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PERSPECTIVE

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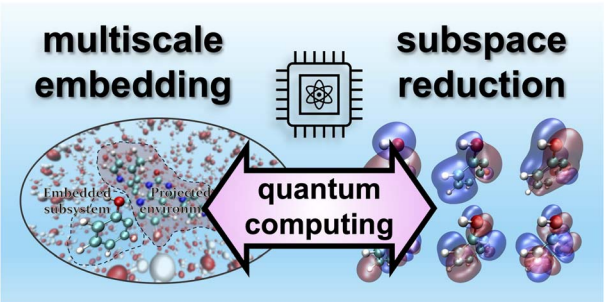
Extending quantum computing through subspace, embedding and classical molecular dynamics techniques

Thomas M. Bickley, Angus Mingare, Tim Weaving, Michael Williams de la Bastida, Shunzhou Wan, Martina Nibbi, Philipp Seitz, Alexis Ralli, Peter J. Love, Minh Chung, Mario Hernández Vera, Laura Schulz and Peter V. Coveney*

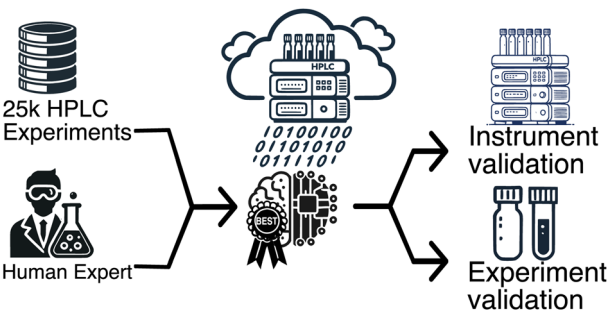


PAPERS

3445

Machine learning anomaly detection of automated HPLC experiments in the cloud laboratory

Filipp Gusev, Benjamin C. Kline, Ryan Quinn, Anqin Xu, Ben Smith, Brian Frezza and Olexandr Isayev*



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PAPERS

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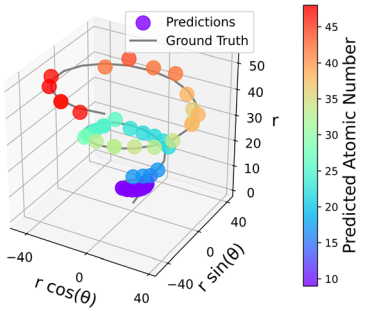
Do Llamas understand the periodic table?

Ge Lei* and Samuel J. Cooper



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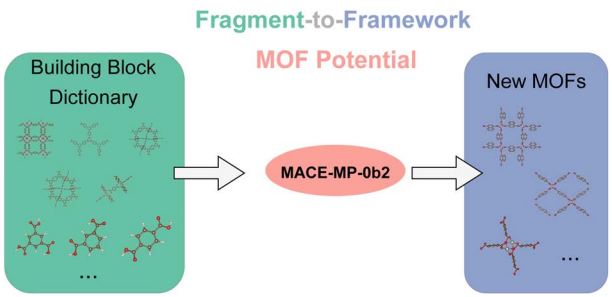
FFLAME: a fragment-to-framework learning approach for MOF potentials

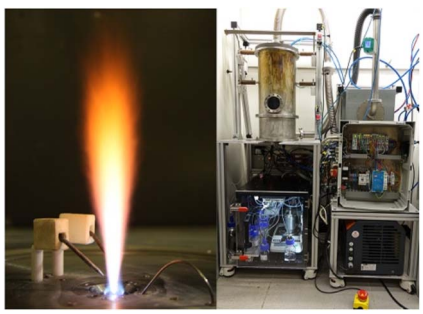
Xiaoqi Zhang, Yutao Li, Xin Jin and Berend Smit*



3478

An automated platform for "on-demand" high-speed catalyst synthesis by flame spray pyrolysis

