Training a Foundation Model for Materials on a Budget

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Abstract

Foundation models for materials modeling are advancing quickly, but their training remains expensive, often placing state-of-the-art methods out of reach for many research groups. We introduce Nequix, a compact E(3)-equivariant potential that pairs a simplified NequIP design with modern training practices, including equivariant root-mean-square layer normalization and the Muon optimizer, to retain accuracy while substantially reducing compute requirements. Built in JAX, Nequix has 700K parameters and was trained in 500 A100-GPU hours. On the Matbench-Discovery and MDR Phonon benchmarks, Nequix ranks third overall while requiring less than one quarter of the training cost of most other methods, and it delivers an order-of-magnitude faster inference speed than the current top-ranked model. We release model weights and a JAX codebase at `https://github.com/atomicarchitects/nequix`.

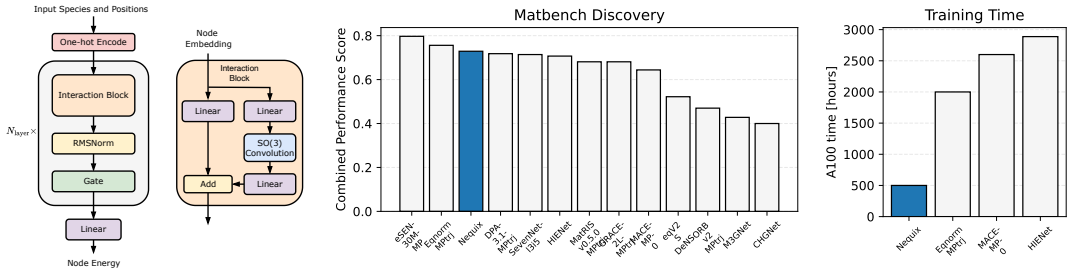
1 Introduction

Machine learned inter-atomic potentials (MLIPs) are rapidly improving in capability and scope, with foundation models trained on broad datasets of atomistic materials offering the promise of augmenting or replacing expensive *ab initio* density functional theory (DFT) calculations [Batatia et al., 2023]. While performance on community benchmarks such as Matbench-Discovery [Riebesell et al., 2025] is rising, the computational costs of both data generation and curation as well as the training of MLIP models on these datasets remain prohibitively expensive for many labs.

We pursue an orthogonal goal to scaling: a lower computational cost recipe that preserves strong downstream accuracy. Concretely, we revisit a simplified E(3)-equivariant architecture based on NequIP [Batzner et al., 2022] with modern training practices: root-mean-square layer normalization for stability, dynamic batching to maximize GPU utilization, and optimizer choices inspired by "speedrunning" deep learning workflows [Jordan et al., 2024a]. The resulting model, Nequix, is 700K parameters and can be trained in 500 gpu hours, while remaining competitive with larger and more costly to train models on Matbench-Discovery and other phonon prediction tasks.

Our contributions are threefold: 1) a simplified NequIP architecture featuring an equivariant layer-norm, with an efficient JAX implementation. ; 2) a budget-conscious training pipeline leveraging dynamic batching and the Muon optimizer [Jordan et al., 2024b] which achieves fast convergence; and 3) evaluations on the Matbench-Discovery and MDR phonon [Loew et al., 2025] benchmarks. Compared to prior MPtrj-trained models [Chen and Ong, 2022, Deng et al., 2023, Batatia et al., 2023, Bochkarev et al., 2024, Neumann et al., 2024, Barroso-Luque et al., 2024, Fu et al., 2025, Zhang et al., 2025, Yan et al., 2025] , we rank in *third* on both benchmarks with one fourth the training cost of any other published training cost and a $10\times$ faster inference speed over current top ranking model.

Preprint.



(a) (b) (c)

Figure 1: (a) Nequix architecture, a simplified version of NequIP [Batzner et al., 2022], with a species-independent residual connection and layer normalization. (b) Combined performance scores of compliant models on the Matbench Discovery, unique prototypes. (c) Available published training times of current compliant models.

