An artificial neural network for surrogate modeling of stress fields in viscoplastic polycrystalline materials

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The purpose of this work is the development of a trained artificial neural network for surrogate modeling of the mechanical response of elasto-viscoplastic grain microstructures. To this end, a U-Net-based convolutional neural network (CNN) is trained using results for the von Mises stress field from the numerical solution of initial-boundary-value problems (IBVPs) for mechanical equilibrium in such microstructures subject to quasi-static uniaxial extension. The resulting trained CNN (tCNN) accurately reproduces the von Mises stress field about 500 times faster than numerical solutions of the corresponding IBVP based on spectral methods. Application of the tCNN to test cases based on microstructure morphologies and boundary conditions not contained in the training dataset is also investigated and discussed.

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INTRODUCTION

In material science and engineering, material modeling is of central importance for gaining insight into the interplay between material properties, microstructure, and behavior essential for material optimization and design. In the case of polycrystalline structural materials (e.g., steels), for example, a common approach to the modeling of their mechanical behavior is the numerical solution of initial-boundary-value problems (IBVPs) for mechanical equilibrium, for example, via spectral or finite-element methods (e.g., refs. 1-7). Unfortunately, such methods are computationally quite expensive, especially at high resolution or “fidelity” (e.g., refs. 8-10). As such, alternative approaches to the modeling of complex polycrystalline and polyphase materials and their behavior under mechanical loads are the focus of current research.

Among existing alternatives to the currently used solvers, perhaps the most prominent ones are those based on artificial neural networks (ANNs) and machine learning (ML) (e.g., ref. 11). Development of corresponding models is based in particular on the fitting of ANN parameters to data via constrained optimization, i.e., ANN training, yielding a trained ANN (tANN). Except for a few early works (e.g., refs. 12,13), most such “surrogate” models based on tANNS have been introduced in the last 4 years (e.g., refs. 14-18). The data employed for training and testing can be experimental, empirical, or synthetic in nature. An example of the latter are results from the numerical solution of IVBPs based on physical models. In the “data-driven” case, training is based on such data alone. Going beyond this, one can employ physical relations (e.g., constitutive relations) on which the data are based as additional training constraints. In the case of physics-informed neural networks (PINN) (e.g., ref. 19), the data are the initial and boundary conditions of an IBVP, and the ANN is used to approximate a least-squares-based (numerical) solution of the IBVP (e.g., ref. 20). Recent reviews of applications in the field of continuum mechanics and material modeling include, for example, ref. 21. In the current work, attention is focused on the data-driven approach. The data are obtained from the numerical solution of a BVP for quasi-static mechanical equilibrium based on viscoplastic material modeling of grains in a heterogeneous polycrystalline ensemble.

A number of data-driven approaches have been proposed for applications in solid mechanics. For example, Yang et al.22 trained a conditional generative adversarial network (cGAN) to reproduce stress and strain fields in strained isotropic two-phase composites. Mianroodi et al.23 trained a U-Net-based convolutional neural network (CNN) using results for the von Mises stress for grain microstructures consisting of isotropic elastic and ideal elastoplastic grains subject to uniaxial extension. More recently, Rashid et al.24 introduced a neural-operator-based approach, the Fourier Neural Operator (FNO)25,26, in particular for the surrogate modeling of stress and strain in heterogeneous composites. The U-Net-based CNN and FNO-based ANN have been compared recently by Kapoor et al.27 in the context of surrogate modeling of stress fields in heterogeneous elastoplastic solids.

In the current work, a U-Net-based CNN is trained to output the von Mises stress field \(\sigma_{vM}\) in heterogeneous periodic microstructures consisting of inelastic grains subject to uniaxial extension. More specifically, the constitutive behavior of each grain is modeled via \(J_2\) elasto-viscoplasticity with linear isotropic hardening. Results from the numerical solution of BVPs for quasi-static mechanical equilibrium in periodic unit cells based on this grain behavior and spectral numerical solution methods (e.g., refs. 12,17) are employed to train the CNN. For brevity, this is referred to as the “reference model” (RM) in this work. In the “Results” section, results for \(\sigma_{vM}\) and its average \(\overline{\sigma}_{vM}\) over the unit cell \(U\) of the grain microstructure obtained from the trained CNN (tCNN) are presented and compared with corresponding results from the RM. In particular, these include the dependence of the tCNN on details of the training dataset and training such as (i) the number of grains in the microstructure, and (ii) the range of material properties chosen for each grain. In addition, the performance of the tCNN for (i) microstructure morphologies (e.g., matrix-inclusion), and (ii) extension levels, not

