

Color Refinement for Relational Structures

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AlMoTh'25

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What the next 20 minutes will be about

Color Refinement

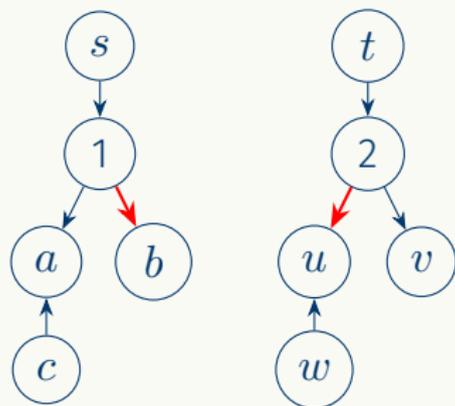
How to lift CR to Relational Structures

Doing it Efficiently

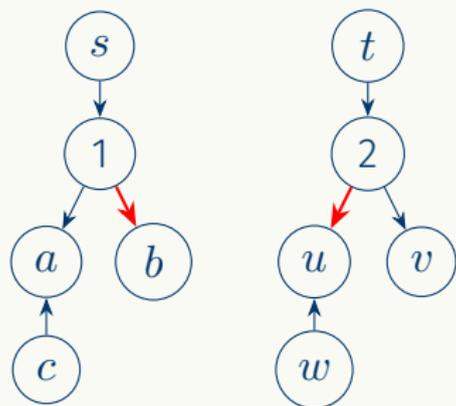
Simpler Approaches?

Outlook

Color Refinement: Iteratively compute stable coloring γ_∞



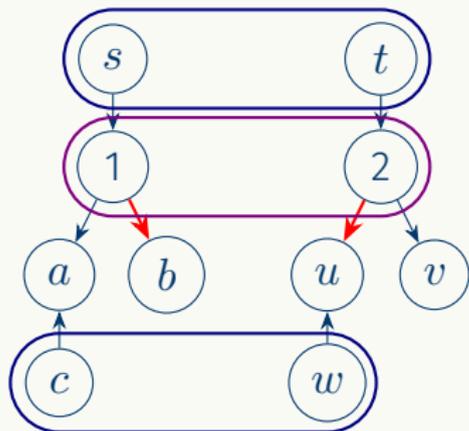
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ex. edge and vertex colors λ, α s.t.
num. of λ -neighbors of color α differ.

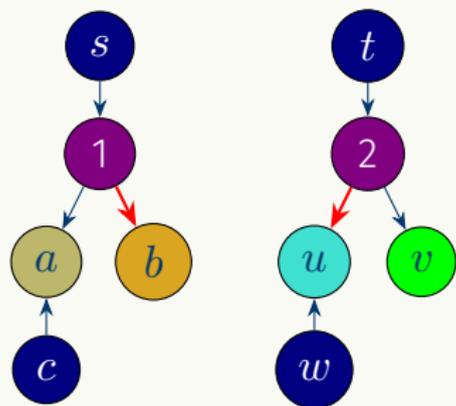
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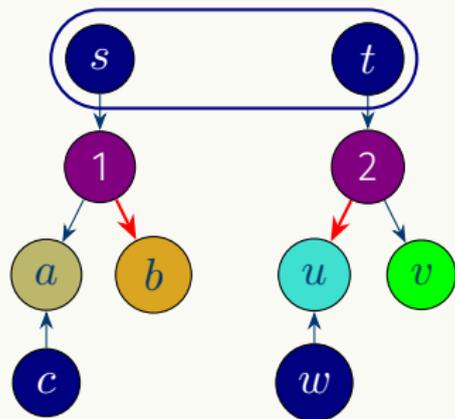
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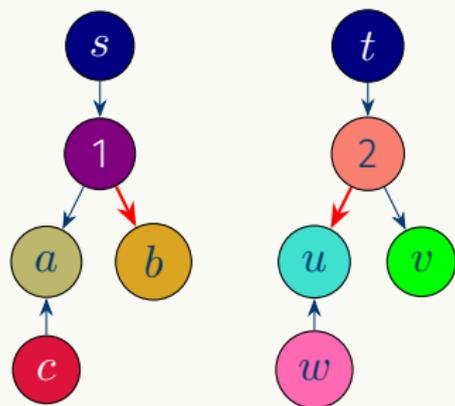
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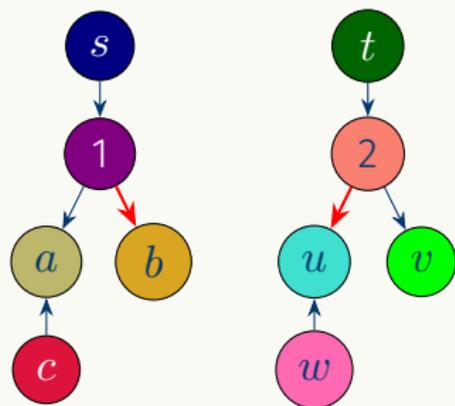
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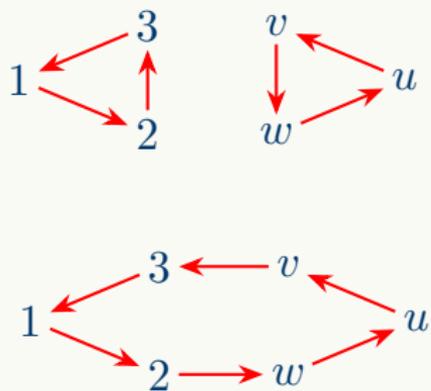
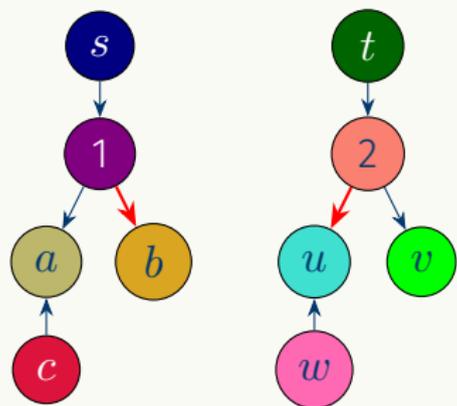
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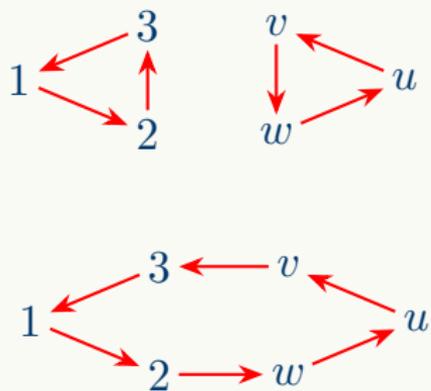
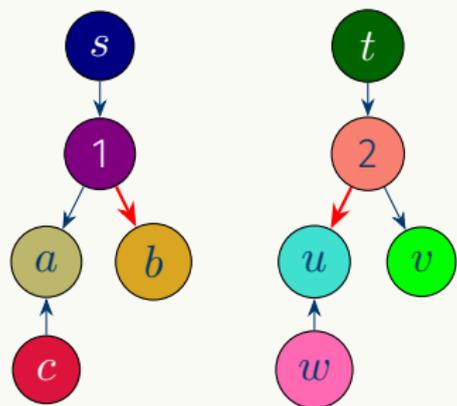
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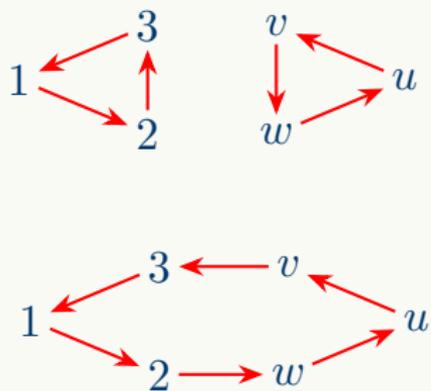
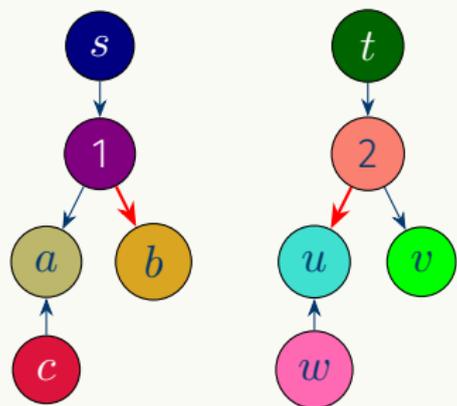


