

A bit about me



From New Orleans, LA, USA

B.S. ChemE at Auburn University (2014)



Ph.D. ChemE at University of Colorado (2018)

- thermochemistry, symbolic ML, water splitting

Postdoc MSE at Berkeley (2019-2022)

- batteries, synthesis science, deep learning
- started working with Prof. Miura!



Prof at UMN in ChemE + MSE (2022-present)

- DFT, ML, solid-state chemistry, materials for energy



First trip to Japan (3 days ago – now!)

Objectives of this course

- understand where **ML fits into the broader** Solid-State Chemistry and Materials Science **disciplines**
- understand how to **apply ML algorithms** to problems in these disciplines using the **Python** programming language
- understand how to **formulate, execute, and interpret** ML projects

ML in design of chemicals and materials



FACEBOOK AI Carnegie Mellon University

Open Catalyst Project

Using AI to model and discover new catalysts to address the energy challenges posed by climate change.

Kitchin, John R. "Machine learning in catalysis." *Nature Catalysis* (2018).

opencatalystproject.org

BiopharmaTrend.com

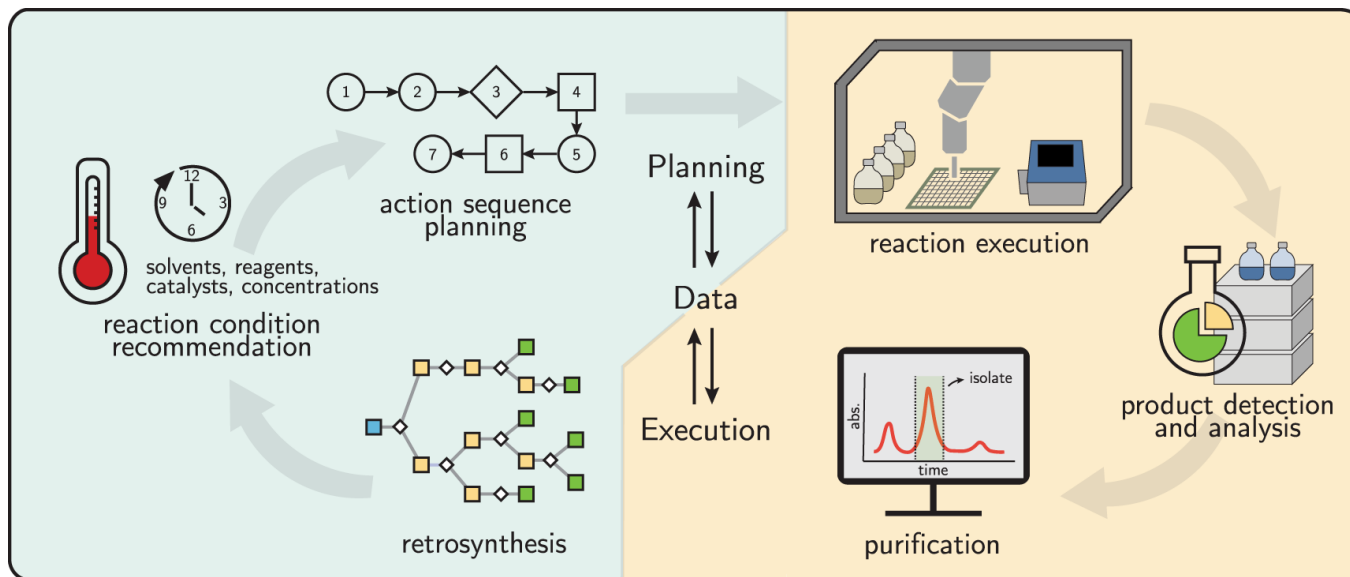
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9 Notable AI-powered Biotech Companies
Founded in 2021



