

Inria

Decentralized Workflow Execution Through Molecular Composition

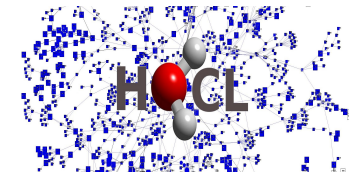
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ICSOC. Paphos, December 5, 2011



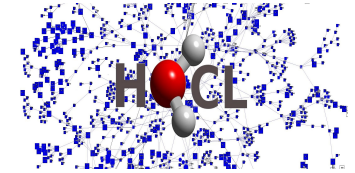
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Autonomic Service Computing?

- The *Service* abstraction
 - Autonomous
 - Network-enabled
- Service-Oriented applications
 - Built as temporal compositions of services (workflows)
 - Loosely-Coupled, dynamic composition
- Need for autonomic execution
 - Self-adaptation, self-healing, self-organization...
 - → **Decentralized Workflow execution**
 - → **Self-coordination**



Service Coordination: background

- Workflow engines
 - One coordinator
 - Managing all data and control flows
 - Scalability?
 - Reliability?
- Workflow execution languages
 - Data / Control driven
 - Implicit / Explicit parallelism
 - BPEL, XPDL, SCUFL, MoML, DAX
 - Designed for centralized, static coordination

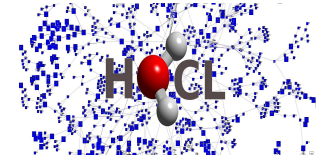


BPEL

Business Process Execution Language



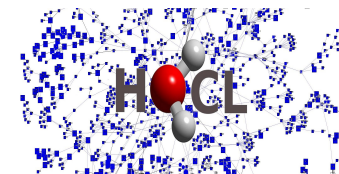
Towards autonomous coordination



- High-level abstractions...
 - Self-adaptation
 - Self-healing
 - Self-...
 - **Self-coordination**
- ...To help designing autonomic service-oriented application

“Nature-inspired metaphors have been shown to be of high interest for service coordination.” [Viroli et al., 2009]

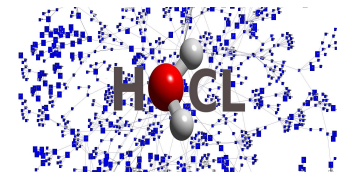
- → **Workflow execution as a succession of chemical reaction**



Chemical Programming Model

- An execution model
 - A program's runtime envisioned as a succession of chemical reactions
 - Reactions taking place (implicitly)
 - In parallel, autonomously, non-deterministically
 - Until no more reactions can take place (state of *inertia*)
- Writing a chemical program
 - Data: define the solution as a multiset
 - Program: define a set of rules rewriting the multiset
- Languages
 - GAMMA [Banâtre *et al.*, 1990]
 - HOCL (Higher-Order Chemical Languages) [Radenac, 2007]
 - Rules can apply on rules → program is able to change dynamically

