

Enforcing Analytic Constraints in Neural Networks Emulating Physical Systems

Tom Beucler^{1,2,*} Michael Pritchard¹ Stephan Rasp³ Jordan Ott⁴ Pierre Baldi⁴ and Pierre Gentine²¹Department of Earth System Science, University of California, Irvine, California 92697-3100, USA²Department of Earth and Environmental Engineering, Columbia University, New York, New York 10027, USA³Technical University of Munich, Boltzmannstr. 3 85748 Garching, Munich, Germany⁴Department of Computer Science, University of California, Irvine, California 92697, USA

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Neural networks can emulate nonlinear physical systems with high accuracy, yet they may produce physically inconsistent results when violating fundamental constraints. Here, we introduce a systematic way of enforcing nonlinear analytic constraints in neural networks via constraints in the architecture or the loss function. Applied to convective processes for climate modeling, architectural constraints enforce conservation laws to within machine precision without degrading performance. Enforcing constraints also reduces errors in the subsets of the outputs most impacted by the constraints.

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Introduction.—Many fields of science and engineering (e.g., fluid dynamics, hydrology, solid mechanics, chemistry kinetics) have exact, often *analytic*, closed-form constraints, i.e., constraints that can be explicitly written using analytic functions of the system's variables. Examples include translational or rotational invariance, conservation laws, or equations of state. While physically consistent models should enforce constraints to within machine precision, data-driven algorithms often fail to satisfy well-known constraints that are not explicitly enforced. In particular, neural networks (NNs) [1], powerful regression tools for nonlinear systems, may severely violate constraints on individual samples while optimizing overall performance.

Despite the need for physically informed NNs for complex physical systems [2–5], enforcing *hard* constraints [6] has been limited to physical systems governed by specific equations, such as advection equations [7–9], Reynolds-averaged Navier-Stokes equations [10,11], boundary conditions of idealized flows [12], or quasigeostrophic equations [13]. To address this gap, we introduce a systematic method to enforce analytic constraints arising in more general physical systems to within machine precision, namely, the architecture-constrained NN or ACnet. We then compare ACnets to unconstrained (UCnets) and loss-constrained NNs (LCnets, in which soft constraints are added through a penalization term in the loss function, e.g., Refs. [14–16]) in the particular case of climate modeling, where the system is high dimensional and the constraints (such as mass and energy conservation) are few but crucial [17].

Theory.—Formulating the constraints: Consider a NN mapping an input vector $\mathbf{x} \in \mathbb{R}^m$ to an output vector $\mathbf{y} \in \mathbb{R}^p$. Enforcing constraints is easiest for linearly-constrained NNs, i.e., NNs for which the constraints ($\mathcal{C}$) can be written as a linear system of rank n :

$$(\mathcal{C}) \stackrel{\text{def}}{=} \left\{ \mathbf{C} \begin{bmatrix} \mathbf{x} \\ \mathbf{y} \end{bmatrix} = \mathbf{0} \right\}. \quad (1)$$

We call $\mathbf{C} \in \mathbb{R}^n \times \mathbb{R}^{m+p}$ the constraints matrix, and use bold font for vectors and tensors to distinguish them from scalars. For the regression problem to have nonunique solutions, the number of independent constraints n has to be strictly less than $m + p$.

In Fig. 1, we consider a generic regression problem subject to analytic constraints ($\mathcal{C}$) that may be nonlinear, and propose how to formulate a linearly constrained NN. First, define the regression's inputs $\mathbf{x}_0$ and outputs $\mathbf{y}_0$, which, respectively, become the *temporary* NN's features and targets. Then (*formulation 1*), write the constraints ($\mathcal{C}$) as an identically zero function $\mathbf{c}$ of the inputs, the outputs, and additional parameters $\mathbf{z}$ the constraints may involve. We recommend nondimensionalizing all variables to facilitate the design, interpretation, and performance of the loss function. While the function $\mathbf{c}$ may be nonlinear, it can always be written as the sum of (i) terms $\mathbf{x}$ that *only* depend on inputs and (ii) terms $\mathbf{y}$ that depend on inputs, outputs, and additional parameters. Thus the constraints can be written as

