

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 ^1H NMR spectrum of a 250 mM Brucine sample in CDCl_3 measured in a single scan taking 10 seconds to acquire.

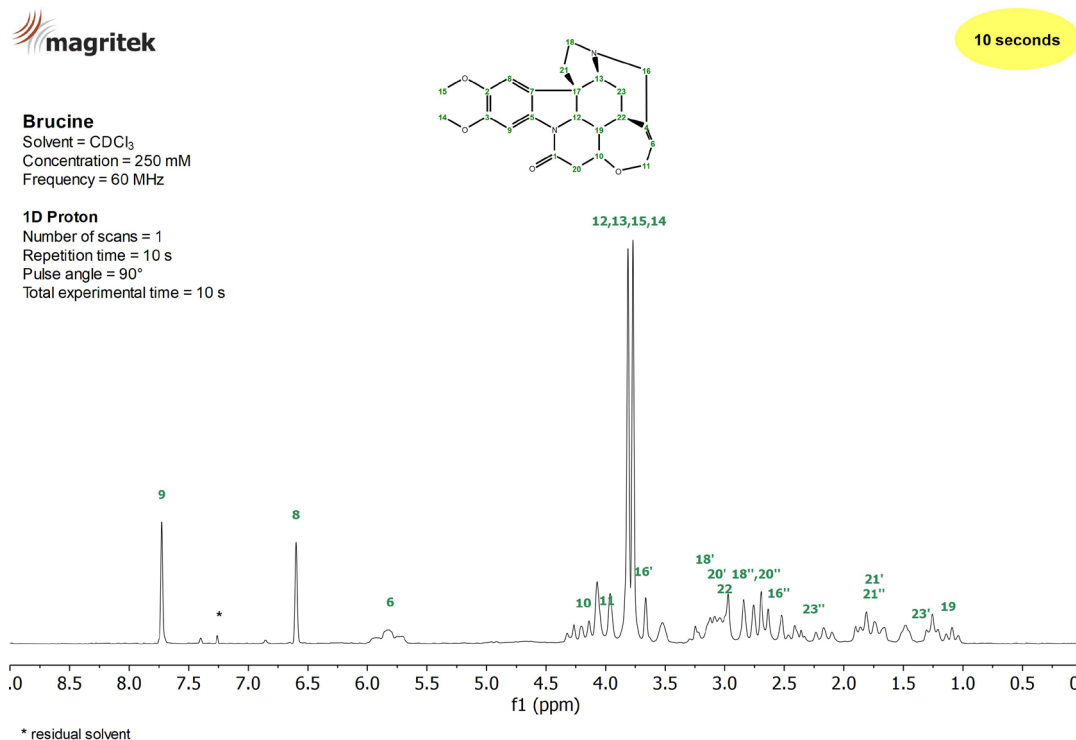


Figure 1: ^1H NMR spectrum of a 250 mM Brucine sample in CDCl_3 measured on a Spinsolve 60 MHz system in a single scan.

^{13}C Spectrum

Figure 3 shows the ^{13}C NMR spectrum of 250 mM Brucine in CDCl_3 acquired using NOE polarization transfer from ^1H to ^{13}C and ^1H decoupling. The 1D Carbon experiment using NOE is sensitive to all ^{13}C nuclei in the sample. It clearly resolves all the expected resonances.

