MULTI-OBJECTIVE CALIBRATION OF THE DOUBLE-ELLIPSOID HEAT SOURCE MODEL FOR GMAW PROCESS SIMULATION

by

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> Original scientific paper https://doi.org/10.2298/TSCI210131181B

The scope of application of simulation models in welding is limited by the accuracy of their output results. This paper presents a calibration procedure for a 3-D quasi-stationary model of heat transfer for gas metal arc welding. The double-ellipsoid heat source used in this model has five input parameters whose value cannot be specified accurately. To estimate these values, we employed a multi-objective calibration procedure with two objective functions using the paretosearch optimization algorithm. Objective functions represented the error between simulated and experimentally observed values of penetration depth and weld bead width during gas metal arc welding of P355GH steel plates. All input parameters were assumed to be a power function of line energy. To reduce computational time, we replaced the numerical model with a response surface methodology metamodel based on an optimal set of simulation results from the numerical model. The results of the simulations based on calculated values of input parameters for the heat source model showed excellent matching with the experimental results.

Key words: welding simulation, multi-objective calibration, double-ellipsoid heat source, RSM

Introduction

Although simulation modelling of different production processes has a growing trend in recent years, the application of simulation models of welding processes is still limited. The main reason is a lack of knowledge of correct values of some input parameters, which in turn provides insufficiently reliable output results regardless of the complexity of the model. In the case of the welding heat transfer model, the parameters that most affect the accuracy of the output results are the parameters related to the model of a heat source. Widely used Goldak's double-ellipsoid heat source [1], is defined by five input parameters neglecting the power of the heat source, fig. 1. The first of those five parameters is arc efficiency. For the GMAW process, the value of arc efficiency varies from 0.80 to 0.88 in the case of Ar shielding [2]. In the case of 90% Ar to 10% CO₂ shielding, arc efficiency was found to be in the range from 0.68-0.86 [3]. Some authors have reported values between 0.675 and 0.722 [4]. It is obvious that such wide limits for arc efficiency significantly affect the results of the simulation.

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The other four parameters a_f , a_r , b_h and c_h are the semiaxes of the front and rear ellipsoid, fig 1. Power of the heat source is distributed, eqs. (1)-(3), between the front and rear ellipsoid in a ratio which corresponds to values of parameters $f_f = 2a_f/(a_f + a_r)$ and $f_r = 2a_r/(a_f + a_r)$ [5]:

$$q_{\rm f}(x, y, z) = \frac{6\sqrt{3}f_{\rm f}Q}{a_{\rm f}b_{\rm h}c_{\rm h}\pi\sqrt{\pi}} e^{-3\frac{x^2}{a_{\rm f}^2}} e^{-3\frac{y^2}{b_{\rm h}^2}} e^{-3\frac{z^2}{c_{\rm h}^2}}$$
(1)

$$q_r(x, y, z) = \frac{6\sqrt{3}f_rQ}{a_rb_hc_h\pi\sqrt{\pi}} e^{-3\frac{x^2}{a_r^2}} e^{-3\frac{y^2}{b_h^2}} e^{-3\frac{z^2}{c_h^2}}$$
(2)

$$Q = \eta U I \tag{3}$$

Goldak *et al.* [1] suggested that the value of parameters b_h and c_h should be measured directly from an experimental cross-section of the weld. They also suggested that in absence of experimental data Christensen's [6] expressions could be used. For f_f and f_r they proposed values of 0.6 and 1.4, respectively. Joshi *et al.* [7] assumed that $f_f = 0.48$ and $f_r = 1.52$ while Chen *et al.* [8] as well as Jia *et al.* [9] suggested that values of parameters f_f and f_r should be 0.5 and 1.5, respectively. Guided by Goldak's [10] recommendation, Nasiri and Enzinger [11] have used parameters of double-ellipsoid heat source about 10% smaller than actual weld geometry.

