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Deep neural network method for predicting the iron concentration in silicon solar cells by current-voltage characteristics

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Corresponding Author:	Oleg Olikh, Dr. hab. Taras Shevchenko National University of Kyiv: Kiivs'kij nacional'nij universitet imeni Tarasa Sevchenka Kyiv, UKRAINE
First Author:	Oleg Olikh, Dr. hab.
Order of Authors:	Oleg Olikh, Dr. hab. Oleg Lozitsky Oleksii Zavhorodnii
Abstract:	<p>Defect-assisted recombination processes frequently limit the photovoltaic device performance. The low-cost and express methods of impurity contamination control are in demand at solar cell manufacturing. In this paper, we applied deep learning-based approach to extract the iron concentration in silicon solar cell from an ideality factor values.</p> <p>SCAPS-1D was the software of choice for simulation of solar cell with the back surface field design and for generation of labeled training and test datasets. Our results demonstrated the deep neural network ability to predict iron concentration with mean squared relative error up to 0.005 by using ideality factor values, SC base thickness and doping level, and temperature. The simulation shows that the prediction error is reduced for high doping level, low temperature, and using of two values of ideality factor (for structure with interstitial iron atoms only as well as for structure with coexistence of Fei and iron-boron pair).</p>
Suggested Reviewers:	<p>Marcelo Cappelletti Universidad Nacional Arturo Jauretche mcappelletti@unaj.edu.ar He has a great skill in solar cell simulation (Superlattices Microstruct. 123 (2018) 338-348, Superlattices Microstruct. 107 (2017) 136-143, Semicond. Sci. Technol. 31 (2016) 115020)</p> <p>Ziv Hameiri National University of Singapore ziv.hameiri@gmail.com He has a great skill in evaluation of recombination processes by using ideality factor (Mater. Sol. Cells 117 (2013) 251-258)</p> <p>Ashraf Uddin, Prof. University of New South Wales a.uddin@unsw.edu.au He has investigated the using of ideality factor to estimate the recombination in solar cell (IEEE Journal of Photovoltaics 8 (2018) 1701-1709)</p> <p>Rachel Kurchin, Dr Massachusetts Institute of Technology rkurchin@alum.mit.edu Her papers are devoted to extract defect properties from inexpensive IV measurements of completed devices</p>

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Junichiro Shiomi, Prof

The University of Tokyo: Tokyo Daigaku

shiomi@photon.t.u-tokyo.ac.jp

His research interests focus on materials informatics, which combines material property calculations/measurements and informatics algorithms.

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To: Physics Letters A Editorial Board
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Dear Editors,

Enclosed with this letter you will find the electronic submission of manuscript entitled “Deep neural network method for predicting the iron concentration in silicon solar cells by current-voltage characteristics” O. Olikh, O. Lozitsky, and O. Zavhorodnii.

This is an original paper which has not been simultaneously submitted as a whole or in part anywhere else. No conflict of interest exists in the submission of this manuscript.

It is well known that impurities are crucial for the solar cells performance. There are many experimental methods for impurity evaluation, such as the infrared spectroscopy, deep level transient spectroscopy, photoluminescence, thermally stimulated capacitance and current, secondary ion mass spectrometry etc. These methods are complicated enough and demand a special setup. At the same time, there is a simpler and commonly used technique, which is the analysis of the solar cell current–voltage characteristics. The present manuscript describes the method of contaminant concentration evaluation by using the ideality factor value, which extracted from current-voltage curve. The method is based on results of numerical simulation of solar cells and use a deep neural network. We believe that such way of defect characterisation would be of interest to the readers.

We would very much appreciate if you would consider the manuscript for publication in the *Physics Letters A*.

Sincerely yours,
Oleg Olikh and co-authors
Taras Shevchenko National University of Kyiv
Kyiv 01601, Ukraine
E-mail: olegolikh@knu.ua

Highlights

Deep neural network method for predicting the iron concentration in silicon solar cells by current-voltage characteristics

Oleg Olikh, Oleg Lozitsky, Oleksii Zavhorodnii

- Proposed deep learning-based method to predict iron contamination in Si-SC by using IV curve.
- The simulated IV characteristics are used to create training and test datasets.
- The DNN's configurations are proposed.
- The mean squared relative error of prediction is up to 0.005.

Deep neural network method for predicting the iron concentration in silicon solar cells by current-voltage characteristics

Oleg Olikh^{a,*}, Oleg Lozitsky^a and Oleksii Zavhorodnii^a

^aTaras Shevchenko National University of Kyiv, 64/13, Volodymyrska Street, City of Kyiv, Ukraine, 01601

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ABSTRACT

Defect-assisted recombination processes frequently limit the photovoltaic device performance. The low-cost and express methods of impurity contamination control are in demand at solar cell manufacturing. In this paper, we applied deep learning-based approach to extract the iron concentration in silicon solar cell from an ideality factor values. SCAPS-1D was the software of choice for simulation of solar cell with the back surface field design and for generation of labeled training and test datasets. Our results demonstrated the deep neural network ability to predict iron concentration with mean squared relative error up to 0.005 by using ideality factor values, SC base thickness and doping level, and temperature. The simulation shows that the prediction error is reduced for high doping level, low temperature, and using of two values of ideality factor (for structure with interstitial iron atoms only as well as for structure with coexistence of Fe_i and iron-boron pair).