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included in the training dataset, is also investigated and discussed. The conclusions and outlook are presented in the “Discussion” section. Methods employed in this work are presented in the “Methods” section. In particular, this includes a brief summary of the model for isotropic J2 viscoplasticity with linear isotropic constitutive response assumed for each grain in the polycrystalline ensemble. Data generation, the U-Net-based network architecture employed, and network training are also explained and discussed.

RESULTS

The purpose of this section is to compare the output of the tCNN for the von Mises stress field $\sigma_{VM}$ with corresponding results from the RM for selected test cases. As explained in more detail in the “Methods” section below, the current CNN is trained with results for $\sigma_{VM}$ on unit cells $U$ with 10-grain microstructures and a resolution of $64 \times 64$ pixels subject to uniaxial extension. Training (80%) and validation (20%) datasets are based on data from 1000 such microstructures generated randomly. For simplicity, data used for training and testing are based on a single (quasistatic) extension rate of $1 \times 10^{-3}$ s$^{-1}$. Test cases here include microstructures with (i) different numbers of grains, (ii) material property contrasts in neighboring grains, and (iii) grain morphologies, which differ from those in the training and validation datasets. The tCNN is also tested for extensions larger than those included in the training dataset, is also investigated and discussed. Note the decrease in property contrast in going from Case 1 to Case 4. This is reflected in the MAEs of the corresponding test datasets shown in Table 2. Note the decrease in MAE with decreasing contrast in material properties. For example, the contrast in $E$ is at least 215 GPa in Case 1, and 145 GPa in Case 4.

Example results for Case 1 and Case 3 are shown in Figs. 4 and 5. The slight difference in the grain shape distributions in Figs. 4 and 5.

As noted above, since the material properties in each grain are homogeneous, the contrast in the property distributions is also related to the contrast in material properties at grain boundaries and triple junctions. In comparison to the results in Figs. 1–3 (b, bottom) for $|\sigma_{CNN} - \sigma_{RM}|$ in the case that material properties values are chosen from the ranges in Table 3, note the increase in $|\sigma_{CNN} - \sigma_{RM}|$ in Fig. 4b (bottom: up to 100 MPa) and Fig. 5b (bottom: up to 80 MPa) due to the larger contrast, especially in the former case. This is also true for the relative error shown in Figs. 4c and 5c. The figure indicates that the results from the tCNN in Case 3 (with a maximum error of 2%) are more accurate than Case 1 in which the maximum error is around 5% as shown in Fig. 4c. Training with a larger dataset in this case would result in a reduction of such errors and better agreement with the RM.

Test microstructures with different morphologies

In this subsection, the tCNN is applied to grain and microstructure morphologies not included in the training dataset. These take the form of a single inclusion embedded in a matrix material. Inclusions of circular and square form are investigated. For each, 50 test results are generated based on a random choice of material properties from the ranges in Table 3. Examples of these are shown in Figs. 6 and 7. Analogous to the case of grain boundaries and triple junctions in the polycrystalline cases in Figs. 1–3 (b, bottom), $|\sigma_{CNN} - \sigma_{RM}|$ is largest at the sharp matrix–inclusion (MI) interface where the contrast in material properties is greatest and pixel resolution is the lowest.

Table 1. MAE for test datasets based on 5- and 20-grain microstructures and for the training dataset based on 10-grain microstructures.

| Number of grains | 5   | 10  | 20  |
|------------------|-----|-----|-----|
| MAE (MPa)        | 1.578 | 1.733 | 1.869 |

Note the maximum in $|\sigma_{CNN} - \sigma_{RM}|$ at the MI interface in the circular case (Fig. 6b (bottom)) at $F_{11} = 1.001$ not present at the MI interface in the square case (Fig. 7b (bottom)). This is due to the fact that (in contrast to the latter system) the former system is still elastic at $F_{11} = 1.001$, as implied by the unit cell stress-deformation results in Fig. 6c and the larger initial resistance stress $\xi_0$ of the circular inclusion and corresponding contrast at the MI interface (Fig. 6a). For comparable contrasts in material properties at the (in particular sharp) interfaces of the circular and square inclusions with the matrix, one expects the largest stress concentration at the corners of the latter, and so the largest values of $|\sigma_{CNN} - \sigma_{RM}|$ over the extension history. This is also reflected by the fact that the MAE of the test dataset for the circular inclusion case (1.42 MPa) is slightly lower than in the square inclusion case (1.58 MPa).
Extension histories not in the training dataset

