67	30.33	0.31	0.00037604
2	550	38.33	226.67	57	23.67	0.31	0.00373747
3	616.67	45	173.33	67	27	0.31	0.0420268
4	483.33	38.33	226.67	67	27	0.35	0.08506395
5	616.67	51.67	200	57	27	0.35	0.02255316
6	483.33	38.33	226.67	67	27	0.31	0.02850679
7	550	51.67	200	57	23.67	0.31	0.04083019
8	483.33	45	200	77	23.67	0.31	0.00034579
9	550	38.33	173.33	57	27	0.35	0.00852373
10	550	38.33	226.67	77	30.33	0.31	0.01906443
11	616.67	38.33	200	67	27	0.35	0.01527427
12	483.33	51.67	226.67	67	27	0.31	0.06371478
13	483.33	38.33	200	57	30.33	0.31	0.00348125
14	550	38.33	200	57	30.33	0.35	0.05410506
15	550	38.33	173.33	67	23.67	0.35	0.00800637
16	483.33	45	226.67	57	30.33	0.31	0.03427427
17	483.33	45	173.33	67	30.33	0.31	0.07724744
18	550	51.67	226.67	67	30.33	0.31	0.00908735
19	616.67	45	200	57	23.67	0.35	0.01946102
20	550	45	200	57	23.67	0.35	0.06582813

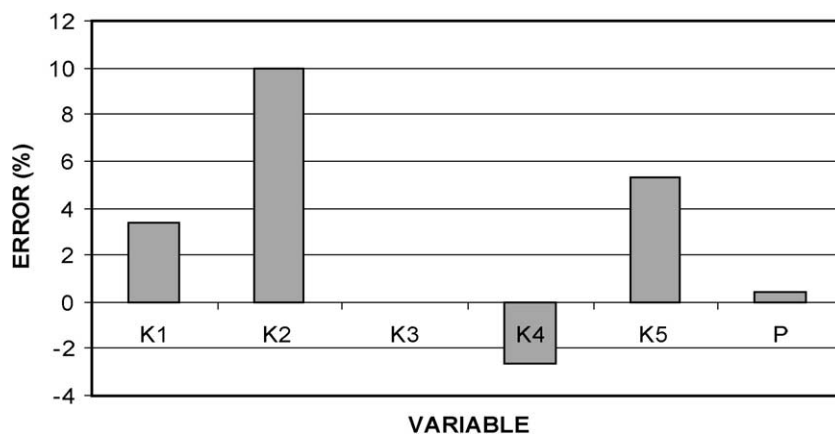


Fig. 10. Distribution of the error in parameter estimations (Case Study No. 1).

tional points were added in the search of the “correct” permeability and porosity parameters.

5. Results and discussion

With reference to Case study No. 1, the neural network was constructed with an $6 \times 2 \times 1$ architecture with a sum of square errors of $5.82e-002$, and $9.68e-002$, respectively. Additionally, all the points

in the DACE model cross-validated within three times of the standard deviation. The minimum objective function value found within the initial sample (60 points) was $2.17e-03$. Additional points (20) were added so that the figure of merit was maximized; from those points the best solution found (8th additional sampled point) observed an objective function value of $3.46e-04$, that is an order of magnitude lower than the best found in the initial sample. The parameter values and the objective function value for the addi-

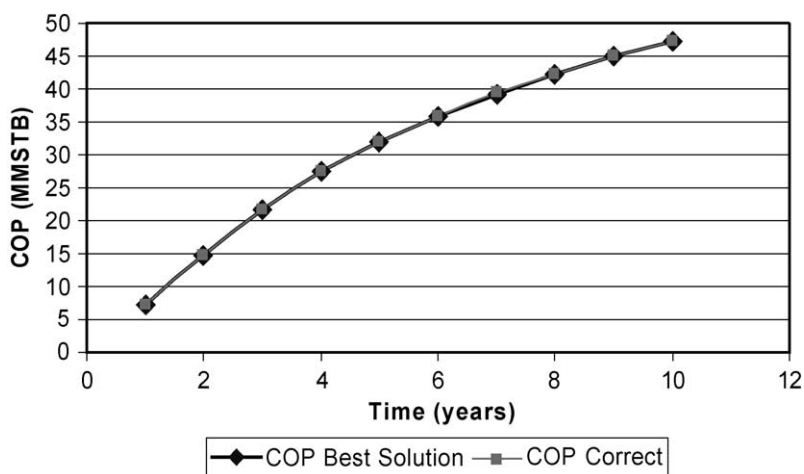


Fig. 11. Cumulative oil production obtained using the best solution and the “correct” solution (Case Study No. 1).