1. Introduction

Metal contamination control remains an important challenge for silicon processing both for microelectronics, logic technologies and solar cells (SCs) [1, 2, 3, 4]. Typically, metal related defect characterization is performed by Fourier-transform infrared spectroscopy, electron-paramagnetic resonance, minority carrier lifetime measurements, deep level transient spectroscopy (DLTS), Laplace DLTS etc [5, 6, 7]. However, these techniques are time-consuming, require special equipment or/and sample preparing. At the same time, the current-voltage (IV) measurements are the standard rapid industrial SC characterization technique. IV characteristics contain important information about electrically active defects [6, 8]. And several methods are proposed for diagnosing of defects with using of the IV characteristic. [6, 8, 9, 10, 11]. The temperature dependencies of current components [10, 11] or IV differential parameters [8, 9] are under consideration. But the numerous and high accuracy IV measurements are required in the first and second cases, respectively.

In our previous work [12], we have shown that the SC ideality factor value (n) can be used to estimate the iron concentration (N_{Fe}). It should be noted that the ideality factor is quite often used to characterize the various semiconductor barrier structures [13, 14, 15, 16, 17]. However, a defect's signature in an ideality factor is convoluted with those from so many other physical processes. As a result, obtained analytic expressions $N_{\text{Fe}} = f(n)$ are not general and the numerous grading curves are required to determine N_{Fe} ; besides the IV measurements over a temperature range are necessary [12]. On the other hand, in the last decade, the deep learning, which enables to solve problems without clear algorithmization, have been successfully used in various fields of theoretical and applied physics [18, 19, 20]. Furthermore, materials informatics (combination of material property calculations/measurements and informatics algorithms) has been asserted [21] to become the fourth (along with theory, simulations, and experiments) paradigm of science. The aim of this work is to apply the deep learning approach for predicting of the iron concentration from ideality factor (so to say "deep learning for deep levels"). Further, unlike in previous work [12], the back surface field (BSF) n^+-p-p^+ structure was under consideration and the influence of the base thickness on ideality factor was taken into account as well.

As the approximation to the practical using, the paper considers a fairly simple system which consists of crystalline silicon (c-Si) SC and iron impurity. However, the system is important in practice. Silicon solar cells constitute 90% of current global production capacity [22] and BSF is one of popular designs used for industrial mass production of c-Si SCs [23]. Iron is a major as well as one of the most detrimental metallic impurities in c-Si SCs [2, 3, 4]. The flowchart

*Corresponding author

 olegolikh@knu.ua (Oleg Olikh)

ORCID(s):

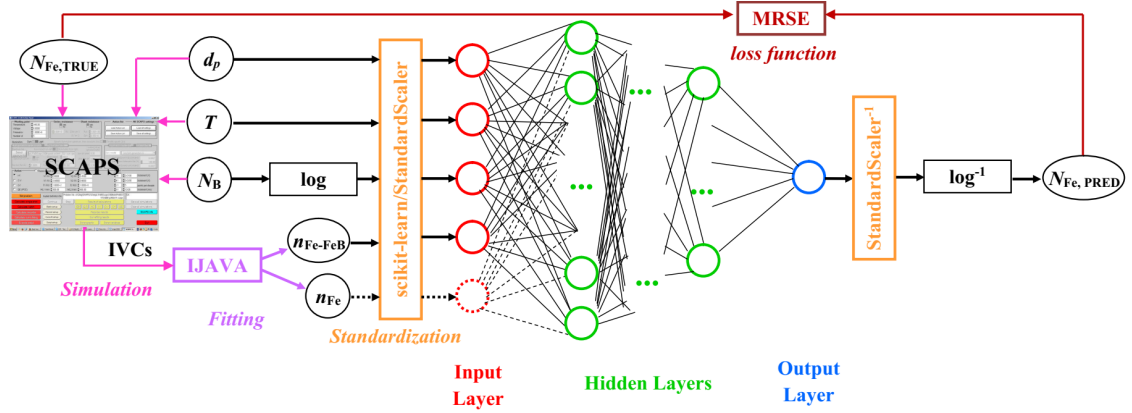


Figure 1: Schematic of deep learning based approach for predicting the iron concentration. Additional details are discussed in the body of the article.

of the used heuristic approach is shown in Fig. 1. The following constituents can be distinguished. First, the dark IV characteristics are simulated for SCs with both known contaminant composition and various parameters. In our numerical simulation we applied SCAPS-1D [24, 25], which widely used to model solar cells [26, 27, 28, 29, 30, 31]. Second, the obtained characteristic is fitted according to the double-diode model and the ideality factor is estimated. As a result of aforesaid steps, the labeled datasets were produced. Obviously, the labeled dataset from experimental IVs would be preferable, but it is practically difficult to find the thousands of samples with the required parameters. Third, the training of deep neural network (DNN) to estimate an iron contamination by using SC's base thickness, doping level, temperature, and ideality factor value. Fours, the DNN testing.

2. Simulation Details

The presented calculation uses n^+p-p^+ structure: the emitter layer n^+ with the donor concentration $N_D = 10^{19} \text{ cm}^{-3}$ and the thickness $0.5 \mu\text{m}$; p and p^+ are uniformly doped with boron; the base p with the thickness $d_p = 150\text{--}240 \mu\text{m}$ is doped with concentration $N_B = 10^{15}\text{--}10^{17} \text{ cm}^{-3}$ and the BSF-layer p^+ with the thickness d_{BSF} ($1 \mu\text{m}$) and the acceptor concentration N_{BSF} ($5 \times 10^{18} \text{ cm}^{-3}$).

