IDENTIFICATION OF THE THERMOPHYSICAL PROPERTIES OF ORTHOTROPIC SEMI-TRANSPARENT MATERIALS

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SUMMARY

- 1. Introduction
- 2. Physical Problem
- 3. Mathematical Formulation
- 4. Direct Problem and Inverse Problem
- 5. Validation of the Solution of the Direct Problem

6. Inverse Problem

7. Conclusions



1. INTRODUCTION

- Identification of thermophysical properties at high temperatures;
- Flash method for thermal diffusivity;
- Semi-transparent materials;
- Coupled conduction-radiation heat transfer.





2. PHYSICAL PROBLEM





3. MATHEMATICAL FORMULATION





3. MATHEMATICAL FORMULATION

Equation of Radiative Transfer

$$\xi \frac{\partial I^{l}}{\partial x} + \eta \frac{\partial I^{l}}{\partial y} + \mu \frac{\partial I^{l}}{\partial z} = -(\kappa_{a} + \sigma_{s})I^{l} + S^{l} \quad \text{in } 0 < x < a, 0 < y < b, 0 < z < c$$

where
$$S^{l} = \kappa_{a} n_{r}^{2} I_{b} (T) + \frac{\sigma_{s}}{4\pi} \int_{\Omega'=4\pi} I^{l'} p(\vec{s}' \rightarrow \vec{s}) d\Omega'$$

Boundary Conditions

$$I\left(\xi,\eta,\mu\right) = I\left(-\xi,\eta,\mu\right) \qquad \text{at } \Gamma_{1}:\begin{cases} x=0\\ 0 < y < b\\ 0 < z < c \end{cases}$$

$$I\left(-\xi,\eta,\mu\right) = \varepsilon n_{r}^{2}I_{b} + \frac{1-\varepsilon}{\pi} \int_{\xi>0} I\left(\xi',\eta',\mu'\right) \xi' d\Omega' \qquad \text{at } \Gamma_{2}:\begin{cases} x=a\\ 0 < y < b\\ 0 < z < c \end{cases}$$

$$I\left(\xi,\eta,\mu\right) = I\left(\xi,-\eta,\mu\right) \qquad \text{at } \Gamma_{3}:\begin{cases} 0 < x < a\\ y=0\\ 0 < z < c \end{cases}$$

$$I\left(\xi,-\eta,\mu\right) = \varepsilon n_{r}^{2}I_{b} + \frac{1-\varepsilon}{\pi} \int_{\eta>0} I\left(\xi',\eta',\mu'\right) \eta' d\Omega' \qquad \text{at } \Gamma_{4}:\begin{cases} 0 < x < a\\ y=b\\ 0 < z < c \end{cases}$$

$$I\left(\xi,\eta,\mu\right) = \varepsilon n_{r}^{2}I_{b} + \frac{1-\varepsilon}{\pi} \int_{\mu<0} I\left(\xi',\eta',\mu'\right) \mu' d\Omega' \qquad \text{at } \Gamma_{5}:\begin{cases} 0 < x < a\\ 0 < y < b\\ 0 < z < c \end{cases}$$

$$I\left(\xi,\eta,\mu\right) = \varepsilon n_{r}^{2}I_{b} + \frac{1-\varepsilon}{\pi} \int_{\mu<0} I\left(\xi',\eta',\mu'\right) \mu' d\Omega' \qquad \text{at } \Gamma_{5}:\begin{cases} 0 < x < a\\ 0 < y < b\\ z=0 \end{cases}$$

$$I\left(\xi,\eta,-\mu\right) = \varepsilon n_{r}^{2}I_{b} + \frac{1-\varepsilon}{\pi} \int_{\mu>0} I\left(\xi',\eta',\mu'\right) \mu' d\Omega' \qquad \text{at } \Gamma_{6}:\begin{cases} 0 < x < a\\ 0 < y < b\\ z=c \end{cases}$$