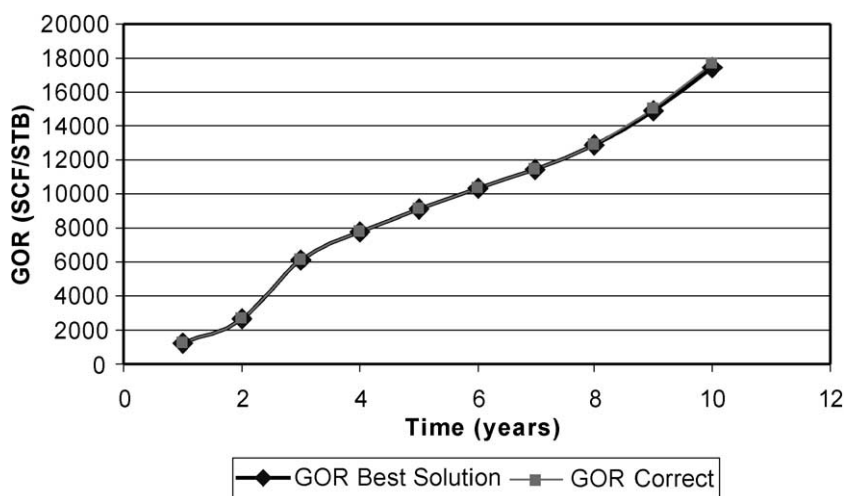


Fig. 12. Cumulative gas oil obtained using the best solution and the “correct” solution (Case Study No. 1).

tional points are shown in Table 5. The maximum percentage error in the parameters estimation (K_1 , K_2 , K_3 , K_4 , K_5 , P) is 10%, as illustrated in Fig. 10. The maximum percentage errors in the estimation of COP and GOR were 0.18% and 0.45%, respectively. Figs. 11 and 12 shows the excellent agreement between the values of COP and GOR obtained using the “correct” parameters and those found by the NEGO algorithm. Note that the results were obtained using only 80 computationally expensive objective function evaluations.

With reference to Case study No. 2, the neural network was constructed with a $23 \times 3 \times 1$ architecture with a sum of square errors of $4.97e-002$ and $9.05e-002$, respectively. A 98% of the points in the DACE model cross-validated within three times of the standard deviation. The minimum objective function value found within the initial sample (187 points) was $2.16e-02$. Additional points (80) were added so that the figure of merit was maximized; from those points the best solution found (56th additional sampled point) observed an objective function value of $1.37e-02$, that is approximately 50% lower than the best found in the initial sample. The parameter values and the objective function value for the additional points are shown in Table 6. The maximum percentage error in the parameters estimation is approximately 20%, as illustrated in Fig. 13. The maximum percentage errors in the estimation of COP

and CGOR, and COV were 0.79%, 6.81%, and 8.09%, respectively. Figs. 14 and 15 shows the excellent agreement between the values of COP and CGOR

Table 6
NEGO estimates for the permeability and porosity parameters (Case Study No. 2)

Variable	“Correct” value	Bis	Error (%)	Best solution	Error (%)
K_{11}	480	474.47	1.152	592	23.333
K_{12}	450	402.384	10.581	435	3.3333
K_{13}	500	472	5.6	500	0
K_{14}	400	361.882	9.529	363.333	9.166
K_{15}	470	495.345	5.392	403.417	14.1667
K_{21}	48	46.577	2.964	46.4	3.3333
K_{22}	45	42.730	5.044	40.5	10
K_{23}	50	45.972	8.056	50	0
K_{24}	40	46.412	16.031	41	2.5
K_{25}	47	49.709	5.764	45.825	2.5
K_{31}	18	20.038	11.32	19.8	10
K_{32}	16	12.418	22.384	14.4	10
K_{33}	20	23.607	18.036	20	0
K_{34}	15	17.219	14.799	15.375	2.5
K_{35}	17	15.959	6.122	16.575	2.5
K_{41}	9	8.455	6.06	9.9	10
K_{42}	7	6.905	1.348	6.3	10
K_{43}	10	10.416	4.168	10	0
K_{44}	5	5.831	16.619	5.125	2.5
K_{45}	8	6.935	13.311	7.8	2.5
P_1	0.3	0.304	1.472	0.33	10
$P_{2,3}$	0.2	0.1904	4.78	0.18	10
P_4	0.1	0.0815	18.408	0.1	0
f		0.0216	—	0.0137	—