For reliable output results of the simulation model, it is necessary to determine the exact values of these input parameters. In the case of individual simulations, we can achieve this goal, minimizing the error between the simulation and experimental results by combining the simulation model with some multivariable optimization algorithm [12-18]. The complexity of the problem arises when it is necessary to determine the optimal values of these parameters for several different simulations [19]. In this case, it is possible to assume that there is a functional dependence between the heat source parameters and some of the process parameters [9, 19, 20]. This way, we can determine the values of input parameters for the whole interval of different simulation conditions. The use of multi-objective [21, 22] instead of single-objective optimization algorithm [14, 23-25] allows the error estimation for each parameter included in the objective functions separately. That gives us the possibility to choose the appropriate functional relations depending on the need for the greater or lesser error of individual parameters concerning experimental results.

The application of numerical models in simulations of the welding process is practically impossible without the use of computers. Despite the development of computer technology, welding simulations still require a large amount of time to perform. However, in a situation such as model calibration or some form of optimization, the number of simulation runs can be large, which in turn leads to a very long time required to obtain the final results. Therefore, it makes sense to use metamodels as an efficient replacement for numerical models. The most commonly used metamodelling techniques are regression analysis [23, 26-28] and neural networks [29-31].

In this article, we proposed an RSM metamodel based calibration methodology for the reduction of the error between simulated and experimental weld penetration depth and weld bead width. This calibration methodology implied the determination of heat source parameters as functions of line energy. The RSM metamodel was used as a replacement for the numerical heat transfer model to reduce the computational time. This way, we were able to use a multi-objective paretosearch optimization algorithm to calculate the values of double-ellipsoid heat source parameters for multiple simulation conditions.

Heat transfer analysis

The heat transfer during welding, can be described with a non-stationary, partial differential equation, eq. (4) [22], where ρ , c_{ρ} , and λ are density, specific heat capacity and thermal conductivity of material while L and q_1 are latent heat of melting/solidification and volumetric heat source described by Goldak's double-ellipsoid model. An analytical solution for this type of equation is connected with difficulties related to non-linearities of material physical properties, the complexity of boundary conditions, and the heat source model:

$$\rho c_p \frac{\partial T}{\partial t} = \lambda \left(\frac{\partial^2 T}{\partial x^2} + \frac{\partial^2 T}{\partial y^2} + \frac{\partial^2 T}{\partial z^2} \right) - \rho L \frac{\partial f_{\text{liq}}}{\partial t} + q_l \tag{4}$$

Latent heat of melting/solidification is described using liquid phase fraction in mushy zone between solidus T_{sol} , and liquidus temperature T_{liq} , eq. (5):

$$f_{\rm liq} = \begin{cases} 0 & \text{for} \quad T \le T_{\rm sol} \\ \frac{T - T_{\rm sol}}{T_{\rm liq} - T_{\rm sol}} & \text{for} \quad T_{\rm sol} < T < T_{\rm liq} \\ 1 & \text{for} \quad T \ge T_{\rm liq} \end{cases}$$
(5)

In case of constant welding speed, v_w , it is possible to transform eq. (4) to the quasisteady-state form, eq. (7). This kind of transformation requires the application of a moving co-ordinate system ζyz , fig. 2. The connection between co-ordinate systems xyz and ζyz is defined:

$$\xi = x - v_w t \tag{6}$$

Considering eqs. (4) and (6), heat transfer eq. (7) in moving co-ordinate system can be written:

$$-v_{\rm w}c_{\rm eff}\frac{\partial T}{\partial\xi} = \frac{\lambda}{\rho c_p} \left(\frac{\partial^2 T}{\partial\xi^2} + \frac{\partial^2 T}{\partial y^2} + \frac{\partial^2 T}{\partial z^2}\right) + q_l \quad (7)$$

where c_{eff} is effective heat capacity:



Figure 2. Moving co-ordinate system

 $c_{eff} = \left(1 + \frac{L}{c_p} \frac{\partial f_{liq}}{\partial T}\right)$ (8)

Equation (7) is solved in MATLAB iteratively, using the multigrid and the successive over-relaxation finite differences method.

Experimental procedure

The calibration procedure was performed using experimental data from specimens with dimensions $300 \text{ mm} \times 150 \text{ mm} \times 5.3 \text{ mm}$. The base material for specimens was P355GH steel with the chemical composition given in tab. 1. The filler material used was OK Autrod

Mn

1.45

for model validation, tab. 3.