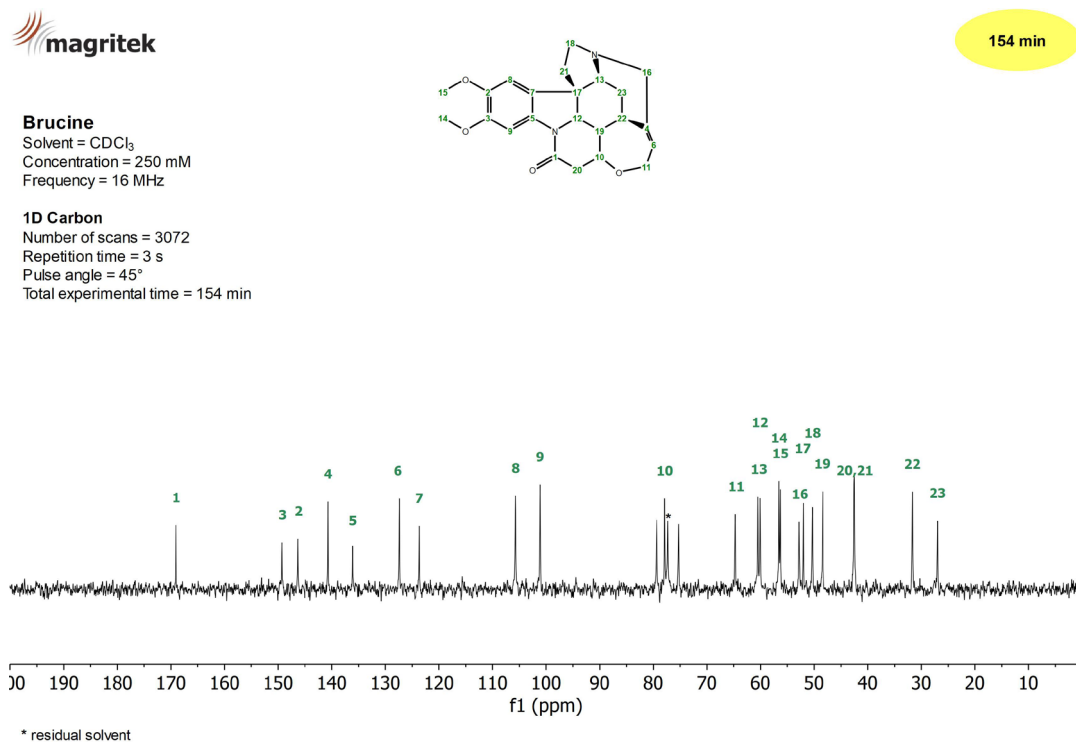


Figure 3: ^{13}C NMR spectrum of a 250 mM Brucine sample in CDCl_3 measured on a Spinsolve 60 MHz system in 154 minutes.

2D COSY

Spinsolve™ 60

The 2D COSY experiment allows one to identify coupled ^1H 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 observed nicely. 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 observed nicely.



Brucine

Solvent = CDCl_3
Concentration = 250 mM
Frequency = 60 MHz

COSY

Number of scans = 1
Total experimental time = 17 min

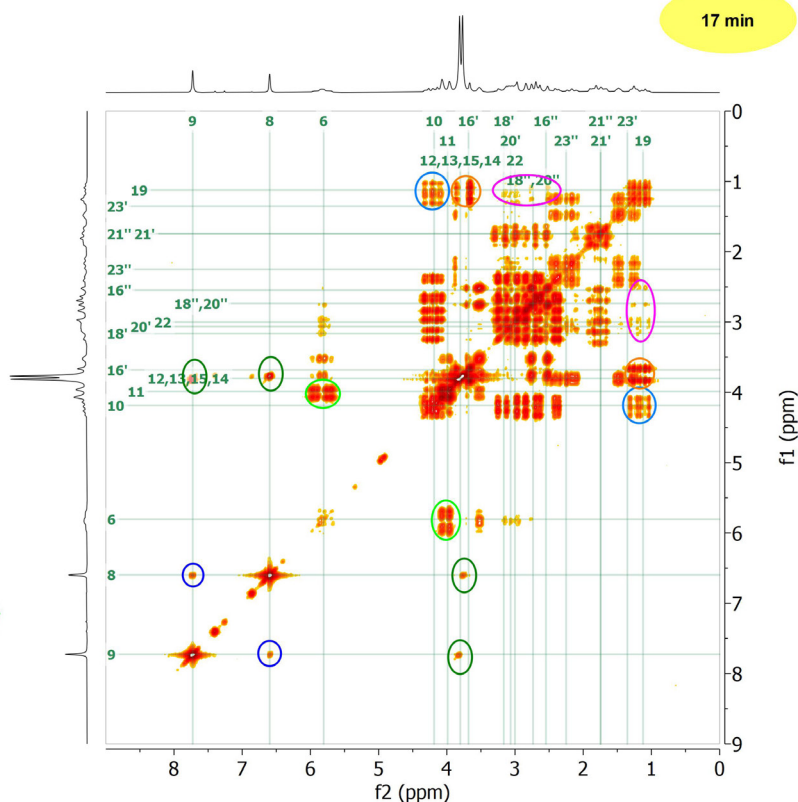
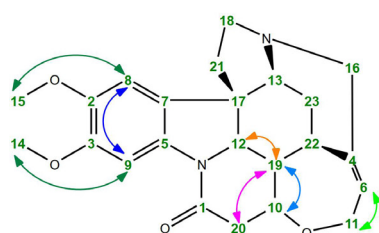


Figure 2: ^1H 2D COSY experiment of a 250 mM Brucine sample in CDCl_3 acquired in 17 minutes on a Spinsolve 60 MHz system.

2D HSQC-ME

The HSQC is a powerful sequence widely used to correlate ^1H with the one-bond coupled ^{13}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 CH_2 groups (blue) from CH and CH_3 groups (red). Figure 4 shows the HSQC-ME spectrum of a 250 mM Brucine in CDCl_3 acquired in 8 minutes.