3. MATHEMATICAL FORMULATION

Energy Conservation Equation

$$C\frac{\partial T}{\partial t} = \frac{\partial}{\partial x} \left(k_x \frac{\partial T}{\partial x} \right) + \frac{\partial}{\partial y} \left(k_y \frac{\partial T}{\partial y} \right) + \frac{\partial}{\partial z} \left(k_z \frac{\partial T}{\partial z} \right) - \nabla \cdot q^{rad}$$

in 0 < x < a, 0 < y < b, 0 < z < c, for t > 0

where:
$$\nabla \cdot q^{rad} = \frac{\kappa_a \tau_0}{N_{pl}} \left[4\pi n_r^2 I_b - \int_{\Omega = 4\pi} I^l d\Omega \right]$$

$$\begin{split} \frac{\partial T}{\partial x} &= 0 & \text{at } \Gamma_1 \text{ for } t > 0 \\ k_x \frac{\partial T}{\partial x} &+ Bi^{rad} T = \frac{\varepsilon \tau_0}{N_{\text{pl}}} \bigg[\int_{\xi > 0} I^l . \xi . d\Omega - n_r^2 \pi I_b \bigg] + Bi^{rad} T_\infty & \text{at } \Gamma_2 \text{ for } t > 0 \\ \frac{\partial T}{\partial y} &= 0 & \text{at } \Gamma_3 \text{ for } t > 0 \\ k_y \frac{\partial T}{\partial y} + Bi^{rad} T &= \frac{\varepsilon \tau_0}{N_{\text{pl}}} \bigg[\int_{\eta > 0} I^l . \eta . d\Omega - n_r^2 \pi I_b \bigg] + Bi^{rad} T_\infty & \text{at } \Gamma_4 \text{ for } t > 0 \\ -k_z \frac{\partial T}{\partial z} + Bi^{rad} T &= \frac{\varepsilon \tau_0}{N_{\text{pl}}} \bigg[\int_{\mu < 0} I^l . \mu . d\Omega - n_r^2 \pi I_b \bigg] + Bi^{rad} T_\infty & \text{at } \Gamma_5 \text{ for } t > 0 \\ k_z \frac{\partial T}{\partial z} + Bi^{rad} T &= \frac{\varepsilon \tau_0}{N_{\text{pl}}} \bigg[\int_{\mu < 0} I^l . \mu . d\Omega - n_r^2 \pi I_b \bigg] + Bi^{rad} T_\infty & \text{at } \Gamma_5 \text{ for } t > 0 \\ + Bi^{rad} T_\infty + \varepsilon_{10.6 \, \mu m} q_{laser} (x, y, t) \end{split}$$

Initial Condition

T = 0 in 0 < x < a, 0 < y < b, 0 < z < c, for t = 0

Dimensionless Variables



$$\sigma_{s\lambda} = \sigma_{s\lambda}^* d_{ref}^*$$
, $N_{pl\lambda} = \frac{\rho_\lambda \kappa_{ref} \Delta I_{max}}{4\sigma T_0^{*4}}$



4. DIRECT AND INVERSE PROBLEMS

DIRECT PROBLEM

Known:

- Boundary and initial conditions
 - C, k_x , k_y , k_z , Bi^{rad} , κ_a and σ_s



Determine:

- Temperature distribution T(x, y, z, t)
- Intensity distribution $I^{l}(x, y, z, \xi^{l}, \eta^{l}, \mu^{l}, t)$

INVERSE PROBLEM

Known:

Boundary and initial conditions *Temperature measurements* Y_m(t_i)
taken at locations (x_m, y_m) m=1,...,M
at the boundary z = 0 and times t_i,
i=1,...,I



Estimate:

• C, k_x, k_y, k_z and Bi^{rad}



Finite-volumes for the Equation of Radiative Transfer and for the Energy Conservation Equation





Conduction





Comparison with analytical solution

















<u>Conduction-Radiation (1D)</u> <u>Gray medium</u>









<u>Conduction-Radiation (1D)</u> <u>Non-gray medium</u>

λ (µm)	n_r	$ ho_{\lambda}$	$\beta_{\lambda} (\mathrm{m}^{-1})$
0.5-1.0	1.5	0.04	10
1.0-2.7	1.5	0.04	100
2.7-4.3	1.5	0.04	1000
4.3-10.3	1.5	0.06	10000
10.3-50	1.8	0.15	10000







Conduction – Cylindrical coordinates





Comparison with analytical solution





Radiation – Cylindrical coordinates









<u>Conduction-Radiation (1D)</u> <u>Gray medium – Cylindrical coordinates</u>









Non-gray medium – Cylindrical coordinates





The inverse problem of interest is concerned with the estimation of the vector of unknown parameters

$$\mathbf{P} = [k_x, k_y, k_z, C, Bi^{rad}]$$

by using transient temperature measurements taken at the non-heated surface Γ_5 at z = 0.





