

Digital Discovery

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See Shu Huang and Jacqueline M. Cole, pp. 1710–1720. Image reproduced by permission of Shu Huang and Nan Tian, who used imagery from rawpixel.com from Freepik, from *Digital Discovery*, 2023, 2, 1710.

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Phillip M. Maffettone,* Pascal Friederich,* Sterling G. Baird, Ben Blaiszik, Keith A. Brown, Stuart I. Campbell, Orion A. Cohen, Rebecca L. Davis, Ian T. Foster, Navid Haghmoradi, Mark Hereld, Howie Joress, Nicole Jung, Ha-Kyung Kwon, Gabriella Pizzuto, Jacob Rintamaki, Casper Steinmann, Luca Torresi and Shijing Sun

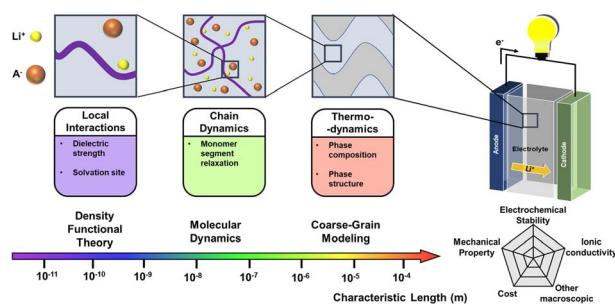


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Computational and data-driven modelling of solid polymer electrolytes

Kaiyang Wang, Haoyuan Shi, Tianjiao Li, Liming Zhao, Hanfeng Zhai, Deepa Korani and Jingjie Yeo*



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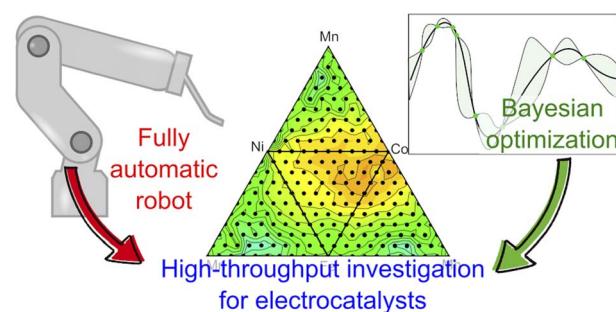
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An automatic robot system for machine learning-assisted high-throughput screening of composite electrocatalysts

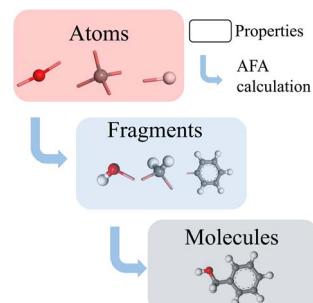
Masanori Kodera* and Kazuhiro Sayama*



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Haoxiang Lin and Xi Zhu*

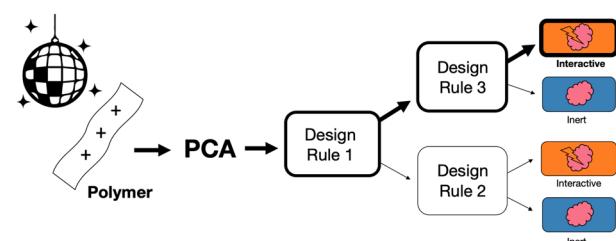


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An interpretable machine learning framework for modelling macromolecular interaction mechanisms with nuclear magnetic resonance

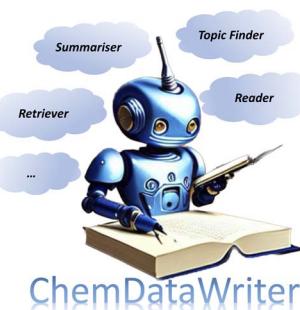
Samantha Stuart, Jeffrey Watchorn and Frank X. Gu*