$$\mathbf{c}(\mathbf{x}_0, \mathbf{y}_0, \mathbf{z}) = \mathbf{C} \begin{bmatrix} \mathbf{x}(\mathbf{x}_0) \\ \mathbf{y}(\mathbf{x}_0, \mathbf{y}_0, \mathbf{z}) \end{bmatrix}, \quad (2)$$

where $\mathbf{C}$ is a matrix. Finally (*formulation 2*), choose $\mathbf{x}$ and $\mathbf{y}$ as the NN's new inputs and outputs. If $\mathbf{x}$ and $\mathbf{y}$ are not bijective functions of $(\mathbf{x}_0, \mathbf{y}_0)$, add variables to the NN's inputs and outputs to recover $\mathbf{x}_0$ and $\mathbf{y}_0$ after optimization (e.g., we add $x_{0,t}$ and $y_{0,t-1}$ to $\mathbf{x}$ in example 2). We are now in a position to build a computationally efficient NN that satisfies the linear constraints ($\mathcal{C}$).

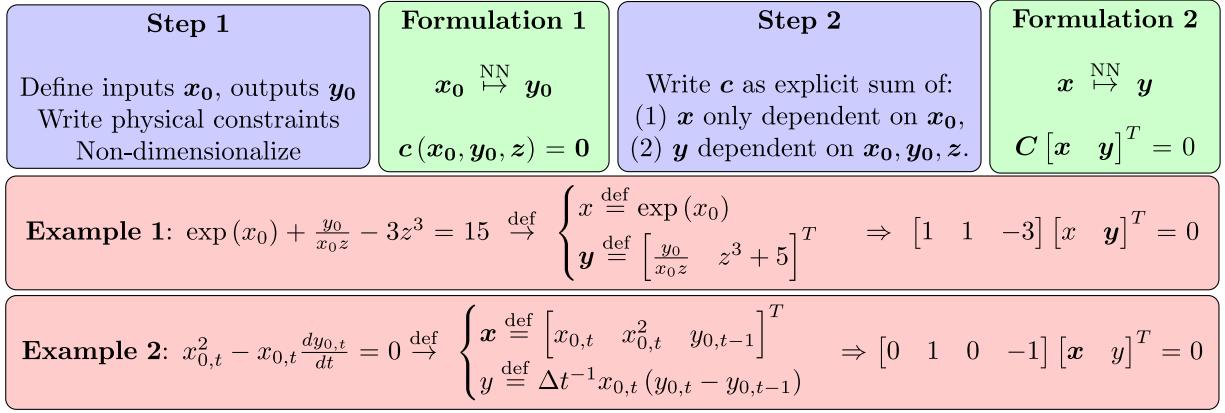


FIG. 1. Framework to treat constrained regression problems using linearly constrained NNs, with two examples: (i) A regression problem with one nonlinear constraint, and (ii) a time-prediction problem with one differential nonlinear constraint that we discretize using a forward Euler method of timestep Δt . Note that the choice of $\mathbf{x}$, $\mathbf{y}$, and $\mathbf{C}$ is not unique.

Enforcing the constraints: Consider a NN trained on preexisting measurements of $\mathbf{x}$ and $\mathbf{y}$. For simplicity's sake, we measure the quality of its output $\mathbf{y}_{\text{NN}}$ using a standard mean-squared error (MSE) misfit:

$$\text{MSE}(\mathbf{y}_{\text{Truth}}, \mathbf{y}_{\text{NN}}) \stackrel{\text{def}}{=} \|\mathbf{y}_{\text{Err}}\|_2 \stackrel{\text{def}}{=} \frac{1}{p} \sum_{k=1}^p y_{\text{Err},k}^2, \quad (3)$$

where we have introduced the error vector, defined as the difference between the NN's output and the "truth":

$$\mathbf{y}_{\text{Err}} \stackrel{\text{def}}{=} \mathbf{y}_{\text{NN}} - \mathbf{y}_{\text{Truth}}. \quad (4)$$

