

## Brucine (2,3-Dimethoxystrychnidin-10-one)

Brucine (2,3-Dimethoxystrychnidin-10-one) is an alkaloid, structurally related to strychnine, but less toxic. Figure 1 shows the  $^1\text{H}$  NMR spectrum of a 250 mM Brucine sample in  $\text{CDCl}_3$  measured in a single scan taking 10 seconds to acquire.

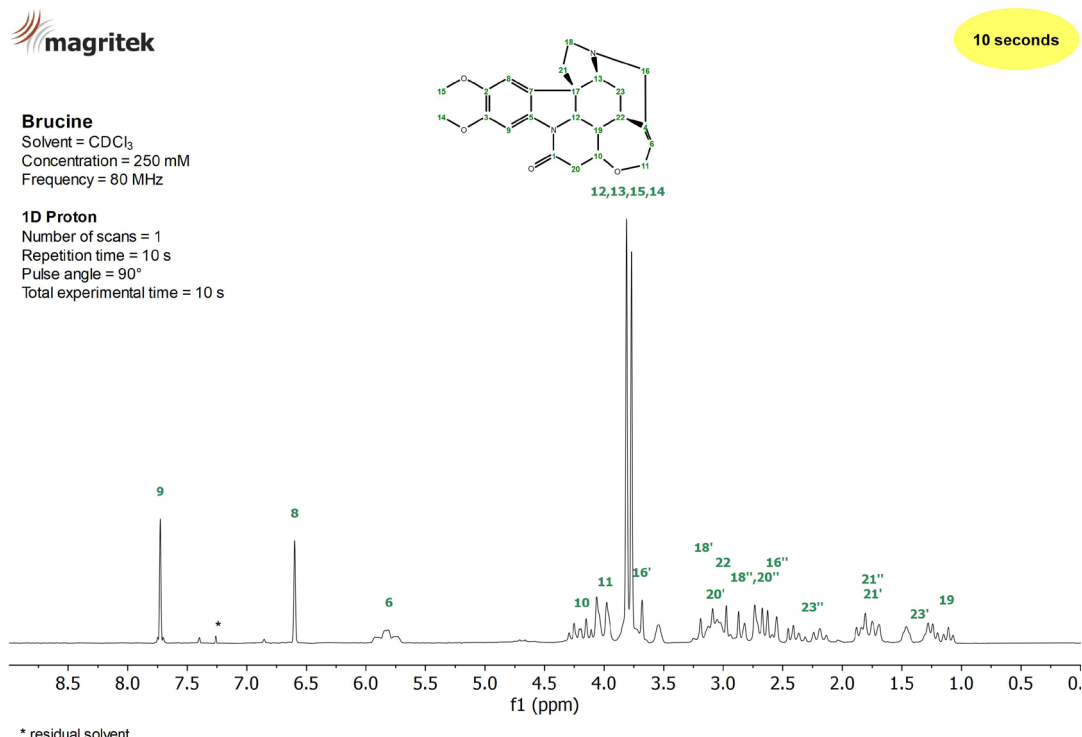


Figure 1:  $^1\text{H}$  NMR spectrum of a 250 mM Brucine sample in  $\text{CDCl}_3$  measured on a Spinsolve 80 MHz system in a single scan.

## $^{13}\text{C}$ Spectrum

Figure 3 shows the  $^{13}\text{C}$  NMR spectrum of 250 mM Brucine in  $\text{CDCl}_3$  acquired using NOE polarization transfer from  $^1\text{H}$  to  $^{13}\text{C}$  and  $^1\text{H}$  decoupling. The 1D Carbon experiment using NOE is sensitive to all  $^{13}\text{C}$  nuclei in the sample. It clearly resolves all the expected resonances.