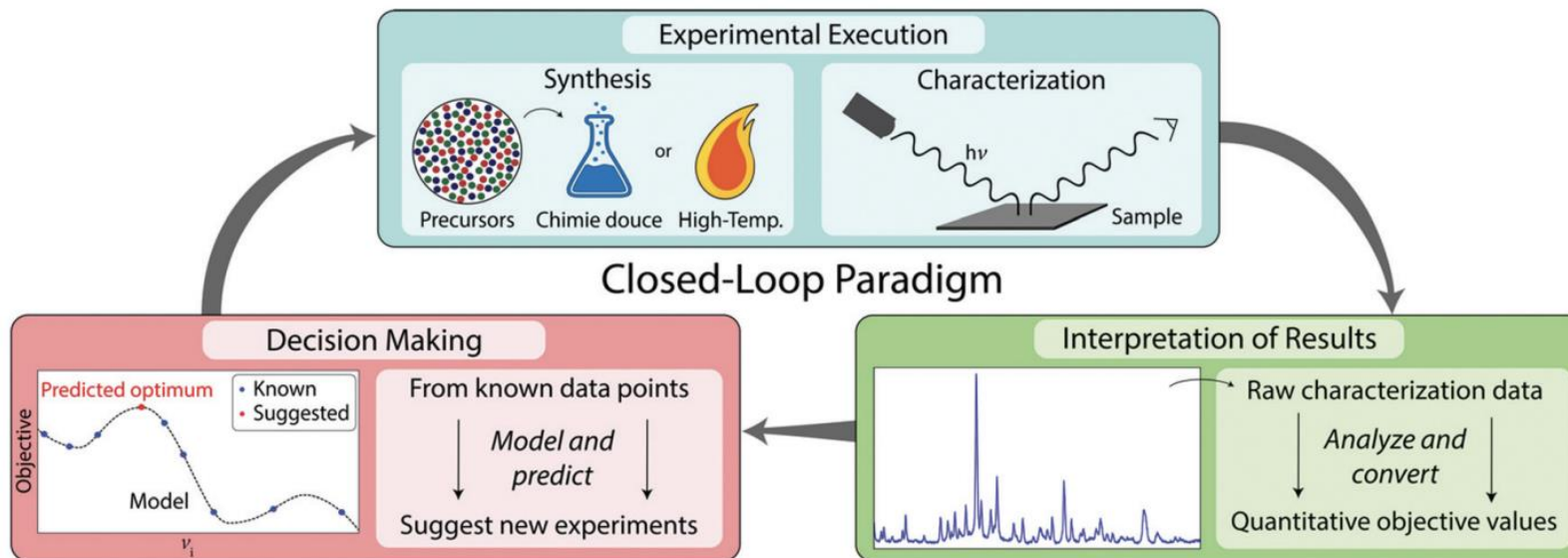
Chen, Hongming, *et al.* "The rise of deep learning in drug discovery." *Drug Discovery Today* (2018).

ML in design of chemicals and materials



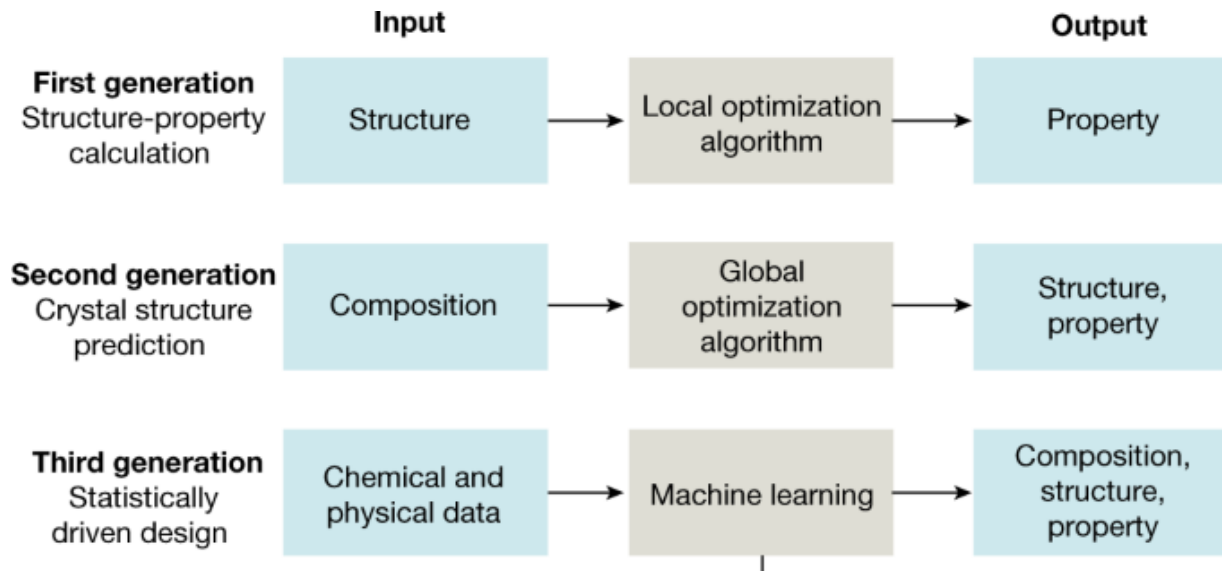
Gao, Raghavan, Coley. "Autonomous platforms for data-driven organic synthesis." *Nature Communications* (2022).