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Wanted: similar method for more complex objects.

In this talk: Relational structures.



But how to identify the “right” approach?

Key Properties of Color Refinement

Theorem (Dvořák 2010; Dell, Grohe, Rattan 2018; Immerman, Lander 1990; and others)

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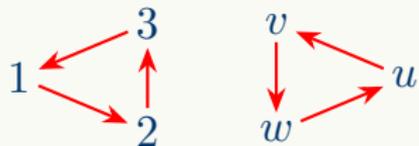
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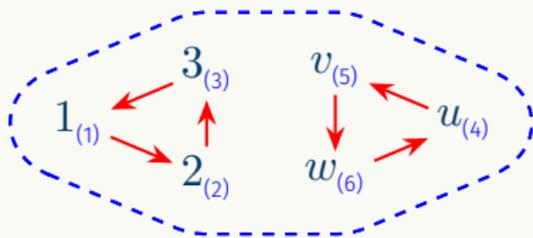
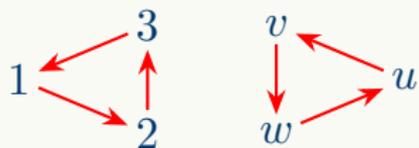
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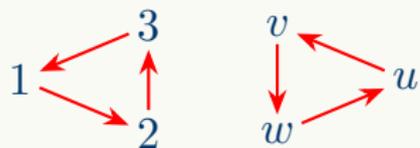
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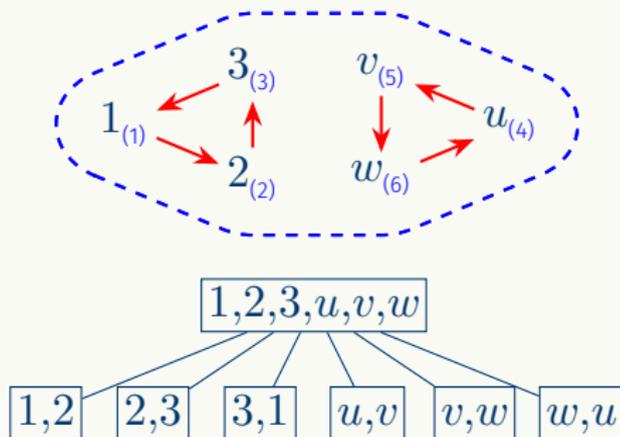
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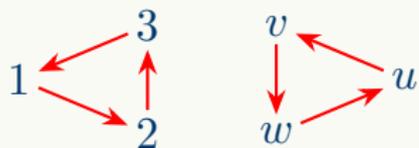


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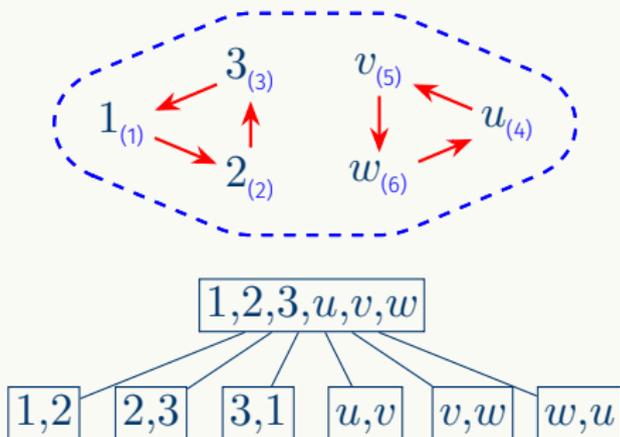


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- **Standard, most general notion from DB theory**, (Abiteboul, Hull, Vianu 1995).

Atomic formulas and Boolean operators as usual:

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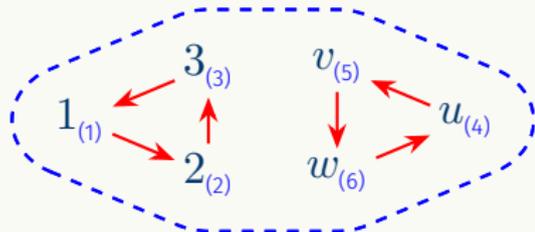
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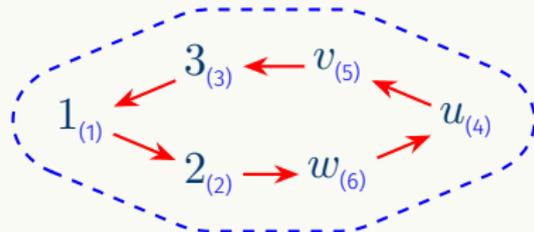
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GF(C) — Example

 $R^{\mathcal{A}} \quad 1, 2, 3, u, v, w$

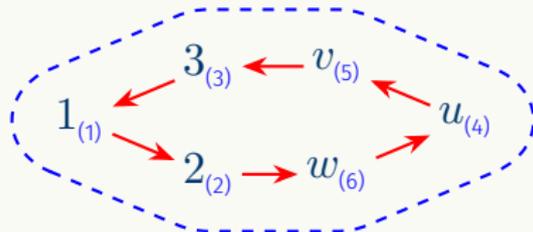
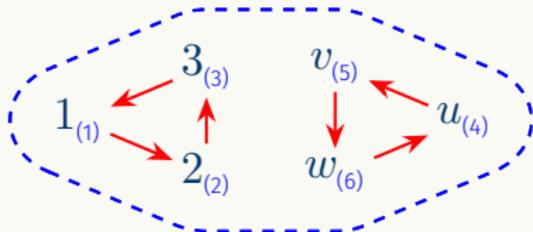
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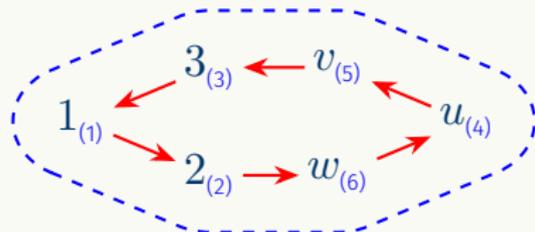
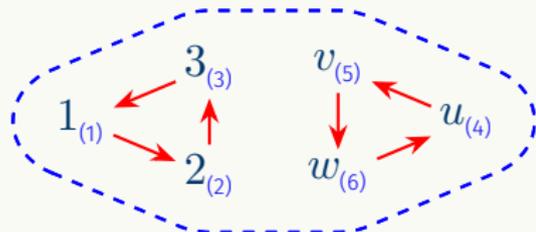