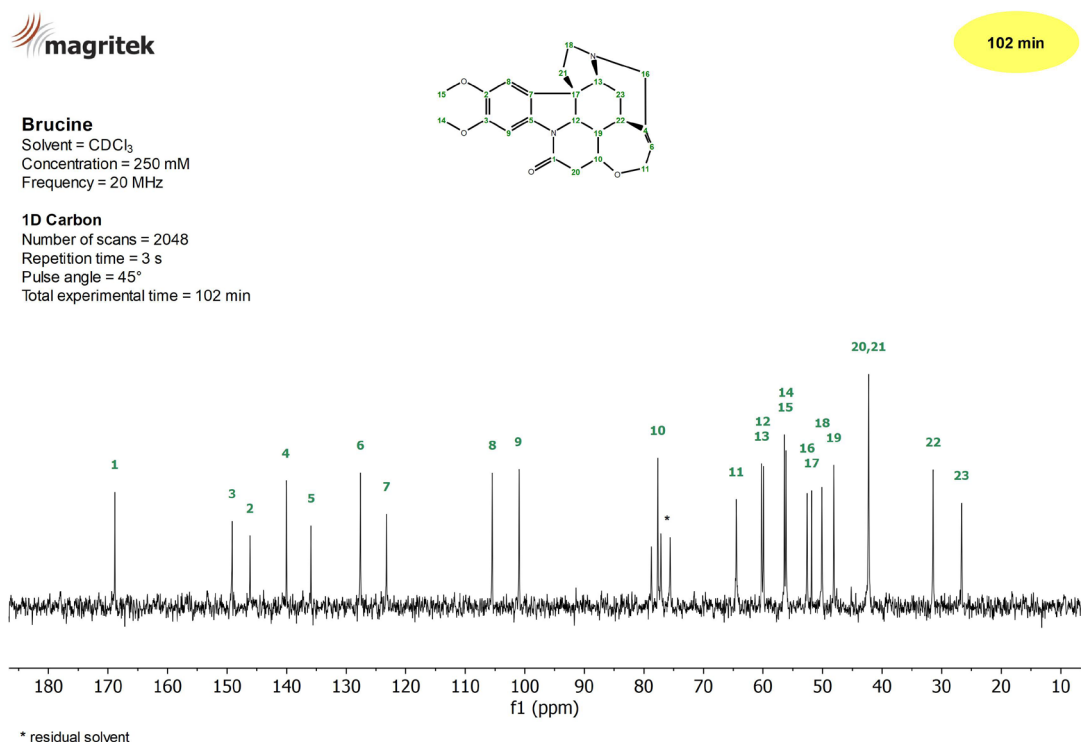


Figure 3:  $^{13}\text{C}$  NMR spectrum of a 250 mM Brucine sample in  $\text{CDCl}_3$  measured on a Spinsolve 80 MHz system in 102 minutes.

## 2D COSY

The 2D COSY experiment allows one to identify coupled  $^1\text{H}$  nuclei as they generate cross peaks out of the diagonal of the 2D data set. In Figure 2 a large number of cross peaks can be nicely observed. For example, the protons at position 6 and 11 (light green) couple with each other. Furthermore, proton 19 couples with proton 10 (light blue), 12 (orange) and 20 (pink). In addition, the couplings between protons 8 and 9 (dark blue) as well as the couplings of protons 8 and 9 with protons 14 and 15 (dark green) can be nicely observed.