The simulations were carried out over the temperature range $290\text{--}340 \text{ K}$. The SCAPS setting file was created for each temperature using the following material parameters. The bandgap E_G and bandgap narrowing ΔE_G models are, respectively, from Pässler [32] and Yan and Cuevas [33]:

$$E_G = E_{G0} - \alpha \Theta \left\{ \frac{1 - 3\Delta^2}{e^{\frac{\Theta}{T}} - 1} + \frac{3\Delta^2}{2} \left(\sqrt{1 + \frac{\pi^2}{3(1 + \Delta^2)} \left(\frac{2T}{\Theta} \right)^2} + \frac{3\Delta^2 - 1}{4} \left(\frac{2T}{\Theta} \right)^3 + \frac{8}{3} \left(\frac{2T}{\Theta} \right)^4 + \left(\frac{2T}{\Theta} \right)^6 - 1 \right) \right\}, \quad (1)$$

$$\Delta E_G = 4.20 \times 10^{-5} \left[\ln \left(\frac{N_D}{10^{14}} \right) \right]^3; \quad \Delta E_G = 4.72 \times 10^{-5} \left[\ln \left(\frac{N_{B,BSF}}{10^{14}} \right) \right]^3, \quad (2)$$

where $E_{G0} = 1.1701 \text{ eV}$, $\alpha = 3.23 \times 10^{-4} \text{ eV/K}$, $\Theta = 446 \text{ K}$, $\Delta = 0.51$. The carrier thermal velocities are calculated from models by Green [34]:

$$v_{th,n} = \sqrt{\frac{8qkT}{0.28m_0\pi}}; \quad v_{th,p} = \sqrt{\frac{8qkT}{0.41m_0\pi}}, \quad (3)$$

where m_0 is the free electron mass. The effective states density masses in the conduction band m_{dC}^* and the valence band m_{dV}^* are calculated according to models from Couderc et al. [35]:

$$\left(\frac{m_{dC}^*}{m_0} \right)^{1.5} = 1.094 - 1.312 \times 10^{-5}T + 6.753 \times 10^{-7}T^2 + 4.609 \times 10^{-10}T^3, \quad (4)$$

$$\left(\frac{m_{dV}^*}{m_0}\right)^{1.5} = 0.3426 + 3.376 \times 10^{-3}T - 4.689 \times 10^{-6}T^2 + 2.525 \times 10^{-9}T^3. \quad (5)$$

The carrier mobilities and the free carrier effective masses were taken from Klaassen [36] and O'Mara et al. [37], respectively. The temperature and doping dependencies of Auger recombination coefficients are calculated from models by Altermatt et al. [38]:

$$C_p(T) = (7.91 \times 10^{-32} - 4.13 \times 10^{-35}T + 3.59 \times 10^{-37}T^2) \times \left(1 + (564812T^{-1.6545} - 1) \left(1 - \tanh \left[\left\{\frac{p}{5 \times 10^{16}}\right\}^{0.29}\right]\right)\right), \quad (6)$$

$$C_n(T) = 2.8 \times 10^{-31} \times \left(1 + (235548T^{-1.5013} - 1) \left(1 - \tanh \left[\left\{\frac{n}{5 \times 10^{16}}\right\}^{0.34}\right]\right)\right). \quad (7)$$

The Auger recombination coefficient was taken from Nguyen et al. [39].

The outside surface recombination with electron and hole velocities 10^3 cm/s was taken into account.

The simulations are carried out under the assumption that the defect-assisted recombination corresponds to the iron-related deep levels only. As the base and the SBF-layer uniform contaminant, iron is assumed to be in concentration $N_{Fe} = 10^{10}$ – 10^{13} cm $^{-3}$. The simulations have been performed for the following two cases. In the first one, the concentration of totally dissolved iron is given by a sum of concentrations of the interstitial iron Fe_i and the trigonal iron-boron pair Fe_iB_s :

$$N_{Fe} = N_{Fe_i} + N_{Fe_iB_s}. \quad (8)$$

The defect distributions are inhomogeneous, and depend on the Fermi level F position, and are given by [40, 41]:

$$\frac{N_{FeB}}{N_{Fe}} = \frac{N_B 10^{-23} \exp\left(-\frac{E_b}{kT}\right)}{\left[1 + \frac{N_B}{10^{23}} \exp\left(-\frac{E_b}{kT}\right)\right] \left[1 + \exp\left(-\frac{F-E_{Fe_i}}{kT}\right)\right]}, \quad N_{Fe_i} = N_{Fe} - N_{FeB}, \quad (9)$$

where $E_b = 0.582$ eV is the binding energy of the Fe_iB_s pairs, E_{Fe_i} is the donor level, associated with Fe_i . This case corresponds to the equilibrium condition and in this article will be referred as “Fe-FeB”.

In the second one, the Fe_i is suggested to be presented only with homogeneous distribution ($N_{Fe_i} = N_{Fe}$). This case can be realized by heat treatment (210°C, 3 min) [42] or intense illumination [43] and will be referred as “Fe-FeB” hereafter.

The donor level $E_{Fe_i} = E_V + 0.394$ eV with electron $\sigma_{n,Fe} = 3.47 \times 10^{-11}T^{-1.48}$ cm 2 and hole $\sigma_{p,Fe} = 4.54 \times 10^{-16} \exp\left(-\frac{0.05}{kT}\right)$ cm 2 capture cross-sections [40, 44] is associated with Fe_i in simulations. The donor level $E_{FeB}^D = E_V + 0.10$ eV, $\sigma_{n,FeB}^D = 4 \times 10^{-13}$ cm 2 , $\sigma_{p,FeB}^D = 2 \times 10^{-14}$ cm 2 and acceptor level $E_{FeB}^A = E_C - 0.26$ eV, $\sigma_{n,FeB}^A = 5.1 \times 10^{-9}T^{-2.5}$ cm 2 , $\sigma_{p,FeB}^A = 3.32 \times 10^{-10} \exp\left(-\frac{0.262}{kT}\right)$ cm 2 [45, 40, 44] are used for Fe_iB_s .