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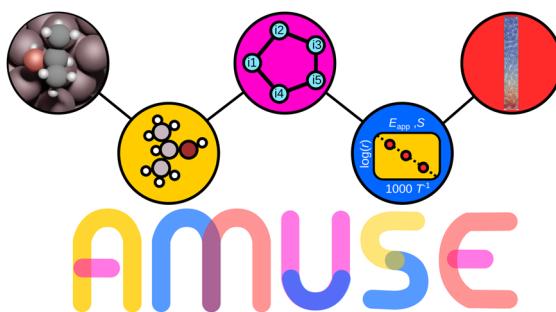
ChemDataWriter: a transformer-based toolkit for auto-generating books that summarise research

Shu Huang and Jacqueline M. Cole*



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Automated MULTiscale simulation environment

Albert Sabadell-Rendón,* Kamila Kaźmierczak, Santiago Morandi, Florian Euzenat, Daniel Curulla-Ferré and Núria López*

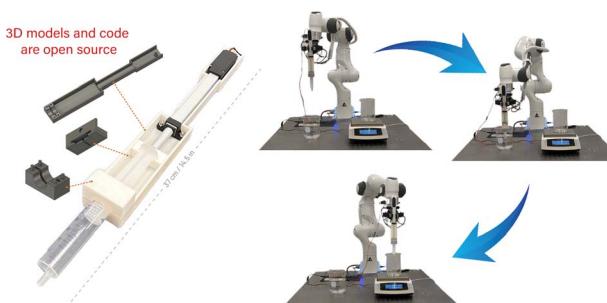
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Autonomous biomimetic solid dispensing using a dual-arm robotic manipulator

Ying Jiang, Hatem Fakhruldeen, Gabriella Pizzuto, Louis Longley, Ai He, Tianwei Dai, Rob Clowes, Nicola Rankin and Andrew I. Cooper*

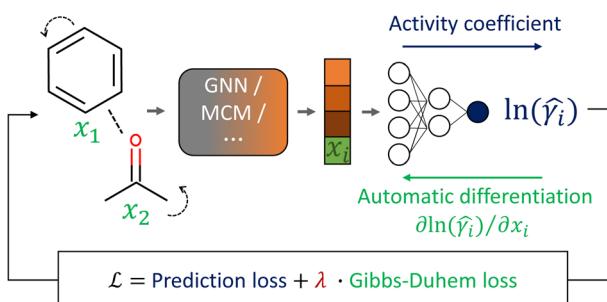
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Digital pipette: open hardware for liquid transfer in self-driving laboratories

Naruki Yoshikawa,* Kourosh Darvish, Mohammad Ghazi Vakili, Animesh Garg* and Alán Aspuru-Guzik*

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Gibbs–Duhem-informed neural networks for binary activity coefficient prediction

Jan G. Rittig, Kobi C. Felton, Alexei A. Lapkin and Alexander Mitsos*

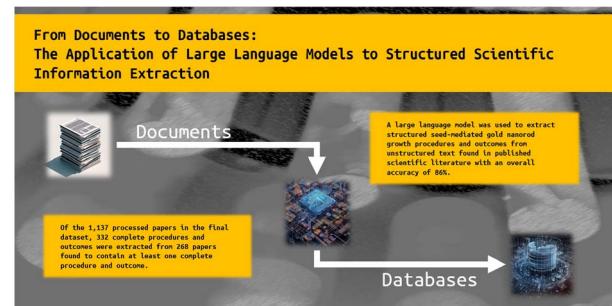


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Extracting structured seed-mediated gold nanorod growth procedures from scientific text with LLMs

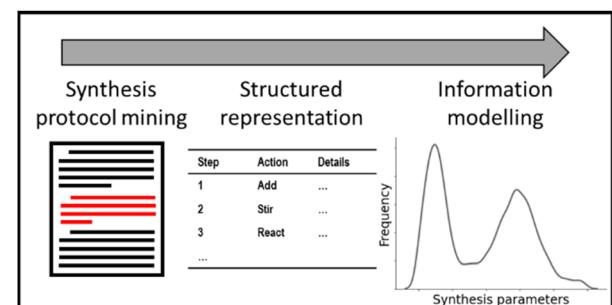
Nicholas Walker,* Sanghoon Lee, John Dagdelen, Kevin Cruse, Samuel Gleason, Alexander Dunn, Gerbrand Ceder, A. Paul Alivisatos, Kristin A. Persson and Anubhav Jain*