ML in design of chemicals and materials



Szymanski et al. "Toward autonomous design and synthesis of novel inorganic materials." *Materials Horizons* (2021).

ML in design of chemicals and materials



Butler et al. "Machine learning for molecules and materials science." *Nature* (2021).

What is the objective with ML?

Defining some basic terminology

Challenges of ML for science?

Given what you know now, discuss w/ one another some challenges of applying ML for science and engineering applications

Letter | [Published: 11 September 2019](#)

Anthropogenic biases in chemical reaction data hinder exploratory inorganic synthesis

[Xiwen Jia](#), [Allyson Lynch](#), [Yuheng Huang](#), [Matthew Danielson](#), [Immaculate Lang'at](#), [Alexander Milder](#), [Aaron E. Ruby](#), [Hao Wang](#), [Sorelle A. Friedler](#) ✉, [Alexander J. Norquist](#) ✉ & [Joshua Schrier](#) ✉

[Nature](#) **573**, 251–255 (2019) | [Cite this article](#)

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Some reactants are popular, but they aren't as good as random reactants!

Machine Learning May Sometimes Simply Capture Literature Popularity Trends: A Case Study of Heterocyclic Suzuki–Miyaura Coupling

Wiktor Beker, Rafał Roszak, Agnieszka Wołos, Nicholas H. Angello, Vandana Rathore, Martin D. Burke*, and Bartosz A. Grzybowski*

✓ **Cite this:** *J. Am. Chem. Soc.* 2022, 144, 11, 4819–4827

Publication Date: March 8, 2022 ✓

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Similar story in organic chemistry...