For the solution of the present parameter estimation problem, different minimization techniques were used:

• the <u>Levenberg-Marquardt method</u> applied to the minimization of the *ordinary least squares norm (OLS)*,

 $S_{OLS}(\mathbf{P}) = [\mathbf{Y} - \mathbf{T}(\mathbf{P})]^T [\mathbf{Y} - \mathbf{T}(\mathbf{P})]$

• the <u>Gauss method</u> applied to the minimization of the *maximum a posteriori objective function (MAP)*,

 $S_{MAP}(\mathbf{P}) = [\mathbf{Y} - \mathbf{T}(\mathbf{P})]^T \mathbf{W} [\mathbf{Y} - \mathbf{T}(\mathbf{P})] + (\boldsymbol{\mu} - \mathbf{P})^T \mathbf{V}^{-1} (\boldsymbol{\mu} - \mathbf{P})$

• and the <u>Hybrid method</u> applied to the minimization of the *ordinary least squares norm (OLS)*, which combines deterministic (BFGS method) and evolutionary/stochastic methods (Particle Swarm and Differential Evolution methods).



MAXIMUM LIKELIHOOD OBJECTIVE FUNCTION

$$S_{ML}(\mathbf{P}) = [\mathbf{Y} - \mathbf{T}(\mathbf{P})]^T \mathbf{W} [\mathbf{Y} - \mathbf{T}(\mathbf{P})]$$

where P = vector of unknown parameters Y = vector of measured temperatures T(P) = vector of estimated temperatures





For uncorrelated measurements:

$$\mathbf{W} = \begin{vmatrix} 1/\sigma_1^2 & & 0 \\ & 1/\sigma_2^2 & \\ & \ddots & \\ 0 & & 1/\sigma_I^2 \end{vmatrix}$$



THE LEVENBERG-MARQUARDT METHOD

$$\mathbf{P}^{k+1} = \mathbf{P}^{k} + [\mathbf{J}^{T}\mathbf{W}\mathbf{J} + \lambda^{k}\mathbf{\Omega}^{k}]^{-1}\mathbf{J}^{T}\mathbf{W}[\mathbf{Y} - \mathbf{T}(\mathbf{P}^{k})]$$

where λ^k is the *damping parameter* and Ω^k is a *diagonal matrix*.

- The Levenberg-Marquardt Method is related to *Tikhonov's regularization* approach.
- Compromise between steepest-descent method and Gauss' method.
- Simple, powerful and straightforward iterative procedure.
- Capable of treating complex physical situations.
- Easy to program.
- Stable and converges fast.



<u>Remark</u>: With the statistical hypotheses described above, the minimization of the least-squares norm yields *maximum likelihood* estimates, that is, the values estimated for the unknown parameters \mathbf{P} are those most likely to produce the measured data \mathbf{Y} .

<u>Remark</u>: Although very popular and useful in many situations, the minimization of the least-squares norm is a non-Bayesian estimator. A Bayesian estimator is basically concerned with the analysis of the *posterior probability density*, which is the conditional probability of the parameters **P** given the measurements **Y**.



The **statistical inversion approach** is based on the following principles:

- 1. All variables included in the model are modeled as random variables.
- 2. The randomness describes our degree of information concerning their realizations.
- 3. The degree of information concerning these values is coded in the probability distributions.
- 4. The solution of the inverse problem is the posterior probability distribution.

Jari P. Kaipio and Erkki Somersalo, *Computational and Statistical Methods for Inverse Problems*, Springer, 2004.



BAYES' FORMULA

$$\pi_{posterior}(\mathbf{P}) = \pi(\mathbf{P} | \mathbf{Y}) = \frac{\pi_{prior}(\mathbf{P})\pi(\mathbf{Y} | \mathbf{P})}{\pi(\mathbf{Y})}$$

Where: $\pi_{\text{posterior}}(\mathbf{P}) = \text{posterior probability density (conditional probability of the parameters$ **P**given the measurements**Y**) $<math>\pi_{\text{prior}}(\mathbf{P}) = \text{prior density (information about the parameters prior to the measurements)}$ $\pi(\mathbf{Y}|\mathbf{P}) = \text{likelihood function (expresses the likelihood of different measurement outcomes$ **Y**with**P** $given)}$ $\pi(\mathbf{Y}) = \text{probability density of the measurements (normalizing constant)}$

posterior \propto prior x likelihood



Maximum a Posteriori Objective Function

 $S_{MAP}(\mathbf{P}) = [\mathbf{Y} - \mathbf{T}(\mathbf{P})]^T \mathbf{W} [\mathbf{Y} - \mathbf{T}(\mathbf{P})] + (\boldsymbol{\mu} - \mathbf{P})^T \mathbf{V}^{-1} (\boldsymbol{\mu} - \mathbf{P})$