Konstantin M. Engel, Patrik O. Willi, Robert N. Grass and Wendelin J. Stark*



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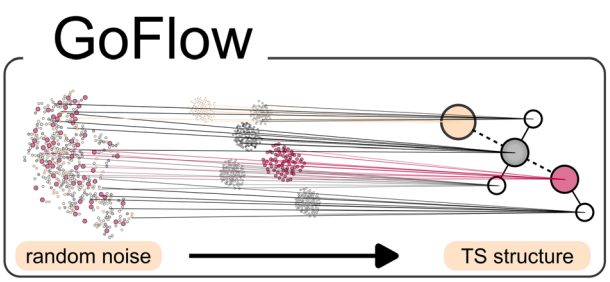
GoFlow: efficient transition state geometry prediction with flow matching and E(3)-equivariant neural networks

Leonard Galustian, Konstantin Mark, Johannes Karwounopoulos, Maximilian P.-P. Kovar and Esther Heid*



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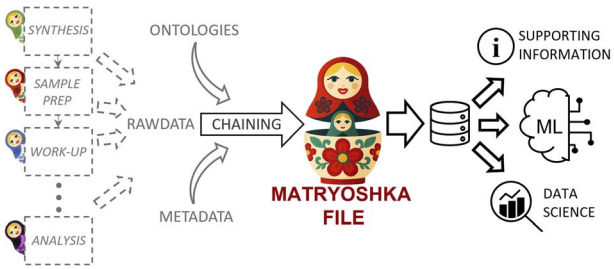

A FAIR research data infrastructure for high-throughput digital chemistry

Alice Gauthier, Laure Vancauwenberghe, Jean-Charles Cousty,* Cyril Matthey-Doret, Robin Franken, Sabine Maennel, Pascal Miéville and Oksana Riba Grognuz

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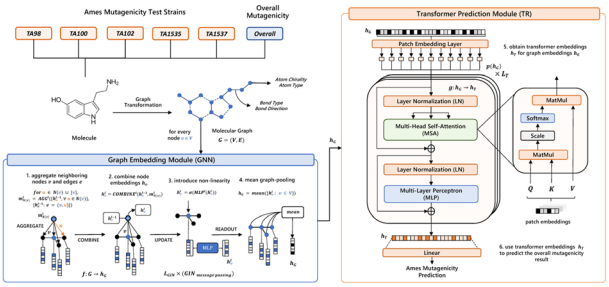

Advancing mutagenicity predictions in drug discovery with an explainable few-shot deep learning framework

Luis H. M. Torres,* Sofia M. da Silva, Joel P. Arrais, Catarina Pimentel and Bernardete Ribeiro

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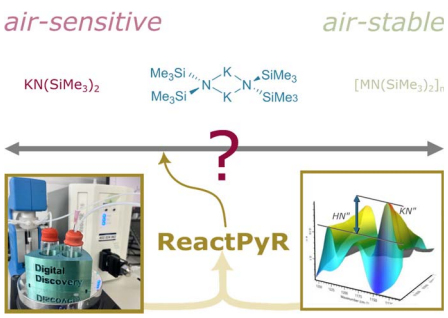

ReactPyR: a python workflow for ReactIR allows for quantification of the stability of sensitive compounds in air

Nicola L. Bell,* Emanuele Berardi, Marina Gladkikh, Richard Drummond Turnbull and Freya Turton

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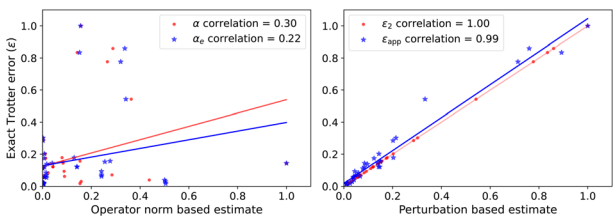

Estimating Trotter approximation errors to optimize Hamiltonian partitioning for lower eigenvalue errors

Shashank G. Mehendale, Luis A. Martínez-Martínez, Prathami Divakar Kamath and Artur F. Izmaylov*

