Algebraic Machine Learning



Fundação Champalimaud

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What is Algebraic Machine Learning?

Algebraic Machine Learning (Martin-Maroto, de Polavieja, arXiv:1803.05252)

- Learning system inspired by Model Theory and Universal Algebra
- Data and prior knowledge are formally embedded as generating elements and relations in an algebraic structure
- Learning from data via iterative modification of the algebraic structure
 - No error minimization
 - Mathematically transparent
 - Capable of finding formal knowledge
 - No hyperparameter tuning

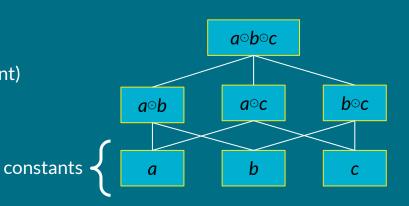
Let's give a short introduction!

- Algebraic structure: semilattices
 - o Operation: (associative, commutative, idempotent)
 - Order relation: $a \le b \Leftrightarrow a \circ b = b$
- Generating elements: constants
 - Components of data



Datapoints can be represented using the operation





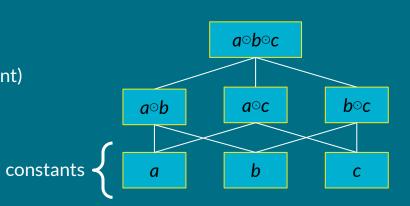
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 - Concepts (e.g. labels)

v: vertical bar



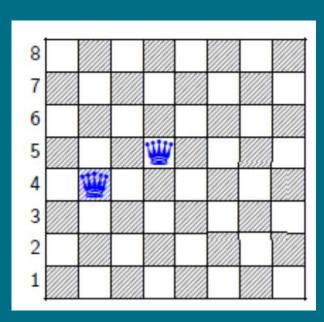


"v is not in the image"



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N-queens problem

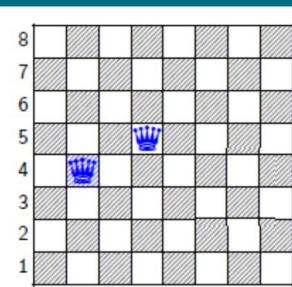
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 - Components of data
 - Concepts (e.g. labels)
- Prior knowledge and data: relations

Martin-Maroto - Polavieja, Semantic embeddings in Semilattices, arXiv:2205.12618

$$\forall x \forall y (R_x \circ C_y < Q_{xy})$$

N-queens problem

But how are they enforced?



• Atomization: a set theoretical based representation of a semilattice

Martin-Maroto - de Polavieja, Finite Atomized Semilattices, arXiv:2102.08050

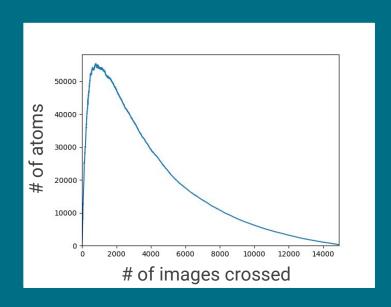
- Atoms φ are determined by a set of constants
- $\phi < a \Leftrightarrow \exists c \text{ constant}, \phi < c \le a$
- \circ $a \le b \Leftrightarrow \exists \phi, \phi < a \text{ and } \phi \not < b \text{ (ϕ discriminates (a,b))}$

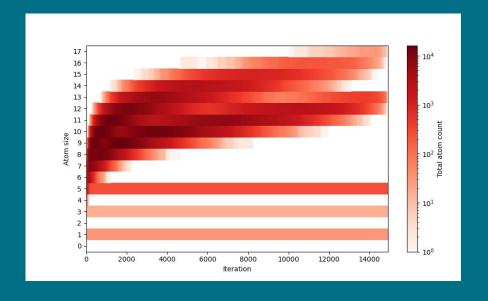
Proposition. Every finite semilattice can be represented through an atomization.