The training data is based in particular on results for \( \sigma_{RM} \) corresponding to uniaxial extension of the unit cell up to \( F_{11} = 1.004 \). Since \( \sigma_{RM} \) is an input to the network, the tCNN can be used to calculate \( \sigma_{VM} \) for values larger than those in the training data, corresponding to unit cell uniaxial extension beyond that represented by the training data. A typical maximum is just above 500 MPa (see, for example, Fig. 11, bottom right). Results for this are shown in Fig. 8.

Note the significant increase in \( |\sigma_{CNN} - \sigma_{RM}| \) in Fig. 8b (bottom) and in the corresponding relative error in Fig. 8c (red triangles) above \( F_{11} = 1.0065 \).

Figure 8c shows the corresponding stress–strain curve for further loading. According to this curve, the error notably
increases for values of \( \sigma_{vM} \) greater than 500 MPa. One reason for this is very little data in the training data for such values. This can be seen for example in Fig. 11 (below right), which displays the distribution of pixel values for \( \sigma_{vM} \) at different \( F_{11} \). As shown, only 0.084% of the pixels in the whole training dataset have a von Mises stress level above 500 MPa.

### Computational efficiency

We consider next the time required to calculate the von Mises stress field based on the RM and the tCNN. Recall that the former is based on the numerical solution of BVPs for quasi-static mechanical equilibrium in periodic unit cells employing spectral numerical solution methods (e.g., refs. 1, 2, 4, 7). The corresponding results for the RM are obtained using the DAMASK simulation package5 (see next section). The calculations are carried out on a single core of an Intel® Core™ i9900K clocked at 3.60 GHz. On this basis, the run time of DAMASK for a single training simulation averaged over 50 training simulations is approximately 75 s, and the corresponding run time for the tCNN is about 0.15 s. As such, the tCNN is about 500 faster than the RM. Increasing the resolution would result in an even larger difference in computational time. On the other hand, in contrast to the RM, the accuracy of the tCNN is limited to (i) the range of the training data as well as (ii) uniaxial extension.

The current approach can be extended and further developed in a number of directions. These include for example (i) output of all components of the stress field, (ii) training for different deformation and loading conditions, or (iii) training for multiple rates of deformation and/or loading. As well, training can be based on more sophisticated “physics-informed” loss functions (accounting, for example, for mechanical equilibrium), or on more sophisticated “physics-encoded” network architectures. This represents work in progress to be reported on in the future.

### METHODS

#### RM: J2 viscoplasticity with isotropic hardening

As discussed in the introduction, data for the training and testing of the ANN are obtained from the numerical solution of a boundary-value problem for quasi-static mechanical equilibrium based on viscoplastic material modeling of grain behavior in a grain microstructure. In the current context of isothermal and quasi-static conditions, these include in particular mechanical equilibrium \( \text{div}\mathbf{P} = 0 \) in terms of the first Piola–Kirchhoff stress \( \mathbf{P} \). In the context of the viscoplastic decomposition \( \mathbf{F} = \mathbf{F}_e \mathbf{F}_p \) of the deformation gradient \( \mathbf{F} \), the linear elastic relation \( \mathbf{S}_e = \lambda (I \cdot \mathbf{E}_e) + 2\mu \mathbf{E}_e \), \( \mathbf{E}_e \) is

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**Table 2.** MAE for test datasets based on the material property distributions in Eq. (1).

| Case | 1    | 2    | 3    | 4    |
|------|------|------|------|------|
| MAE (MPa) | 3.662 | 1.925 | 1.519 | 1.265 |