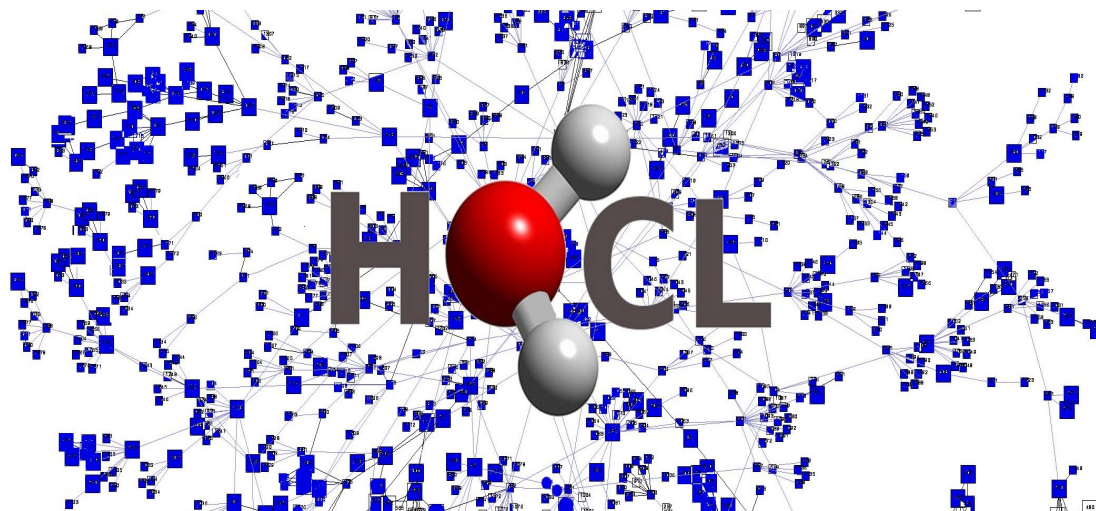
A Basic Chemical Program



- A specification
- A possible execution (amongst others)

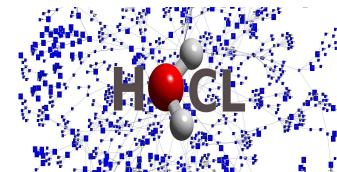
replace x, y by x if $x \geq y$ in $\langle 2, 3, 5, 8, 9 \rangle$

$\langle 2, 3, 5, 8, 9 \rangle \rightarrow^* \langle 3, 5, 9 \rangle \rightarrow^* \langle 9 \rangle$

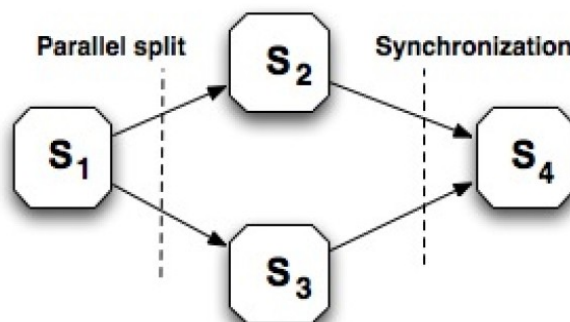


Workflow as a molecular composition

Defining a *molecular* workflow



- Express all data and control flow (reaction rules and molecules).



```
< // Multiset (Solution)
```

```
WS1:<CALL:S1, PARAM:<in1>,DEST:WS2, DEST:WS3>, // WS1 Sub-solution
```

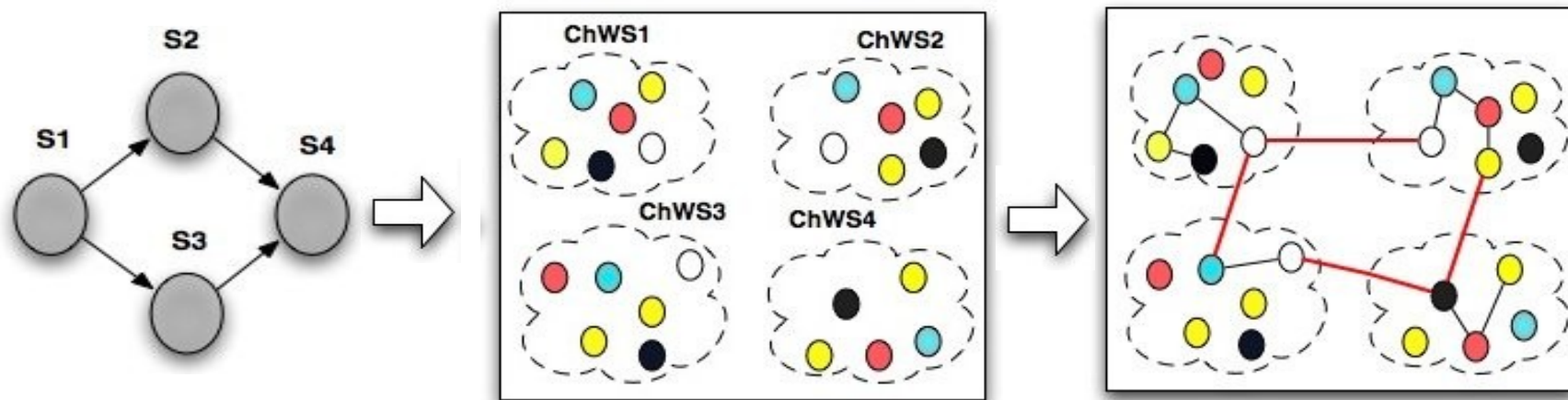
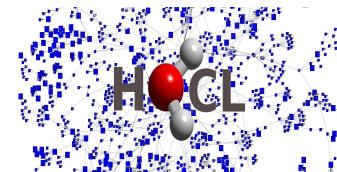
```
WS2:<DEST:WS4, replace RESULT:WS1:value1 by CALL:S2, PARAM:<(value1)> > ,
```

```
WS3:<DEST:WS4, replace RESULT:WS1:value1 by CALL:S3, PARAM:<(value1)> > ,
```

```
WS4:<replace RESULT:WS2:value2, RESULT:WS3:value3 by CALL:S4, PARAM:<(value2)> >
```