The dark forward IV characteristic were generated by SCAPS over a voltage range up to 0.45 V. According to the two-diode model, the dark SC current is given by [46]

$$I = I_{01} \left[\exp\left(-\frac{q(V - R_s I)}{kT}\right) - 1 \right] + I_{02} \left[\exp\left(-\frac{q(V - R_s I)}{nkT}\right) - 1 \right] + \frac{V - R_s I}{R_{sh}}, \quad (10)$$

where I_{01} and I_{02} are the saturation currents, R_{sh} and R_s are the shunt and series resistances. The two-diode model is often applied for description of real Si SCs and we used Eq. 10 to fit the simulated data by taking n , I_{01} , I_{02} , R_{sh} , and R_s as fitting parameters. The fitting was performed by using of the meta-heuristic method IJAVA [47]. It should be noted that influence of both R_s (obtained values $< 10^{-2}$ Ω) and R_{sh} (obtained values $> 10^{18}$ Ω) can be neglected in simulated IV.

It is the ideality factor value n which is used in our further calculation. The ideality factors, which are obtained in Fe-case and Fe-FeB-case, are referred as n_{Fe} and n_{Fe-FeB} hereafter. The typical simulated dependencies of the ideality factor are shown in Fig. 2. The detailed discussion about n_{Fe} and n_{Fe-FeB} values are presented elsewhere [48], however it should be noted that (i) n can take equal values for different values of SC parameters; (ii) dependencies of n_{Fe} and n_{Fe-FeB} vary not only in absolute values but although in behavior slightly.

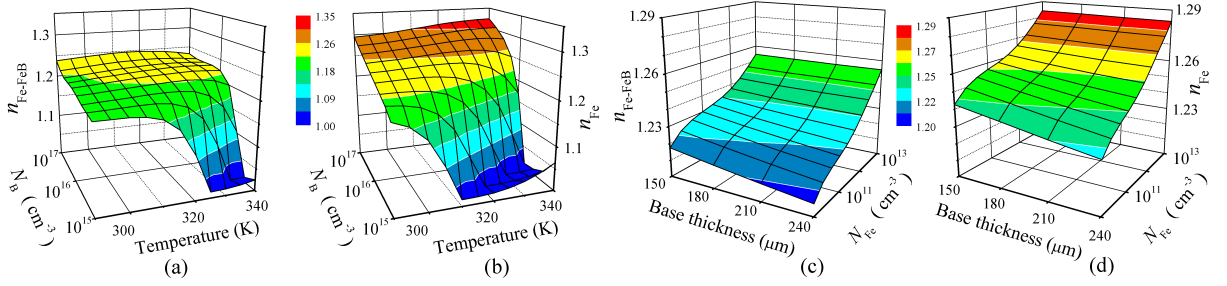


Figure 2: Ideality factor versus temperature and boron concentration (a, b) or base thickness and iron concentration (c, d). The Fe-FeB-case (a, c) and Fe-case (b, d). $N_{\text{Fe}} = 10^{10} \text{ cm}^{-3}$ (a, b), $d_p = 180 \text{ μm}$ (a, b), $N_B = 10^{16} \text{ cm}^{-3}$ (c, d), $T = 320 \text{ K}$ (c, d).

3. Deep neural network models

Training a deep neural network requires a large number of samples. In order to build a training dataset, we used IV characteristics, which are simulated with using of 4 d_p values, 9 N_B values, 11 T values and 19 N_{Fe} values. These base thickness, doping level, temperature, and iron concentration values are regularly (for T and d_p in linear scale, for N_{Fe} and N_B in logarithmic scale) distributed over the ranges 150–240 μm, 10¹⁵–10¹⁷ cm⁻³, 290 – 340 K, and 10¹⁰–10¹³ cm⁻³, respectively. Thus, 7524 IV characteristics are simulated in Fe-case as well as in Fe-FeB-case to build a training dataset.

Besides, several test datasets are prepared. The d_p , N_B , and N_{Fe} values, which equal to values from training dataset, and T values, which is divergent from training dataset, are used to build the test dataset, labeled “T-varied”. These dataset is based on 894 pairs of IV characteristics. The similar approach was used to prepare “d-varied” (1189 samples), “Fe-varied” (856 samples), and “B-varied” (514 samples) test datasets. The base thickness, doping level, temperature, and iron concentration values, which are divergent from training dataset values, are used to prepare “All-varied” (684 samples).

The precise values of parameters are listed in Supplementary Material.

We have tried to construct the DNN, which is able to estimate iron contamination by using SC parameters (d_p and N_B), measurement temperature, and result of IV fitting (ideality factor value). As it is shown in Fig. 1 two DNNs with different input parameters were under consideration. The input sample of the first one consist of $\{d_p, \log N_B, T, n_{\text{Fe-FeB}}\}$. In practice, these input set can be obtained from one dark IV measurement. These neural network referred as DNN_{FeFeB} hereafter. The second one uses $\{d_p, \log N_B, T, n_{\text{Fe-FeB}}, n_{\text{Fe}}\}$ in input layer. In practice, the obtaining of such a set requires additional SC processing (e.g. intense illumination) and two IV measurements. The label DNN_{FeFeB-Fe} is used below.

The dense deep neural network was implemented through the high-level Keras API provided by TensorFlow [49]. The input layers consist of 4 or 5 nodes — see Fig. 1. 1 node and linear activation were used in output layer. The five configurations of hidden layers were under consideration: (i) “pipe”: each hidden layer contains equal number of nodes; (ii) “trapezium”: six hidden layers, number of neurons linearly decreases from 100% (first layer) to 50% (last layer); (iii) “triangle”: ten layers, number of neurons linearly decreases from 100% (first layer) to 10% (last layer); (iv) “butterfly”: two serial reflected trapezium configurations; (v) “fir”: two serial trapezium configurations.