2 Methods

Architecture Nequix follows a simplified version of the NequIP [Batzner et al., 2022] architecture, as shown in figure 1a. We adopt two modifications suggested by Park et al. [2024]; the species-specific self-connection layer within the interaction block is replaced with a single linear layer, and unused non-scalar representations are discarded from the final layer. Lastly, we add an equivariant root-mean-square layer normalization (RMSNorm) Liao et al. [2023], which we find improves performance in our optimization setting. We document the full architecture hyper-parameters and the rationale behind each decision in Table A1.

Implementation Nequix is implemented in JAX [Bradbury et al., 2018] with Equinox [Kidger and Garcia, 2021], taking advantage of just-in-time compilation and efficient automatic differentiation. Following standard energy-conserving MLIP practice [Fu et al., 2025], forces are obtained as the negative energy gradient with respect to atomic positions, $-\nabla_{\mathbf{r}} E$, and stresses as the energy derivative with respect to strain, normalized by volume, $\sigma = V^{-1}\,\partial E/\partial \varepsilon$, where E is the predicted total energy of the system, $\mathbf{r}$ is an atom position, ε is the strain tensor, V is the simulation cell volume, and σ denotes the stress tensor.

Dynamic batching Materials radius graphs vary widely in their numbers of nodes and edges due to differences in unit-cell size and atom count. With fixed-size batching, the largest graphs dictate memory usage, leaving GPUs underutilized for most batches. To keep batch workloads more uniform while respecting memory limits, we use dynamic batching [Speckhard et al., 2025]: each batch is filled up to caps on the total nodes and edges. We set these caps to $1.1\times$ (ideal batch size) $\times$ (dataset-average nodes-per-graph and edges-per-graph), respectively. Batches are then padded to these caps to fulfil the static shape requirements of JAX [Godwin* et al., 2020].

Optimization and normalization We compare the widely used Adam optimizer Kingma and Ba [2014] with the recently proposed Muon optimizer Jordan et al. [2024b]. Using a smaller version of Nequix on MPtrj [Jain et al., 2013, Deng et al., 2023] with hidden irreps of `128x0e + 64x1o`, we sweep learning rates of $\{0.03, 0.01, 0.003, 0.001\}$ for Adam and Muon, each with and without RMSNorm. The lowest validation error runs for each optimizer are shown in Fig. 2. We find that the Muon configuration achieves comparable energy/force errors to Adam in 60-70% of the epochs, and results in a 7% reduction in energy MAE. We also find a significant reduction in the variance of stress error, which we notice in runs that use RMSNorm. Notably, the presence of the RMSNorm layer generally resulted in lower validation error for Muon-based training configurations, and higher for those using Adam.

Training procedure The final Nequix model is trained for 100 epochs on MPtrj [Jain et al., 2013, Deng et al., 2023], which we leave 5% out of training for validation. More details on the training settings are provided in Sec. A.1. The model was trained on 4 NVIDIA A100 GPUs in 125 hours, for a total cost of 500 GPU hours.

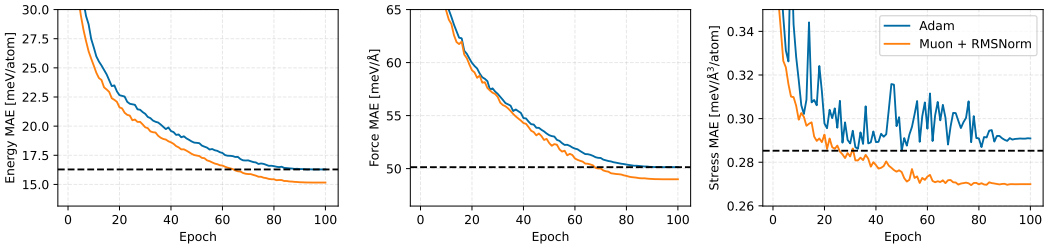


Figure 2: Validation metrics during training a smaller version of Nequix configuration with Adam and Muon, trying learning rates in $\{0.03, 0.01, 0.003, 0.001\}$ and with/without RMSNorm. This model configuration uses the same hyperparameters as the final model, except with hidden irreps of `128x0e + 64x1o`. The dotted horizontal line shows the best validation performance reached during the Adam training.