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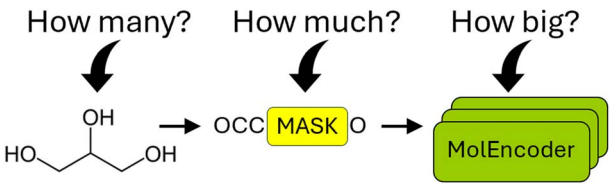
MolEncoder: towards optimal masked language modeling for molecules

Fabian P. Krüger,* Nicklas Österbacka, Mikhail Kabeshov, Ola Engkvist and Igor Tetko



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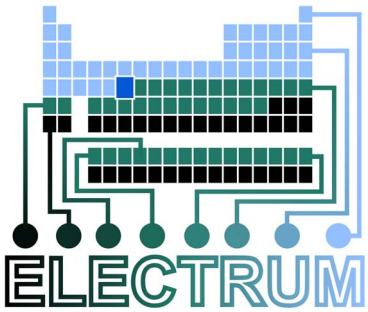
ELECTRUM: an electron configuration-based universal metal fingerprint for transition metal compounds

Markus Orsi* and Angelo Frei*



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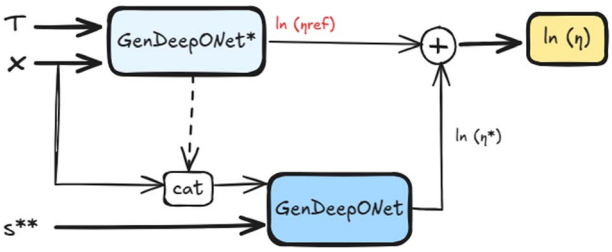
Generalized DeepONets for viscosity prediction using learned entropy scaling references

Maximiliam Fleck,* Marcelle B. M. Spera, Samir Darouich, Timo Klenk and Niels Hansen*



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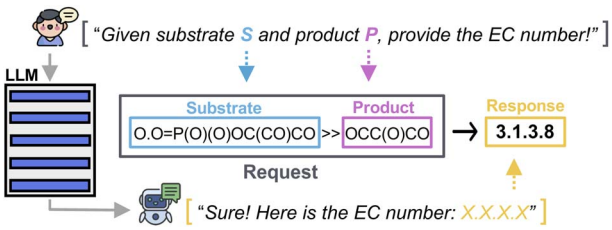
Leveraging large language models for enzymatic reaction prediction and characterization

Lorenzo Di Fruscia and Jana M. Weber*



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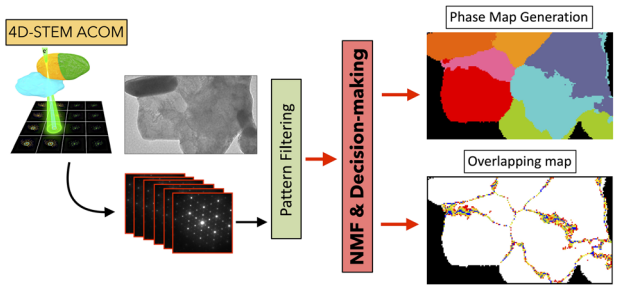

Unsupervised multi-clustering and decision-making strategies for 4D-STEM orientation mapping

Junhao Cao, Nicolas Folastre, Gozde Oney, Edgar Rauch, Stavros Nicolopoulos, Partha Pratim Das and Arnaud Demortière*

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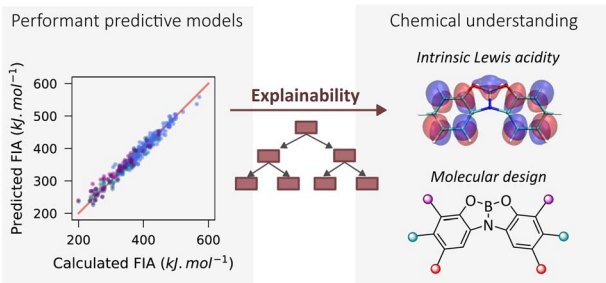