Nb

0.014 For each of four specimens, single pass

12.50 uncoated wire with a diameter of 1.2 mm. As a shielding gas, a two-component mixture of 82% Ar and 18% CO2 was used. Four specimens were welded in a flat position using ARC Mate 100iC welding robot with Migatronic Sigma Galaxy 400 power source.

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0.062

Table 1. Chemic	al composition	n of base materi	al - P355GH
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Figure 3. Measured weld geometry

Doromotor	Specimen				
Parameter	Ι	II	III	IV	
Voltage [V]	20.2	20.3	21.2	21.8	
Current [A]	185	190	208	222	
Welding speed [mms ⁻¹]	8	8	8	8	
Wire feed rate [m per minute]	4.3	4.5	5.2	5.7	
Wire diameter [mm]	1.2	1.2	1.2	1.2	
Gas-flow [Lpm]	12	12	12	12	

Table 2. Welding parameters

Table 3. Measured weld dimensions	Table 3	. Measured	weld	dimensions
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Parameter		Spec	imen	
Farameter	Ι	II	III	IV
W [mm]	6.5	6.6	7.2	7.3
<i>D</i> [mm]	2.1	2.1	2.4	2.3

Metamodelling

Metamodelling represents a technique by which a particular model is replaced by another that appropriately imitates the original model, and at the same time, has better characteristics related to computational efficiency [32].

Solving eq. (7) iteratively using the finite differences method is associated with a large number of iterations, which results in a large computational time. Bearing in mind that the calibration process is based on the multi-objective optimization method, this further increases the

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time required to reach a solution. In this case, an RSM-based metamodel is employed to reduce the computational time.

Based on the welding parameters shown in tab. 2, for specimens I and IV, and using the Design-Expert software package, an I-optimal plan of experiments was created. As already mentioned, Goldak's model has five unknown input parameters: η , $a_{\rm f}$, $a_{\rm r}$, $b_{\rm h}$, and $c_{\rm h}$. To reduce the number of experimental runs it is assumed that parameters $a_{\rm f}$ and $b_{\rm h}$ are equal. Therefore, the following variables were used as input ones:

$$x_1 = \eta, x_2 = b_h/b_{hexp}, x_3 = c_h/c_{hexp}, x_4 = a_r/a_f$$

Levels of input parameters are listed in tab. 4.

Table 4. Levels of input parameters

Levels		Fac	tors	
Levels	x_1	x_2	<i>x</i> ₃	x_4
Level 1	0.6	0.9	0.9	2
Level 2	0.9	1.1	1.1	4

We have used numerical model based on eq. (7) to calculate values of weld bead width and penetration depth for specimens I and IV using design matrix, tab. 5, and values of welding parameters listed in tab. 2.

Table 5.	Optimal	design	matrix
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Due	Factors]	Respons	es [mm]]	
Run	x_1	<i>x</i> ₂	<i>x</i> ₃	x_4	WI	$D_{\rm I}$	W _{IV}	$D_{\rm IV}$
1	0.9	0.954	1.1	2.52	6.2	2.7	6.9	3.1
2	0.6	0.9	0.9	4	4.5	1.7	5.4	2.2
3	0.6	1.1	0.9	2	5.6	1.9	6.5	2.2
4	0.6	0.9	1.1	2	5.1	2.3	5.9	2.5
5	0.723	0.9	0.955	2	5.8	2.3	6.4	2.7
6	0.759	1.01891	0.9	3.2	5.9	2.2	6.6	2.4
7	0.615	1.1	1.1	2.6	4.8	1.7	6.1	2.4
8	0.9	1.1	0.951	2.54	6.6	2.4	7.3	2.6
9	0.616395	1.1	0.959	4	4.7	1.6	5.5	1.9
10	0.9	0.9	0.9	2	6.3	2.5	7.3	3.1
11	0.6	1.1	1.1	4	4	1.5	5.3	2
12	0.615	0.96	1.1	4	4.4	1.8	5.4	2.2
13	0.834	1.023	1.023	3.24329	5.9	2.4	6.8	2.8
14	0.68871	0.982	1.1	2.8	5.3	2.2	6.2	2.6
15	0.9	0.952	0.952901	4	5.9	2.5	6.6	2.7
16	0.84	1.1	1.1	4	5.6	2.2	6.5	2.7
17	0.760449	1.021	1.02	2	5.9	2.3	6.9	2.8
18	0.6	0.971	0.972	2.7	5.1	1.9	5.8	2.3
19	0.9	1.1	0.921	4	6.2	2.3	6.9	2.6
20	0.7605	0.9	1.022	3.2	5.5	2.4	6.2	2.6

