



eRA09: Kevin Buckley: A Grid-Based Facility for Large-Scale Cross-Correlation of Continuous Seismic Data
Outside the boundary: Workflows

Workflows provide a mechanism for knowledge transfer about processes within an automated representation of the process as a whole.

Two seemingly opposite drivers at work:

Hiding the raw computation in the sub-processes from the end user, eg

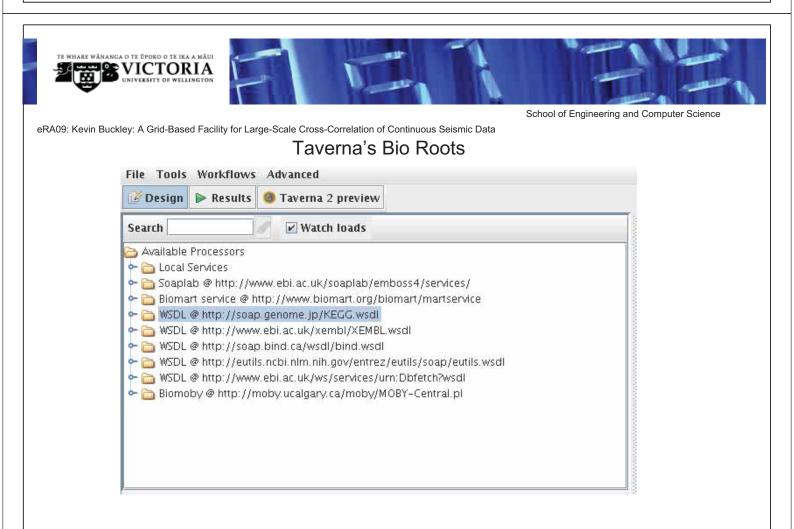
■ Interact with a specified web service via a WSDL scraper

Making the sub-processes explicit to the end user

- End user might specify a different web service
- End user might not use a web service at all

Workflow tool-kits come with a lot of pre-defined process units for "doing things"

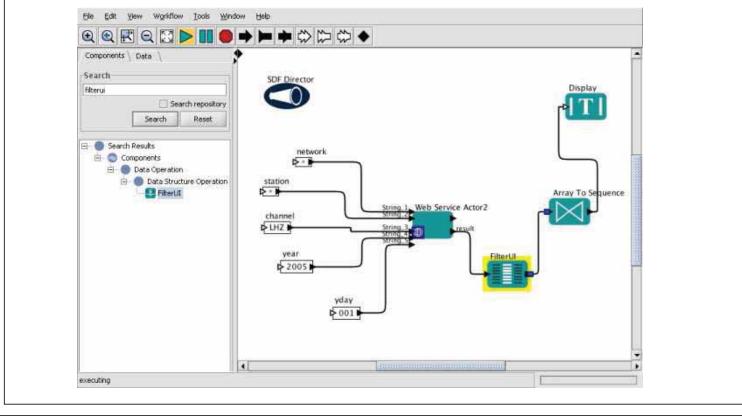
Does your domain science, let alone your sub-processes, have units in the workflow tool-kit you want to use, that let you do what you want?





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Kepler WSDL





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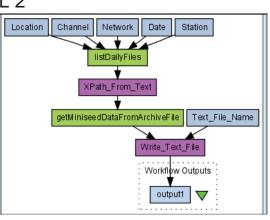
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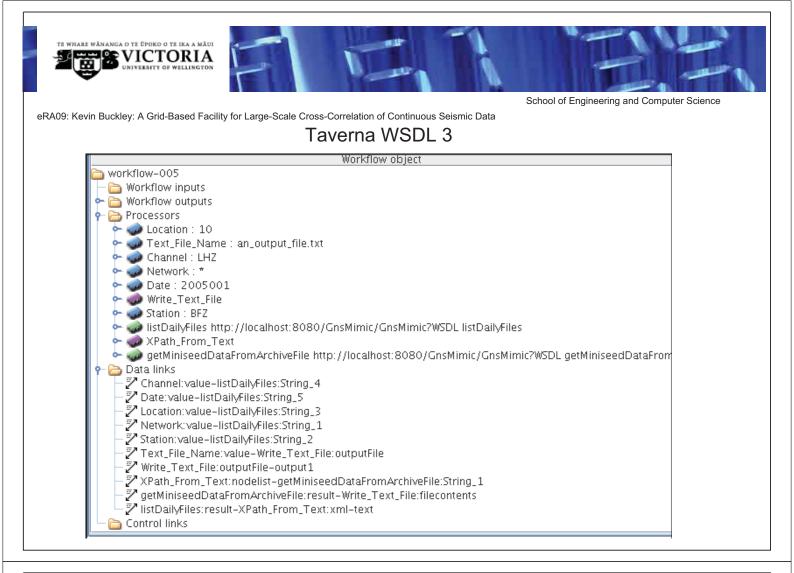
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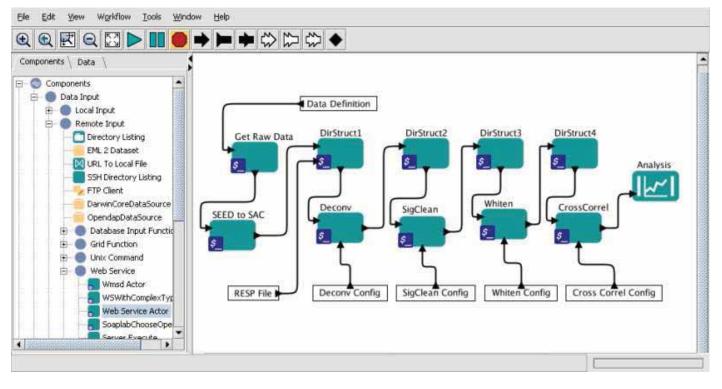