18 Nov 2022 in **Research & Technology**

Cuprate superconductivity mechanism may be coming into focus

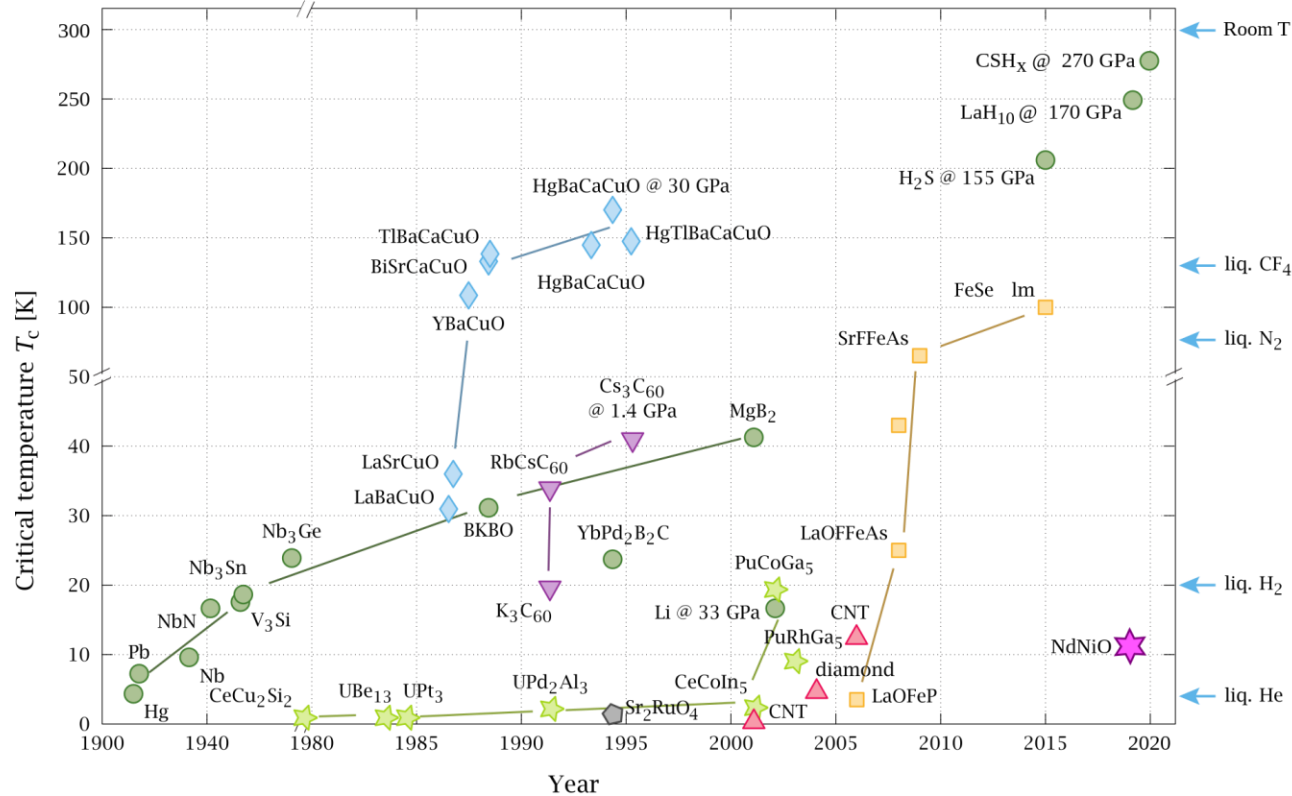
Theory and experiment have pinpointed factors that determine the magnetic attraction between electron pairs.

Daniel Garisto

Since their dramatic debut in 1986, cuprate superconductors have been some of the best-studied materials in existence. Nonetheless, many mysteries about the materials have

“Even if we are correct, in 30 years people will still say that the theory is not understood,” Tremblay says.

Challenges of ML for science?



Challenges of ML for science?

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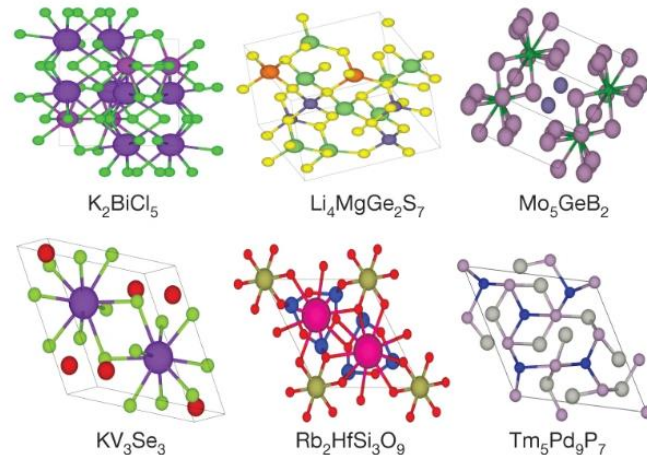
Article | [Open access](#) | [Published: 29 November 2023](#)

Scaling deep learning for materials discovery

[Amil Merchant](#) ✉, [Simon Batzner](#), [Samuel S. Schoenholz](#), [Muratahan Aykol](#), [Gowoon Cheon](#) & [Ekin Dogus Cubuk](#) ✉ ← PhD Physics

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nature

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An autonomous laboratory for the accelerated synthesis of novel materials

[Nathan J. Szymanski](#), [Bernardus Rendy](#), [Yuxing Fei](#), [Rishi E. Kumar](#), [Tanjin He](#), [David Milsted](#), [Matthew J. McDermott](#), [Max Gallant](#), [Ekin Dogus Cubuk](#), [Amil Merchant](#), [Haegyeom Kim](#), [Anubhav Jain](#), [Christopher J. Bartel](#), [Kristin Persson](#), [Yan Zeng](#) ✉ & [Gerbrand Ceder](#) ✉

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PhD Materials Science

UMN CEMS postdoc!



