C8953
NMR strukturní analýza
seminář
NOESY

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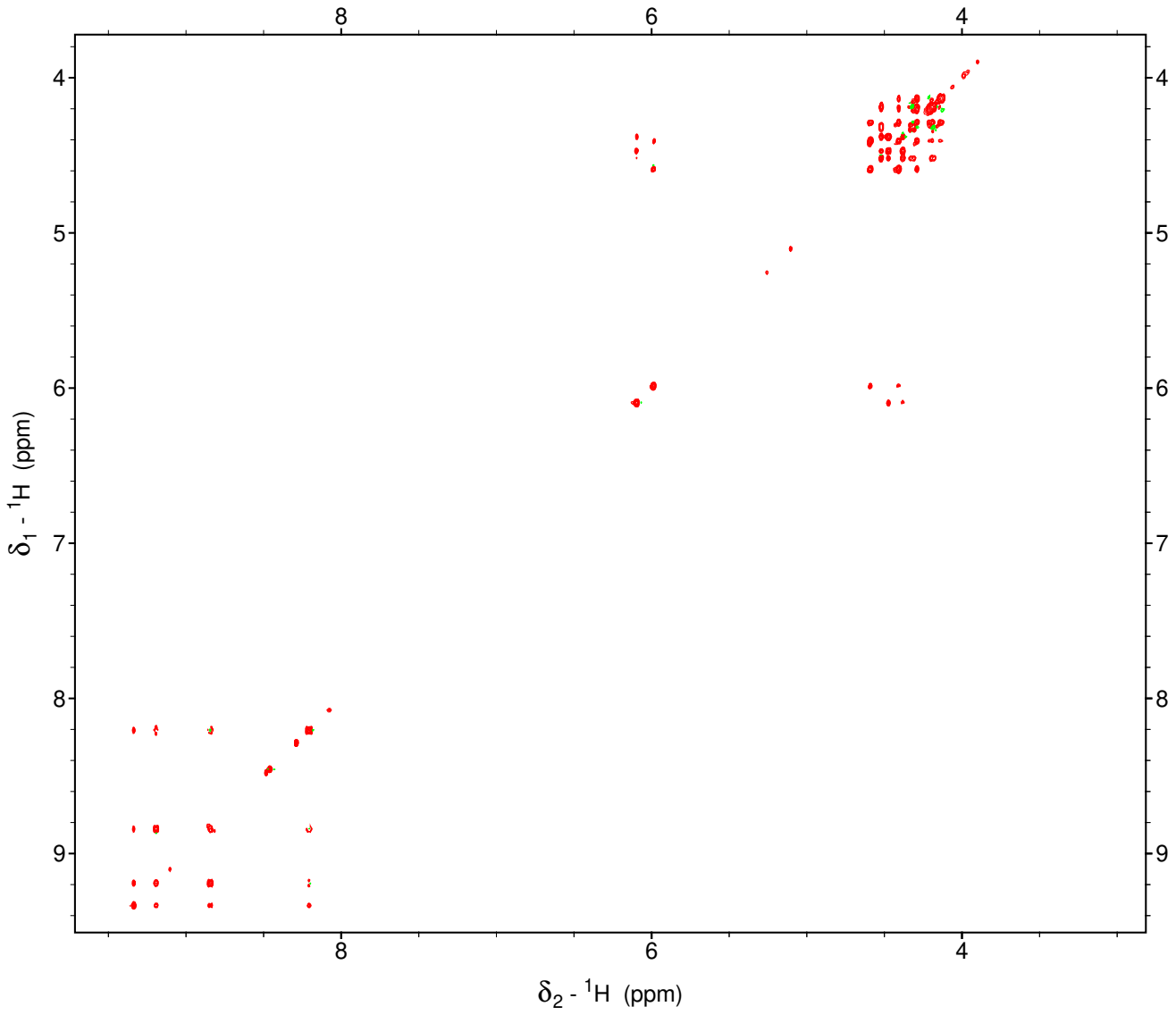