In the reference case of an "unconstrained network" (UCnet), we optimize a multilayer perceptron [18,19] using MSE as its loss function $\mathcal{L}$. To enforce the constraints ($\mathcal{C}$) within NNs, we consider two options: (i) *Constraining the loss function (LCnet, soft constraints)*.—We first test a soft penalization of the NN for violating physical constraints using a penalty $\mathcal{P}$, defined as the mean-squared residual from the constraints:

$$\begin{aligned} \mathcal{P}(\mathbf{x}, \mathbf{y}_{\text{NN}}) &\stackrel{\text{def}}{=} \left\| \mathbf{C} \begin{bmatrix} \mathbf{x} \\ \mathbf{y}_{\text{NN}} \end{bmatrix} \right\|_2, \\ &= \frac{1}{n} \sum_{i=1}^n \left(\sum_{j=1}^m C_{ij} x_j + \sum_{k=1}^p C_{i(k+m)} y_{\text{NN},k} \right)^2, \end{aligned} \quad (5)$$

and given a weight $\alpha \in [0, 1]$ in the loss function $\mathcal{L}$:

$$\mathcal{L}(\alpha) = \alpha \mathcal{P}(\mathbf{x}, \mathbf{y}_{\text{NN}}) + (1 - \alpha) \text{MSE}(\mathbf{y}_{\text{Truth}}, \mathbf{y}_{\text{NN}}). \quad (6)$$

(ii) *Constraining the architecture (ACnet, hard constraints)*.—Alternatively, we treat the constraints as *hard* and augment a standard, optimizable NN with n fixed conservation layers that sequentially enforce the constraints ($\mathcal{C}$)

to within machine precision (Fig. 2), while keeping the MSE as the loss function:

$$(\text{ACnet}) \Rightarrow \{\min \text{MSE} \text{ s.t. } \mathbf{C} [\mathbf{x} \ \mathbf{y}_{\text{NN}}]^T = \mathbf{0}\}. \quad (7)$$

The optimizable NN calculates a "direct" output whose size is $p - n$. We then calculate the remaining output's components of size n as exact "residuals" from the constraints. Concatenating the direct and residual vectors results in the full output $\mathbf{y}_{\text{NN}}$ that satisfies the constraints to within machine precision. Since our loss uses the full output $\mathbf{y}_{\text{NN}}$, the gradients of the loss function are passed through the constraints layers during optimization, meaning that the final NN's weights and biases depend on the constraints ($\mathcal{C}$). ACnet improves upon the common approach of calculating residual outputs *after* training because ACnet exposes the NN to residual output data *during* training (see Supplemental Material [20] C.3). A possible implementation of the constraints layer uses custom (Tensorflow in our case) layers with fixed parameters that solve the system of equations ($\mathcal{C}$), in row-echelon form, from the bottom to the top row (Supplemental Material [20] B.1). Note that we are free to choose which outputs to calculate as residuals, which introduces n new hyperparameters (Supplemental Material [20] B.2).

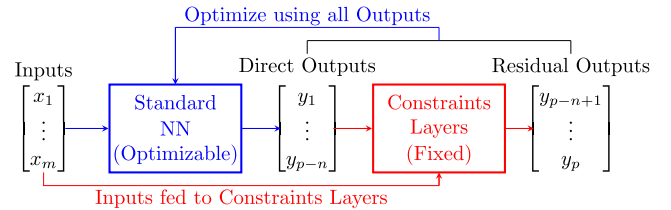


FIG. 2. ACnet: Direct outputs are calculated using a standard NN, while the remaining outputs are calculated as residuals from the fixed constraints layers.

Linking constraints to performance: Intuitively, we might expect the NN performance to improve once we enforce constraints arising in physical systems with few degrees of freedom, but this may not hold true with many degrees of freedom. We formalize the link between constraints and performance by (i) decomposing the NN's prediction into the truth and error vectors following Eq. (4); and (ii) assuming that constraints exactly hold for the truth (no errors in measurement). This yields

$$\mathbf{C} \begin{bmatrix} \mathbf{x} \\ \mathbf{y}_{\text{NN}} \end{bmatrix} \stackrel{\text{def}}{=} \overbrace{\mathbf{C} \begin{bmatrix} \mathbf{x} \\ \mathbf{y}_{\text{Truth}} \end{bmatrix}}^{\mathbf{0}} + \mathbf{C} \begin{bmatrix} \mathbf{0} \\ \mathbf{y}_{\text{Err}} \end{bmatrix}. \quad (8)$$