- The errors are additive, with zero mean and normally distributed.
- The statistical parameters describing the errors are known.
- There are no errors in the independent variables.
- **P** is a random vector with known mean μ and known covariance matrix **V**.
- **P** is distributed normally and is independent of **Y**.

Hypotheses:



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For uncorrelated measurements:

$$\mathbf{W} = \begin{bmatrix} 1/\sigma_1^2 & & 0 \\ & 1/\sigma_2^2 & & \\ & & \ddots & \\ 0 & & & 1/\sigma_I^2 \end{bmatrix}$$

For the minimization of $S_{MAP}(\mathbf{P})$: $\frac{\partial S_{MAP}(\mathbf{P})}{\partial P_1} = \frac{\partial S_{MAP}(\mathbf{P})}{\partial P_2} = \dots = \frac{\partial S_{MAP}(\mathbf{P})}{\partial P_N} = 0$

$$-2\mathbf{J}^T\mathbf{W}[\mathbf{Y}-\mathbf{T}(\mathbf{P})] - 2\mathbf{V}^{-1}[\boldsymbol{\mu}-\mathbf{P}] = 0$$

where **J** is the <u>Sensitivity Matrix</u>.



$$-2\mathbf{J}^T\mathbf{W}[\mathbf{Y}-\mathbf{T}(\mathbf{P})] - 2\mathbf{V}^{-1}[\boldsymbol{\mu}-\mathbf{P}] = 0$$

Linear Problems: J does not depend on P \implies T(P) = J P

$$\mathbf{P} = [\mathbf{J}^T \mathbf{W} \mathbf{J} + \mathbf{V}^{-1}]^{-1} [\mathbf{J}^T \mathbf{W} \mathbf{Y} + \mathbf{V}^{-1} \boldsymbol{\mu}]$$

Nonlinear Problems: $\mathbf{J} \equiv \mathbf{J}(\mathbf{P}) \implies \mathbf{T}(\mathbf{P}) = \mathbf{T}(\mathbf{P}^k) + \mathbf{J}^k (\mathbf{P} - \mathbf{P}^k)$

$$\mathbf{P}^{k+1} = \mathbf{P}^k + [\mathbf{J}^T \mathbf{W} \mathbf{J} + \mathbf{V}^{-1}]^{-1} \{\mathbf{J}^T \mathbf{W} [\mathbf{Y} - \mathbf{T} (\mathbf{P}^k)] + \mathbf{V}^{-1} (\boldsymbol{\mu} - \mathbf{P}^k)\}$$



Hybrid Method – Minimization of OLS



• **<u>DE Method:</u>**

- Alternative to the Genetic Algorithm method.
- Proposed in 1995 by Kenneth Price and Rainer Storn from Berkeley.
- The method initializes with a random generated random matrix **P** which

contains N vector parameters \mathbf{x}

• From the initial population matrix, generations are created until the best generation (optimum) is found.

• The next generation is created as:



where

 α , β and γ are three randomly chosen members of the population matrix **P**.

F is a weighting function which defines the mutation (0.5 < F < 1).

k is the generation counter.

 δ_1 and δ_2 are delta Dirac functions that defines the crossover.

If $f(\mathbf{x}^{k+1}) < f(\mathbf{x}^k)$ \longrightarrow \mathbf{x}^{k+1} replaces \mathbf{x}^k in the population matrix \mathbf{P} If $f(\mathbf{x}^{k+1}) > f(\mathbf{x}^k)$ \longrightarrow \mathbf{x}^k is kept in the population matrix \mathbf{P} and \mathbf{x}^{k+1} is discarded • The crossover is obtained as: $\mathbf{x}_i^{k+1} = \delta_1 \mathbf{x}_i^k + \delta_2 [\alpha + F(\beta - \gamma)]$

$$\delta_1 = \bigoplus \begin{array}{c} 0, \text{ if } R < CR \\ 1, \text{ if } R > CR \end{array} \qquad \qquad \delta_2 = \bigoplus \begin{array}{c} 1, \text{ if } R < CR \\ 0, \text{ if } R > CR \end{array}$$