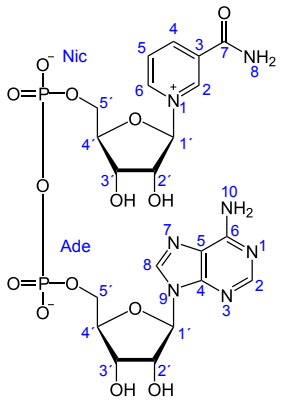
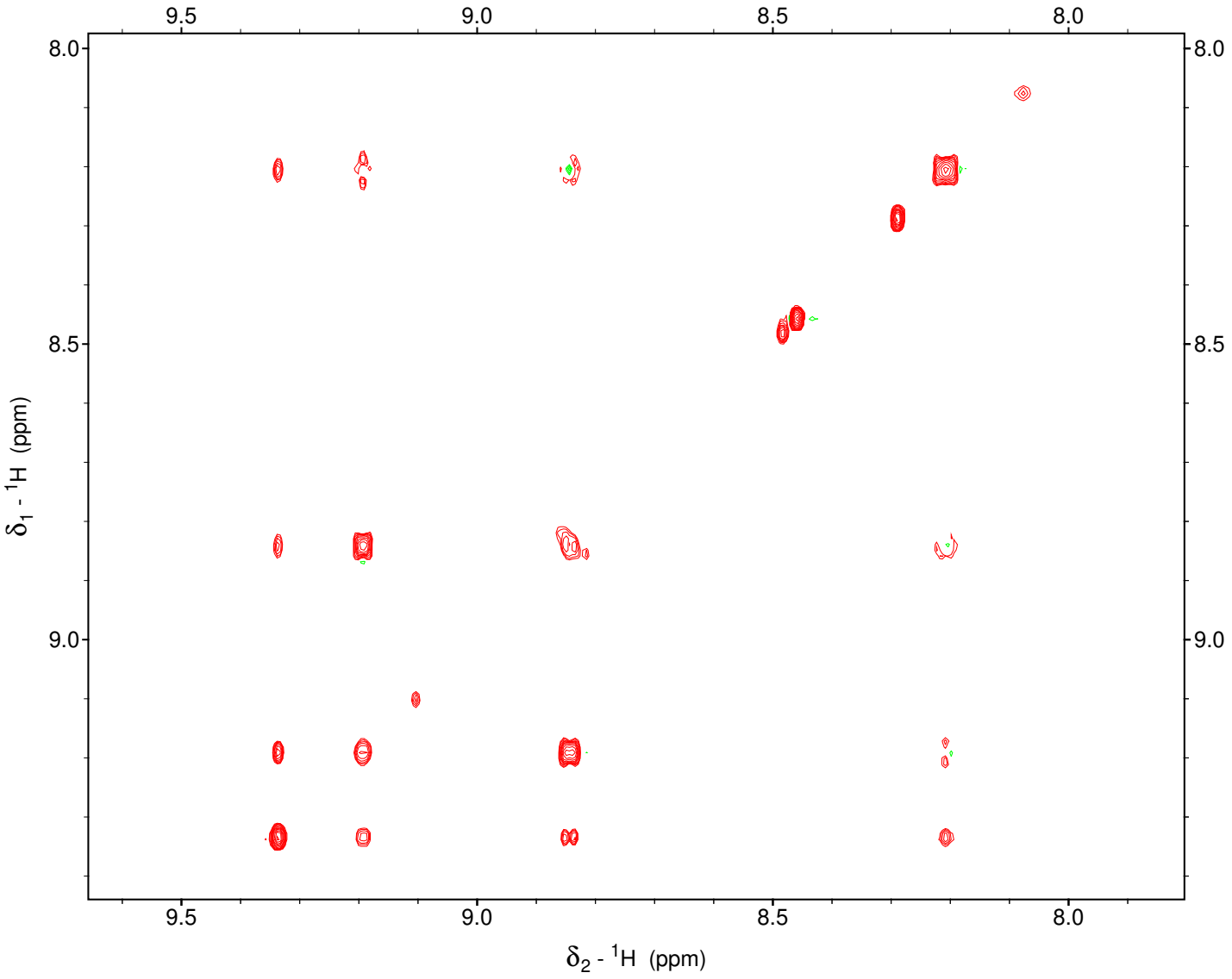
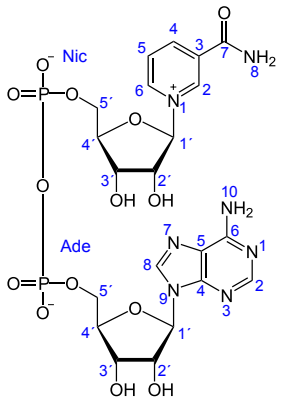
NAD^+: TOCSY (40ms)



