Some quick commands from the MATLAB command line

• nplot [filenames]

Plot one or several (if can be combined) scans

Example: nplot 0123[45, 46, 50:53] (Combine and plot 012345, 012346, 012350-012353)

nplot 012345 fit gauss1 (Plot, and fit one Gaussian peak)

nplot 012345 calc pal1+pal2-2*pal3 (Plot linear combination of polarization states)

(For a list of available options, type help nplot, and see the nplot manual)

plotmultiple [filenames]

Plot one or several FLATCONE scans

Note: nplot and plotmultiple use local files if found, and automatically try to download from the instrument computer otherwise.

• transfer [filenames]

Copy one or more files from spectrometer to current local directory

Example: transfer 0555[10:20] (Copy files 055510 – 055520) transfer 055520:end (Copy all files from 055520 till most recent one)

printscanlist [filenames]

Print a list of scans

Example: printscanlist all (list all data files in current directory)

printscanlist 0555[10:20] var kf (show also the value of kf for each scan)

datafile ['find'|'show'] [filename]

Search datafiles on the instrument computer (works also from the linux command line on the instrument computer)

• powxbu [...]

Create a list of powder reflections for a2/a4 alignment

Example: powxbu (opens an interactive window) powxbu create silicon 2.662 (calculate silicon bragg peaks for ki=2.662) powxbu writexbu powder.xbu 2:5 (write xbu-file with reflections No. 2,3,4,5)

powcal [...]

Read scan data from a4 powder scans and fit a2 and a4 zeros

Example: powcal (opens an interactive window)
powcal readfiles 0123[45:56] (read and fit data files 012345-012356)
powcal fitzeros (determine za2, za4 from fitted data)

For all commands: Type help [command] for more information and examples.

Type edit options to change a number of options. (You should copy the options.m file to your current own directory before editing.)