Constructing and explaining machine learning models for the exploration and design of boron-based Lewis acids

Juliette Fenogli,* Laurence Grimaud* and Rodolphe Vuilleumier*

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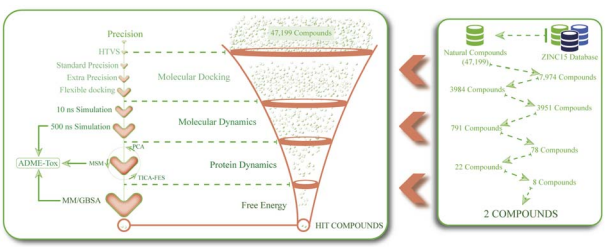

Database mining of ZINC15 natural compounds reveals potential thyroid receptor β agonists for NAFLD management: an *in silico* study

Ahmet Buğra Ortaakarsu,* Michel Hosny, Mansour Sobeh and Mohamed A. O. Abdelfattah

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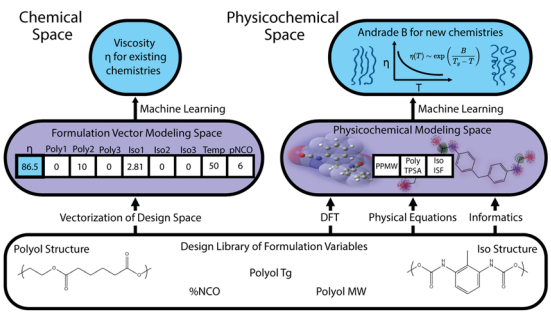

Machine learning of polyurethane prepolymer viscosity: a comparison of chemical and physicochemical approaches

Joseph A. Pugar, Calvin Gang, Isabelle Millan, Karl Haider and Newell R. Washburn*

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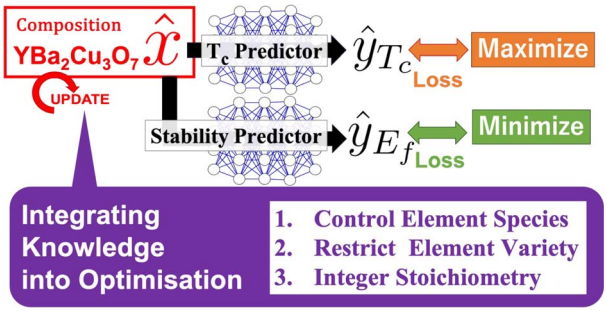
A straightforward gradient-based approach for designing superconductors with high critical temperature: exploiting domain knowledge *via* adaptive constraints

Akihiro Fujii,* Anh Khoa Augustin Lu, Koji Shimizu and Satoshi Watanabe



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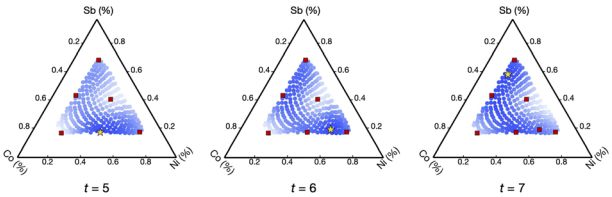
Active learning path-dependent properties using a cloud-based materials acceleration platform

Dan Guevarra, Michael J. Statt,* Kostiantyn Popovich, Brian A. Rohr,* John M. Gregoire, Kevin Tran, Santosh K. Suram, Joel A. Haber* and Willie Neiswanger*



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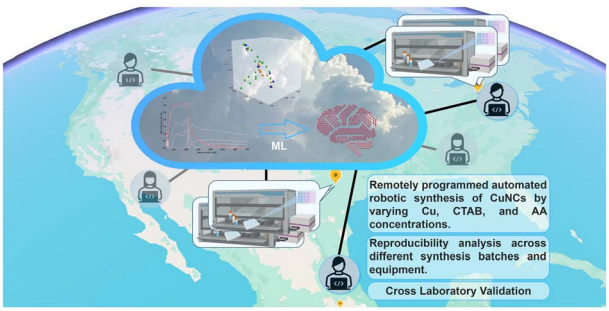
Cross-laboratory validation of machine learning models for copper nanocluster synthesis using cloud-based automated platforms