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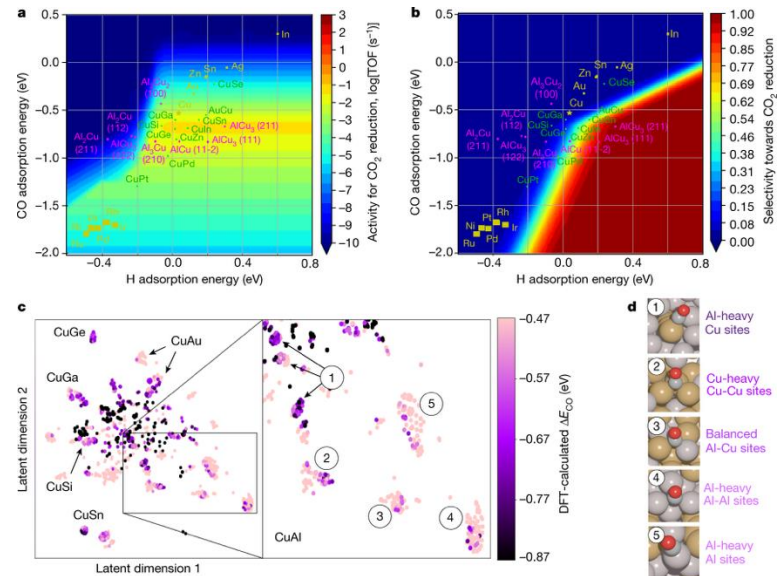
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Accelerated discovery of CO₂ electrocatalysts using active machine learning

[Miao Zhong](#), [Kevin Tran](#), [Yimeng Min](#), [Chuanhao Wang](#), [Ziyun Wang](#), [Cao-Thang Dinh](#), [Phil De Luna](#), [Zongqian Yu](#), [Armin Sedighian Rasouli](#), [Peter Brodersen](#), [Song Sun](#), [Oleksandr Voznyy](#), [Chih-Shan Tan](#), [Mikhail Askerka](#), [Fanglin Che](#), [Min Liu](#), [Ali Seifitokaldani](#), [Yuanjie Pang](#), [Shen-Chuan Lo](#), [Alexander Ip](#), [Zachary Ulissi](#) ✉ & [Edward H. Sargent](#) ✉

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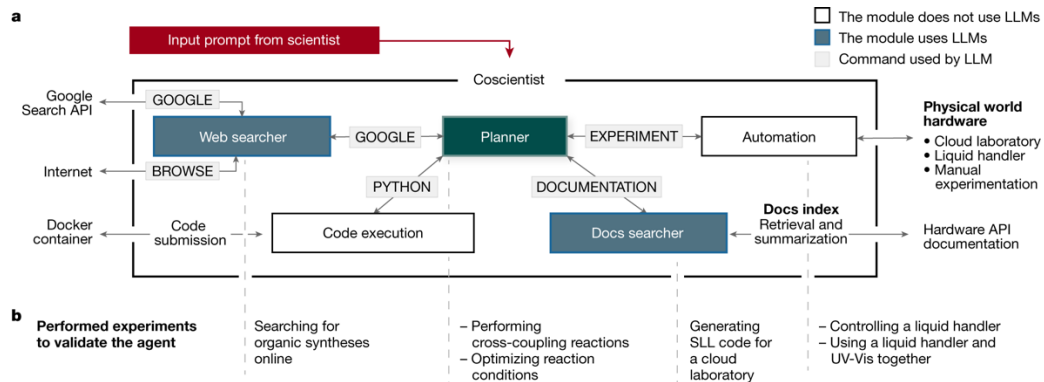
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Autonomous chemical research with large language models

