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Cover
See Matthew D. Witman and Peter Schindler, pp. 625–635. Image reproduced by permission of David Witman and Peter Schindler from *Digital Discovery*, 2025, **4**, 625.

Inside cover
See Linjiang Chen, Fei Zhang, Weiwei Shang, Jun Jiang *et al.*, pp. 636–652. Image reproduced by permission of Jun Jiang from *Digital Discovery*, 2025, **4**, 636.

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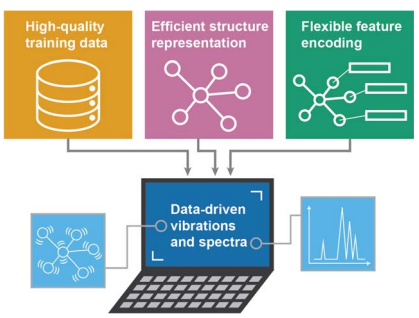
AI-powered exploration of molecular vibrations, phonons, and spectroscopy

Bowen Han, Ryotaro Okabe, Abhijatmedhi Chotrattanapituk, Mouyang Cheng, Mingda Li* and Yongqiang Cheng*



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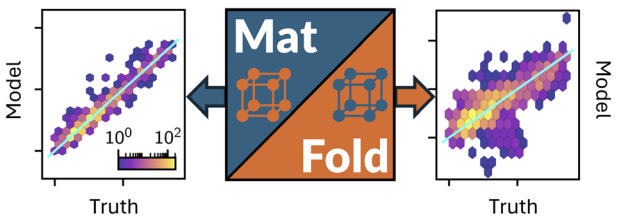
MatFold: systematic insights into materials discovery models' performance through standardized cross-validation protocols

Matthew D. Witman* and Peter Schindler*



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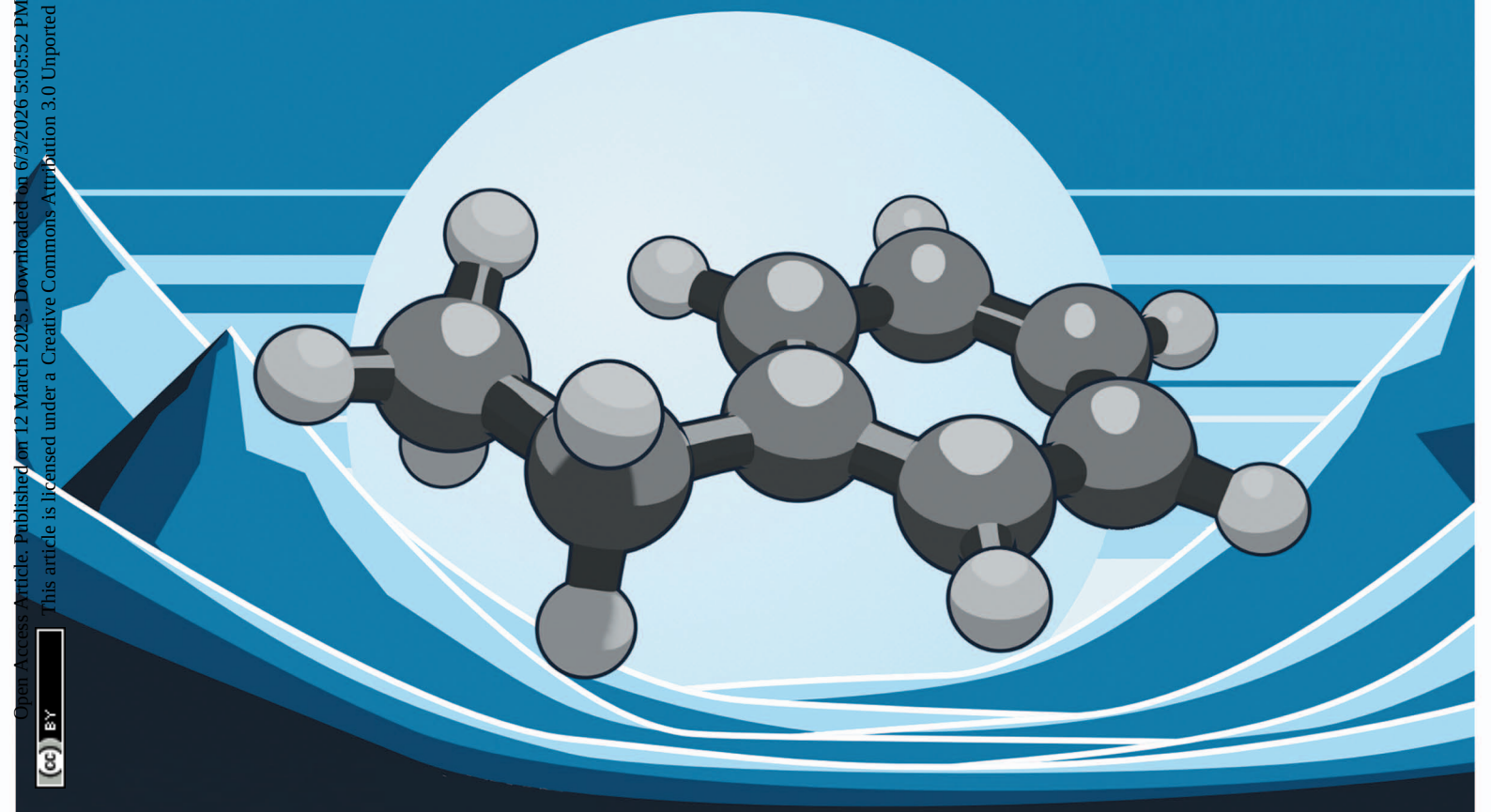
Cheminformatics, Automation and Machine Learning in Chemistry (CAMLC micro-credential)