Ricardo Montoya-Gonzalez, Rosa de Guadalupe González-Huerta, Martha Leticia Hernández-Pichardo and Subha R. Das*



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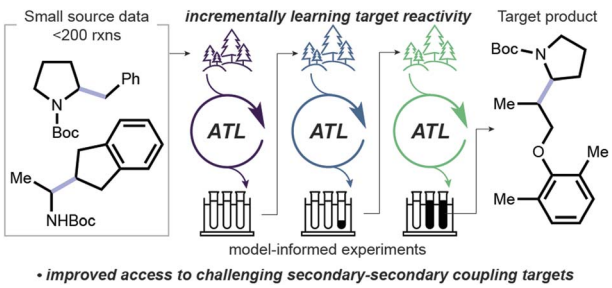
Prospective active transfer learning on the formal coupling of amines and carboxylic acids to form secondary alkyl bonds

Eunjae Shim, Ambuj Tewari, Paul M. Zimmerman* and Tim Cernak*



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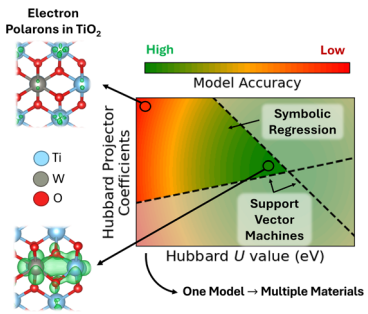

Machine learning generalised DFT+*U* projectors in a numerical atom-centred orbital framework

Amit Chaudhari, Kushagra Agrawal and Andrew J. Logsdail*

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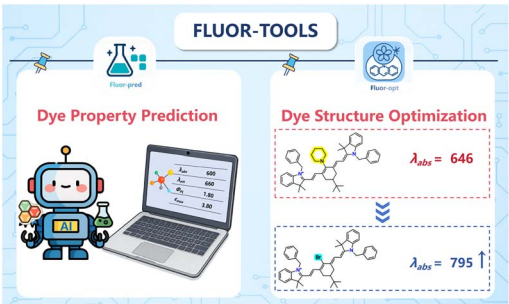

Fluor-tools: an integrated platform for dye property prediction and structure optimization

Wenxiang Song, Yuyang Zhang, Le Xiong, Xinmin Li, Jingwei Zhang, Guixia Liu, Weihua Li, Youjun Yang* and Yun Tang*

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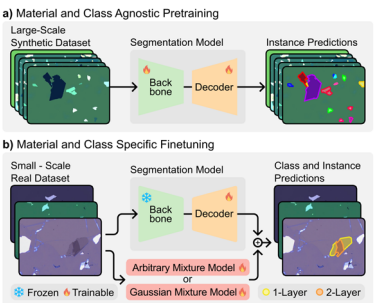

MaskTerial: a foundation model for automated 2D material flake detection

Jan-Lucas Uslu,* Alexey Nekrasov, Alexander Hermans, Bernd Beschoten, Bastian Leibe, Lutz Waldecker and Christoph Stampfer

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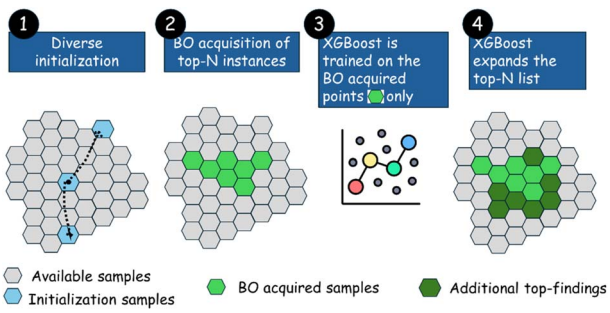