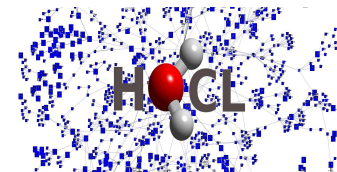
```
>
```

Molecular workflow: design & runtime



- At runtime
 - Interactions between services modeled by reactions
 - Molecules react in chain following the patterns of the workflow
- At design time
 - Workflow defined using specific reaction rules and molecules (**multiset**)
 - Each Service is a sub-solution with specific molecules
 - Each Service is provided with a library of generic rules

Chemical rules for distributed execution

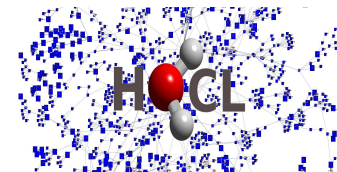


- Independent from any workflow
- Chemical rules to ensure the execution

```
let invokeServ = replace ChWSi:⟨CALL:Si, PARAM:⟨in1,...,inn⟩, FLAG_INVOKE:1, ω⟩,  
by ChWSi:⟨RESULT:ChWSi:⟨value⟩, ω⟩  
let preparePass = replace ChWSi:⟨RESULT:ChWSi:⟨value⟩, DEST:ChWSj, ω⟩  
by ChWSi:⟨PASS:ChWSj:⟨COMPLETED:ChWSi:⟨value⟩⟩⟩
```

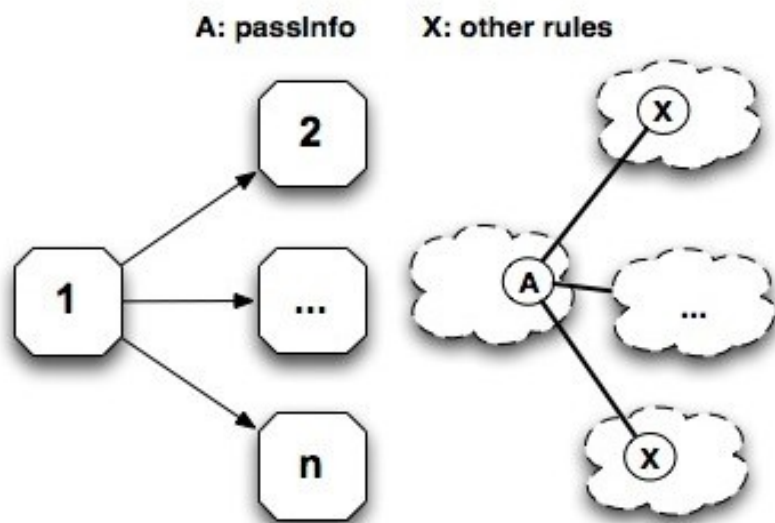
- Control and data transfer among ChWSes.

```
let passInfo = replace ChWSj:⟨PASS:ChWSi:⟨ω1⟩, ω2⟩, ChWSi:⟨ω3⟩  
by ChWSj:⟨ω2⟩, ChWSi:⟨ω1, ω3⟩
```