The mean squared relative error (MSRE) was chosen as the loss function:

$$\text{MSRE} = \frac{1}{N_s} \sum_{i=1}^{N_s} \frac{(N_{\text{Fe,TRUE},i} - N_{\text{Fe,PRED},i})^2}{N_{\text{Fe,TRUE},i} \cdot N_{\text{Fe,PRED},i}}, \quad (11)$$

where N_s is the number of samples in dataset, $N_{\text{Fe,TRUE},i}$ is the iron concentration, which was used in simulation of i -th sample, $N_{\text{Fe,PRED},i}$ is the DNN prediction for i -th sample.

Hyperparameters include the number of nodes for first hidden layer, the number of hidden layers (in pipe configuration), the batch size, the activation function, the optimizer, the learning rate, the preprocessing method, the dropout rate, the regularization function, the regularization rate, and the weight initializer. A grid search (coarse tuning to limit one hyperparameter) and random search (fine tuning) were performed over the predefined hyperparameter space,

shown in Table 1, and the best hyperparameter combination is chosen.

Table 1

Hyperparameter space for DNNs.

Hyperparameter	Values
# nodes for first hidden layer	30, 40, 50, 75, 100, 120, 150
# hidden layers	4, 5, 6, 8, 10, 15
batch size	8, 16, 32, 64, 128
activation function	ReLU, sigmoid, tanh, SELU, ELU
optimizer	SGD, RMSprop, Adam, Adadelta, Adagrad, Adamax, Nadam, Ftrl
learning rate	10^{-5} , 10^{-4} , 10^{-3} , 10^{-2}
# epochs	100, 300, 400, 600, 1000, 1500
preprocessing method	StandartScaler, MinMaxScaler
regularization function	None, L2, L1, Dropout
regularization rate	10^{-5} , 10^{-4} , 10^{-3} , 10^{-2}
dropout rate	0.2, 0.3, 0.4, 0.5
weight initializer	Xavier Normal or Uniform, He Normal or Uniform, Random Normal or Uniform, Ones

10-fold cross-validation was used to estimate DNN training. The MSRE, coefficient of determination R^2 , and coefficient of correlation R were three metrics used to evaluate the performance of the DNN models on test datasets. Finally, in order to increase a DNNs performance, the full dataset, which consists of training dataset and all test datasets, was used for the training models.

4. Results and Discussion

The results of hyperparameter search are listed in Table 2. In particular, the trapezium and pipe configurations are chosen for $\text{DNN}_{\text{FeFeB}}$ and $\text{DNN}_{\text{FeFeB-Fe}}$ respectively.

Table 2

Chosen hyperparameter combinations.

Hyperparameter	$\text{DNN}_{\text{FeFeB}}$	$\text{DNN}_{\text{FeFeB-Fe}}$
# nodes for hidden layers	120, 108, 96, 84, 72, 60	100, 100, 100, 100
batch size	32	32
activation function	ReLU	ELU
optimizer	Adamax	Adamax
learning rate	10^{-3}	10^{-3}
# epochs	400	1500
preprocessing method	StandartScaler	StandartScaler
regularization function	None	None
weight initializer	Xavier Normal	Xavier Normal

The training and test results of $\text{DNN}_{\text{FeFeB}}$ are presented in Table 3, Table 4, and Fig. 3. As we can see, that MSRE of $\text{DNN}_{\text{FeFeB}}$ prediction is sufficiently large. But it should be noted, that the fraction of prediction with a great difference between $N_{\text{Fe,TRUE},i}$ and $N_{\text{Fe,PRED},i}$ is not big in most cases. Thus squared relative error (SRE) does not exceed 0.05 for 87%, 88%, and 96% samples from T-varied, d-varied and Fe-varid datasets respectively — see bars in Fig. 3. For B-varied dataset (with doping level value, which was not used in the training dataset) the biggest MSRE = 1.06 connects to occurrence of some samples with a really big SRE (>20). While SRE is less than 0.05 in 54% of samples from B-varied test dataset. The worst predictions is quite expectedly to be observed for the All-varied dataset: R^2 equals to 0.813 and SRE is less than 0.05 for 18% only. On the other hand, the Fe-varied dataset is most similar to real demand. And determination and correlation coefficients are high enough (0.991 and 0.996) in this case.

The dependencies on DNN prediction error on SC parameters values are under consideration too — see Figs. 4–7. These figures represent data for training dataset, the results for test datasets are quite similar. Thus Fig. 4(a) shows that

Table 3

Results of 10-fold cross-validation

Dataset	MSRE	
	$\text{DNN}_{\text{FeFeB}}$	$\text{DNN}_{\text{FeFeB-Fe}}$
training	0.31 ± 0.07	0.03 ± 0.01
full	0.28 ± 0.05	0.03 ± 0.01

Table 4

DNN's testing results

Dataset	$\text{DNN}_{\text{FeFeB}}$			$\text{DNN}_{\text{FeFeB-Fe}}$		
	MSRE	R^2	R	MSRE	R^2	R
T-varied	0.41	0.936	0.967	0.020	0.994	0.997
d-varied	0.37	0.961	0.980	0.018	0.996	0.998
B-varied	1.06	0.881	0.939	0.084	0.991	0.995
Fe-varied	0.06	0.991	0.996	0.005	0.999	0.999
All-varied	0.54	0.813	0.901	0.138	0.948	0.974

