

Commands

Setup

- A window to create an new experiment **edc /new**
- To increase the number of experiments and to copy over the parameters from the experiment before **iexpno**
- To look at the library **list_pp**
- Copy parameters and to which directory and number they will go (mf) **wrpa^{expno}**
- To move over to another exp.no (jexp) **re^{expno}**

Temp

- To open a temp window **edte**
- Temperature in K **te**
- Change probe temperature to **te** **teset**
- To measure the temperature of the shims **coiltemp**
- To measure the actual temperature (reference samples MeOH or EtylenGlycol) **calctemp**

Prepare for acquisition

- To check/change nuclei (F1 active in tuning) **edasp**
- To open a probe window **edhead**
- To tune and match automatically **atma**
- To tune and match manually **atmm**
- To tune and match (31P-probe) **wobb (wobb-sw)**
- **acqu**
- **stop**
- To choose solvent and auto lock **lock**
- To open a lock window **lockdisp**
- To store shimvalues **wsh**
- To read old shimsfiles **rsh**
- To wiew a shimfile **vish**
- To shim automatically **topshim / tshim, topshim^{gui}**
- To get solvent and probe specific values from prosol **getprosol**
- Update pulses related to measured p1 **getprosol 1H p1 p11**
- Choose experiment (copy all) **rpar**
- To see the current exp. pulsprogram graficly **showpp**
- To check or change number of scans **ns**
- To check or change sweepwitdt **sw**
- To check or change transmitter offset **o1p**
- Experiment time consuming **expt**
- Programme to calibrate d1, o1p etc. **gs**

Acquisition

- To rga and start an exp **xaua**
- To let the machine count the appropriate receiver gain **rga**
- Check the set receiver gain value **rg**
- Start an acquisition **zg**
- Go fouriertransform phase and linebroadning **zgefp**
- To restart and add ns to an already run exp **go**

- To put up a que of experiment (in the same folder) **multizg**
- To manually put experiments or commands to the spooler **qu[^]<x>**
- auto-spool (marked in Options)
- put zg, go, rga, atma to the que

- A au program that performs a single scan 90° pulse calibration based on nutation. **pulsecal**
- To calibrate the powerlevel for the ¹H 90° pulse for a special length of the pulse (e.g spin lock) **pulse**
- To array and the 90° pulse **paropt**

- Transmitted already run scans **tr, tr[^]number of scans**

- To save and stop the acquisition **halt, halt[^] number of scans**
- To stop the acquisition **stop**
- To stop the acquisition in an more rough way **kill**

Processing

- Opens a dialog box in which you can set all processing parameters **edp**
- To autoprocess **xaup**
- read row from 2D data and store as 1D **rser[^]number in array**
- read column from 2D data and store as 1D **rsc**

- Open a interactive window multiplication window **.winf**
- Linebrodning, to set values **lb**
- Makes linebroadning **em**
- gaussian broadening parameter used along with lb for gm **gb**
- gaussian muliplication **gm**
- Fourier transform **ft**
- Linebroadning and fouriertransform **ef**
- Automatic phasing **apk**
- Fouriertransform with chosen line broadning and already set phasing **efp**
- Fouriertransform with chosen gaussian function and already set phasing **gfp**
- To process an array in paropt **multiefp**
- Fourier transform 2D exp **xfb**
- Fourier transform 2D, T1/T2 exp **xf2**

• Basline correction (give you a window)	bas
• Basline correction 1D and auto integration	abs
• Basline correction first dimension 2D	abs1
• Basline correction second dimension 2D	abs2
• Basline correction without integration	abs n
• Symmetrization 2D e.g cosy	sym
• To measure the half width, (zoom in the desired peak) chose Calculate width of current peak.	pp
• To measure the half width	hwcal
• To measure the half width and the hump (expand the area around the peak, save displayed region, must be well-defined, calibrate it)	humpcal
• Find the internal zero-reference	sref
• Zero fill	si
• Calculate the signal to noise	sinocal
• Serial integration	intser
• To make a list	edlist
• Write miscellaneous lists	wmisc
• Read miscellaneous lists	rmisc
• Edit miscellaneous lists	emisc
<u>Various</u>	
• Setup	ii
• open BSMS display	bsmsdisp
• Autoprogram for selective exp	bnmr
• To make macron	edmac
• To edit au programs	edau
• To delete the processing on selected dataset	delp[^]dataset
• Edit shim files	edtune
• Edit solvent and probe specific parameters	edprosol
• To get info about the current exp	edcpul
• To get information about pulseprograms	edpul[^]*.info
	edpul[^]pulprog.info
	edpul[^]*noesy*