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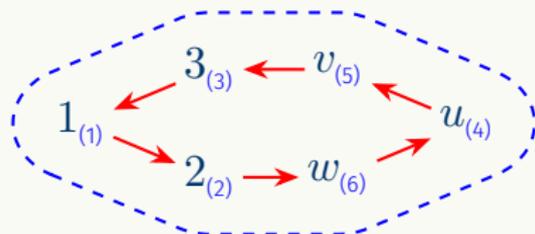
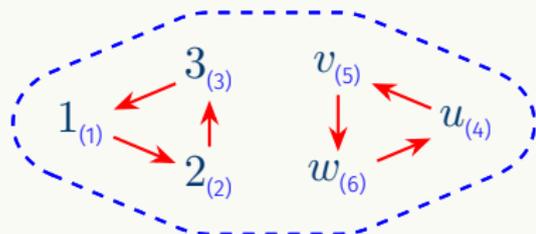
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RCR should also distinguish \mathcal{A} and \mathcal{B} !

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Still based on “**message passing**”: should be able to encode \mathcal{A} as a graph $G_{\mathcal{A}}$ and use Color Refinement on $G_{\mathcal{A}}$.

Relational Color Refinement

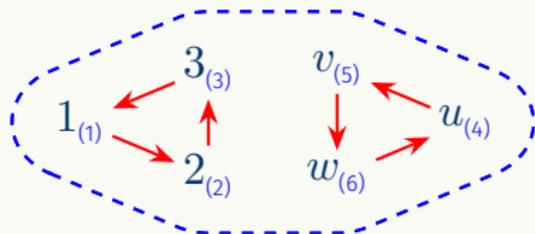
Idea: Color *tuples* instead of *elements*:

$$\varrho_{i+1}(\mathbf{a}) := (\varrho_i(\mathbf{a}), \{ \{ (stp(\mathbf{a}, \mathbf{b}), \varrho_i(\mathbf{b})) : \mathbf{a} \cap \mathbf{b} \neq \emptyset \} \})$$

Still based on “message passing”: should be able to encode \mathcal{A} as a graph $G_{\mathcal{A}}$ and use Color Refinement on $G_{\mathcal{A}}$.

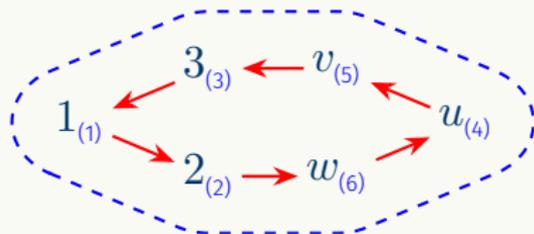
$$R^{\mathcal{A}} \quad 1,2,3,u,v,w$$

$$E^{\mathcal{A}} \quad \begin{array}{ccc} 1,2 & 2,3 & 3,1 \\ u,v & v,w & w,u \end{array}$$



$$R^{\mathcal{B}} \quad 1,2,3,u,v,w$$

$$E^{\mathcal{B}} \quad \begin{array}{ccc} 1,2 & 2,3 & 3,1 \\ u,v & v,w & w,u \end{array}$$

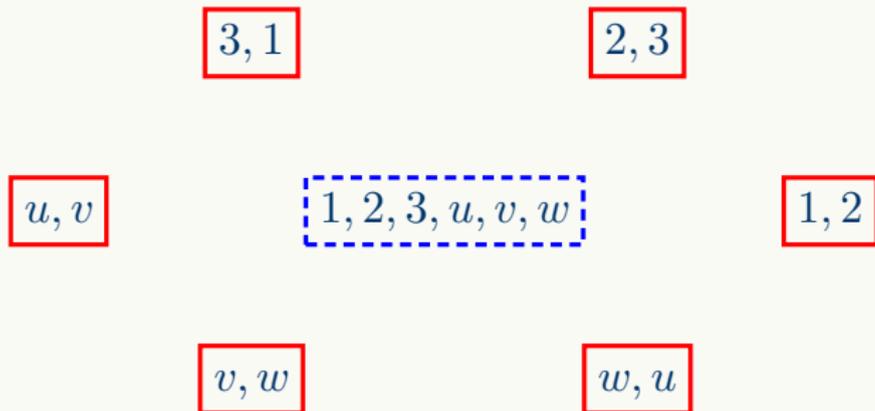


Representing a Structure as a Graph — $G_{\mathcal{A}}$

$R^{\mathcal{A}}$	1,2,3,u,v,w		
$E^{\mathcal{A}}$	1,2	2,3	3,1
	u,v	v,w	w,u

Representing a Structure as a Graph — $G_{\mathcal{A}}$

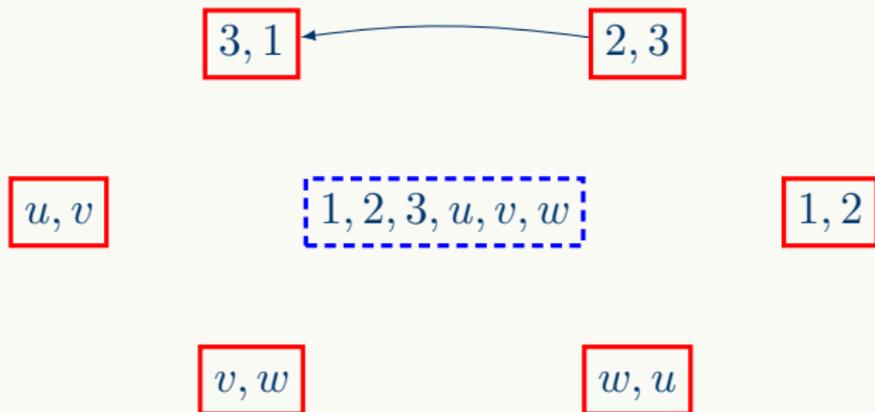
$R^{\mathcal{A}}$	$1, 2, 3, u, v, w$		
$E^{\mathcal{A}}$	$1, 2$	$2, 3$	$3, 1$
	u, v	v, w	w, u



Representing a Structure as a Graph — $G_{\mathcal{A}}$

Draw edge if tuples “intersect”