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**Fig. 3** Twenty-grain test microstructure subject to uniaxial extension along the horizontal axis. a Young’s modulus \( E \), Poisson’s ratio \( \nu \), initial flow resistance \( \xi_0 \), linear isotropic hardening modulus \( h_0 \). b Results for \( \sigma_{vM} \) (above), \( \sigma_{vM} \) (middle), \( \sigma_{vM} \) (below) (all MPa), at different \( F_{11} \). c \( \sigma_{vM} \) (blue curve), \( \sigma_{vM} \) (green points), and relative error (red triangles) given by 100 \( (\sigma_{vM} - \sigma_{vM})/\sigma_{vM} \) versus \( F_{11} \).
assumed between the elastic second Piola-Kirchhoff stress $S$ and the elastic Green strain $E$:

$$E = \frac{1}{2} \left( F^T F - I \right); \quad \text{note that} \quad F = F_p S F_p^{-T}.$$  

In the following, the Lame constants with $\lambda = E\nu / ((1 + \nu)(1 - 2\nu))$ and $\mu = E/(2(1 + \nu))$ are determined in terms of the Young’s modulus $E$ and the Poisson ratio $\nu$. The evolution of $F_p$.

Fig. 4 Case 1 (high-contrast) test microstructure subject to uniaxial extension along the horizontal axis. a Young’s modulus $E$, Poisson’s ratio $\nu$, initial flow resistance $\xi_0$, linear isotropic hardening modulus $h_0$. b Results for $\sigma_{RM}$ (above), $\sigma_{CNN}$ (middle), $|\sigma_{CNN} - \sigma_{RM}|$ (below) (all MPa), at different $F_{11}$. c $\sigma_{RM}$ (blue curve), $\sigma_{CNN}$ (green points), and relative error (red triangles) given by $100 \left( \sigma_{CNN} / \sigma_{RM} \right)$ versus $F_{11}$.

Fig. 5 Case 3 (low-contrast) test microstructure subject to uniaxial extension along the horizontal axis. a Young’s modulus $E$, Poisson’s ratio $\nu$, initial flow resistance $\xi_0$, linear isotropic hardening modulus $h_0$. b Results for $\sigma_{RM}$ (above), $\sigma_{CNN}$ (middle), $|\sigma_{CNN} - \sigma_{RM}|$ (below) (all MPa), at different $F_{11}$. c $\sigma_{RM}$ (blue curve), $\sigma_{CNN}$ (green points), and relative error (red triangles) given by $100 \left( \sigma_{CNN} / \sigma_{RM} \right)$ versus $F_{11}$.

Table 3. Material property ranges used for generating the dataset.

|          | $E$ (GPa) | $\nu$ | $\xi_0$ (MPa) | $h_0$ (GPa) |
|----------|-----------|-------|----------------|-------------|
| Minimum  | 50        | 0.2   | 50             | 0           |
| Maximum  | 300       | 0.4   | 300            | 50          |
is determined by the $J_2$ flow rule $F_p F_p^{-1} = \dot{\gamma}_p S_{dev}^p / |S_{dev}^p|$, where $\dot{\gamma}_p$ is the rate of equivalent plastic shear, and $S_{dev}$ the deviatoric part of $S_e$. The viscoplastic (i.e., rate-dependent) form $\dot{\gamma}_p = \dot{\gamma}_0 (|S_{dev}^p|/\xi)^{n_0}$ for the evolution of $\dot{\gamma}_p$ is determined by the typical material inelastic shear rate $\dot{\gamma}_0$, the power-law exponent $n_0$, and the flow resistance $\xi_0 \dot{\gamma}_p = \xi_0 + h_0 \dot{\gamma}_p$ for linear isotropic hardening, with $\xi_0$ the initial flow resistance, and $h_0$ the isotropic hardening modulus.

The viscoplastic model is implemented in the simulation software toolkit DAMASK\(^6\). This toolkit is also used for numerical solution of the corresponding quasi-static mechanical boundary-value problem on periodic grain microstructures based on spectral (i.e., Fourier) methods.
Material parameters for each grain in the microstructure include $E$, $\nu$, $\xi_0$, $h_0$, $\gamma_0$, and $n_0$. For simplicity, $\gamma_0 = 10^{-3}$ s$^{-1}$ and $n_0 = 20$ are assumed the same for all grains. Values for the remaining material properties of each grain are chosen randomly from a range of values shown in Table 3. The training dataset consists of 1000 grain microstructures based on 10 grains and random material property distributions. Grain morphologies and microstructures are generated randomly via Voronoi tessellation. An example is shown in Fig. 9.
As done, for example, in ref. 23, the ANN input consists of the distribution of these material properties in the unit cell/microstructure as well as results for the scalar von Mises stress field \( \sigma_{\text{vM}} \) during extension. Given the material property distribution \( (E, \nu, h_0, \xi_0) \) and \( \sigma_{\text{vM}} \) at time step \( t \) (i.e. \( \sigma_{\text{vM}}^t \)), then, the trained ANN (tANN) outputs \( \sigma_{\text{vM}} \) at time step \( t + \Delta t \) (i.e., \( \sigma_{\text{vM}}^{t+\Delta t} \)). This is depicted in Fig. 10.