• Atomization: a set theoretical based representation of a semilattice

Martin-Maroto - de Polavieja, Finite Atomized Semilattices, arXiv:2102.08050

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- Crossing: obtain a new semilattice verifying a new relation
 - Full crossing
 Freest semilattice → guaranteed to find rule in data, slower

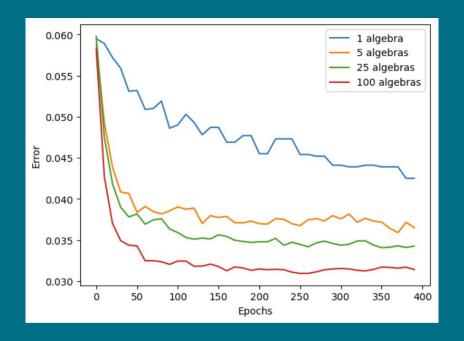




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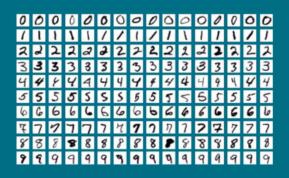
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- Crossing: obtain a new semilattice verifying a new relation
 - Full crossing
 Freest semilattice → guaranteed to find rule in data, slower
 - Sparse crossing
 Less free, more logical consequences → better generalization
 Retain only some of the atoms (stochastic) → faster learning from data
 No overfitting
 Very parallelizable



But does it work?

Some benchmarks

MNIST as a benchmark: Handwritten digit classification



Binary classification: is it or is it not a given number?

- → AML with no prior knowledge: < 0.5% error
- → Multi-layer perceptron: < 0.6% error

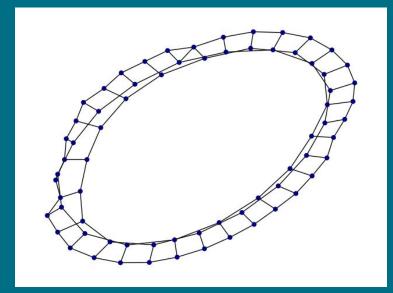
General classification task

- → AML with simple thresholding: ~97.6% accuracy
- → AML with optimization: ~98.2% accuracy
- → Multi-layer perceptron: ~98.3% accuracy
- → Convolutional neural network: 99%+ accuracy

Open question: how can we effectively inject prior knowledge?

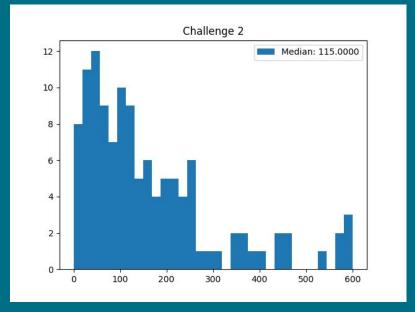
Some benchmarks

Hamiltonian cycles

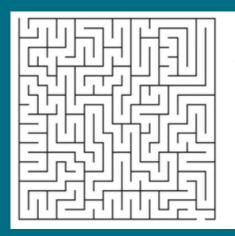


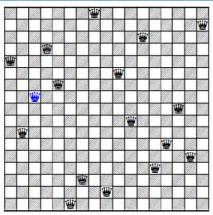
Flinders Hamiltonian Cycle Project, graph 2

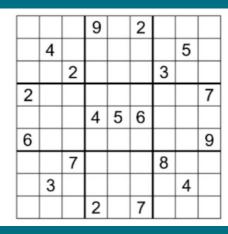
- Snakes and ladders: 5078 transforms
- Algebraic Machine Learning

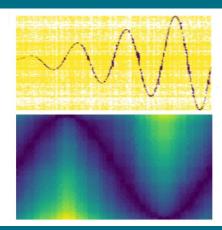


Researching further applications









Further research

- Real values/inference
 - o Infinite Atomized Semilattices (F. Martin-Maroto, A. Ricciardo, D. Mendez, G. de Polavieja, arXiv:2311.01389)
 - Extended semilattices
- Automatic discovery of constants
- Statistical study of atom weights
- Training performance analysis/distances between models
- Automatic tools for model explanation
- Development of more user-friendly software



Thank you for your attention

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