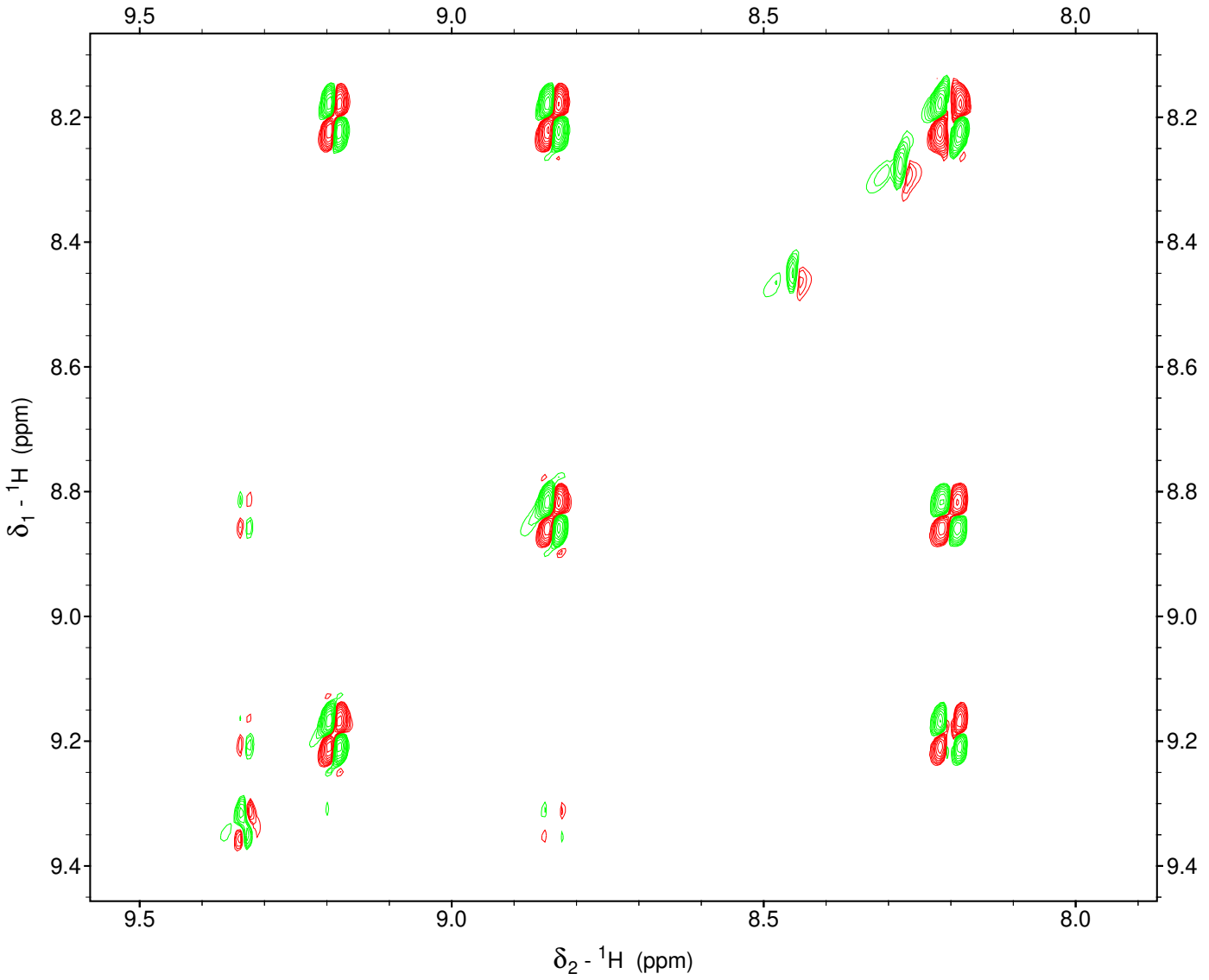


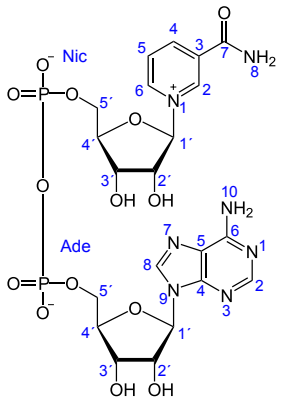
NAD^+: TOCSY (40ms), detail of aromatics