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Unveiling the synthesis patterns of nanomaterials: a text mining and meta-analysis approach with ZIF-8 as a case study

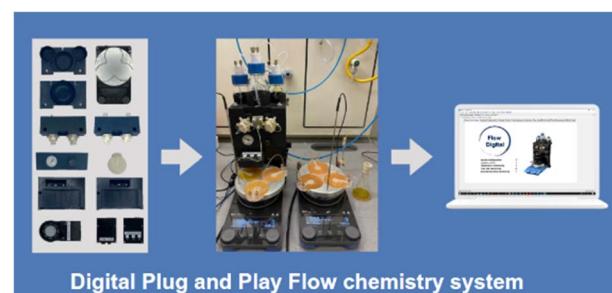
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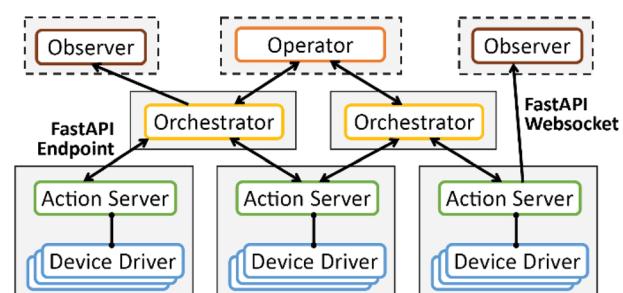
Mireia Benito Montaner, Matthew R. Penny and Stephen T. Hilton*



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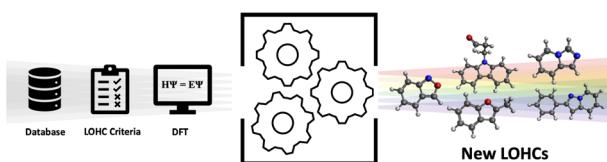
Orchestrating nimble experiments across interconnected labs

Dan Guevarra,* Kevin Kan, Yungchieh Lai, Ryan J. R. Jones, Lan Zhou, Phillip Donnelly, Matthias Richter, Helge S. Stein and John M. Gregoire*



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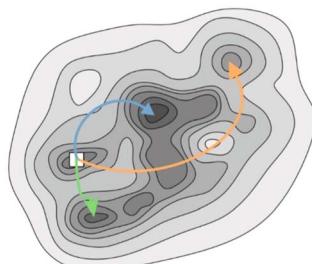
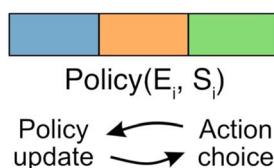
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Uncovering novel liquid organic hydrogen carriers: a systematic exploration of chemical compound space using cheminformatics and quantum chemical methods

Hassan Harb, Sarah N. Elliott, Logan Ward, Ian T. Foster, Stephen J. Klippenstein, Larry A. Curtiss and Rajeev Surendran Assary*

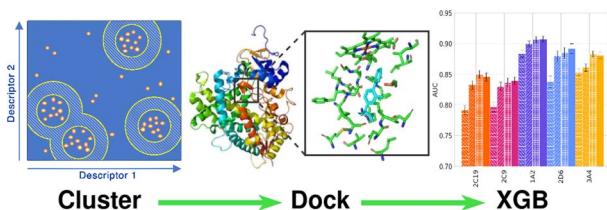
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RL-CSP

Reinforcement learning in crystal structure prediction

Elena Zamaraeva, Christopher M. Collins, Dmytro Antypov, Vladimir V. Gusev, Rahul Savani,* Matthew S. Dyer, George R. Darling, Igor Potapov, Matthew J. Rosseinsky* and Paul G. Spirakis

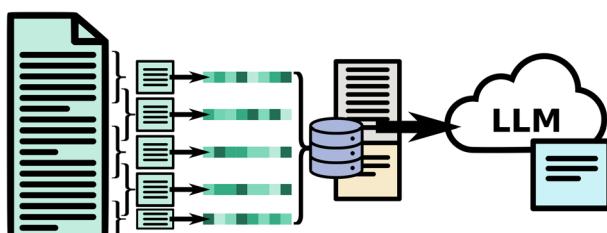
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Machine learning-augmented docking. 1. CYP inhibition prediction