Mathematical metamodels of functional relations between weld geometry and parameters x_1 to x_4 for specimens I and IV were obtained using the Design-Expert software package. For specimen I, weld bead width W_1 and penetration depth D_1 are described by 2FI models. Models are power transformed:

$$W_{\rm I} = \left(-506.919 - 294.903x_1 + 780.403x_2 + 735.052x_3 - 29.1603x_4 + 583.514x_1x_2 + 133.67x_1x_3 - 4.88747x_1x_4 - 1058.41x_2x_3 - 22.4403x_2x_4 + 26.0028x_3x_4\right)^{1/3}$$
(9)
$$D_{\rm I} = \left(-45.3105 + 1.86122x_1 + 50.8888x_2 + 50.098x_3 - 1.25343x_4 - 9.2795x_1x_2 + 128.4888x_2 + 50.098x_3 - 1.25343x_4 + 50.2795x_1x_2 + 128.4888x_3 + 50.098x_3 - 1.25343x_4 + 50.2795x_1x_2 + 128.4888x_3 + 50.098x_3 - 1.25343x_4 + 50.2795x_1x_2 + 128.4888x_3 + 50.098x_3 + 50.09$$

$$+16.3006x_{1}x_{3}+1.83625x_{1}x_{4}-54.3195x_{2}x_{3}+0.486464x_{2}x_{4}-1.16573x_{3}x_{4}\Big)^{1/2.32}$$
(10)

Comparation of experimental and predicted values of parameters W_1 and D_1 are shown in figs. 4. and 5, respectively. Based on these figures, it is clear that both metamodels represent a good replacement for the numerical model.



Similarly, for specimen IV, weld bead width W_{IV} and penetration depth D_{IV} can be described by power transformed linear models:

$$W_{\rm IV} = \left(9.14992 + 96.5959x_1 + 36.8124x_2 - 24.4857x_3 - 8.51831x_4\right)^{1/2.27}$$
(11)

$$D_{\rm IV} = \left(1.52052 + 2.58036x_1 - 1.27186x_2 + 1.12647x_3 - 0.174858x_4\right)^{1/1.09}$$
(12)

Comparison between actual and predicted values for parameters W_{IV} and D_{IV} is shown in figs. 6. and 7, respectively. The adequacy for all models is tested using the statistical analysis of variance (ANOVA). Results of ANOVA are shown in tab. 6. The *F*-values of models imply that all models are significant while low *P*-values less than 0.05 indicate model terms are significant. The predicted R^2 is in reasonable agreement with the adjusted R^2 for all models.

Model	SS	df	MS	F	р	R^2	Adj. R ²	Pred. R^2	Adeq P
WI	70766.92	10	7076.69	229.39	< 0.0001	0.9961	0.9917	0.9687	53.8446
DI	81.49	10	8.15	97.45	< 0.0001	0.9908	0.9807	0.9335	36.9678
W _{IV}	3924.10	4	981.03	171.90	< 0.0001	0.9787	0.9730	0.9628	40.8285
$D_{\rm IV}$	2.62	4	0.6551	57.91	< 0.0001	0.9392	0.9230	0.8835	25.1054

Table 6. The ANOVA analysis for metamodels W_1 , D_1 , W_{1V} , and D_{1V}



Optimization methodology

Multi-objective optimization methods involve simultaneous optimization of two or more objective functions instead of optimization of the single objective as a linear combination of multiple objective functions. This way, it is possible to estimate each objective function separately, which gives us the possibility to choose the appropriate solution. In this case, that means that we can choose the solution with smaller or larger error either for one or multiple objective functions.

The optimization procedure was performed using the MATLAB software package and paretosearch algorithm. This algorithm uses pattern search on a set of points to search iteratively for non-dominated solutions [33].

Objective O_1 and O_2 are defined using least-squares method, tab. 7. Here, W_j^{exp} and D_j^{exp} are measured values of weld bead width and penetration depth for specimens I and IV, tab. 3, and *j* is number of specimens used for calibration, in this case, j = 2.