- R is a random number with uniform distribution between 0 and 1
- CR is the crossover factor (0.5 < CR < 1)

- <u>PS (Particle Swarm) method</u>:
 - Created in 1995 by an Electric Engineer (Russel Eberhart) and a Social-Psychologist (James Kennedy) as an alternative to Genetic Algorithm.
 - Based on the social behavior of various species (including humans).
 - Balances the individuality and sociability of individuals in order to find a optimum.
- Individuality
 Chances to find alternatives places
 Convergence

Sociability

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Learning process among the individuals

Chances to find alternatives places. Individuals can find a local minima

- **PS method:**
- Update process

$$\mathbf{x}_{i}^{k+1} = \mathbf{x}_{i}^{k} + \mathbf{v}_{i}^{k+1}$$
$$\mathbf{v}_{i}^{k+1} = \alpha \mathbf{v}_{i}^{k} + \beta \mathbf{r}_{1i} (\mathbf{p}_{i} - \mathbf{x}_{i}^{k}) + \beta \mathbf{r}_{2i} (\mathbf{p}_{g} - \mathbf{x}_{i}^{k})$$
Individuality Sociability

where

 \mathbf{x}_{i} is i-th individual of the vector of parameters

 \mathbf{r}_{1i} and \mathbf{r}_{2i} are are random numbers with uniform distribution between 0 and 1

 \mathbf{p}_i is the best value found for the vector \mathbf{x}_i

 \mathbf{p}_{g} is the vest value found for the entire population

 $0 < \alpha < 1; 1 < \beta < 2$







Simulated measurements ($\sigma = 0.8$ K) $C^* = 2.5 \times 10^6 \text{ Jm}^{-3} \text{K}^{-1}$ $k_x^* = 5 \text{ Wm}^{-1} \text{K}^{-1}$ $k_y^* = 5 \text{ Wm}^{-1} \text{K}^{-1}$ $k_z^* = 5 \text{ Wm}^{-1} \text{K}^{-1}$ $h_{rad}^* = 1372 \text{ Wm}^{-2} \text{K}^{-1}$. $\kappa_a^* = 10 \text{ m}^{-1}$ $\sigma_s^* = 10^4 \text{ m}^{-1}$



The sample was assumed to be a parallelepiped with dimensions $2a^* = 2b^* = 0.01$ m and $c^*=0.001$ m, heated by a laser with a power of 23 W and a Gaussian distribution. For the heat flux imposed by the laser, 99% of its power was assumed to be delivered within a circle with radius of 2 mm centered at the sample. The sample is assumed to be initially at the uniform temperature of 1800K, which is the same temperature of the surrounding environment.









Model for the Gradient	Surrogate	Complete	Surrogate	Complete	<mark>1st step:</mark> Surrogate 2 nd step: Surrogate	<mark>1st step:</mark> Surrogate 2 nd step: Complete	<mark>1st step:</mark> Surrogate 2 nd step: Surrogate	<u>1st step:</u> Surrogate 2 nd step: Complete
Model for the Direct Problem	Complete	Complete	Complete	Complete	<u>1st step:</u> Surrogate 2 nd step: Complete	<u>1st step:</u> Surrogate 2 nd step: Complete	<u>1st step:</u> Surrogate 2 nd step: Complete	<u>1ªt step:</u> Surrogat <mark>e</mark> 2 nd step: Complete
Method	Levenberg-Marquardt	Levenberg-Marquardt	Gauss	Gauss	<u>1st step:</u> Hybrid 2 nd step: Levenberg- Marquardt	<u>1st step:</u> Hybrid 2 nd step: Levenberg- Marquardt	<u>1st step:</u> Hybrid 2 nd step: Gauss	<u>1st step:</u> Hybrid 2 nd step: Gauss
Objective Function	Least-squares	Least-squares	Maximum a Posteriori	Maximum a Posteriori	Least-squares	Least-squares	<u>1^ª step:</u> Least-squares 2 nd step:MAP	<u>1^ª step:</u> Least-squares 2 nd step:MAP
Technique	↽	2	ю	4	Ð	Q	7	ω