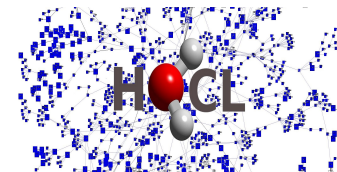


Workflow patterns: Parallel Split

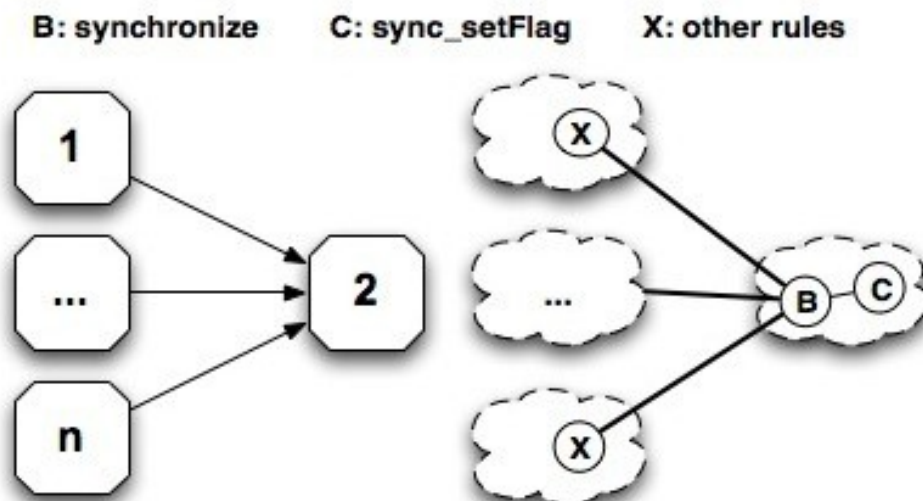
```
let passInfo = replace ChWSj:⟨PASS:ChWSi:⟨ $\omega_1$ ⟩,  $\omega_2$ ⟩, ChWSi:⟨ $\omega_3$ ⟩  
by ChWSj:⟨ $\omega_2$ ⟩, ChWSi:⟨ $\omega_1, \omega_3$ ⟩
```



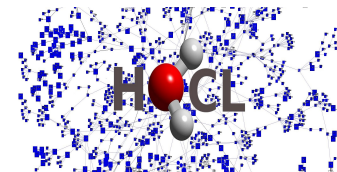
Sorkflow pattern: Synchronization



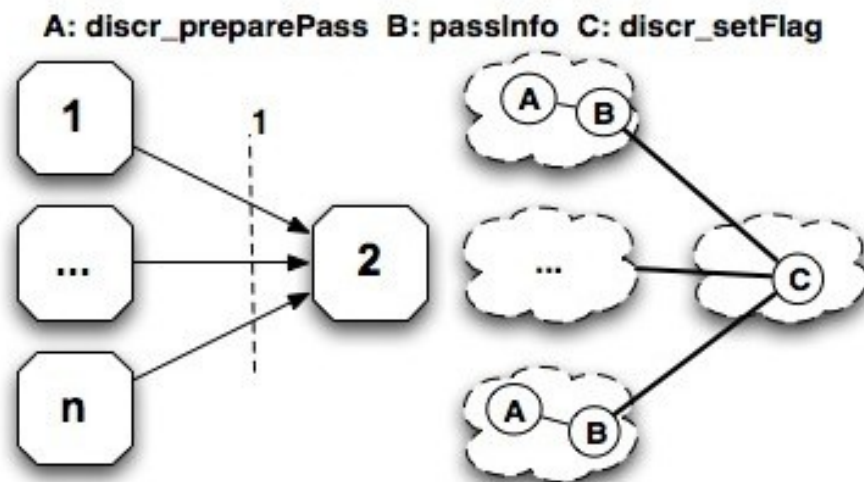
```
let synchronize = replace SYNC_SRC:<ChWSi,  $\omega_1$  >, COMPLETED:ChWSi:<value>,
                    SYNC_INBOX:<  $\omega_2$  >
                    by SYNC_INBOX:< COMPLETED:ChWSi:<value>,  $\omega_2$  >, SYNC_SRC:<  $\omega_1$  >
let sync_setFlag = replace-one SYNC_SRC:< > by FLAG_INVOKE:1
```