Table 7. Objective functions				
$O_1(x_1, x_2, x_3, x_4)$	$O_2(x_1, x_2, x_3, x_4)$			
$\sum_{j=1}^{2} \left(\frac{W_{j}^{\exp} - W_{j}^{\min}}{W_{j}^{\exp}} \right)^{2}$	$\sum_{j=1}^{2} \left(\frac{D_{j}^{\exp} - D_{j}^{\sin}}{D_{j}^{\exp}} \right)^{2}$			

As the unknown parameters, variables x_1, x_2, x_3 , and x_4 have been chosen. We intended to determine if there is a functional relations between variables x_1, x_2, x_3, x_4 , and welding parameters. So we assumed the following relations between them, tab. 8.

Table 8. Assumed functional relations



All functional relations were supposed to be a power function of line energy, where U is arc voltage, I – the arc current, v_w – the welding speed, and a_i , b_i , and c_i are unknown con-

stants. The main goal of the optimization was to determine values of constants a_i , b_i , and c_i so that differences between experimental and computed values defined by objective functions, be minimal. The constraints were supposed as follows, tab. 9.

Variables				
$0.6 \le x_1 \le 0.9$	$0.9 \le x_2 \le 1.1$	$0.9 \le x_3 \le 1.1$	$2 \le x_4 \le 4$	

The optimization problem can be stated:

$$\min \left[O_1 \left(x_1, x_2, x_3, x_4 \right), \ O_2 \left(x_1, x_2, x_3, x_4 \right) \right]$$

$$x_i = a_i + b_i \left(\frac{U_1}{v_w} \right)^{c_i}$$

$$x_{i\min} \le x_i \le x_{i\max}$$
(13)

The result of multi-objective optimization is shown in the form of the Pareto front, fig. 8. It contains a set of 1000 non-dominated solutions. It is not possible to directly choose one solution from this set, which is superior to other solutions considering both objective functions. To overcome this problem we have used the Taguchi signal-to-noise ratio ratio. The signal-to-noise ratio implies three categories of performance characteristics: the lower-the-better, the higher-the-better, and the nominal-the-better. The lower-the-better characteristic is used in the case when it is necessary to minimize the response. It can be expressed:

$$\frac{S}{N} = -10\log\left(\frac{1}{n}\sum_{i=1}^{n}Y_{i}^{2}\right)$$
(14)

where *n* is the number of observed values, in this case, the number of objective functions, and Y_i is the value of the observed characteristic, *i.e.*, the value of the objective function. For all three performance characteristics, a higher *S*/*N* ratio matches to better performance characteristics. In our case, this means that the non-dominated solution with the highest *S*/*N* ratio is superior to other solutions. The *S*/*N* ratio for a set of non-dominated solutions from Pareto front is shown at fig. 9.

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Figure 8. Pareto front of the optimization problem

Figure 9. The *S*/*N* ratio for all non-dominated solutions

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The solution with the highest S/N ratio of 47.7764 [dB] was chosen as optimal one, from the set of 1000 non-dominated solutions, fig. 9. Values of constants a_i , b_i , and c_i which correspond to the optimal solution are listed in tab. 9.

Table 9. Value	s of constants	a_i, b_i , and c_i
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Constants	<i>i</i> = 1	<i>i</i> = 2	<i>i</i> = 3	<i>i</i> = 4
a_i	0.0790	1.0999	0.9001	0.5908
b_i	4.4783	6.0708 · 10 ⁻⁶	-3.1099.10-7	1.4094
C _i	-0.3069	0.3564	0.7648	5.7131.10-6

Functional relations, tab. 8, between parameters

 $x_1 = \eta, x_2 = b_h/b_{hexp}, x_3 = c_h/c_{hexp}, x_4 = a_r/a_f$

and line energy are shown on figs. 10 and 11. As can be seen, parameters x_2 , x_3 , and x_4 have nearly constant values over the whole interval of line energy. The value of parameter x_3 , which is 10% smaller than actual weld geometry, is in good agreement with the proposition of Nasiri and Enzinger [11], while the parameter x_2 has a value which is 10% higher than actual geometry. The parameter x_4 has a value of 2, which corresponds to $f_f = 0.67$ and $f_r = 1.33$, and agrees with Nguyen [34]. Only parameter x_1 changes decreasing from the value of 0.76 for specimen I to 0.71 for specimen IV. This corresponds to values of arc efficiency found by Haelsig *et al.* [3].