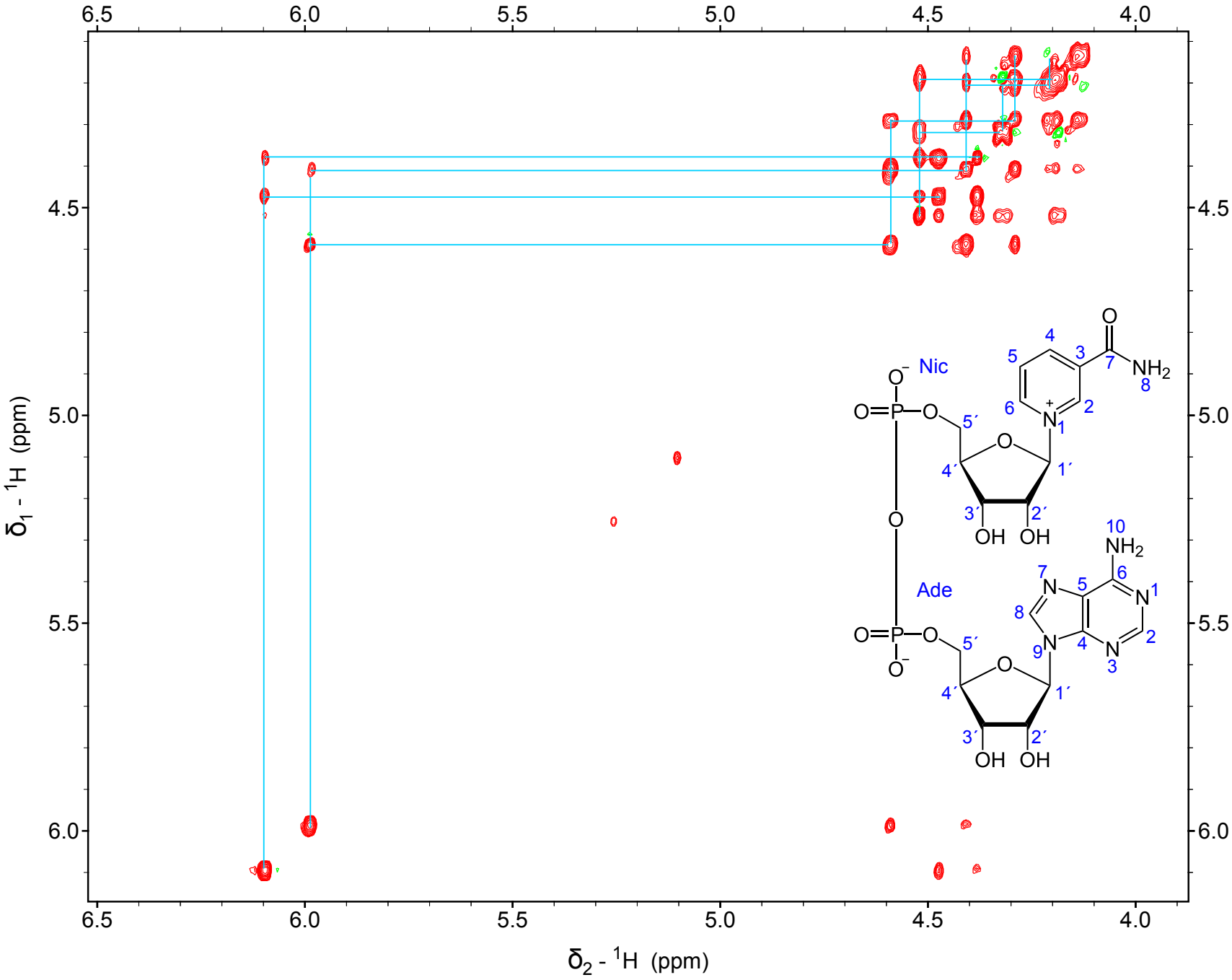


NAD^+: DQF-COSY, detail of aromatics





NAD^+: TOCSY (40ms), detail of aliphatics



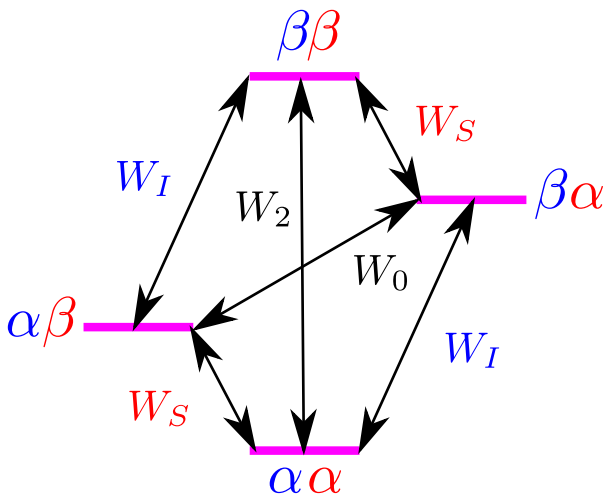
NOESY - introduction

Nuclear Overhauser effect

- dipol-dipol interaction