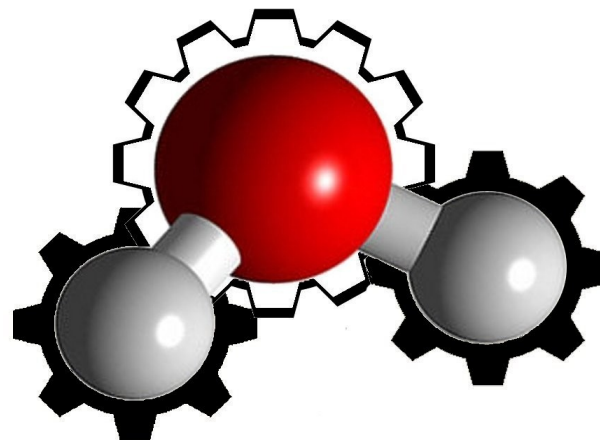


Workflow pattern: Discriminator



```
let discr_preparePass = replace DEST:ChWSj, RESULT:ChWSi:<value>  
    by PASS:ChWSj:<COMPLETED:ChWSi:<value>, DISCRIMINATOR:Yes>  
let discr_setFlag = replace-one DISCRIMINATOR:Yes by FLAG_INVOKE:1
```



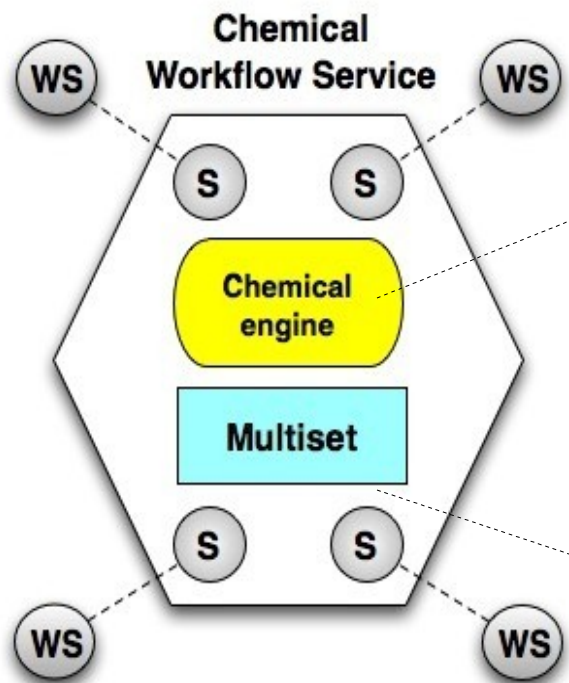


Architectures

- Coordination mechanism built upon HOCL
- Two architectures, two software prototypes
 - Centralized
 - Decentralized

Centralized Architecture

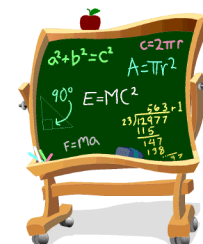
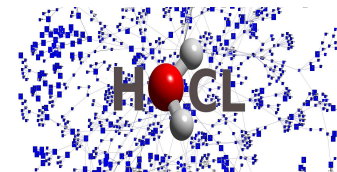
- Central node coordinating all data and control flows
 - A chemical encapsulation per Web service participating in the workflow
 - **Multiset** containing the workflow definition
 - **Chemical engine** processing the content of the multiset