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A multi-robot–multi-task scheduling system for autonomous chemistry laboratories

Junyi Zhou, Man Luo, Linjiang Chen,* Qing Zhu, Shan Jiang, Fei Zhang,* Weiwei Shang* and Jun Jiang*

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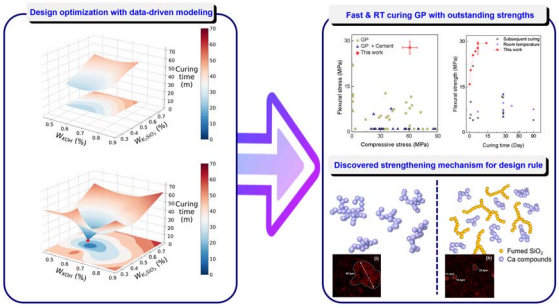
Accentuating the ambient curing behavior of geopolymers: metamodel-guided optimization for fast-curing geopolymers with high flexural strength

Kyungwon Kim, Hyejeong Song, Sanghun Lee, Hyeongkyu Cho, Hyung Mi Lim and Hyunseok Ko*



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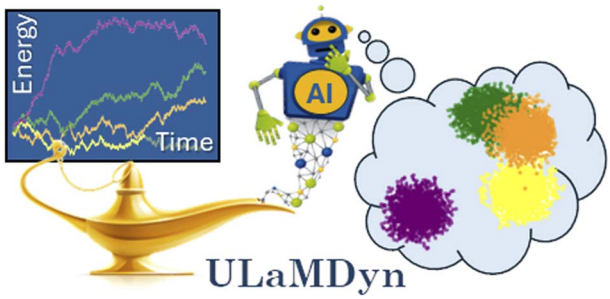
ULaMDyn: enhancing excited-state dynamics analysis through streamlined unsupervised learning

Max Pinheiro Jr,* Matheus de Oliveira Bispo, Rafael S. Mattos, Mariana Telles do Casal, Bidhan Chandra Garain,* Josene M. Toldo, Saikat Mukherjee and Mario Barbatti*



