Algorithms in Cheminformatics – DM840 (autumn 2022)

Daniel Merkle

This is the slide set used for a small video presentation which you can find here: <u>https://www.imada.sdu.dk/~daniel/2022-DM840-pizza.mp4</u>

LETS START WITH A GAME

(CATALAN)

https://www.imada.sdu.dk/~daniel/DM840-2021/assignment1/assign1-2021.html

https://www.imada.sdu.dk/~daniel/DM840-2022/ (For video and slide set)

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DM840 – Algorithms in Cheminformatics



A "solution" is a shortest path in a directed graph



$\begin{array}{l} \mbox{Reactions} - \mbox{Application of Transformation Rules} \\ \mbox{1,2-ethenediol} + \mbox{formaldehyde} \xrightarrow{\mbox{aldol addition}} \mbox{glyceraldehyde} \end{array}$



The Formose Chemistry

4 reaction patterns:













Generation 4

Shortest Path(s)



A "solution" is a shortest path in a directed graph

http://cheminf.imada.sdu.dk/mod

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Shortest Path(s)



A "solution" is a shortest path in a directed graph



How to find the k best solutions?

In hypergraphs?

Algorithms?

Computational Complexity?

How chemists do it



Are these 2 colored graphs isomorph?



Algorithms? Computational Complexity? State-of-the-Art?

(Sub-)Graph Isomorphism / Canoncial Representaions of Graphs

Are these 2 colored graphs isomorph?



A Quasipolynomial Time Algorithm for Graph Isomorphism: The Details

Posted on November 12, 2015 by j2kun

Update 2015-11-21: Ken Regan and Dick Lipton posted an article with some more details, and a high level overview of how the techniques fit into the larger picture of CS theory.

Update 2015-11-16: Laci has posted the talk on his website. It's an hour and a half long, and I encourage you to watch it if you have the time \bigcirc

Laszlo Babai has claimed an astounding theorem, that the Graph Isomorphism problem can be solved in quasipolynomial time. On Tuesday I was at Babai's talk on this topic (he has yet to release a preprint), and I've compiled my notes here. As in Babai's talk, familiarity with basic group theory and graph theory is assumed, and if you're a casual (i.e., math-phobic) reader looking to understand what the fuss is all about, this is probably not the right post for you. This post is research level theoretical computer science. We're here for the juicy, glorious details.

Note: this blog post will receive periodic updates as my understanding of the details improve.



Algorithms? Computational Complexity? State-of-the-Art?

(Sub-)Graph Isomorphism / Canoncial Representaions of Graphs



Canoncial representations of graphs (i.e. molecules):

Algorithms? Computational Complexity? And ... w.t.h. went wrong in Chemistry?

Counting



More Counting

Variations on a theme 1: Trees

Fundamental construct



Variation on the theme: restrict each node to 0 or 2 children



Applications: Sampling, Analysis of Algorithms, Chemistry



Even more counting





ŃН2

Cycle Bases (Scheduling, Distributed Computing, Chemistry)



How to find a minimal cycle basis? Algorithms? Computaional Complexity?

...and w.t.h. went wrong in chemistry?

QSAR / PCA



Petri Nets





...and many more For example: Traverse a tetrahedron, each edge twice.



...and many more For example: Traverse a tetrahedron, each edge twice.



Topics

...

Canonical Representations Structure Descriptors **Graph Isomorphism** Pólya's **Counting Theory** Artificial Chemistries **Graph Grammars** Cycle Bases Metabolic Networks and Metabolic Pathways Concurrency Theory Petrinets Applications in Industry

Formal Model

Algorithms

Computational Complexity

Applications (including Chemistry)

Topics

...

Canonical Representations
Structure Descriptors
Graph Isomorphism
Pólya's Counting Theory
Artificial Chemistries
Graph Grammars
Cycle Bases
Metabolic Networks and Metabolic Pathways
Concurrency Theory
Petrinets
Applications in Industry

- 10 ECTS

- max. 2 programming projects

no chemistry knowledge needed!

Algorithms in Cheminformatics - DM840



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D-BASF

We create chemistry





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Enzyme Design

https://cheminf.imada.sdu.dk/novo-synergy/

Base

Nuc

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