[Daniil A. Boiko](#), [Robert MacKnight](#), [Ben Kline](#) & [Gabe Gomes](#) ✉

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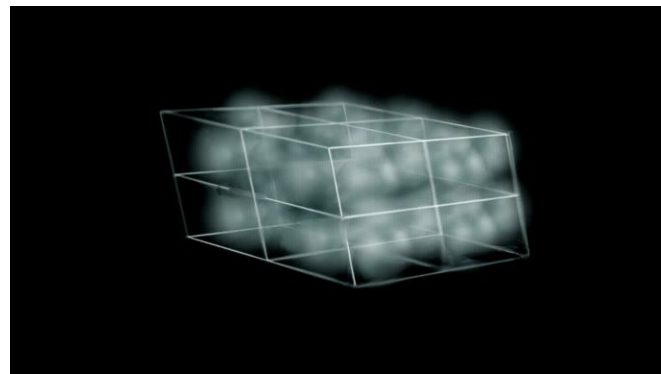
A generative model for inorganic materials design

[Claudio Zeni](#), [Robert Pinsler](#), [Daniel Zügner](#), [Andrew Fowler](#), [Matthew Horton](#), [Xiang Fu](#), [Zilong Wang](#), [Aliaksandra Shysheya](#), [Jonathan Crabbé](#), [Shoko Ueda](#), [Roberto Sordillo](#), [Lixin Sun](#), [Jake Smith](#), [Bichlien Nguyen](#), [Hannes Schulz](#), [Sarah Lewis](#), [Chin-Wei Huang](#), [Ziheng Lu](#), [Yichi Zhou](#), [Han Yang](#), [Hongxia Hao](#), [Jielan Li](#), [Chunlei Yang](#), [Wenjie Li](#), ... [Tian Xie](#)  [+ Show authors](#)

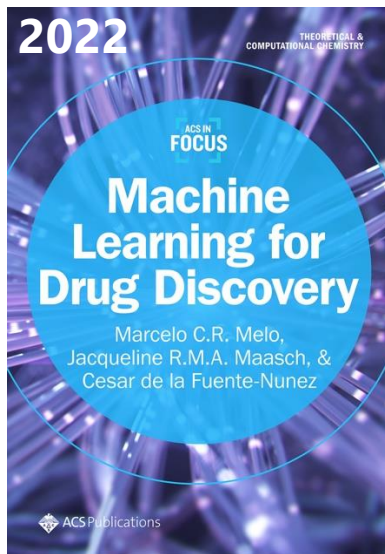
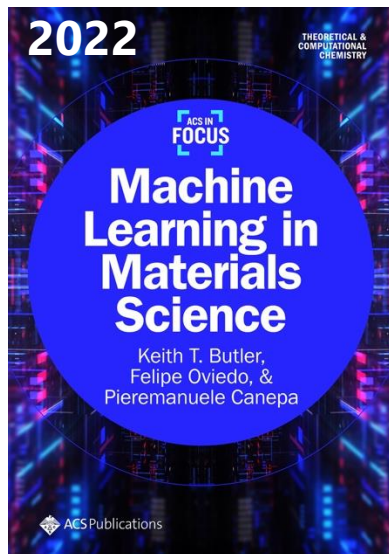
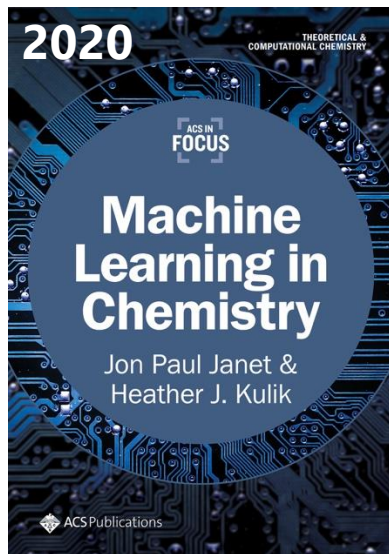
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PhD Materials Science



Machine learning interest on the rise!



You do not need a CS degree to make meaningful contributions with ML!