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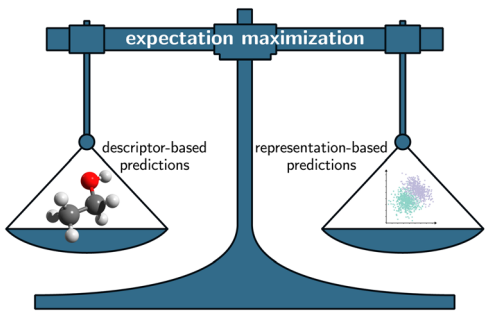
Balancing molecular information and empirical data in the prediction of physico-chemical properties

Johannes Zenn,* Dominik Gond, Fabian Jirasek and Robert Bamler



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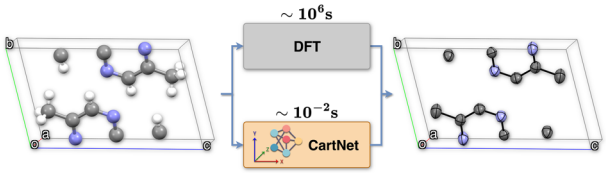

A Cartesian encoding graph neural network for crystal structure property prediction: application to thermal ellipsoid estimation

Àlex Solé, Albert Mosella-Montoro, Joan Cardona, Silvia Gómez-Coca,* Daniel Aravena,* Eliseo Ruiz* and Javier Ruiz-Hidalgo*

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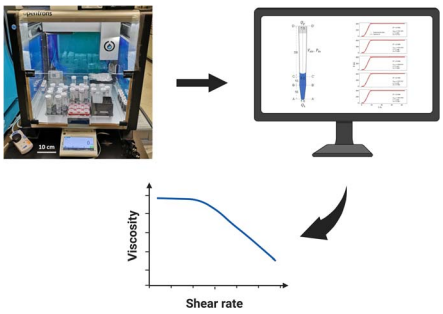

Opentrons for automated and high-throughput viscometry

Beatrice W. Soh,* Aniket Chitre, Shu Zheng Tan, Yuhan Wang, Yinqi Yi, Wendy Soh, Kedar Hippalgaonkar and D. Ian Wilson

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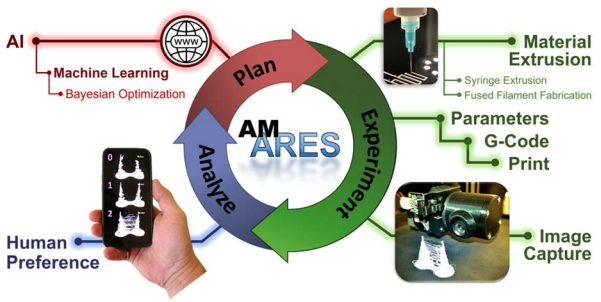

Preferential Bayesian optimization improves the efficiency of printing objects with subjective qualities

James R. Deneault, Woojae Kim, Jiseob Kim, Yuzhe Gu, Jorge Chang, Benji Maruyama, Jay I. Myung and Mark A. Pitt*

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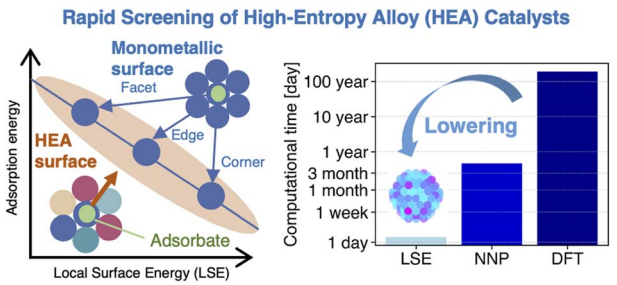

Lowering the exponential wall: accelerating high-entropy alloy catalysts screening using local surface energy descriptors from neural network potentials