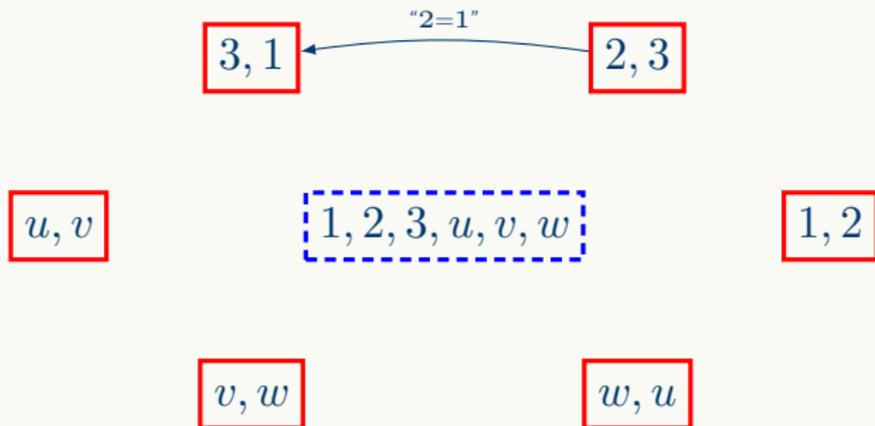
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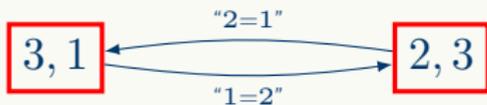
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	u, v	v, w	w, u



u, v

$1, 2, 3, u, v, w$

$1, 2$

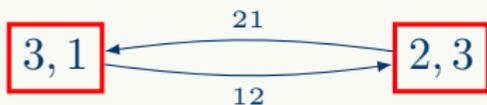
v, w

w, u

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$1, 2$

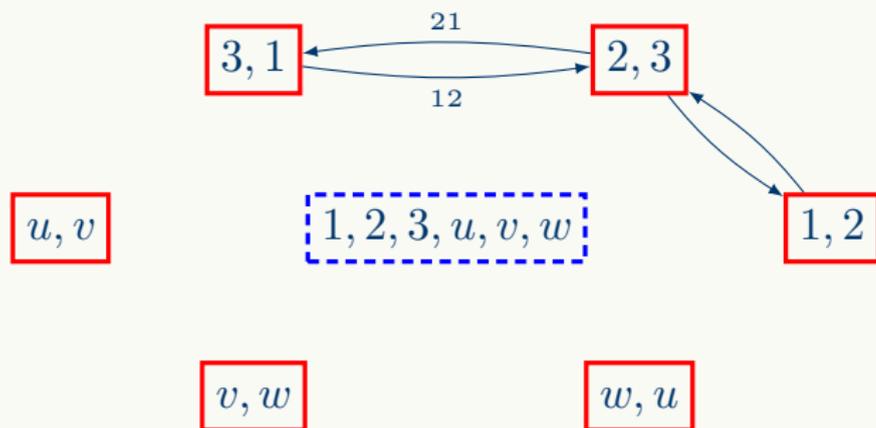
v, w

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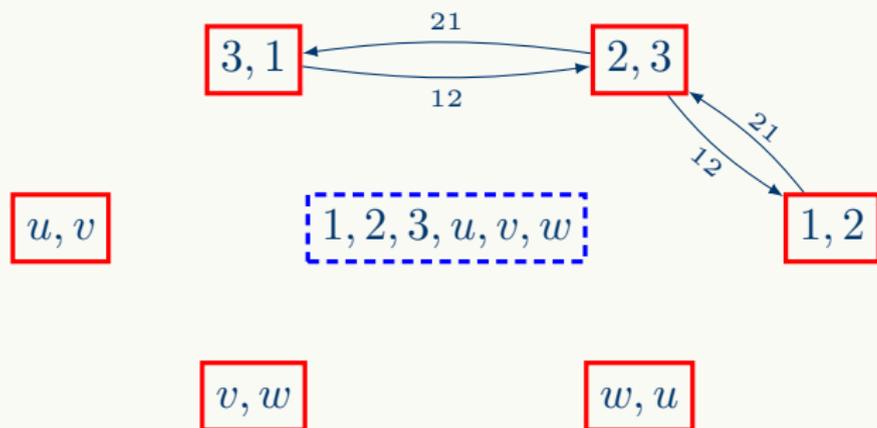
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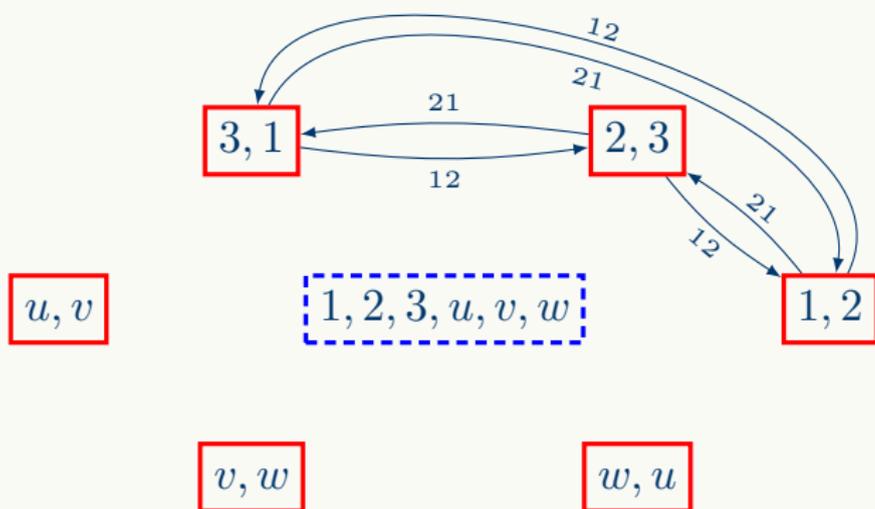
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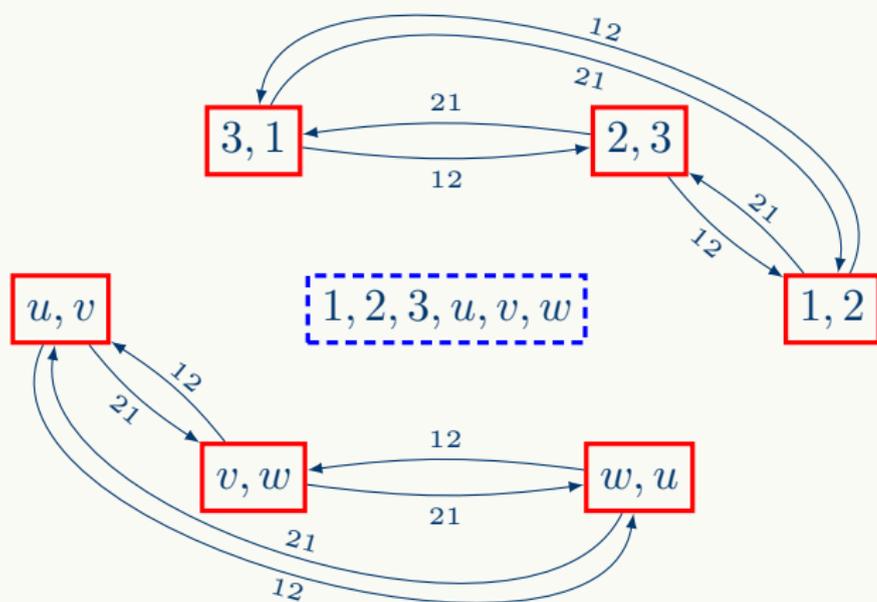
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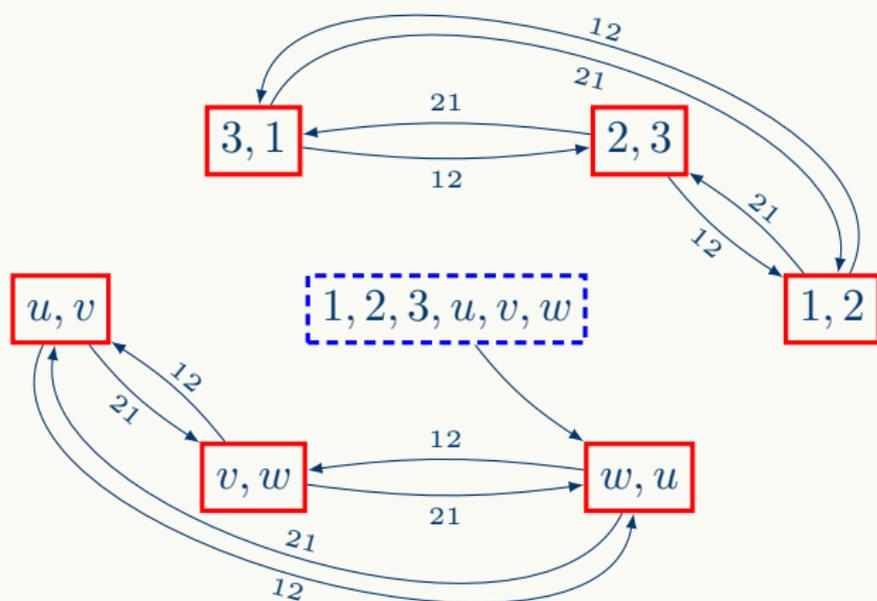
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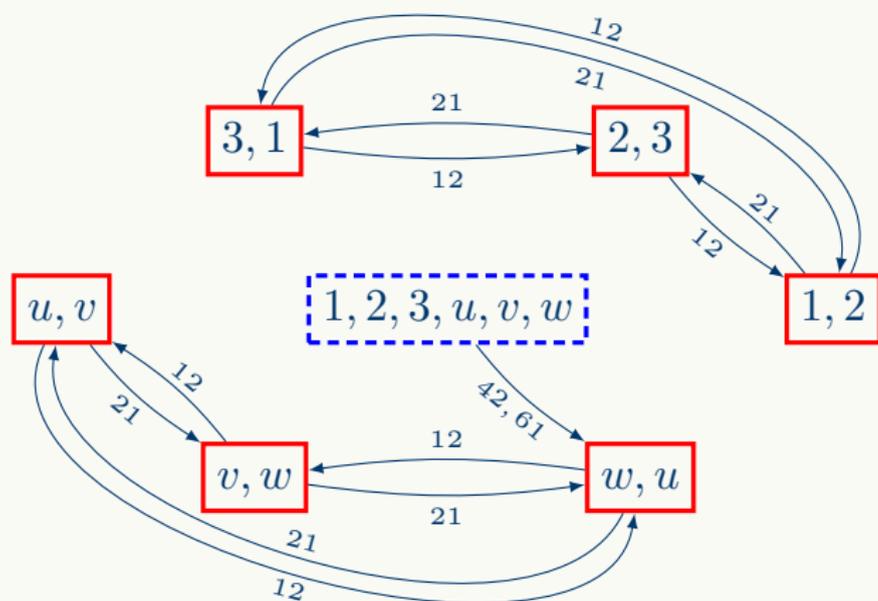