### Brucine

Solvent =  $\text{CDCl}_3$   
 Concentration = 250 mM  
 Frequency = 80 MHz

### COSY

Number of scans = 1  
 Total experimental time = 17 min

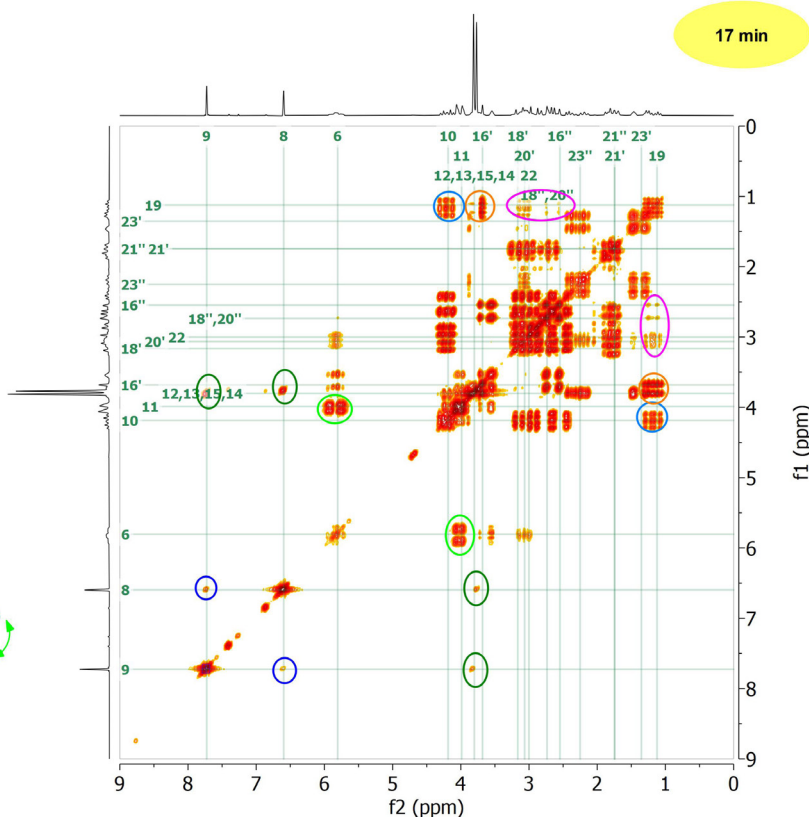
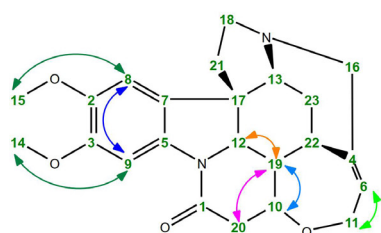


Figure 2:  $^1\text{H}$  2D COSY experiment of a 250 mM Brucine sample in  $\text{CDCl}_3$  acquired in 17 minutes on a Spinsolve 80 MHz system.

## 2D HSQC-ME

The HSQC is a powerful sequence widely used to correlate  $^1\text{H}$  with the one-bond coupled  $^{13}\text{C}$  nuclei. The Spinsolve is equipped with a multiplicity edited version (HSQC-ME) of this method. It provides the editing power of the DEPT-135 sequence, which is useful to differentiate the signals of  $\text{CH}_2$  groups (blue) from  $\text{CH}$  and  $\text{CH}_3$  groups (red). Figure 4 shows the HSQC-ME spectrum of a 250 mM Brucine in  $\text{CDCl}_3$  acquired in 4 minutes. The measurement time was optimized applying NUS (non uniform sampling).

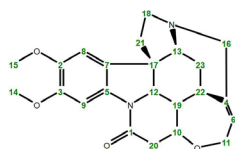


### Brucine

Solvent =  $\text{CDCl}_3$   
 Concentration = 250 mM  
 Frequency  $^1\text{H}$  = 80 MHz

### HSQC-ME

Number of scans = 2  
 Repetition time = 1 s  
 Number of steps = 256  
 NUS = 50%  
 Total experimental time = 4 min



