

Facility

SNS

Instrument

ARCS

Datafiles: Files are selected. See datafiles tab

Workflow algorithms: default (modified)

Inputs: still need some inputs

The other option is "all OK"

The other option is "No files selected"

Options:

"default" Algorithms and parameters from default workflow

"default (modified)" Algorithms and/or parameters changed, but not saved

"from myworkflow.xml" Loaded from a file

"from myworkflow.xml (modified)" Started from a workflow file, with modifications, not saved

Manage User Directories

Save workflow

Load workflow

Generate Script to clipboard

Execute

Help :)

We should provide some standard workflows (powder reduction with Vanadium normalization, single crystal with/without normalization, ...). When instrument is selected, automatically load a standard workflow, that the instrument scientist can change

Real Help -
not link to
webpage

Load Workflow



workflow1.xml
workflow2.xml
powderworkflow.xml

- Load
- ☒ Everything
 - ☐ Data
 - ☐ Algorithms
 - ☐ Algorithms with parameters

Load

Cancel

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Cannot change facility/ instrument on this tab

Datafiles

Run1,Run2, Run3;Run4;Run5,Run6

Browse

WBVanadium

WB Run1, WBRun2

Browse

Mono Van

MVRun1,MVRun2

Browse

Files separated by commas or dashes are added together, files separated by semicolon are processed separately

Manage User Directories

Save workflow

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Execute

Help :)

Datafile(s) should be the only mandatory thing in this whole GUI

We need to define format for add together, say "," and "-", and separately ";" maybe something like "I" for range in which all files are separated

Facility

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- ☒ GetEi
- ☒ Rebin in Energy
- ☐ Filter bad pulses
- ☐ Filter by log value
- ☐ Time independent background subtraction
- ☒ He3Tube Efficiency
- ☒ Ki/Kf correction
- ☐ Detailed balance correction
- ☒ Normalize by current
- ☒ Normalize detector efficiency to white beam Vanadium
- ☒ Load mask from file
- ☐ Solid angle correction
- ☐ Group detectors for powder
- ☐ Absolute normalization using monochromatic Vanadium

The list of algorithms should be instrument specific. (For example no He3 tube efficiency if you use scintillators)

As different algorithms are selected, inputs tab is populated

Manage User Directories

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Help :)

Facility

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GetEi

Incident Energy Guess (meV)

100

☐ Fixe Ei to guess

Rebin in Energy *

White Beam Vanadium

He3TubeEfficiency *

Normalize by Current

Load Mask File *

Goniometer *

Tabs with *
need
additional
input

Add T0 for
CNCS

Note that Algorithms that need no
additional input are not shown on this tab,
even if they are selected (Ki/Kf
correction)

Note that Goniometer tab appeared,
even if there is no algorithm, but
needed for NXSPE single crystal

Should be able to put a first file preview for specific
algorithms (i.e. time independent background should
be able to preview on request sum over all space)

Manage User Directories

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Execute

Help :)

Facility

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Instrument

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GetEi

Rebin in Energy *

White Beam Vanadium

He3TubeEfficiency *

Normalize by Current *

Load Mask File *

Goniometer *

TOF range: Min

2000

Max

9000

☒ Add files together☐ Process separately and compare masks☒ Use/save file

Filename

vanadium.nxs

Browse

Detector diagnostics

Whatever is needed here

Manage User Directories

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Execute

Help :)

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- ☐ SPE
- ☐ PHX
- ☐ PAR
- ☒ NXSPE
- ☒ Mantid Nexus

☒ Single Crystal

If NXSPE is not selected disable Single Crystal in this tab, and if Single Crystal is selected add goniometer in the inputs tab

[Change from default filenames](#)

mat(s) NXSPE, NXS
es: Combinedruns1-3.nxspe, Run4.nxspe, Combinedruns5-
es: Combinedruns1-3.nxs, Run4.nxs, Combinedruns5-6.nxs
CR Axis (0,1,0)
equation: value+23.4