Equation (8) relates how much the constraints are violated to the error vector. More explicitly, if we measure performance using the MSE, we may square each component of Eq. (8). The resulting equation links how much physical constraints are violated to the squared error for each constraint of index $i \in [1, n]$:

$$\underbrace{\left(\mathbf{C} \begin{bmatrix} \mathbf{x} \\ \mathbf{y}_{\text{NN}} \end{bmatrix} \right)_i^2}_{\text{Physical constraints}} = \underbrace{\sum_{k=1}^p C_{i(k+m)}^2 y_{\text{Err},k}^2}_{\text{Squared-error} > 0} + \underbrace{\sum_{k=1}^p \sum_{l \neq k} C_{i(k+m)} C_{i(l+m)} y_{\text{Err},k} y_{\text{Err},l}}_{\text{Cross-term}} \quad (9)$$

In ACnets, we strictly enforce physical constraints, setting the left-hand side of Eq. (9) to 0, within numerical errors. As the squared error is positive definite, the cross term is always negative in ACnets as both terms sum up to 0. It is difficult to predict the cross term before optimization, hence Eq. (9) does not provide *a priori* predictions of performance, even for ACnets. Instead, it links how much the NN violates constraints to how well it predicts outputs that appear in the constraints equations: the more negative the cross term, the larger the squared error for a given violation of physical constraints.

Application.—Convective parametrization for climate modeling: The representation of subgrid-scale processes in coarse-scale, numerical models of the atmosphere, referred to as subgrid *parametrization*, is a large source of error and uncertainty in numerical weather and climate

prediction [29,30]. Machine-learning algorithms trained on fine-scale, process-resolving models can improve subgrid parametrizations by faithfully emulating the effect of fine-scale processes on coarse-scale dynamics [21,31–33] ([see Sec. 2 of Ref. [34] for a detailed review]). The problem is that none of these parametrizations exactly follow conservation laws (e.g., conservation of mass, energy). This is critical for long-term climate projections, as the spurious energy production may both exceed the projected radiative forcing from greenhouse gases and result in large thermodynamic drifts or biases over a long time period. Motivated by this shortcoming, we build a NN parametrization of convection and clouds that we *constrain* to conserve 4 quantities: column-integrated energy, mass, long-wave radiation, and short-wave radiation.

Model and data: We use the Super-Parameterized Community Atmosphere Model 3.0 [35] to simulate the climate for two years in aquaplanet configuration [36], where the surface temperatures are fixed with a realistic equator-to-pole gradient [37]. Following the sensitivity tests of Ref. [32], we use 42M samples from the simulation's first year to train the NN (training set) and 42M samples from the simulation's second year to validate the NN (validation set). Since we use the validation set to adjust the NN's hyperparameters and avoid overfitting, we additionally introduce a test set using 42M different samples from the simulation's second year to provide an unbiased estimator of the NN performances. Note that each sample represents a single atmospheric column at a given time, longitude, and latitude.

Formulating the conservation laws in a neural network: The parametrization's goal is to predict the rate at which subgrid convection vertically redistributes heat and water based on the current large-scale thermodynamic state. We group all variables describing the local climate in an input vector $\mathbf{x}$ of size 304 (5 vertical profiles with 30 levels each, prescribed large-scale conditions $\mathbf{LS}$ for all profiles of size 150, and 4 scalars):

$$\mathbf{x} = [(\mathbf{q}_v, \mathbf{q}_l, \mathbf{q}_i, T, v, \mathbf{LS}, p_s, S_0) \quad \text{SHF} \quad \text{LHF}]^T, \quad (10)$$

where all variables are defined in Supplemental Material [20], Sec. A. We then concatenate the time tendencies from convection and the additional variables involved in the conservation laws to form an output vector $\mathbf{y}$ of size 216 (7 vertical profiles with 30 levels, followed by 6 scalars):

$$\mathbf{y} = [\dot{\mathbf{q}}_v \quad \dot{\mathbf{q}}_l \quad \dot{\mathbf{q}}_i \quad \dot{T} \quad \dot{T}_{KE} \quad \text{lw} \quad \text{sw} \quad \text{LW}_t \quad \text{LW}_s \quad \text{SW}_t \quad \text{SW}_s \quad P \quad P_i]^T. \quad (11)$$

We normalize all variables to the same units before nondimensionalizing them using the constant 1 W m^{-2} (Supplemental Material [20] A.5). Finally, we derive the dimensionless conservation laws (Supplemental Material [20] A.1–A.4) and write them as a sparse matrix of size $4 \times (304 + 218)$:

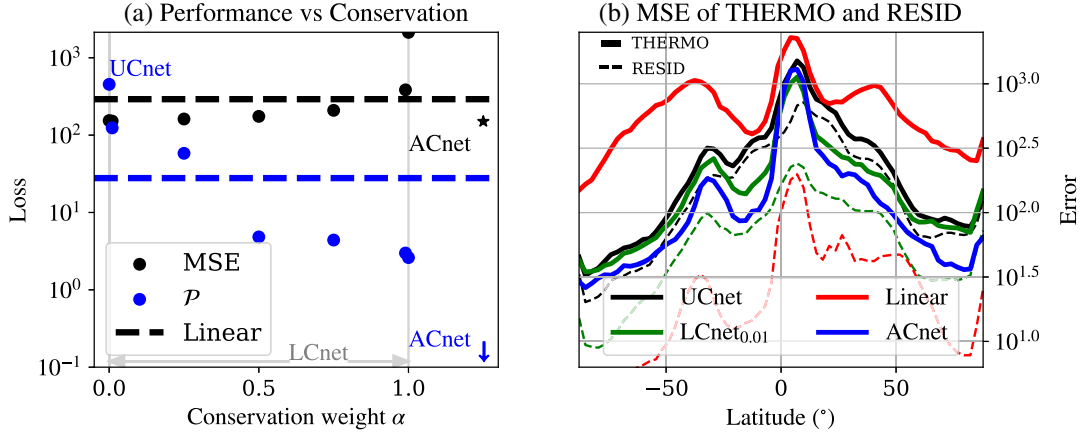


FIG. 3. (a) MSE and $\mathcal{P}$ averaged over all samples of the test dataset for UCNet, LCnets of varying α , and ACNet. The dashed lines indicate MSE and $\mathcal{P}$ for our multilinear regression baseline. (b) Mean-squared error in the thermodynamic term (THERMO) and the enthalpy residual (RESID) versus latitude for our lowest-MSE NN in each category.

$$C = \begin{bmatrix} 0 & 1 & \ell_s & -\ell_s \delta p & -\ell_f \delta p & 0 & -\delta p & \delta p & 0 & 0 & -1 & 1 & 1 & -1 & -\ell_f & \ell_f \\ 0 & 0 & 1 & -\delta p & -\delta p & -\delta p & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & \delta p & 0 & 1 & -1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & \delta p & 0 & 0 & -1 & 1 & 0 & 0 \end{bmatrix}, \quad (12)$$

that acts on $\mathbf{x}$ and $\mathbf{y}$ to yield Eq. (1).

Each row of the constraints matrix C describes a different conservation law: The first row is column-integrated enthalpy conservation (here equivalent to energy conservation), the second row is column-integrated water conservation (here equivalent to mass conservation), the third row is column-integrated long-wave radiation conservation and the last row is column-integrated short-wave radiation conservation.

Implementation: We implement the three NN types and a multilinear regression baseline using the Tensorflow library [38] version 1.13 with Keras [39] version 2.2.4: (i) *LCnets* for which we vary the weight α given to conservation laws from 0 to 1 [Eq. (6)], (ii) our reference *ACnet*, and (iii) *UCnet*, i.e., an unconstrained LCnet of weight $\alpha = 0$. In our reference ACnet, we write the constraints layers in Tensorflow to solve the system of equations (C) from bottom to top, and calculate surface tendencies as residuals of the conservation equations (Supplemental Material [20] B.1); switching the residual outputs to different vertical levels does not significantly change the validation loss nor the constraints penalty (Supplemental Material [20] B.3). After testing multiple architectures and activation functions (Supplemental Material [20] C.2), we chose 5 hidden layers of 512 nodes with leaky rectified linear-unit activations as our standard multilayer perceptron architecture, resulting in ~ 1.3 M trainable parameters. We optimized the NN's weights and biases with the RMSprop optimizer [40] for LCnets

(because it was more stable than the Adam optimizer [41]), used Sherpa for hyperparameter optimizations [42], and saved the NN's state of minimal validation loss over 20 epochs.