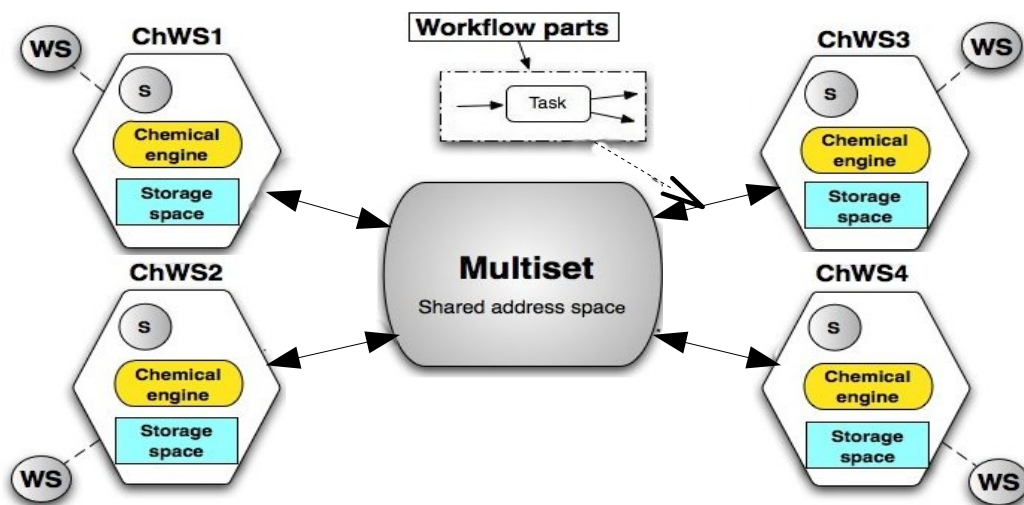
```
let invokeServ = replace ChWSi:⟨CALL:Si, PARAM:⟨in1, ..., inn⟩, FLAG_INVOKE, ω ⟩,
    by ChWSi:⟨RESULT:ChWSi:⟨value⟩, ω ⟩
let sm_preparePass = replace DEST:WSj, RESULT:WSi:⟨value⟩
    by PASS:WSj:⟨RESULT:WSi:⟨value⟩, MERGE:Yes⟩
let sm_setFlag = replace-one MERGE:Yes by FLAG_INVOKE
```

```
< // Multiset (Solution)
WS1:⟨CALL:S1, PARAM:⟨in1⟩,DEST:WS2, DEST:WS3⟩, // WS1 Sub-solution
WS2:⟨DEST:WS4, replace RESULT:WS1:value1 by CALL:S2, PARAM:⟨(value1)⟩ ⟩,
WS3:⟨DEST:WS4, replace RESULT:WS1:value1 by CALL:S3, PARAM:⟨(value1)⟩ ⟩,
WS4:⟨replace RESULT:WS2:value2, RESULT:WS3:value3 by CALL:S4, PARAM:⟨(value2)⟩ ⟩
>
```

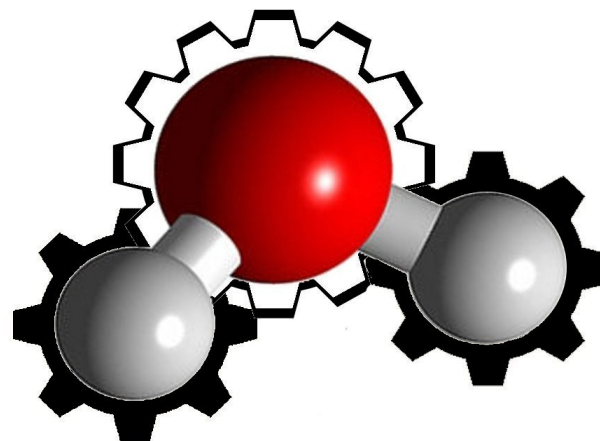
Decentralized Architecture



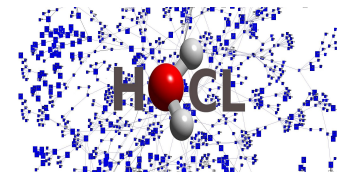
- The multiset becomes a reliable shared address space
 - For nodes to communicate together
 - Reading and writing the multiset for coordination
- Workflow executed in parts by each *Chemical Web Service*
 - Data and control transfer through this shared space
 - Each node co-responsible of the execution