Tomoya Shiota,* Kenji Ishihara and Wataru Mizukami*

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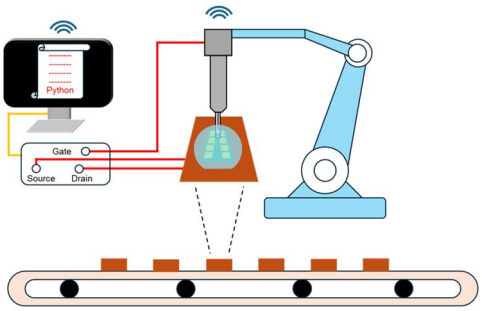
An automated electrolyte-gate field-effect transistor test system for rapid screening of multiple sensors

Zhengru Liu, Long Bian, Wenting Shao, Sean I. Hwang and Alexander Star*



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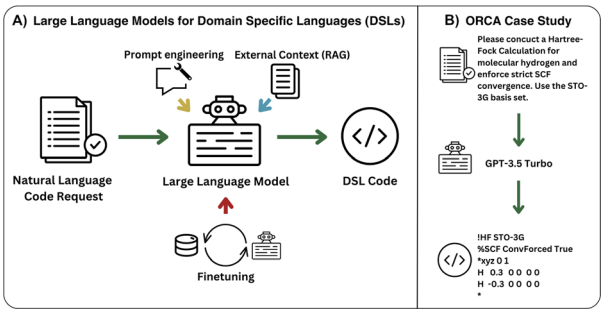
Developing large language models for quantum chemistry simulation input generation

Pieter Floris Jacobs and Robert Pollice*



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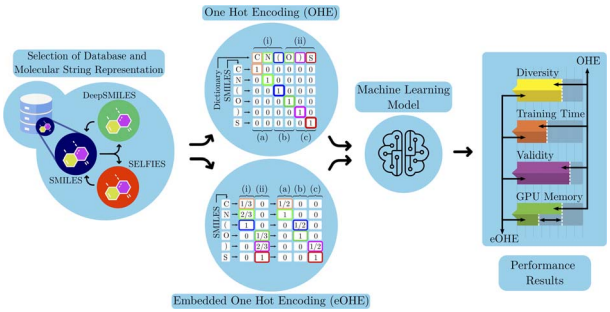
Embedded machine-readable molecular representation for resource-efficient deep learning applications

Emilio Nuñez-Andrade,* Isaac Vidal-Daza, James W. Ryan, Rafael Gómez-Bombarelli and Francisco J. Martin-Martinez*



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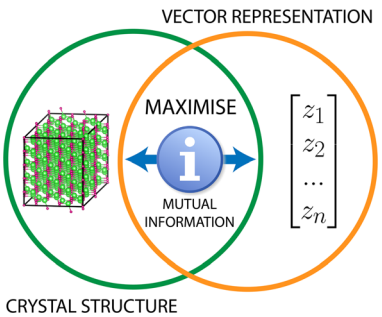
Establishing Deep InfoMax as an effective self-supervised learning methodology in materials informatics

Michael Moran, Michael W. Gaultois, Vladimir V. Gusev,* Dmytro Antypov and Matthew J. Rosseinsky*



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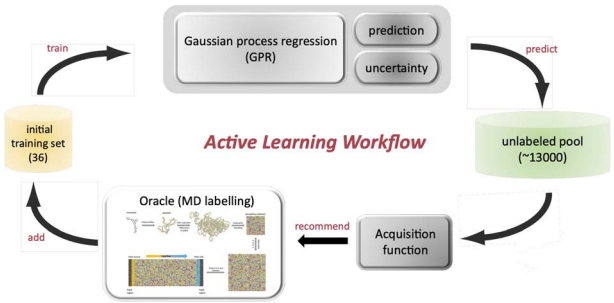