Results: In Fig. 3(a), we compare mean performance (measured by MSE) and by how much physical constraints are violated (measured by $\mathcal{P}$) for the three NN types. As expected, we note a monotonic trade-off between performance and constraints as we increase α from 0 to 1 in the loss function. This trade-off is well measured by MSE and $\mathcal{P}$ across the training, validation, and test sets (Supplemental Material [20] Table V). Interestingly, the physical constraints are easier to satisfy than reducing MSE in our case, likely because it is difficult to deterministically predict precipitation, which is strongly non-Gaussian, inherently stochastic, and whose error contributes to a large portion of MSE. Despite this, UCnet may violate physical constraints more than our multilinear regression baseline.

Our first key result is that *ACnet* performs nearly as well as our lowest-MSE UCnet on average (to within 3%) while satisfying constraints to $\sim (10^{-9}\%)$ (Supplemental Material [20] C.1). This result holds across the training, validation and test sets (Supplemental Material [20] Table IV). In our case, ACnets perform slightly less well than UCnet because they are harder to optimize and the residual outputs exhibit systematically larger errors (Supplemental Material [20] B.2). This systematic, unphysical bias can be remedied by multiplying the weights of these residual outputs in the

loss function (Supplemental Material [20] B.3) by a factor $\beta > 1$ (Supplemental Material [20], Eq. 12 and Fig. 2). β can be objectively chosen alongside the residual outputs via formal hyperparameter optimization (Supplemental Material [20] C.2).

In Fig. 3(b), we compare how much the NNs violate column energy conservation (RESID) to the prediction of a variable that appears in that constraint: the total thermodynamic tendency in the enthalpy conservation equation (THERMO):

$$\overbrace{\left(\mathbf{C} \begin{bmatrix} x \\ y_{\text{NN}} \end{bmatrix} \right)}^{\text{RESID}} = \overbrace{\delta \mathbf{p} \cdot (\dot{\mathbf{T}}_{\text{KE}} - \dot{\mathbf{T}} - \ell_s \dot{\mathbf{q}}_v - \ell_f \dot{\mathbf{q}}_l)}^{\text{THERMO}} + \dots, \quad (13)$$

where the ellipsis includes the surface fluxes, radiation, and precipitation terms. ACnet predicts THERMO more accurately than all NNs (full blue line) by an amount closely related to how much each NN violates enthalpy conservation (dashed lines), followed by LCnet (full green line). This yields our second key result: Enforcing constraints, whether in the architecture or the loss function, can systematically reduce the error of variables that appear in the constraints. This result holds true across the training, validation, and test sets (Supplemental Material [20] Fig. 4). However, possibly since our case has many degrees of freedom, it does not hold true for individual components of THERMO as their cross term in Eq. (9) is more negative for ACnet, nor does it hold for variables that are hard to predict deterministically (e.g., precipitation). Additionally, obeying conservation laws does not guarantee the ability to generalize well far outside of the training set, e.g., in the Tropics of a warmer climate (see Fig. 3 of Ref. [43]). These results nuance the finding that physically constraining NNs systematically improves their generalization ability, which has been documented for machine learning emulation of low-dimensional idealized flows [5,12], and motivate physically constraining machine-learning algorithms capable of stochastic predictions [44] that are consistent across climates [43].

Finally, although the mapping presented in Sec. III has linear constraints, ACnets can also be applied to nonlinearly constrained mappings by using the framework presented in Fig. 1. We give a concrete example in Supplemental Material [20], Sec. D, where we introduce the concept of “conversion layers” that transform nonlinearly constrained mappings into linearly constrained mappings within NNs and without overly degrading performance (Supplemental Material [20] Table IX). Additionally, ACnets can be extended to incorporate inequality constraints on their direct outputs (by using positive-definite activation functions, discussed in Supplemental Material [20] E), making ACnets applicable to a broad range of constrained optimization problems.

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*tom.beucler@gmail.com

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- [20] See Supplemental Material at <http://link.aps.org/supplemental/10.1103/PhysRevLett.126.098302> for more details. Supplemental Material has 5 Sections that can be read independently: (A) introduces the variables and constraints specific to our application, (B) details the implementation of our architecture-constrained network, (C) compares different network types and architectures, (D) presents an example where nonlinear constraints are enforced in neural networks emulating physical systems, and (E) discusses how to enforce inequality constraints, which contains Refs. [13,21–28].
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