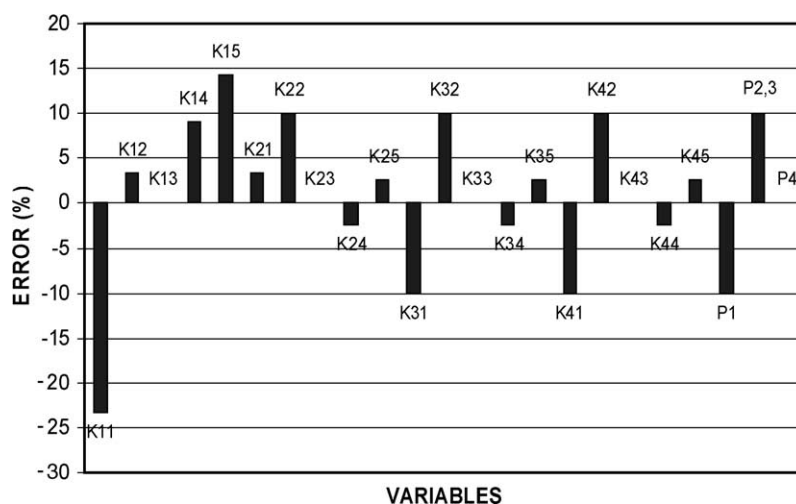


Fig. 13. Distribution of the error in parameters estimations (Case Study No. 2).

obtained using the correct parameters and those found by the proposed solution methodology. Fig. 16 depicts the agreement between the desired variogram and that obtained using the NEGO algorithm. Note that the results were obtained using only 267 computationally expensive objective function evaluations.

Observe that it is possible (e.g., Case study No. 2) to have significant differences in the parameter estimates while having a very small value of the associated objective function. This is a typical pathological feature (non-uniqueness) of the solution of inverse problems constrained by very limited data.

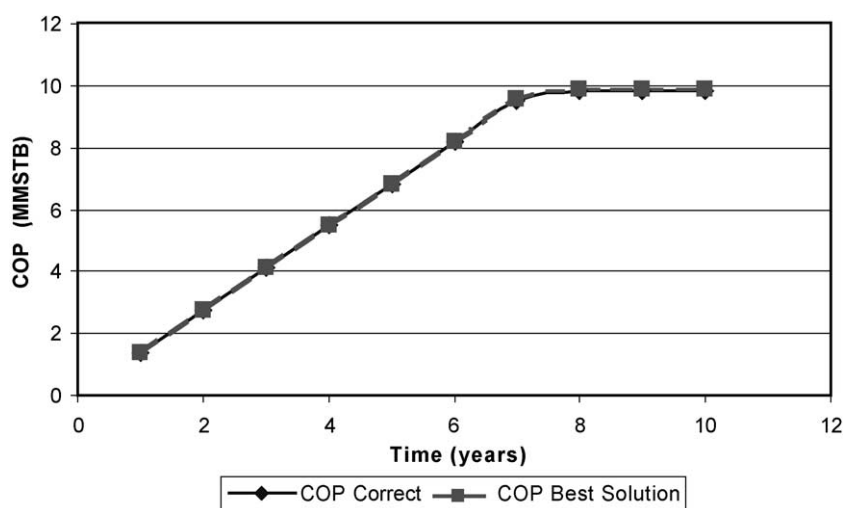


Fig. 14. Cumulative oil production obtained using the best solution and “correct” solution (Case Study No. 2).

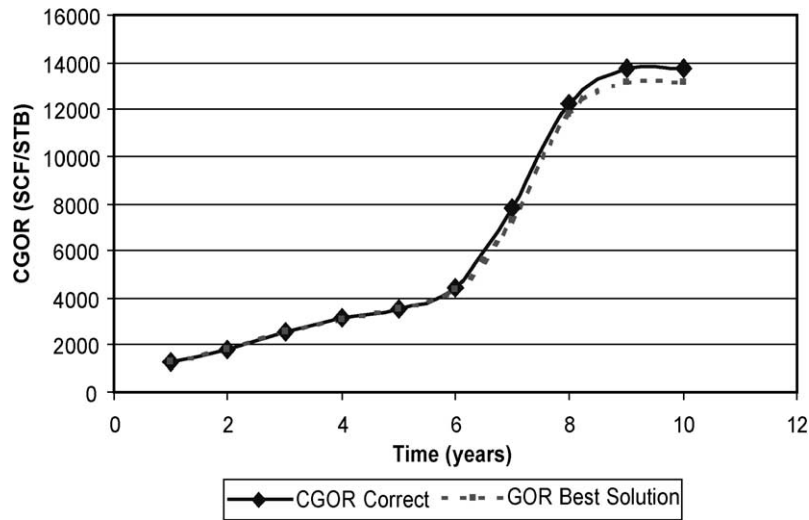


Fig. 15. Cumulative gas oil ratio obtained using the best solution and “correct” solution (Case Study No. 2).