3 Experiments

3.1 Matbench-Discovery benchmark

Matbench-Discovery [Riebesell et al., 2025] provides a standard framework for evaluating interatomic potentials in a high-throughput materials screening task consisting of geometry optimization and energy prediction on a set of 257,487 generated structures, and thermal conductivity prediction on a set of 103 structures. Ground truth is calculated with DFT/PBE level of theory, the same as MPtrj. The primary metrics include: 1) the F1 for stable/unstable classification after relaxation; 2) root mean squared displacement (RMSD) between predicted and reference structures after relaxation; and 3) symmetric relative mean error in predicted phonon mode contributions to thermal conductivity κ (κ_{SRME}). A normalized and weighted combination of these metrics are then used to compute a combined performance score (CPS), which is used for ranking.

Following Riebesell et al. [2025] we integrate our interatomic potential as Atomic Simulation Environment (ASE) calculator, which is then used to perform structure relaxation and phonon calculations with the default settings of the benchmark. For comparison, we consider only models in the compliant subset of the benchmark. This consists only of models that are trained on MPtrj or subsets, which limits data leakage and offers a more fair comparison among methods. Table 1 contains the performance of Nequix along with all current compliant models at the time of writing. We also include the reported training cost for the models when available, visualized in Fig. 1. We find that Nequix ranks third by CPS, outperforming most models at a fraction of the training cost. It is noteworthy that this high ranking is due to high performance in the thermal conductivity task, however the F1 score is still comparable to many of the other methods.

Table 1: Matbench-Discovery compliant leaderboard, sorted by combined performance score (CPS). Metrics are shown for the unique prototypes subset. Train cost is measured in A100 hours.

Model	Params	Train cost	RMSD↓	κ_{SRME} ↓	F1↑	CPS↑
eSEN-30M-MP	30.1M		0.075	0.340	0.831	0.797
Eqnorm MPtrj	1.31M	2000	0.084	0.408	0.786	0.756
Nequix	707K	500	0.085	0.446	0.750	0.729
DPA-3.1-MPtrj	4.81M		0.080	0.650	0.803	0.718
SevenNet-l3i5	1.17M		0.085	0.550	0.760	0.714
HIENet	7.51M	2888	0.080	0.642	0.777	0.707
MatRIS v0.5.0 MPtrj	5.83M		0.077	0.861	0.809	0.681
GRACE-2L-MPtrj	15.3M		0.090	0.525	0.691	0.681
MACE-MP-0	4.69M	2600	0.092	0.647	0.669	0.644
eqV2 S DeNS	31.2M		0.076	1.676	0.815	0.522
ORB v2 MPtrj	25.2M		0.101	1.725	0.765	0.470
M3GNet	228k		0.112	1.412	0.569	0.428
CHGNet	413k		0.095	1.717	0.613	0.400

3.2 MDR phonon benchmark

We also evaluate performance on the MDR phonon benchmark [Loew et al., 2025], a set of 10,000 phonon calculations also done with DFT/PBE level of theory. We follow the identical procedure to Loew et al. [2025], first performing a geometry relaxation, then phonon calculations using displacements of supercell. Mean absolute error (MAE) of properties derived from the phonon calculation: maximum phonon frequency $\omega_{\max}$, vibrational energy S, Helmholtz free energy F, and heat capacity at constant volume C_V. Table 2, demonstrates the performance of Nequix compared to other MPtrj-trained models. Similarly to Matbench-Discovery, we achieve performance within the top three of models, with a fraction of the parameter count of other methods.

Table 2: Model performance of MPtrj-trained models on the MDR phonon benchmark, sourced from Loew et al. [2025] and Fu et al. [2025].