- magnetisation transfer TROUGH SPACE as a consequence of cross-relaxation

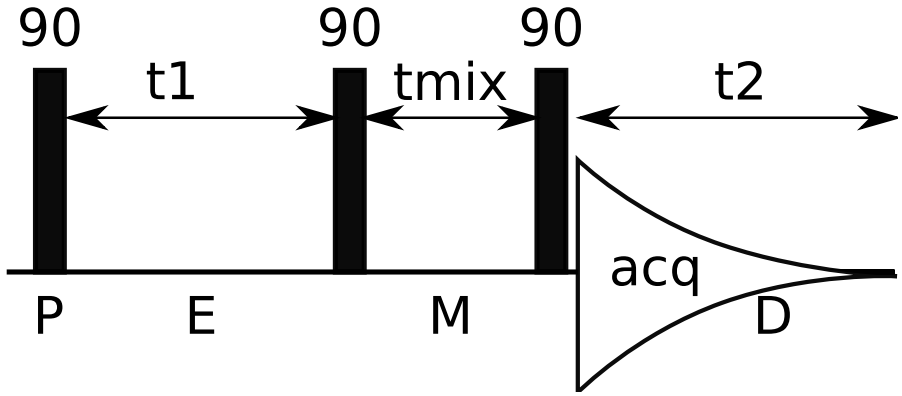
NOESY

- correlates nuclei if their distance is **smaller than 5 Å**



$$\frac{\mathrm{d}\Delta I_z}{\mathrm{d}t} = -\rho_I(I_z - I_z^0) - \sigma_{IS}(S_z - S_z^0)$$