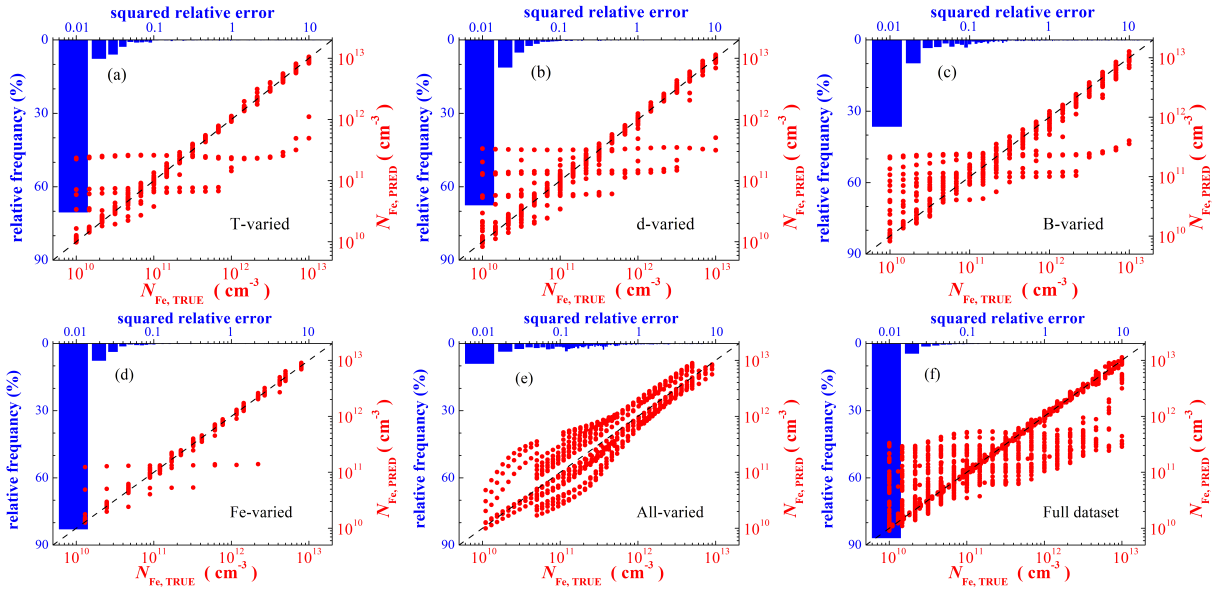


Figure 3: Iron concentrations are plotted against those generated by $\text{DNN}_{\text{FeFeB}}$ on T-varied (a), d-varied (b), B-varied (c), Fe-varied (d), All-varied (e), and full (f) datasets (red points). Bars represent histograms of squared relative error. DNN was learned by training (a)–(e) or full (f) dataset. The black dashed lines are the identify lines servings as the references.

the considerable increase in prediction error value is observed at $T > 320$ K for $\text{DNN}_{\text{FeFeB}}$. So the maximum SRE is about 20 and the SRE is less than 0.01 for 55% of samples at $T = 340$ K (Fig. 4(c)). Whereas those values are equal to 0.02 and 83% at $T = 290$ K (Fig. 4(b)). It has been shown previously [48], that temperature rise causes the increase in the intrinsic recombination's contribution to an ideality factor. As a result, the sign of Shockley-Read-Hall (SRH) recombination in n value become less evident and DNN predictive ability falls.

As it is shown on Fig. 5, the SC base thickness has no influence on the prediction error (mean value as well as relative frequency) practically. But one can see in Fig 2(c,d), that an ideality factor value depend on base thickness at constant N_{Fe} . Therefore d_p is important parameters for DNN training.

The predictive error rises sharply with doping level decrees — see Fig. 6(a). Thus maximum SRE is about 0.05 for $N_B = 10^{17} \text{ cm}^{-3}$ (Fig. 6(c)) whereas squared relative error is less than 0.05 for 56% of samples only for

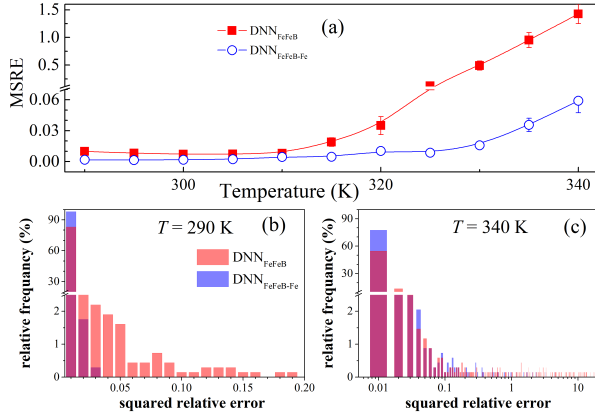


Figure 4: (a) Dependence of the MSRE (training dataset) on the temperature. (b),(c) Histograms of squared relative error for $T = 290$ K and $T = 340$ K. Red: DNN_{FeFeB}; blue: DNN_{FeFeB-Fe}.

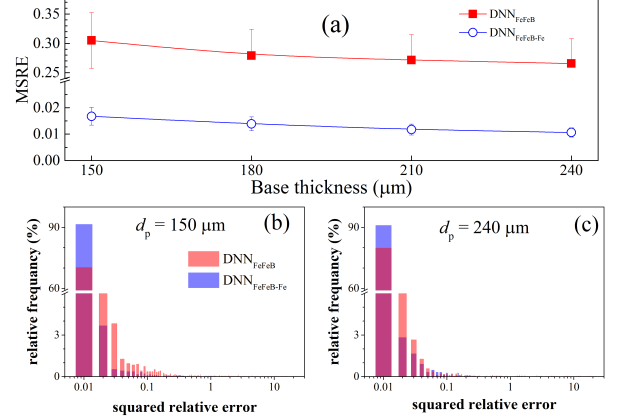


Figure 5: (a) Dependence of the MSRE (training dataset) on the base thickness. (b),(c) Histograms of squared relative error for $d_p = 150$ μm and $d_p = 240$ μm. Red: DNN_{FeFeB}; blue: DNN_{FeFeB-Fe}.