Model	MAE($\omega_{\max}$)	MAE(S)	MAE(F)	MAE(C_V)
eSEN–30M	21	13	5	4
SevenNet–l3i5	26	28	10	5
Nequix	26	33	12	6
SevenNet–0	40	48	19	9
GRACE–2L (r6)	40	25	9	5
MACE	61	60	24	13
CHGNet	89	114	45	21
M3GNet	98	150	56	22

3.3 Inference speed

Lastly, we compare inference speed against existing materials potentials using a script from the MACE [Batatia et al., 2022] repository[1], modified to work with Nequix and eSEN. We run each model in the configuration in which it is used within its ASE calculator, without compilation for the PyTorch-based MACE-MP-0 and eSEN models, and just-in-time compilation for the JAX-based Nequix. Figure 3 compares performance of each model in terms of steps per day vs. number of atoms. In this study, Nequix is about 10× faster than eSEN, and roughly 2× slower than MACE-MP-0, offering a new option in the accuracy vs. speed Pareto frontier at a fraction of the training cost.

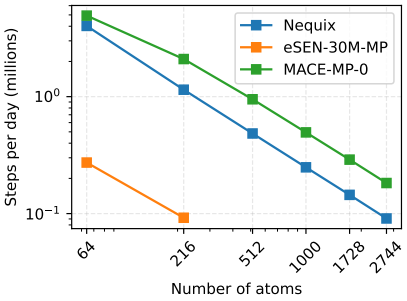


Figure 3: Inference speed of various models in steps per day.

4 Conclusion

We presented Nequix, an E(3)-equivariant interatomic potential that pairs a simplified NequIP architecture with modern training practices. Our results show that Nequix achieves competitive accuracy on Matbench-Discovery and the MDR phonon benchmark at less than one quarter of the reported training cost of many contemporaries. This resource-efficient recipe provides a practical alternative to large-scale foundation models and helps broaden access to high-quality atomistic modeling in settings with more limited compute. We release trained weights and a JAX codebase to streamline reuse and extension.

Looking ahead, we see several promising directions: scaling training duration and data while maintaining budget discipline, exploring pretraining and fine-tuning regimes across broader datasets, and pushing cost even lower through model distillation, pruning, quantization, kernel-implementations, or more data-efficient training. We hope Nequix serves as a strong, efficient baseline for future work on accessible materials foundation models.

[1] `https://github.com/ACEsuit/mace/blob/main/tests/test_benchmark.py`

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A Appendix

A.1 Training and model configuration

Table A1 shows the hyper-parameters used to train Nequix. The model is trained for 100 epochs, using an MAE loss function on energy and stress, and l_2 loss on forces. We use a linear warmup with cosine decay learning rate schedule.

Table A1: Hyper-parameters used and rationale behind selection

Hyper-parameter	Value	Notes/Rational
Radial cutoff	6 Å	Most models, use 5 or 6 Å; 6 performed slightly better in preliminary validation performance.
Hidden irreps	`128x0e + 64x1o + 32x2e + 32x3o`	From SevenNet-l3i5.
L_{max}	3	Consistent with hidden irreps.
N_{layers}	4	Balance of performance and efficiency.
Radial basis size	8	From NequIP and analysis from Sec. 5.2 of Fu et al. [2025]
Radial MLP size	64	From NequIP.
Radial MLP layers	2	From NequIP.
Polynomial cutoff p	6.0	From NequIP.
Radial basis function	Bessel	From NequIP. Also tried Gaussian, which had minimal difference on validation performance.
Learning rate	0.01	Selected from $\{0.03, 0.01, 0.003, 0.001\}$ based on validation performance early in training.
Warmup epochs	0.1	From eSEN.
Warmup factor	0.2	From eSEN.
Optimizer	Muon	See Sec. 2.
Weight decay	0.001	From eSEN. Also tried 0.0, which lead to worse validation performance.
Energy weight	20	From eSEN.
Force weight	20	From eSEN.
Stress weight	5	From eSEN.
Batch size	256 (dynamic)	See Sec. 2
Number of epochs	100	Standard training duration.