Representing a Structure as a Graph — $G_{\mathcal{A}}$

Draw edge if tuples “intersect”

R^A	1,2,3,u,v,w		
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E^A	1,2	2,3	3,1
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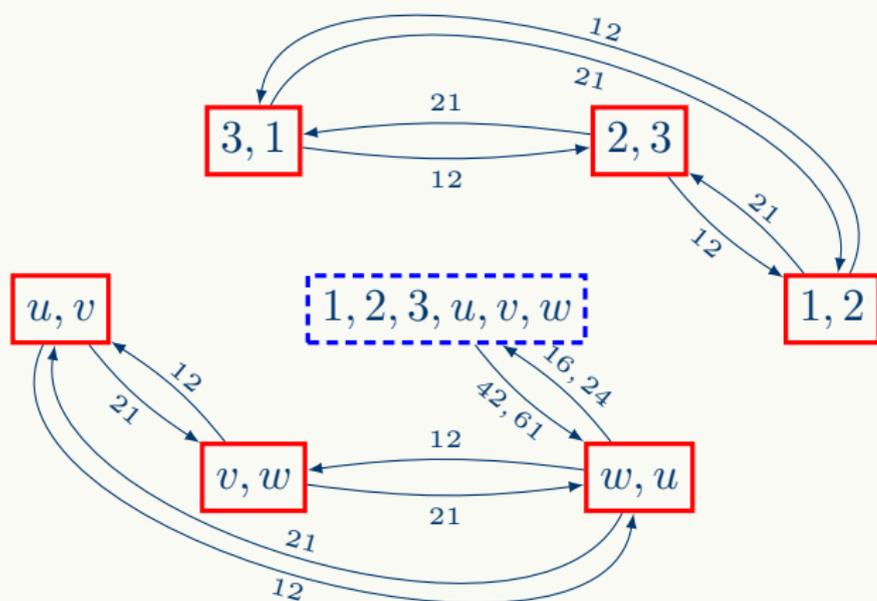


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Draw edge if tuples “intersect”

$R^{\mathcal{A}}$	1,2,3,u,v,w		
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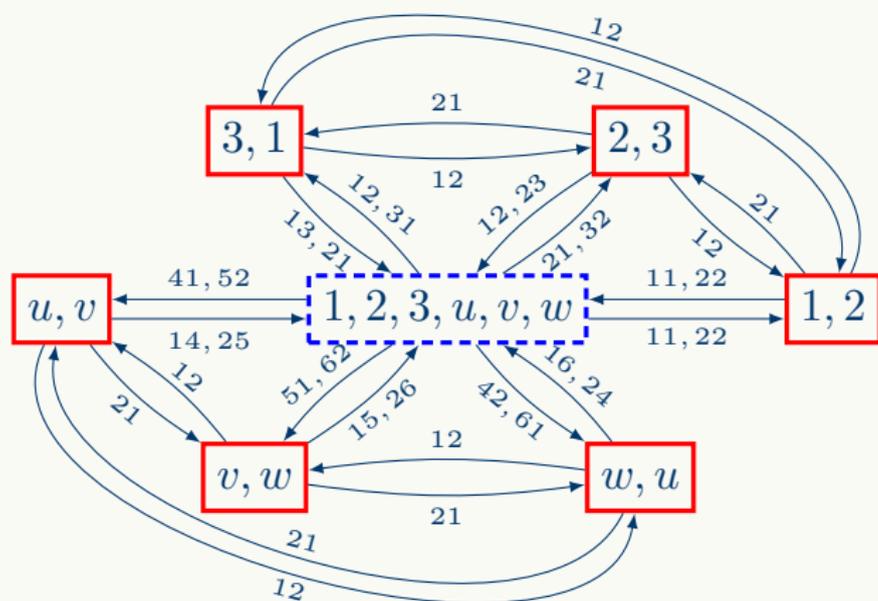
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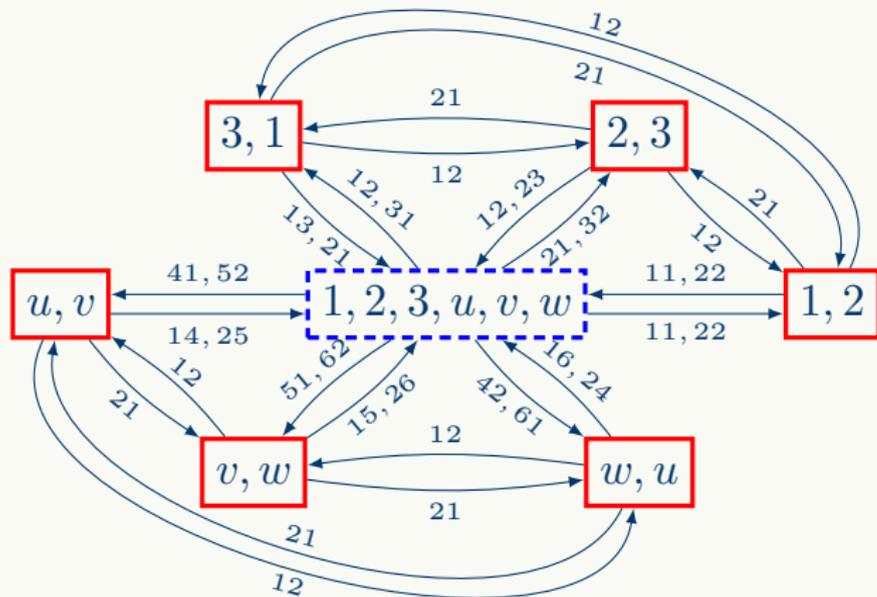
Representing a Structure as a Graph — G_A