[Manage User Directories](#)[Save workflow](#)[Load workflow](#)[Generate Script to clipboard](#)[Execute](#)[Help :\)](#)

Facility

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Instrument

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Datafiles: Data: Run 1, Run2, Run3
Data: Run 4
Data: Run 5, Run 6
Vanadium: WB Run 1, WB Run2
MonoVan: MV Run1, MVRun2

Algorithms: Calculate Ei
He3Tube efficiency
Correct Ki/Kf
Normalize by current
Normalize by WB vanadium
Add mask from file

Inputs: Efixed=100meV (to be calculated)
Energy rebin parameters: -20,1,100
WB vanadium integration range: 2000,9000
He3 Tube parameters: Scale: Default, Temperature: Default,

Save: File format(s) NXSPE, NXS
File names: Combinedruns1-3.nxspe, Run4.nxspe, Combinedruns5-6.nxspe
File names: Combinedruns1-3.nxs, Run4.nxs, Combinedruns5-6.nxs
Store orientation:
Motor CCR Axis (0,1,0)
equation: value+23.4

Manage User Directories

Save workflow

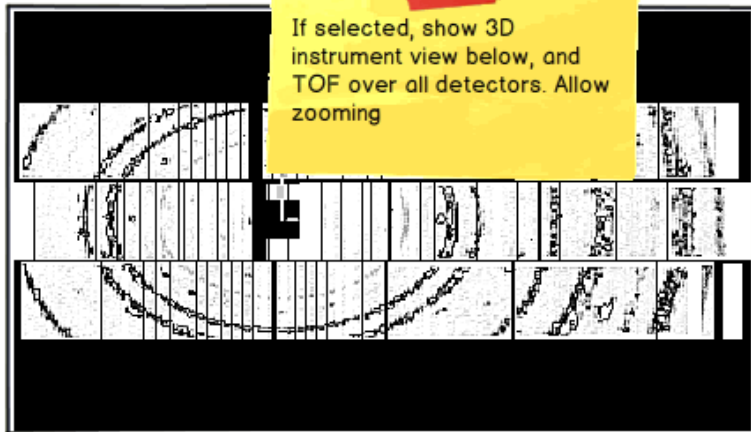
Load workflow

Generate Script to clipboard

Execute

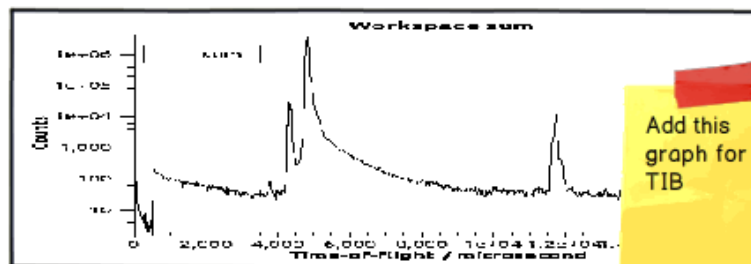
Help :)

This is just a
summary

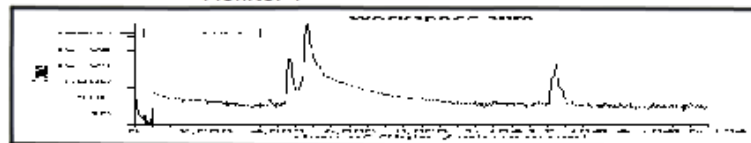
Facility Instrument ☒ Preview first file

Tab is instrument specific, instrument scientist should be able to modify it

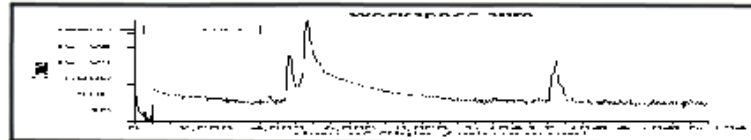
Sum over all detectors



Monitor 1



Monitor 2



Manage User Directories

Save workflow

Load workflow

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Execute

Help :)