Using functional relations, tab. 8, and values of line energy, tab. 2, we have used calculated values of parameters η , b_h , c_h , and a_r/a_f to simulate weld geometry for all specimens. Relative errors for specimens I and IV were 3.1% and 4.1%, respectively. For depth of penetration, relative errors for these two samples were 4.8% and 4.4%. The maximal relative error for the specimens II and III which were used for model validation were 8.3% in case of weld bead for specimen III.

Figures 12 and 13 show simulated weld geometry compared to the experimental one in the case of specimens I and IV. According to tab. 10 and figs. 12 and 13, a calibrated numerical model can reliably predict actual weld geometry.







Table 10. Absolute and relative errors

Figure 13. Comparison between simulated and experimental weld bead geometry for specimen IV

	Parameter									
Spec.	Weld bead width [mm]				Depth of penetration [mm]					
	$W_{\rm sim}$	Wexp	Abs. error	Rel. error	$D_{\rm sim}$	D_{exp}	Abs. error	Rel. error		
Ι	6.3	6.5	0.2	0.031	2.2	2.1	0.1	0.048		
II	6.4	6.6	0.2	0.03	2.2	2.1	0.1	0.048		
III	6.6	7.2	0.6	0.083	2.4	2.4	0	0		
IV	7.0	7.3	0.3	0.041	2.4	2.3	0.1	0.044		

Summary and conclusions

To improve the reliability of the 3-D numerical heat transfer model, we have developed a calibration procedure to determine the input parameters of the double-ellipsoid heat source. The procedure were based on a multi-objective paretosearch optimization algorithm combined with RSM metamodel. This approach proved to be an efficient and reliable way to speed up the optimization process. All input parameters were supposed to be power functions of line energy. The results of simulations based on the heat source parameters calculated using the calibration model show good agreement between simulated and actual weld geometry. It leads to the conclusion that the calibration model based on the functional relations between heat source parameters and welding parameters provides a reliable way to increase the accuracy of the output results of the numerical model.

Acknowledgment

The authors wish to express their gratitude to the Ministry of Education, Science and Technological development of the Republic of Serbia for support through contract No. 451-03-68/2020-14/200108 and to the National CEEPUS Office of the Czech Republic through project CIII-HR-0108-07-1314.

Nomenclature

- $a_{\rm f}$ semiaxis of front half-ellipsoid in x-direction, [m]
- a_r semiaxis of rear half-ellipsoid in x-direction, [m]
- $b_{\rm h}$ semiaxis of front half-ellipsoid in y-direction, [m]
- $c_{\rm h}$ semiaxis of front half-ellipsoid in z-direction, [m]
- $c_{\rm eff}$ effective heat capacity
- c_p specific heat capacity, [Jkg⁻¹K⁻¹]
- \hat{D} depth of penetration [mm]
- $f_{\rm f}$ front proportion coefficient
- $J_{\rm f}$ from proportion coefficient

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- f_{liq} liquid phase fraction, [%]
- $f_{\rm r}$ rear proportion coefficient
- *I* welding current, [A]
- L latent heat, [Jkg⁻¹]
- n number of observed values
- q_l heat density, [Wm⁻³]
- \overline{O}_1 first objective function
- O_2 second objective function
- S/N signal to noise ratio [dB]
- T temperature, [°C]
- Q arc power, [W]
- t time, [s]
- U arc voltage, [V]
- $v_{\rm w}$ welding speed, [ms⁻¹]
- W weld bead width [mm]
- x unknown parameters
- Y value of the observed characteristic

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- thermal conductivity, [Wm⁻¹K⁻¹]
 - density, [kgm⁻³]
 - axis of moving co-ordinate system

- arc efficiency, [%]

Subscripts

Greek symbols

η

λ

ρ

Ĕ

- $i i^{th}$ parameter
- I-IV– specimen number
- i -number of specimens used for calibration
- liq liquidus
- sol solidus

Superscripts

- exp experimental
- sim simulated

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