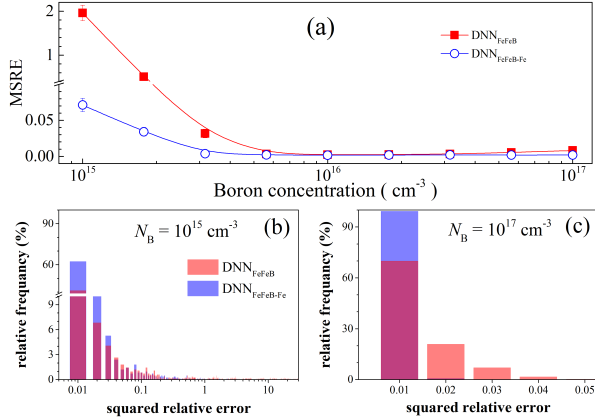


Figure 6: (a) Dependence of the MSRE (training dataset) on the boron concentration. (b),(c) Histograms of squared relative error for $N_B = 10^{15}$ cm⁻³ and $N_B = 10^{17}$ cm⁻³. Red: DNN_{FeFeB}; blue: DNN_{FeFeB-Fe}.

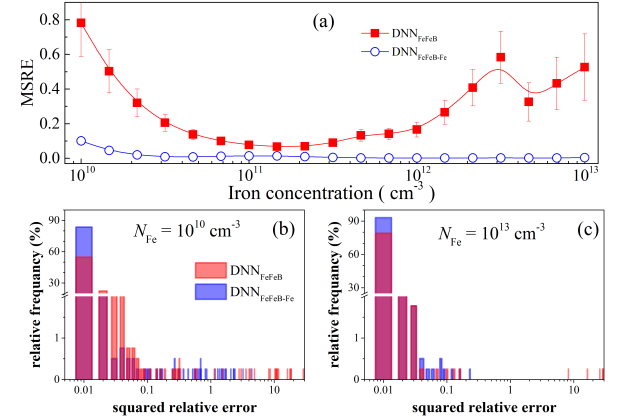


Figure 7: (a) Dependence of the MSRE (training dataset) on the iron concentration. (b),(c) Histograms of squared relative error for $N_{Fe} = 10^{10}$ cm⁻³ and $N_{Fe} = 10^{13}$ cm⁻³. Red: DNN_{FeFeB}; blue: DNN_{FeFeB-Fe}.

$N_B = 10^{15}$ cm⁻³ (Fig. 6(b)). The hole occupation of the Fe-related level determines the SRH recombination efficiency. Accordingly to the Fermi-Dirac statistics, the probability of a hole occupation in a non-degenerate p -type semiconductor with full acceptor depletion can be expressed as

$$f_p = \frac{1}{1 + \frac{N_V}{N_B} \exp\left(\frac{E_V - E_{Fe_i}}{kT}\right)}. \quad (12)$$

If N_B decreases, the level is filled with an electron, the SRH recombination ceases, and the ideality factor value sharply reduces — Fig. 2(a,b). Besides, a weak influence of impurities on ideality factor under low doping condition is a reason of observed MSRE increase. In our opinion, the level filling is additional reason of an error increase at high temperature as well.

Fig. 7(a) shows that MSRE increases at both low and high iron concentrations. First N_{Fe} area of bad DNN accuracy is quite foreseeable, second one is surprising enough. But according to Fig 7(c), the MSRE increasing at $N_{Fe} =$

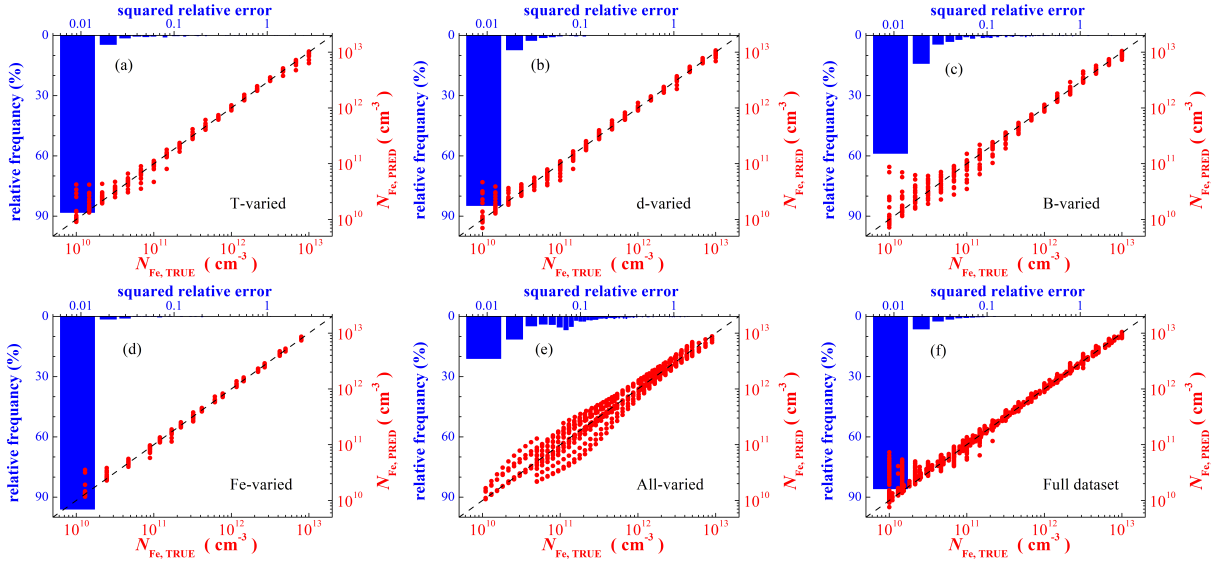


Figure 8: Iron concentrations are plotted against those generated by DNN_{FeFeB-Fe} on T-varied (a), d-varied (b), B-varied (c), Fe-varied (d), All-varied (e), and full (f) datasets (red points). Bars represent histograms of squared relative error. DNN was learned by training (a)–(e) or full (f) dataset. The black dashed lines are the identify lines serving as the references.