Table 1: Estimation techniques

					Estimates		
Technique	Number of	CPU Time	$C^* imes 10^{6}$	³⁵ 4	k_{γ}^{*}	k_z^*	h_{rad}^{*}
	Iterations)	${ m Jm^{-3}K^{-1}}$	Wm ⁻¹ K ⁻¹	$\mathrm{Wm}^{-1}\mathrm{K}^{-1}$	$Wm^{-1}K^{-1}$	$\mathrm{Wm}^{2}\mathrm{K}$
Ļ	16	5h35m18s	2.51 ± 0.03	4.99 ± 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 ± 5
ы	16	6h14m04s	2.51 <u>±</u> 0.03	5.00 ± 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 <u>±</u> 5
3	13	4h33m16s	<u>2.51 ± 0.03</u>	5.00 = 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 ± 5
4	9	2h36m5s	2.51 ± 0.03	2.00 = 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 ± 5
ų	50	1h17m26s	2.19	5.74	5.80	3.6	1246
n	15	5h50m55s	2.51 ± 0.03	4.99 ± 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 ± 5
¢,	50	1h17m26s	2.19	5.74	5.80	3.6	1246
o	16	6h33m02s	2.51 ± 0.03	5.00 ± 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 ± 5
1	20	1h17m26s	2.19	2.74	5.80	3.6	1246
1	11	4h09m59s	2.51 ± 0.03	5.00 ± 0.07	5.01 ± 0.07	5.0 ± 0.2	1373 ± 5
o	50	1h17m26s	2.19	5.74	5.80	3.6	1246
0	4	1h59m11s	2.51 + 0.03	5.00 - 0.07	5.01 + 0.07	5.0 + 0.2	1373 + 5

<u>Table 2</u>: Results obtained with an initial guess close to the exact parameters $(C^{*0} = 2.8 \times 10^6 \text{ J/m}^3 \text{ K}, k_x^{*0} = k_y^{*0} = k_z^{*0} = 8 \text{ W/m.K}, h^{na^{*0}} = 800 \text{ W/m}^2 \text{ K})$

$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	1 NC -	4 NC	3 NC	2 NC	1 NC	Iterations M^{-2} Iterations M^{-1} M^{-1} M^{-1} M^{-1} M^{-1} M^{-1} M^{-1} M^{-1} M^{-2} M^{-2}	$\frac{\text{Number of CPUTTime } C^* \times 10^6 k_x^* k_y^* k_z^* h_{zz}^* h_{zz}^* = h_{zz}^* h_{zz}^* = h_{zz}^* h_{zz}^* = h_{zz}^* h_{zz}^* = h_$	Estimates	⁻¹ Wm ⁻² K Nm ⁻² K 2 1373 ± - 1224 2 1373 ± -	$\begin{array}{c} k_{z}^{*} \\ Wm^{-1}K^{-1} \\ Wm^{-1}K^{-1} \\ - \\ - \\ - \\ - \\ - \\ - \\ - \\ - \\ - \\ $	Estimates k_y^* $W_{\rm In}^{-1} {\rm K}^{-1}$ - - - - - - - -	k_{x}^{*} Wm ⁻¹ K ⁻	$C^* \times 10^6$ $Jm^3 K^{-1}$ $-$ $-$ $-$ 2.04 2.04 2.04 2.04 2.04 2.04 2.04 2.04	CPU Time - - - - 1h19m46s 7h44m03s 6h32m46s 6h32m446s 6h32m446s 3h54m49s 3h54m49s 1h19m46s 3h54m49s	Number of Iterations NC NC NC NC NC S0 50 50 50 50 16 10 10 50 50 50 50 50 50 50 50 50 50 50 50 50	shnique 1 5 6 8
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<u>Table 3:</u> Results obtained with an initial guess far from the exact parameters $\left\{ C^{*1} = 0.1 \times 10^{6} \text{ J/m}^{2}.\text{K}, k_{*}^{*0} = k_{*}^{*0} = k_{*}^{*0} = 50 \text{ W/m.K}, h^{7ad^{*0}} = 5 \text{ W/m}^{2}.\text{K} \right\}$



7. CONCLUSIONS

• The use of a surrogate model for the gradient did not affect the accuracy of the estimated parameters and may cause an increase on the number of iterations and CPU time, due to the loss of computational accuracy.

• The two-step approach was necessary to reach convergence if initial guesses far from the exact parameters were used in the inverse analysis.