Experimental Evaluation

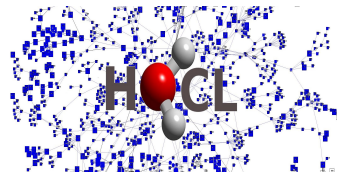


Experiments

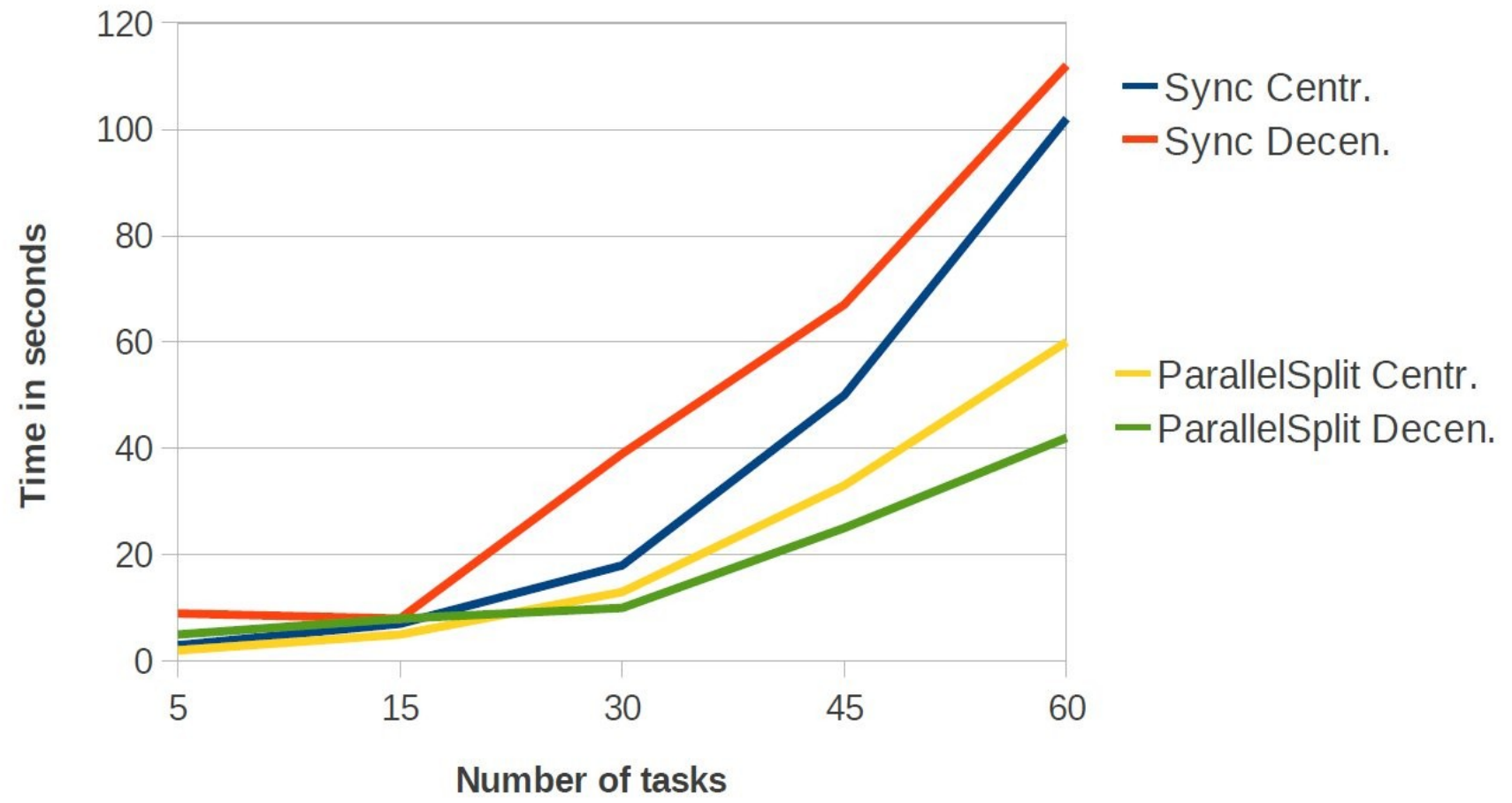


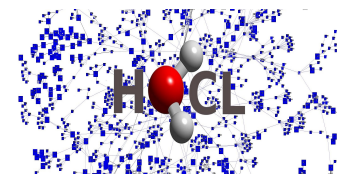
- Both prototypes experimented
- Patterns tested
 - Parallel split
 - Discriminator
 - Synchronization merge
- Size of patterns 5, 15, 30, 45, 60 nodes
- Experiments conducted on the Grid'5000 research infrastructure