Draw edge if tuples “intersect”

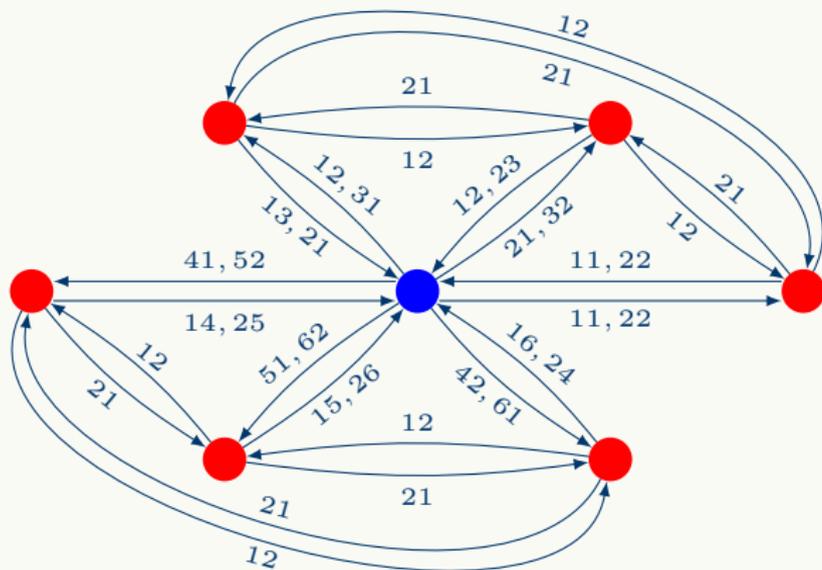
R^A	1,2,3,u,v,w		
E^A	1,2	2,3	3,1
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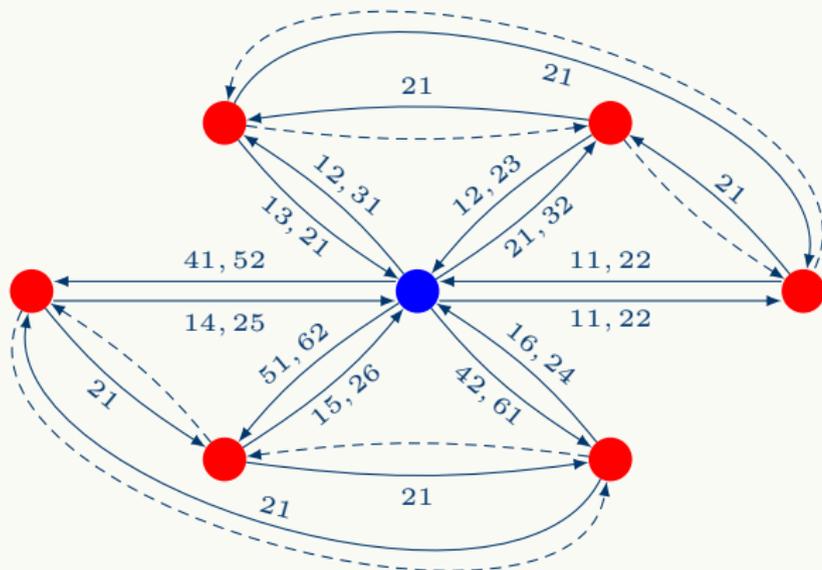
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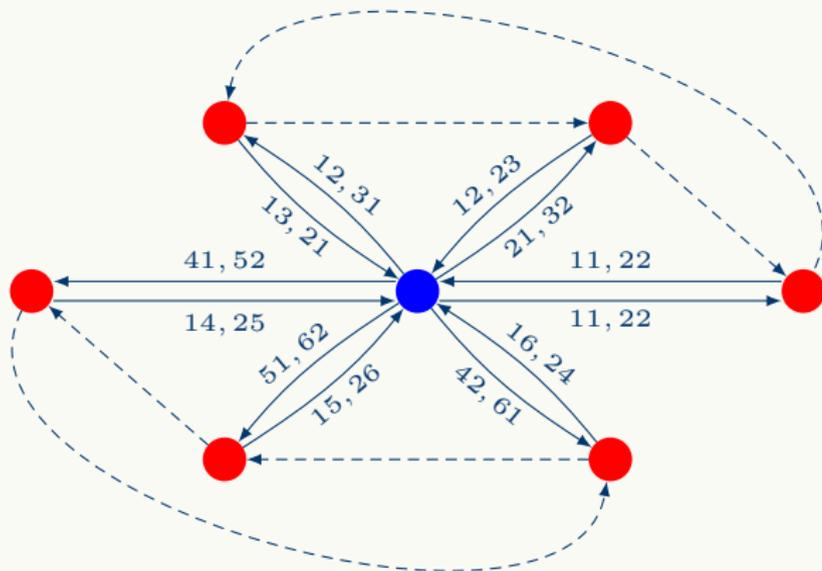
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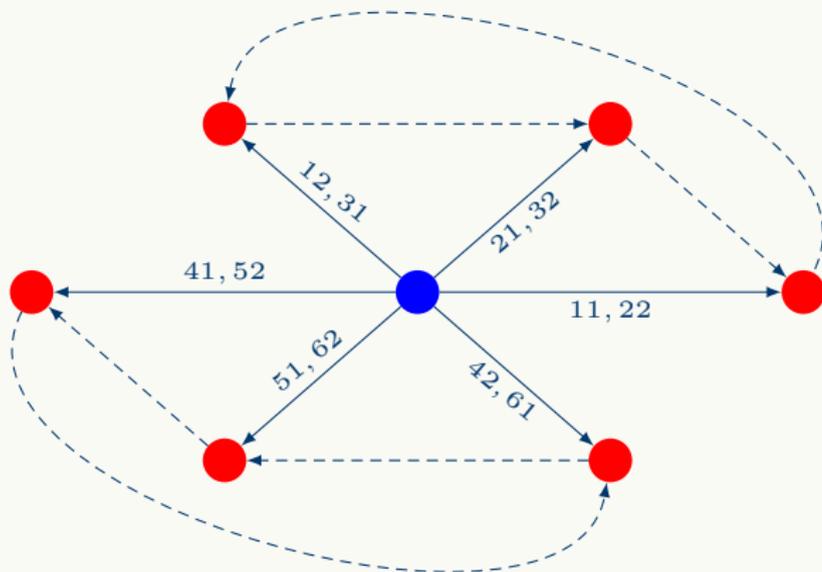
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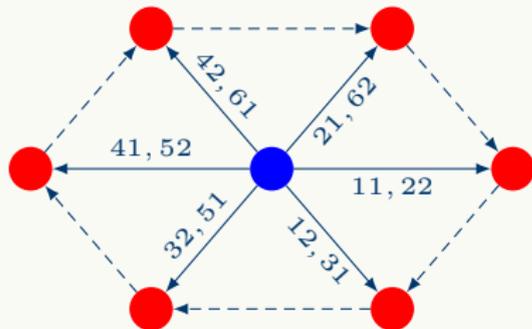
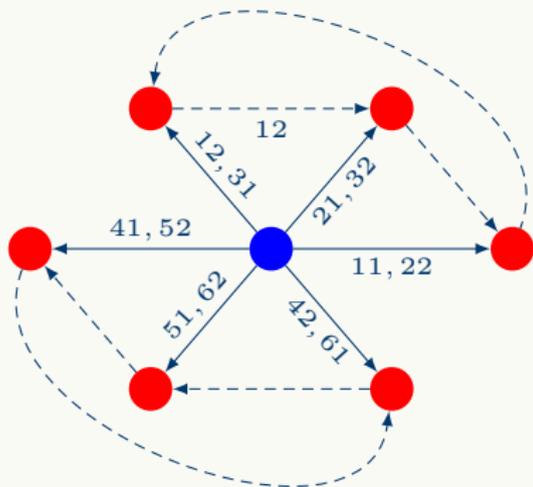
Representing a Structure as a Graph — $G_{\mathcal{A}}$



