

Adosy_ste_BBFO Topspin 3.6.5 600MHz.

Freshly prepared D2O (0.8ml) with 2mg CuSO4.

Lit value for D2O is $1.87 \times 10^{-9} \text{m}^2/\text{s}$

Calibrated at 25C **gradpar** 5.350 G/mm.

This parameter set can be used as exercise or for calibrate the gradient strength (probe dependent).

=====

Check **gradpar** for proper Gradient calibration if needed.

1. Complete the optimization with **Adosy_ste_setup**.
2. Create a new data file #4.
 - **rpar Adosy_ste_BBFO; getprosol.**
 - Import all optimized values: **d1, p1, d20, p30**, gradient shape (SINE or RECT shape), **ns** and **rg**
3. Type **TD**
 - Define **TD(F1)**: number of experiment with different gradient, default is **18**.
 - More points to get higher resolution or for multi-peaks with varied decay rates.
4. Type **dosy** {If defaults are to be used, just press enter four times as follows}.
 - Enter first gradient amplitude: **5** (default)
 - Enter final gradient amplitude: **95** (default)
 - Enter # of points: **TD(F1)** default **18**
 - Ramp type: **"I"** note: **"I"** stands for "linear" option, press "enter".
5. If you want to do only one measurement: Press enter to accept "Do you want to start the measurement?" Click **OK** to start.

=====

Processing:

1. Type the next three commands:
 - Set **SI = TD(F1), xf2; abs2; setdiffparm.**

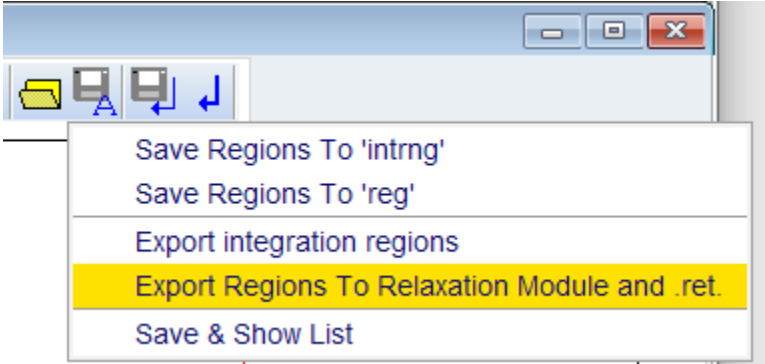
- **Process > Adjust Phase > Select one row (1, 2.....) by right click to phase manually; save and return**

2. Click **Analysis > Dynamics > T1/T2 Relaxation**

- Select **Fid > Spectrum > Slice Number = 1** (check phase correctly)
- Select **Peaks/Ranges > Manual Integration > Define new range** using cursor



> **Export Ranges to Relaxation Module and .ret**



- Select **Relaxation > Intensity** or **Area** (preferred)
- Select **Fitting > Function – lin > Relaxation parameters** – ok
- Select **Calculation**
- Select **Report**
- Select **Publish > Plot layout**