$$\sigma_{IS} = W_2 - W_0$$

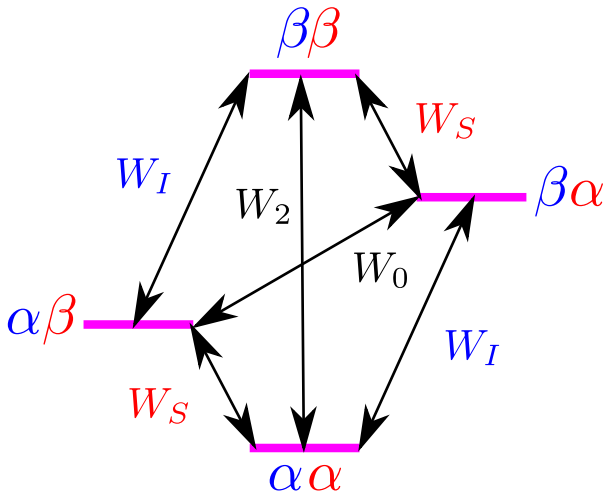


NOE vs. size of a molecule

NOE enhancement
$\frac{M_I^S - M_{0I}}{M_{0I}} \approx \frac{\sigma_{IS}}{\rho_I} = \frac{W_2 - W_0}{W_0 + 2W_1 + W_2}$, where M_I^S is magnetisation of I perturbed by saturation of S

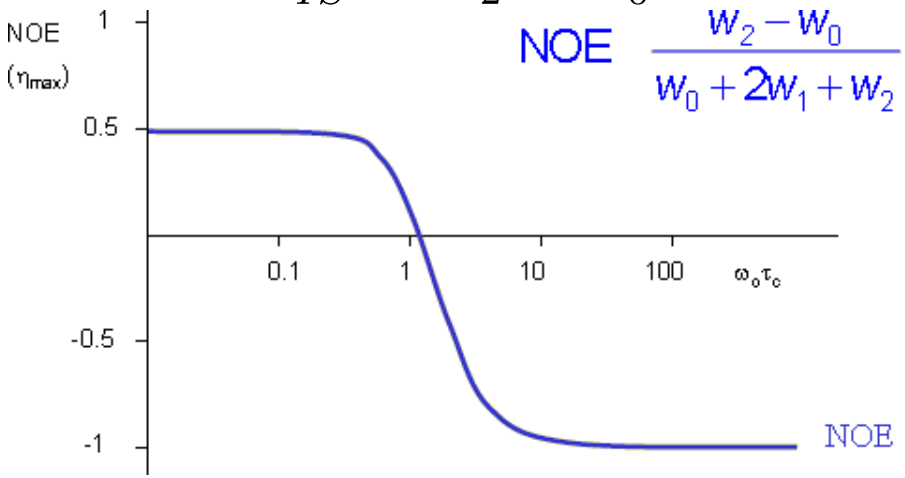
Correlation time τ_c

- $\omega_0 \tau_c < 1 \Leftrightarrow \omega_0 \frac{1}{f} < 1 \Leftrightarrow \omega_0 < f$ (small molecules $\ll$ 1 kDa)
  - **fast molecular motion**, $\beta\beta \rightarrow \alpha\alpha$ dominates $\Rightarrow W_2 > W_0$
  - positive NOE
  - crosspeaks have opposite phase relative to diagonal
- $\omega_0 \tau_c > 1$ (large molecules $\gg$ 1 kDa)
  - **slow molecular motion**, $\alpha\beta \rightarrow \beta\alpha$ dominates $\Rightarrow W_0 > W_2$
  - negative NOE
  - crosspeaks have the same phase
- $\omega_0 \tau_c \approx 1$ (cca 1 kDa)
  - NOE$\approx$0 - no crosspeaks
  - ROESY is an alternative