Active learning-guided exploration of thermally conductive polymers under strain

Renzheng Zhang, Jiaxin Xu, Hanfeng Zhang, Guoyue Xu and Tengfei Luo*

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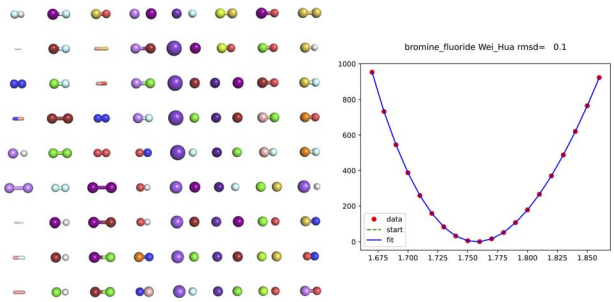

Quantitative evaluation of anharmonic bond potentials for molecular simulations

Paul J. van Maaren and David van der Spoel*

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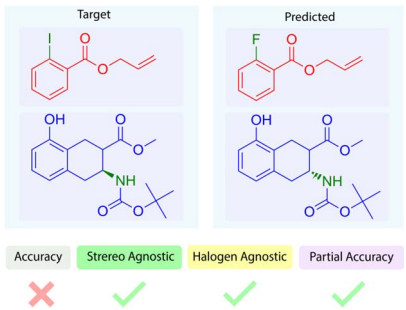

Dissecting errors in machine learning for retrosynthesis: a granular metric framework and a transformer-based model for more informative predictions

Arihanth Srikar Tadanki, H. Surya Prakash Rao and U. Deva Priyakumar*

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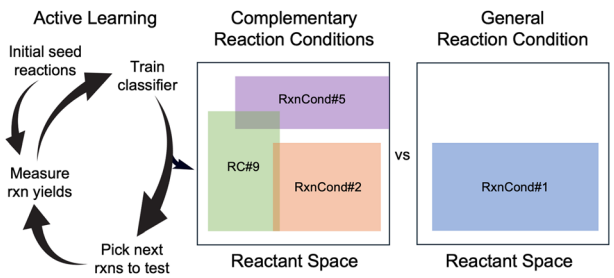

Active learning high coverage sets of complementary reaction conditions

Sofia L. Sivilotti, David M. Friday and Nicholas E. Jackson*

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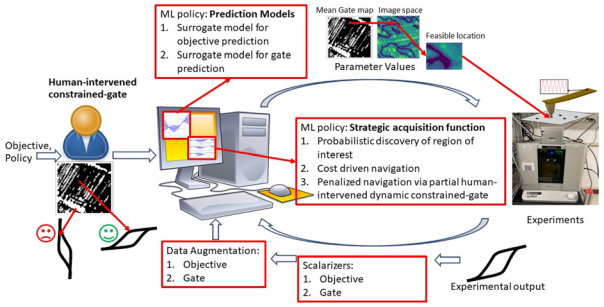
SANE: strategic autonomous non-smooth exploration for multiple optima discovery in multi-modal and non-differentiable black-box functions

Arpan Biswas, Rama Vasudevan, Rohit Pant, Ichiro Takeuchi, Hiroshi Funakubo and Yongtao Liu



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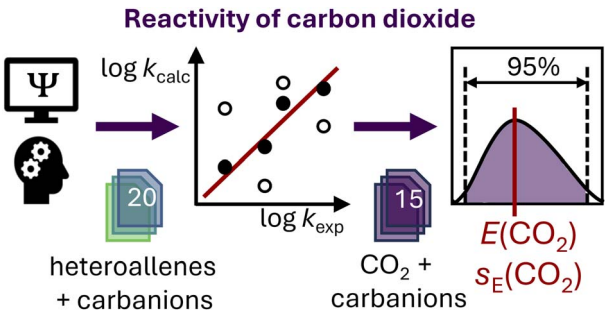
Unveiling CO_2 reactivity with data-driven methods

Maike Eckhoff, Kerstin L. Bublitz and Jonny Proppe*



CORRECTION

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Correction: Distortion/interaction analysis *via* machine learning

Samuel G. Espley, Samuel S. Allsop, David Buttar, Simone Tomasi and Matthew N. Grayson*

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