Benjamin Weiser,* Jérôme Genzling, Mihai Burai-Patrascu, Ophélie Rostaing and Nicolas Moitessier*

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Domain-specific chatbots for science using embeddings

Kevin G. Yager*

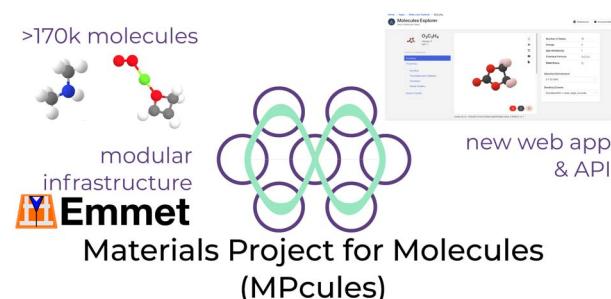


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A database of molecular properties integrated in the Materials Project

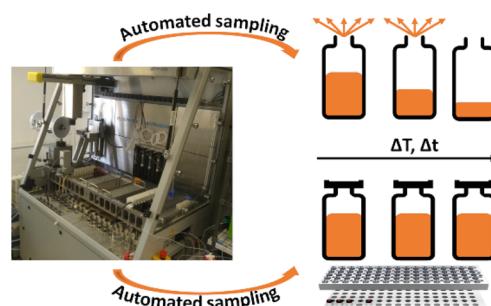
Evan Walter Clark Spotte-Smith,* Orion Archer Cohen, Samuel M. Blau, Jason M. Munro, Ruoxi Yang, Rishabh D. Guha, Hetal D. Patel, Sudarshan Vijay, Patrick Huck, Ryan Kingsbury, Matthew K. Horton and Kristin A. Persson*



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Best practice for sampling in automated parallel synthesizers

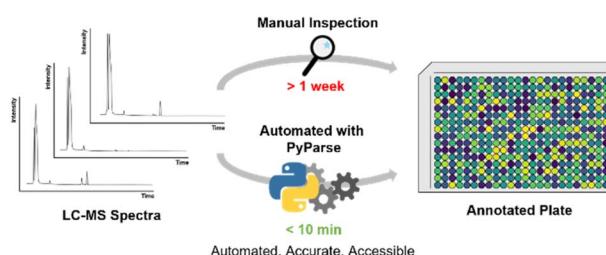
Michael Ringleb, Timo Schuett, Stefan Zechel and Ulrich S. Schubert*



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Automated LC-MS analysis and data extraction for high-throughput chemistry

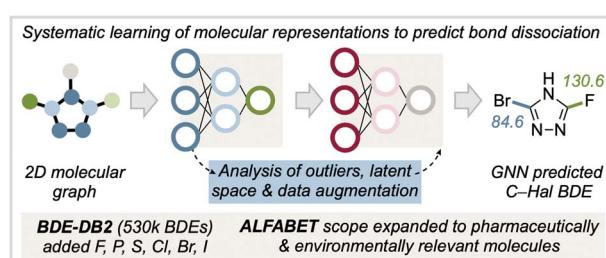
Joseph Mason,* Harry Wilders, David J. Fallon, Ross P. Thomas, Jacob T. Bush, Nicholas C. O. Tomkinson and Francesco Rianjongdee



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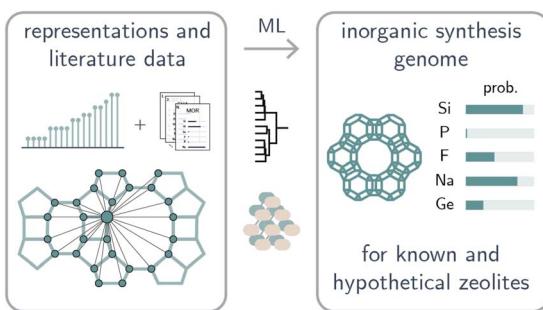
Expansion of bond dissociation prediction with machine learning to medicinally and environmentally relevant chemical space