Red = CH and  $\text{CH}_3$

Blue =  $\text{CH}_2$

\* residual solvent

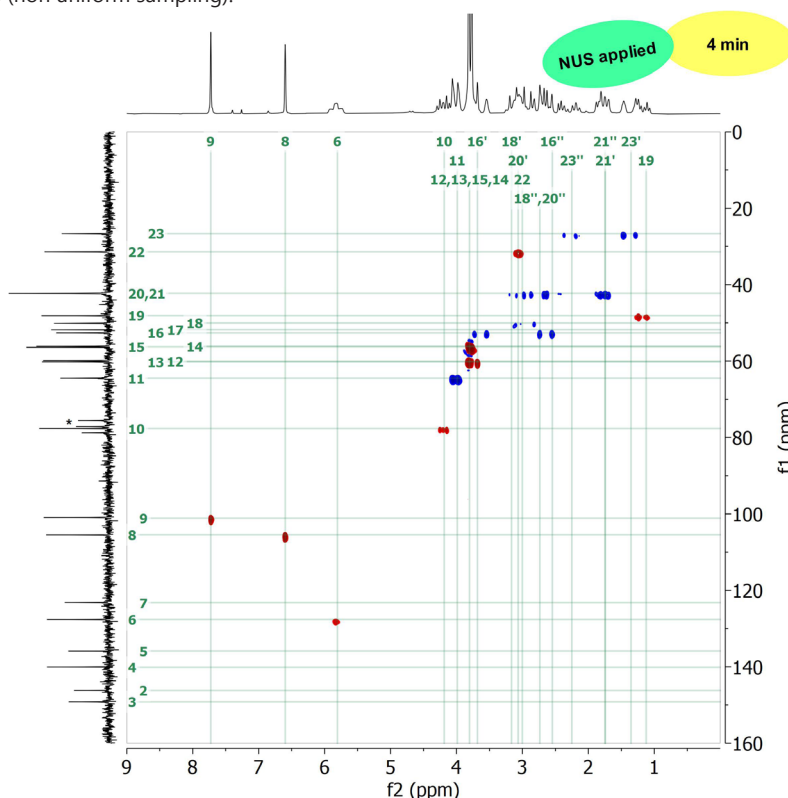


Figure 4: HSQC-ME spectrum of a 250 mM Brucine sample in  $\text{CDCl}_3$  showing the correlation between the  $^1\text{H}$  (horizontal) and  $^{13}\text{C}$  (vertical) signals.

## 2D HMBC

To obtain long-range  $^1\text{H}$ - $^{13}\text{C}$  correlations through two or three bond couplings, the Heteronuclear Multiple Bond Correlation (HMBC) experiment can be used. Figure 5 shows the HMBC spectrum of a 250 mM Brucine sample measured in 34 minutes on our Spinsolve 80 MHz. As an example, the long-range correlations of proton 8 with carbons 17 (light green), 9 (dark green), 7 (dark blue), 5 (light blue), 2 (orange) and 3 (pink) are marked with circles. The experiment shows the correlation with quaternary carbons, too.

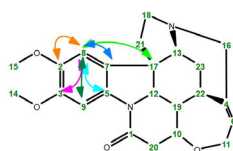


### Brucine

Solvent =  $\text{CDCl}_3$   
 Concentration = 250 mM  
 Frequency  $^1\text{H}$  = 80 MHz

### HMBC

Number of scans = 8  
 Repetition time = 1 s  
 Number of steps = 256  
 Total experimental time = 34 min



\* residual solvent

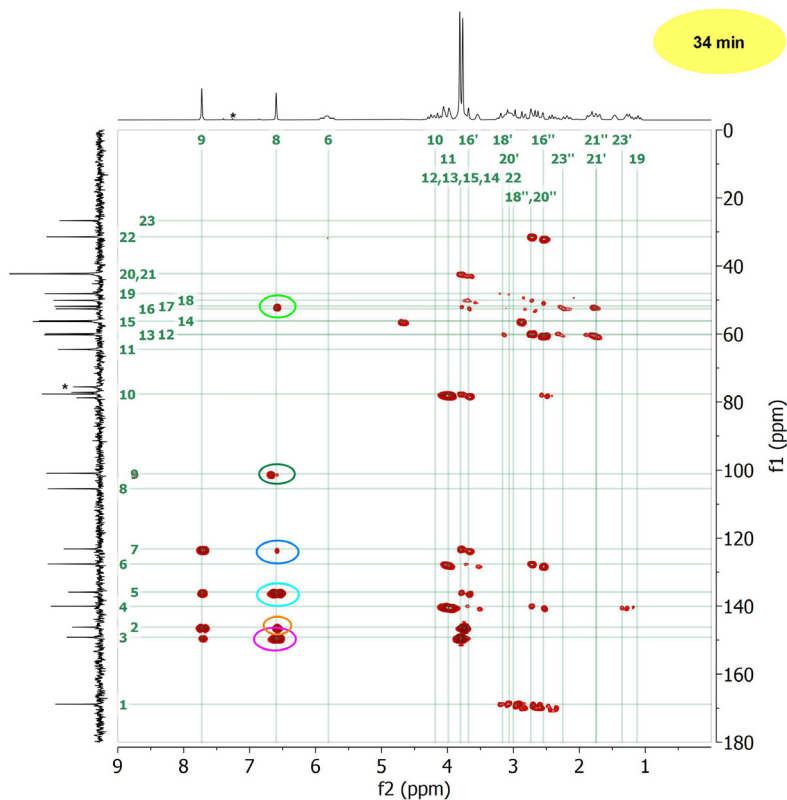


Figure 5: HMBC spectrum of a 250 mM Brucine sample in  $\text{CDCl}_3$  showing the long-range couplings between  $^1\text{H}$  and  $^{13}\text{C}$  nuclei.