Navigating materials design spaces with efficient Bayesian optimization: a case study in functionalized nanoporous materials

Panagiotis Krokidas,* Vassilis Gkatsis, John Theocharis and George Giannakopoulos

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PAPERS

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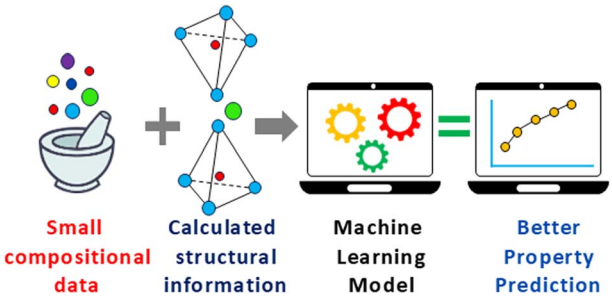
An improved machine learning strategy using structural features to predict the glass transition temperature of oxide glasses

Satwinder Singh Danewalia* and Kulvir Singh



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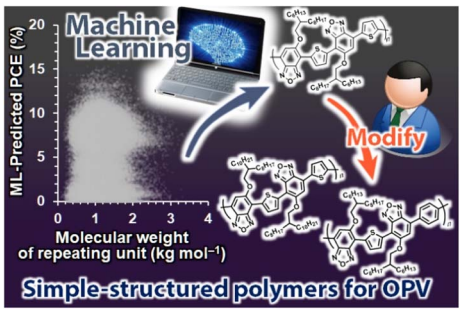
Design of simple-structured conjugated polymers for organic solar cells by machine learning-assisted structural modification and experimental validation

Shogo Tadokoro, Ryosuke Kamimura, Fumitaka Ishiwari and Akinori Saeki*



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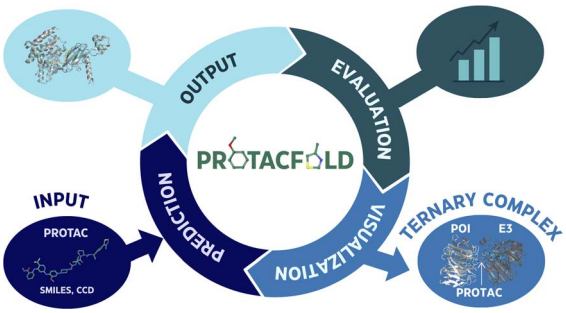
Predicting PROTAC-mediated ternary complexes with AlphaFold3 and Boltz-1

Nils Dunlop, Francisco Erazo, Farzaneh Jalalypour and Rocío Mercado*



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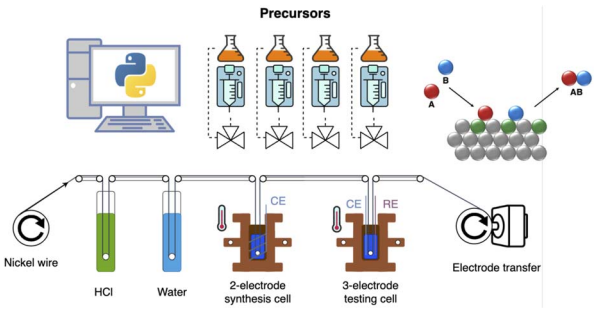
CatBot – a high-throughput catalyst synthesis and testing system with roll to roll transfer

Paolo Vincenzo Freiesleben de Blasio,* Rune Kruger, Nis Fisker-Bødker, Jin Hyun Chang and Christodoulos Chatzichristodoulou



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PAPERS

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Multi-modal contrastive learning for chemical structure elucidation with VibraCLIP

Pau Rocabert-Oriols, Camilla Lo Conte, Núria López and Javier Heras-Domingo*

CORRECTION

3828

Correction: Beyond training data: how elemental features enhance ML-based formation energy predictions

Hamed Mahdavi,* Vasant Honavar and Dane Morgan

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