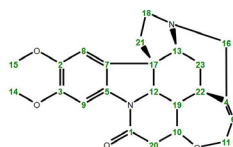


Brucine

Solvent = CDCl_3
Concentration = 250 mM
Frequency ^1H = 60 MHz

HSQC-ME

Number of scans = 2
Repetition time = 1 s
Number of steps = 256
Total experimental time = 8 min



Red = CH and CH_3

Blue = CH_2

* residual solvent

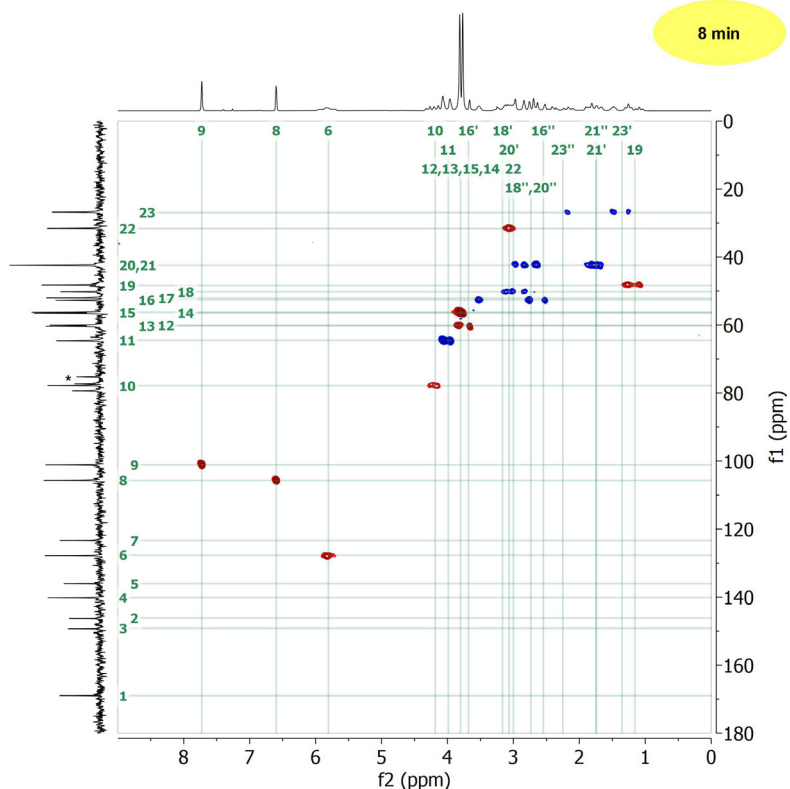


Figure 4: HSQC-ME spectrum of a 250 mM Brucine sample in CDCl_3 showing the correlation between the ^1H (horizontal) and ^{13}C (vertical) signals.

2D HMBC

To obtain long-range ^1H - ^{13}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 60 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.

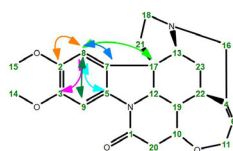


Brucine

Solvent = CDCl_3
Concentration = 250 mM
Frequency ^1H = 60 MHz

HMBC

Number of scans = 16
Repetition time = 1 s
Number of steps = 256
Total experimental time = 68 min



* residual solvent

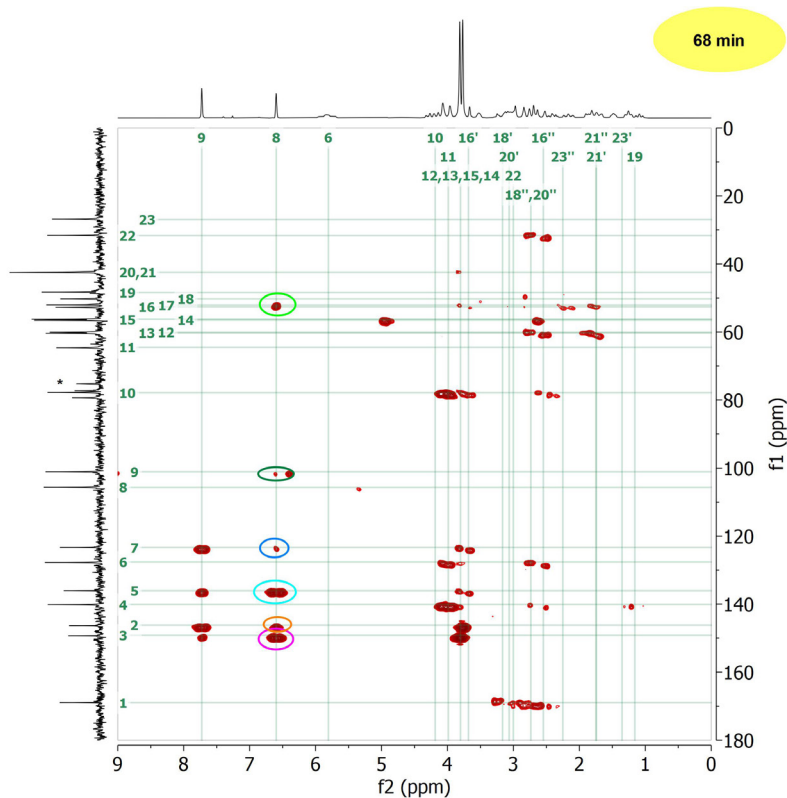


Figure 5: HMBC spectrum of a 250 mM Brucine sample in CDCl_3 showing the long-range couplings between ^1H and ^{13}C nuclei.