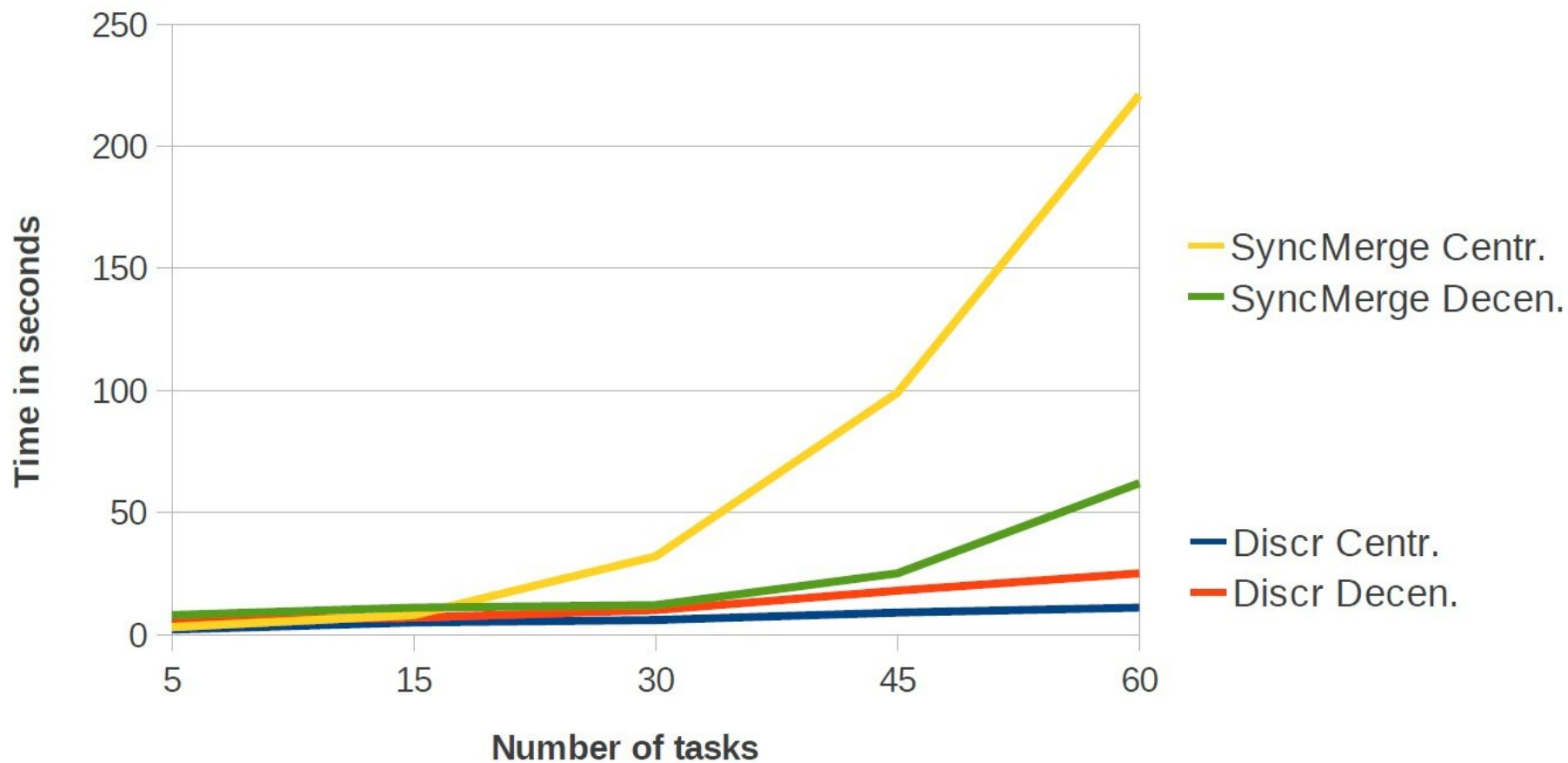


Results (I)

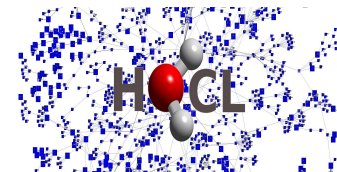




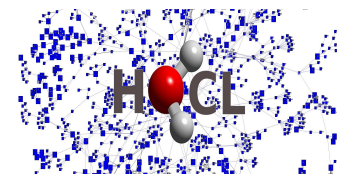
Results (II)



Conclusion & On-going works



- New chemically-inspired abstractions for workflow self-coordination
 - Molecular composition
 - Autonomic behavior
 - Decentralization / Self-coordination
- Proof of concepts
 - Two architectures and software prototypes (centralized, decentralized)
 - highlighting the feasibility of such a concept
- On-going studies
 - Towards a P2P multiset
 - New applications envisioned:
 - Workflow scheduling
 - Broader scope – Social networking, Agile Programming Networks



Thank you!

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