Representing a Structure as a Graph — $G_{\mathcal{A}}$



RCR on \mathcal{A}, \mathcal{B} is equivalent to CR on $G_{\mathcal{A}}, G_{\mathcal{B}}$



This might get large

- RCR can be implemented to run in time

$$\mathcal{O}(\|G_{\mathcal{A}}\| \cdot \log \|G_{\mathcal{A}}\|).$$

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- So RCR runs in time $\mathcal{O}(\|\mathcal{A}\|^2 \cdot \log \|\mathcal{A}\|)$ in the worst case?

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 - $\mathcal{O}(\|\mathcal{A}\|^2)$ edges in the worst case.
- So RCR runs in time $\mathcal{O}(\|\mathcal{A}\|^2 \cdot \log \|\mathcal{A}\|)$ in the worst case?
- Didn't I promise $\mathcal{O}(\|\mathcal{A}\| \cdot \log \|\mathcal{A}\|)$?!

Runtime of Relational Color Refinement

Use a different encoding $H_{\mathcal{A}}$ of \mathcal{A} instead of $G_{\mathcal{A}}$ where

$$\|H_{\mathcal{A}}\| \in \mathcal{O}(\|\mathcal{A}\|).$$

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- The number of cliques a tuple can be involved in depends on its arity, which is a constant.
- The proof that CR on $H_{\mathcal{A}}$ is equivalent to RCR on \mathcal{A} is quite involved.

Isn't this way too complicated?!

Let us look at “straightforward” approaches.

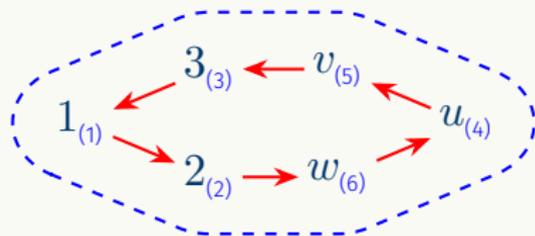
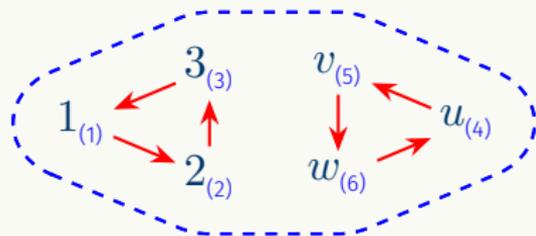
Approach 1:

Run Color Refinement on the Gaifman graph.

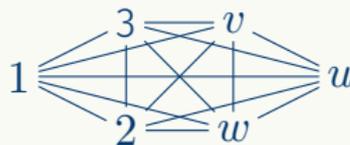
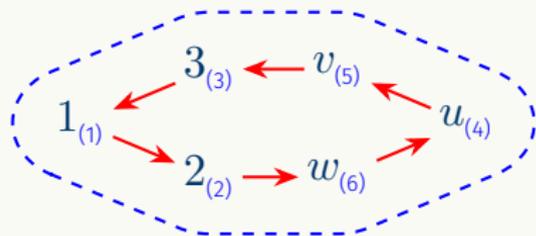
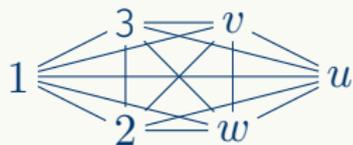
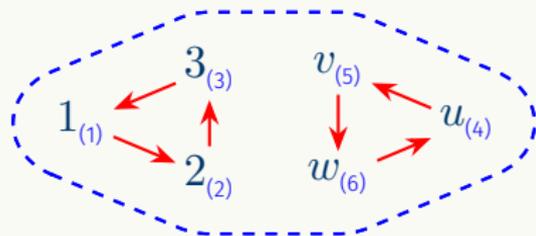
Approach 2:

Run Color Refinement on the incidence graph.

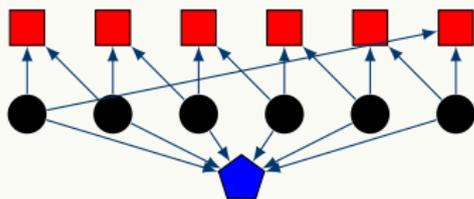
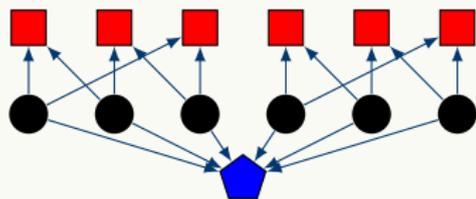
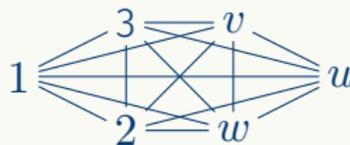
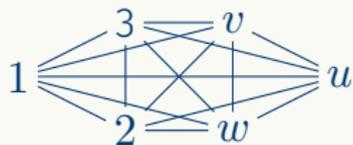
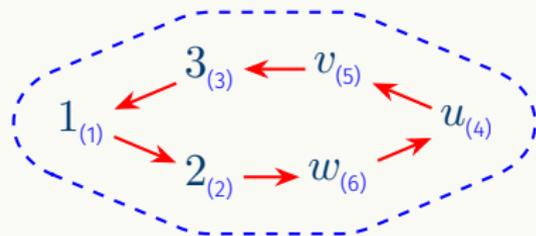
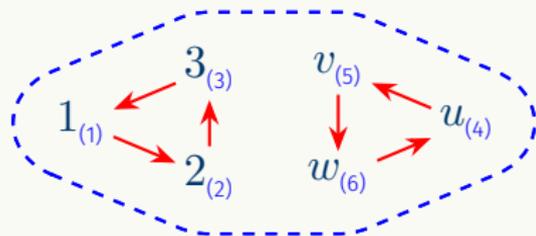
Consider our Examples...