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eRA09: Kevin Buckley: A Grid-Based Facility for Large-Scale Cross-Correlation of Continuous Seismic Data Removing the boundary: Workflows

Once raw data is converted, stage inputs and outputs are the same format - simple pipeline ?

Compare that workflow representation, itself a mere schematic, with the schematics of what the domain scientists think of as what's happening

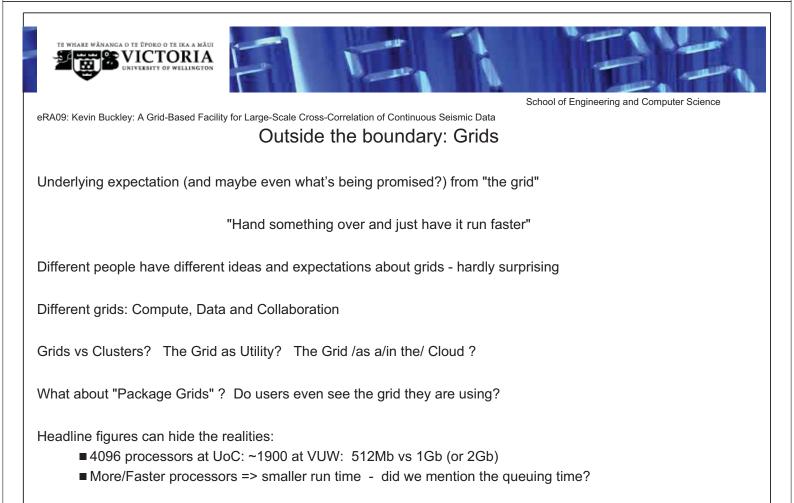
A lot more stuff (local/domain knowledge) going on, eg

- overcoming directory requirement impedences between the components
- handling errors at various stages

and that "stuff" becomes visible in the workflow tools.

Tempting to say "keep things hidden" - but it's not a package, it's a workflow

Other benefits to being explict





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Now remove the boundary

#### Our grids

- VUW actually has two totally seperate grids
- Neither are dedicated resources: cycle-stealing grids
- So neither really allow for message-passing approaches

Our science: where were the boundaries for us now?

- Data constant access to raw data vs storing raw/modified data
- Data three distinct processes: only need the whole at the end
- Compute discretisation gives us resilience to failure
- Collaboration not all users happy at the command line



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The DQS/SGE aware will notice there's an implied step size (of 1) there

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#### Beyond the Boundary

One job was run whilst Yannik was back in Germany - no boundaries indeed !

Reductions of up to 87% (91% and 51%, stage 1 and 2 resp.) in processing times were seen (5 stations over 99 days, 1Hz sampling data, 170Mb of disk)

(Time :	in mins)	Desktop	Grid
Stage1	time	34.2	2.9
Stage2	time	5.2	2.2
Total	time	39.4	5.1

Managed to get a concurrent usage of 180 (out of ~230) otherwise idle processors (10 stations over 360 days, 100Hz sampling data, 114Gb of disk)

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(Time in hrs)	Desktop	Grid
Stage1 time	~120	~3



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## Summary: Seismic Processing

- Moved a desktop process onto the "grid"
- Built in resilience and flexibility across generic grid resources
- Maximised use of local grid resources
- Hidden "the grid" from the user
- Questions raised informed future domain science approaches

As to the future:

- Investigate other approaches upon dedicated grid resources
- Actual implementation of processes in a workflow tool



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### Summary: Here be boundaries Domain Science

- Processes operating in a simple, understandable, pipeline
- Become wrapped, for operational simplicity, into one process

## Workflow Boundary

- Individual processes become hidden (package or workflow?)
- Additional operational artefacts not externally visible

## Grid Computing Boundary

- No visible loops to parallelise
- Data placement/distribution is hard

## Avoiding Boundaries

- Clearly identifiable processes with simple (maybe NO) interfaces between them
- Explicit loops over monolithic processing



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### And finally ... who is supposed to do this "no-boundary" science ?

#### Domain scientists ?

■ Might need to brush up on their BEPL, SCUFL, WSDL, MPI and DRMAA

Workflow/Grid computer scientists ?

■ Might need to read up on SEED, SAC, QuakeML

# Phew !

Might still be a few jobs for people who can straddle the boundaries