From the results it is concluded that the methodology can be used effectively and efficiently for reservoir characterization purposes. In addition, the optimization approach holds promise to be useful in the optimization of objective functions involving the execution of computationally expensive mathematical models (e.g., reservoir numerical simulators), such as those found, not only in reservoir characterization, but

also in other areas of petroleum engineering (e.g., EOR optimization).

6. Conclusions

- A global optimization method for integrating static and dynamic data into a reservoir description

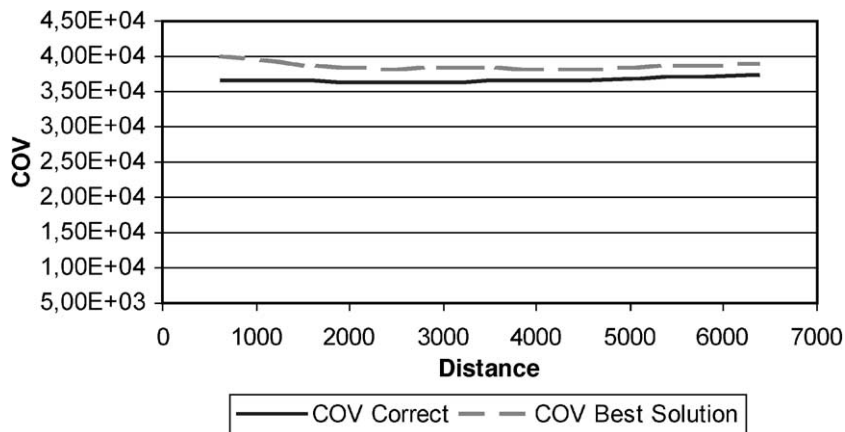


Fig. 16. Covariance values obtained using the best solution and “correct” value of horizontal permeability parameter.

called NEGOT has been proposed. The method includes the construction of a “fast surrogate” of an objective function whose evaluation involves the execution of a time-consuming mathematical model (i.e., reservoir numerical simulator) based on neural networks, DACE modeling, and adaptive sampling. Using adaptive sampling, promising areas are searched considering the information provided by the surrogate model and the expected value of the errors.

- The results suggest that the NEGOT algorithm can be used effectively and efficiently for reservoir characterization purposes. In addition, the optimization approach holds promise to be useful in the solution of inverse problems involving the execution of computationally expensive mathematical models, such as those found, not only in reservoir characterization, but also in other areas of engineering.

- The NEGOT algorithm is expected to outperform competing methods, in terms of computationally expensive objective function evaluations, necessary to meet a stopping criteria. This is because it can identify promising areas without the need of moving step by step along a given trajectory. In addition, by providing estimates of the errors at unsampled points, it is possible to establish a reasonable stopping criterion. Furthermore, provides a fast surrogate model that could be used to visualize the relationship between the sought parameters and the objective function values and to identify the relative significance of each of the parameters.

- Although the proposed approach was not tested using a noisy static and dynamic data, the method holds promise to be effective and efficient under the expected noisy conditions (non-white) present in a reservoir description. This is the subject of current investigation.

Nomenclature

DACE	Design and analysis of computer experiment
\mathbf{x}	Parameters vector
X	Set constraint
f	Objective function
\mathbf{W}	Weighting matrix
d_{obs}	Reference to static and dynamic data available (normalized)
d_{cal}	Denotes the data obtained using a mathematical model
y	DACE response value

μ	Mean of the population
ε	Error in the DACE model
p	Number of dimensions in the vector \mathbf{x}
σ	Standard deviation of the population
θ_h	Correlation parameter
\mathbf{r}'	Correlation vector between the new point and the points used to construct the model
\mathbf{R}	Correlation matrix between the n sample points
$\mathbf{1}$	n -vector of ones
fom	Figure of merit
Φ	Cumulative normal distribution function
ϕ	Density normal distribution function
\bar{y}	DACE predictor
f_{\min}	Current best function value
$s^2(\mathbf{x})$	Mean square error of the predictor
COP	Cumulative oil production
CGOR	Cumulative gas oil ratio
GOR	Gas oil ratio
COV	Covariance
BIS	Best initial solution
NIO	Number of initial observations
OFE	Objective function evaluations
RE:	Relative error
$ \mathbf{x} - \mathbf{x}_{\text{opt}} $	Distance to the optimum value
$ \mathbf{x} - \mathbf{x}_{\text{opt}} _{\eta}$	Distance to the optimum value (test function with noise)
η	Noise factor
randn()	Random number generator normally distributed $N(0,1)$

Subscripts

h	Coordinate directions
obv	Observed
cal	Calculated

Superscript

*	New point
'	Transpose

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