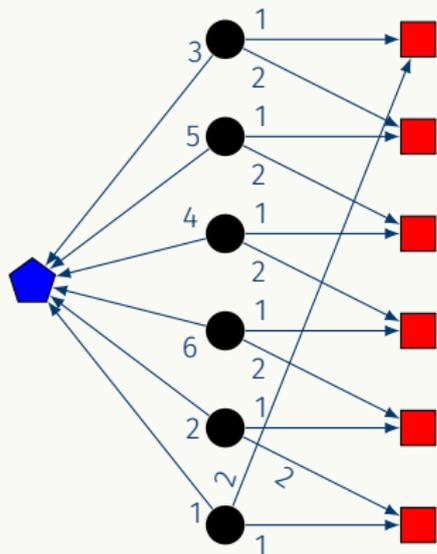
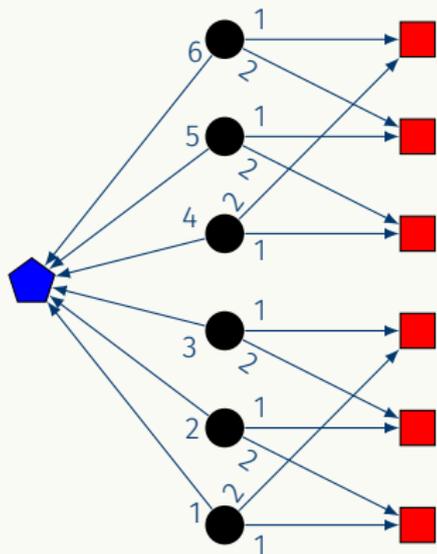
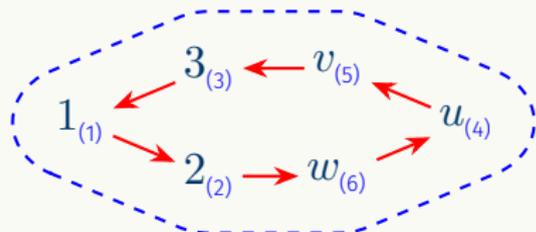
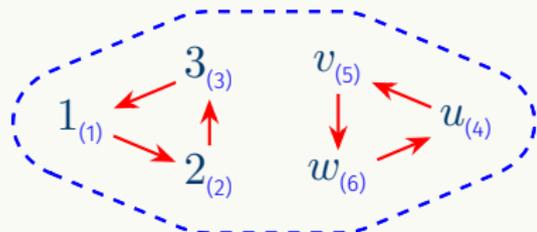
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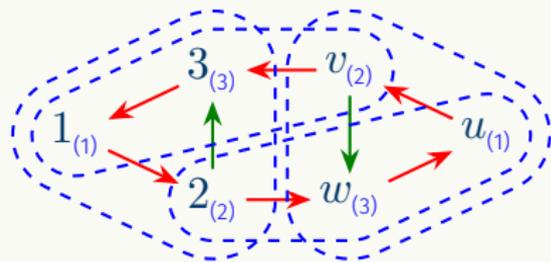
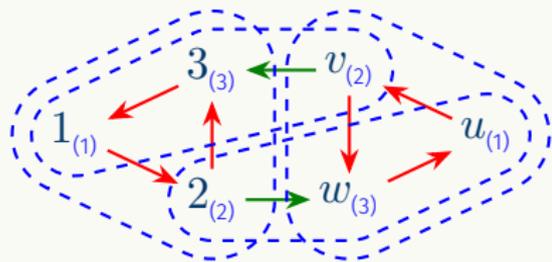
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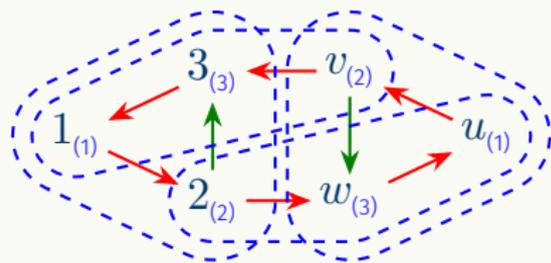
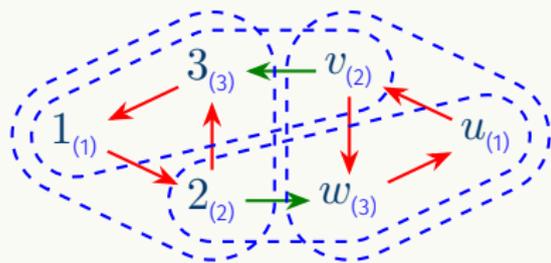
Press more Information into the Representations



Another Counter-Example

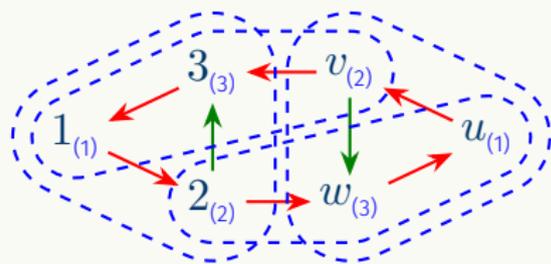
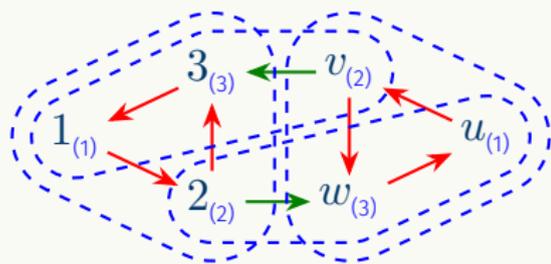


Another Counter-Example



CR does not distinguish their incidence graphs, even with labeled edges.

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$$\exists^{\geq 1}(x_1, x_2, x_3) \cdot (R(x_1, x_2, x_3) \wedge E(x_1, x_2) \wedge F(x_2, x_3) \\ \wedge E(x_3, x_1))$$

Bottom line:

These representations do not seem to capture homomorphisms from acyclic structures.

They somehow lose crucial information.

Main Result & Outlook

Theorem (Scheidt, Schweikardt 2024)

The following are equivalent:

- *Relational Color Refinement distinguishes \mathcal{A} and \mathcal{B} .*
- *There is an acyclic \mathcal{C} s.t. $\text{hom}(\mathcal{C}, \mathcal{A}) \neq \text{hom}(\mathcal{C}, \mathcal{B})$.*
- *There is a $\varphi \in \text{GF}(\mathbb{C})$ s.t. $\mathcal{A} \models \varphi$, and $\mathcal{B} \not\models \varphi$.*
- *Spoiler has a win. strategy for the Guarded-Game on \mathcal{A}, \mathcal{B} .*

RCR can be implemented to run in time $\mathcal{O}(\|\mathcal{A}\| \cdot \log \|\mathcal{A}\|)$.

Outlook: lift results on graphs to relational structures?

- CR can be used to speed up the evaluation of fc-ACQs on binary schemas (Riveros, Scheidt, Schweikardt 2024).
- CR is related to Graph Neural Networks.
- k-dimensional Relational Weisfeiler-Leman?

Further Reading



Benjamin Scheidt, Nicole Schweikardt.

“Color Refinement for Relational Structures”.

2024. DOI: [10.48550/ARXIV.2407.16022](https://doi.org/10.48550/ARXIV.2407.16022).



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