10^{13} cm^{-3} is most likely determined by a few samples with big SRE value. Whereas SRE increasing is more systematic at $N_{\text{Fe}} = 10^{13} \text{ cm}^{-3}$ — Fig 7(b).

The ideality factor value for the case of interstitial iron only presence (n_{Fe}) gives additional information about defects in comparing with $n_{\text{Fe-FeB}}$. It is not surprising that DNN_{FeFeB-Fe} has better operating parameters in comparing with DNN_{FeFeB} — see Table 3, Table 4, Fig. 8. The operation improvement appearances in the MSRE decrease as well as in absence of huge difference between $N_{\text{Fe,TRUE}}$ and $N_{\text{Fe,PRED}}$ values and narrowing of SRE range (Figs. 4–8). Really, it is shown in Fig. 8 that the maximum SRE does not exceed 1 even in the case of All-varied datasets and SRE is less than 0.02 for 93%, 92%, 73%, and 97% of samples of T-varied, d-varied, B-varied, and Fe-varied datasets respectively. The R^2 (0.999) and R (0.999) values for Fe-varied dataset draws attention as well.

Despite the difference in predicting accuracy, DNN_{FeFeB-Fe} features are similar to DNN_{FeFeB} ones. Thus the DNN training with N_{B} values, which is expected in object of future research, is important for prediction accuracy (Fig. 8); the increase in the temperature value (Fig. 4) as well as decrease in doping level (Fig. 6) or iron concentration (Fig. 7) results in error rise. It should be noted that the prediction error gain with N_{Fe} increase not observed in DNN_{FeFeB-Fe} case and SRE range at $N_{\text{Fe}} = 10^{13} \text{ cm}^{-3}$ is more narrow then those at $N_{\text{Fe}} = 10^{10} \text{ cm}^{-3}$ — see Fig. 7(b,c).

The results of training both DNN_{FeFeB} and DNN_{FeFeB-Fe} with full dataset are presented in Table 3, Fig. 3(f), and Fig. 8(f). One can see that the extension of labeled dataset does not practically improve DNN result in our case. In our opinion these is evidence of i) a good DNN configuration tuning; ii) a limited predictive ability of DNN_{FeFeB}, which caused by ambiguity of dependence $n_{\text{Fe-FeB}} = f(N_{\text{Fe}})$.

5. Conclusion and Outlook

In this paper, we extracted the iron concentration in silicon BSF solar cell from an ideality factor value and systematically studied the performance of deep learning in this problem. This is the first attempt of using of the deep learning for deep levels parameter retrieval from current-voltage curve. In this model study, we used simulation to obtain training and test labeled datasets. Our results showed the ability of the deep neural network to predict iron concentration with MSRE up to 0.005 by using ideality factor values, SC base thickness and doping level, and temperature. The simulation has shown that two ideality factor values (for structure with Fe_i only as well as with Fe_iB_s and Fe_i coexistence) would be needed to substantially upgrade a prediction accuracy. It was important to train DNN with a boron concentration value, which agreed with doping level of investigated structures. Moreover the increase of iron

concentration or boron concentration as well as temperature decrease results in a prediction error reducing.

The proposed approach envisages the utilization of a simple and widely applicable setup and does not require a much time. Therefore it could be integrated into manufacturing environment. However, it should be noted that we have simplified the task for our purposes. In our opinion there are two ways to further improve method. First one connects with an refining of labeled datasets and can be realized by using 3D-simulators (e.g. SILVACO TCAD) or an real IV measurements in a broad set of SCs. The improving of DNN operation is the second one and the fine-tuning is like most promising in this case. For example, a not numerous input parameters can be multiplied and transformed to the picture and apply a vision model (e.g. VGG16).

CRedit authorship contribution statement

Oleg Olikh: Conceptualization, Methodology, Formal analysis, Data Curation, Writing - Review & Editing, Visualization, Supervision. **Oleg Lozitsky:** Software, Validation, Investigation, Writing - Original Draft. **Oleksii Zavorodnii:** Software, Validation, Formal analysis, Writing - Original Draft.

Data availability

The simulated IV characteristics, n_{Fe} and $n_{\text{Fe-FeB}}$ values, and trained DNNs are available at <https://github.com/olegolikh/IVcharacteristics.git>.

Declaration of competing interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

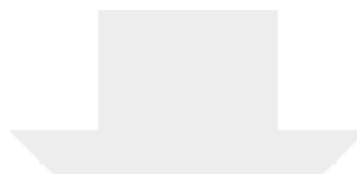
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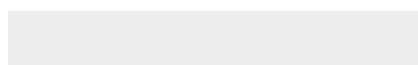
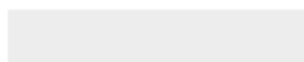
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Declaration of interests

☐The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

☒The authors declare the following financial interests/personal relationships which may be considered as potential competing interests:

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