Shree Sowndarya S. V., Yeonjoon Kim, Seonah Kim,* Peter C. St. John* and Robert S. Paton*



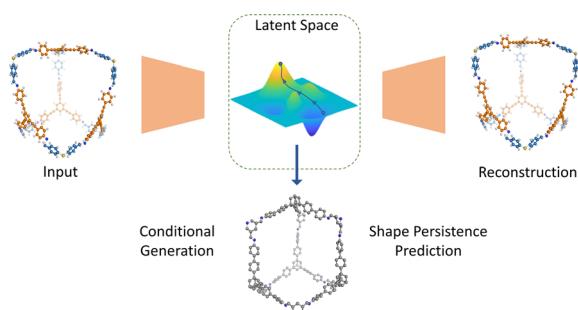
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1911


Inorganic synthesis-structure maps in zeolites with machine learning and crystallographic distances

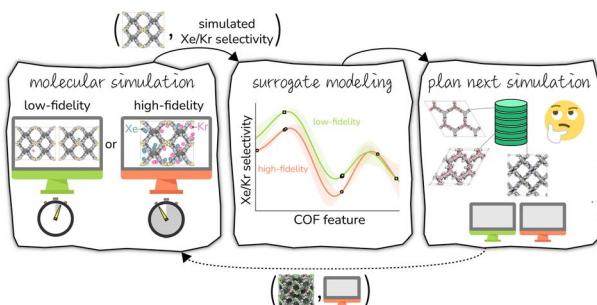
Daniel Schwalbe-Koda,* Daniel E. Widdowson, Tuan Anh Pham and Vitaliy A. Kurlin*

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Deep generative design of porous organic cages via a variational autoencoder

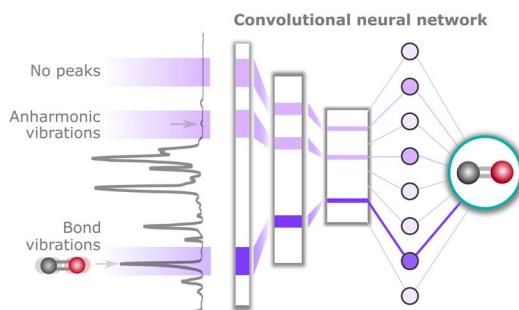
Jiajun Zhou, Austin Mroz and Kim E. Jelfs*

1937


Multi-fidelity Bayesian optimization of covalent organic frameworks for xenon/krypton separations

Nickolas Gantzler, Aryan Deshwal, Janardhan Rao Doppa* and Cory M. Simon*

1957


Understanding the patterns that neural networks learn from chemical spectra

Laura Hannemose Rieger, Max Wilson, Tejs Vegge and Eibar Flores*

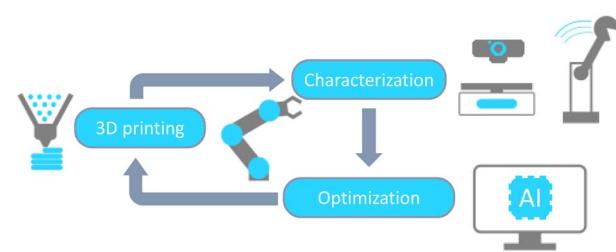


PAPERS

1969

Robotically automated 3D printing and testing of thermoplastic material specimens

Miguel Hernández-del-Valle, Christina Schenk, Lucía Echevarría-Pastrana, Burcu Ozdemir, Enrique Dios-Lázaro, Jorge Ilarraz-Zuazo, De-Yi Wang and Maciej Haranczyk*



1980

Towards a modular architecture for science factories

Rafael Vescovi, Tobias Ginsburg, Kyle Hippe, Doga Ozgulbas, Casey Stone, Abraham Stroka, Rory Butler, Ben Blaiszik, Tom Brettin, Kyle Chard, Mark Hereld, Arvind Ramanathan, Rick Stevens, Aikaterini Vriza, Jie Xu, Qingteng Zhang and Ian Foster*