In the current case of purely bulk behavior of grain microstructures on the unit cell \( U \), fields \( f = \bar{f} + \tilde{f} \) on the unit cell \( U \) are additively split into mean \( \bar{f} \) and fluctuation \( \tilde{f} \) parts, with \( \bar{f} := \nu(U)^{-1} \int_U f \, dv \) and \( \nu(U) := \int_U dv \). In this context, deformation...
“boundary conditions” on $U$ take the form of prescribed values for the mean deformation gradient $\mathbf{F}(t)$. For the current case of uniaxial extension, the Cartesian/matrix form

$$
\mathbf{F}(t) = \begin{bmatrix} \mathbf{F}_{11}(t) & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}, \quad \mathbf{F}_{11}(t) = 1 + \mathbf{F}_{11}(0)t, \quad (2)
$$

of $\mathbf{F}(t)$ applies. For simplicity, data for training and testing are limited to a single extension rate $\mathbf{F}_{11}(0) = 1 \times 10^{-5} \text{s}^{-1}$, i.e., to quasi-static extension. Based on this extension rate, results obtained up to $\mathbf{F}_{11}(4) = 1.004$ are employed as training data, and those between this value and $\mathbf{F}_{11}(8) = 1.008$ are used for testing the tCNN.

The pixel distributions of values for Young's modulus $E$ and $\sigma_{\text{VM}}$ for different values of $\mathbf{F}_{11}$ in the training data are shown in Fig. 11.

The pixel distributions of the other input material properties are similar. As evident, the pixel distribution of $\sigma_{\text{VM}}$ at each value of $\mathbf{F}_{11}$ is quasi-normal in character, as expected for randomly distributed material property values.

**Neural network type, architecture, and training**

The current ANN is based on the U-Net convolutional type and architecture introduced by Ronneberger et al. \(^{29}\). As shown by Mianroodi et al. \(^{23}\), this architecture is suitable for surrogate modeling the stress field in solid mechanics problems, and in particular the von Mises stress $\sigma_{\text{VM}}$. Fig. 12 schematically depicts the U-Net network architecture, (referred to in ref. \(^{29}\) as U-Net due to its shape).

Network input consists of 64 × 64 pixel images for $E$, $\nu$, $\sigma_{\text{VM}}$, and the network outputs one 64 × 64 pixel image for $\sigma_{\text{VM}}$. As usual, both input and output are normalized to 1. In addition, 32 filters capture the main features from the input images. As shown in the figure, four types of operation are performed in the U-Net, namely, separable two-dimensional (2D) convolution with a kernel size of 9, batch normalization, 2D max pooling, and 2D upsampling with bilinear interpolation.

For the training process, Adam optimization is employed with a learning rate of 0.001 and a momentum of 0.9. As discussed above, the loss function for training, validation, and testing is given by the mean absolute error (MAE) based on the difference $\sigma_{\text{CNN}} - \sigma_{\text{VM}}$. TensorFlow\(^{29}\) is used to set up and train the network. As usual, the dataset is divided into training (80%) and (20%) testing subsets. Training is based on 500 epochs and has an MAE of 1.733 MPa; the corresponding MAE for the validation dataset is 1.743 MPa. No sign of overfitting was observed.

**DATA AVAILABILITY**

The data employed in the current work can be obtained from the corresponding author upon request.

**CODE AVAILABILITY**

The code employed in the current work can be obtained from the corresponding author upon request.

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M.S.K. and J.R.M. developed the initial concept and model formulation. M.S.K. carried out the solid mechanics simulations to generate the training and test datasets. N.H.S. designed the neural network, and M.S.K. obtained the neural network parameters. M.S.K., J.R.M., P.G., and B.S. wrote the initial draft. B.S., P.B., and D.R. supervised the work and contributed to the interpretation of the results.

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