$$\frac{d\Delta I_z}{dt} = -\rho_I (I_z - I_z^0) - \sigma_{IS}(S_z - S_z^0)$$

$$\sigma_{IS} = W_2 - W_0$$



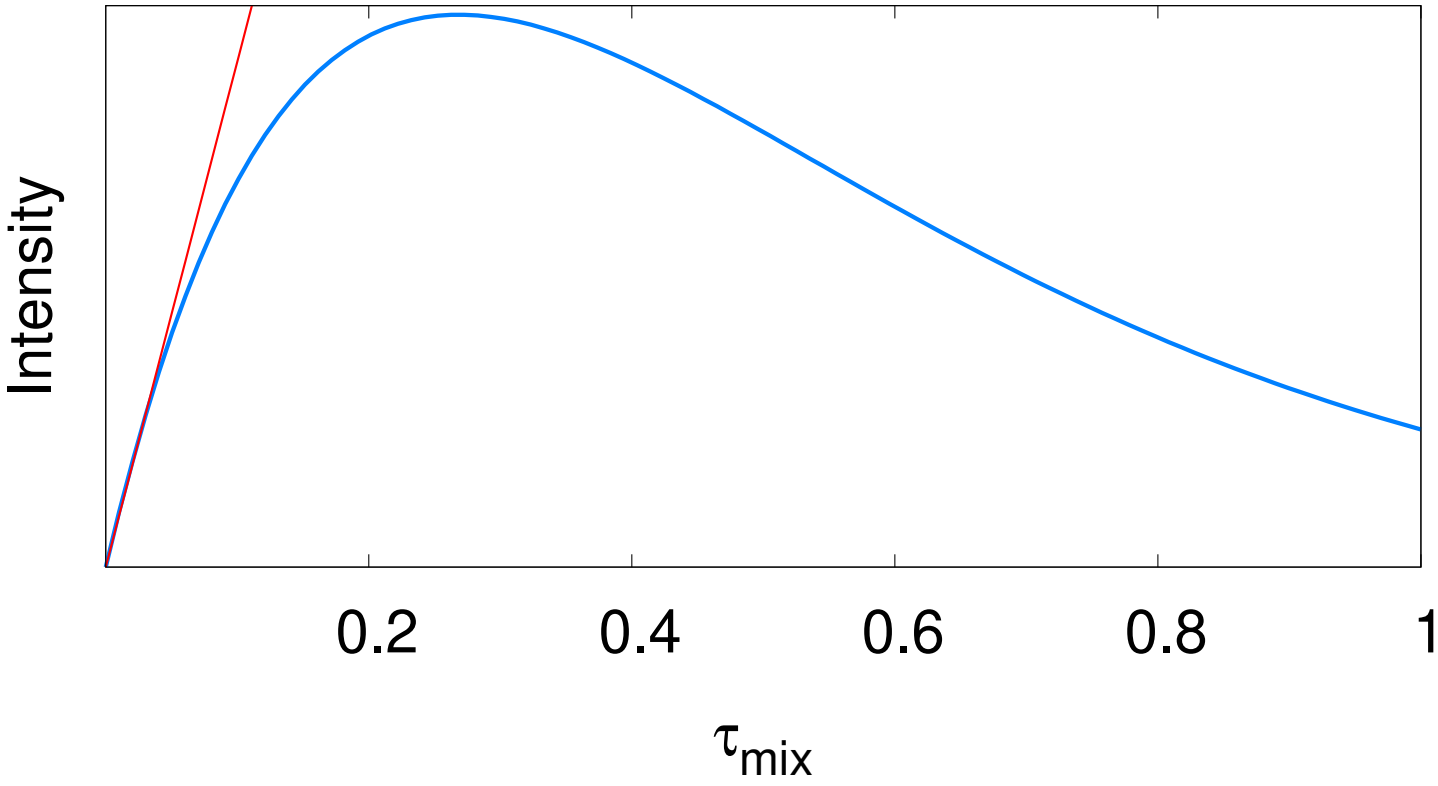
Application of NOESY

Mixing time τ_{mix}

- small molecules $\tau_{mix} \approx 500 - 800\,ms$
- biomolecules $\tau_{mix} \approx 50 - 300\,ms$

approximative determination of interatomic distatces ($< 5\,Å$)

- at short τ_{mix}
- $r_{ij} \approx A \times I_{ij}$

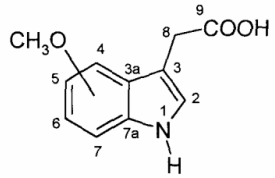


NOE differential experiment

PROBLEM 4

NOE- Difference Spectroscopy

Figure 4.1 shows the ^{1}H NMR and a ^{1}H NOE difference spectrum of a 3-indolylacetic acid derivative **13** bearing a methoxy group at the benzenoic ring.



What is the position of the methoxy group? **(400 MHz ^{1}H)**

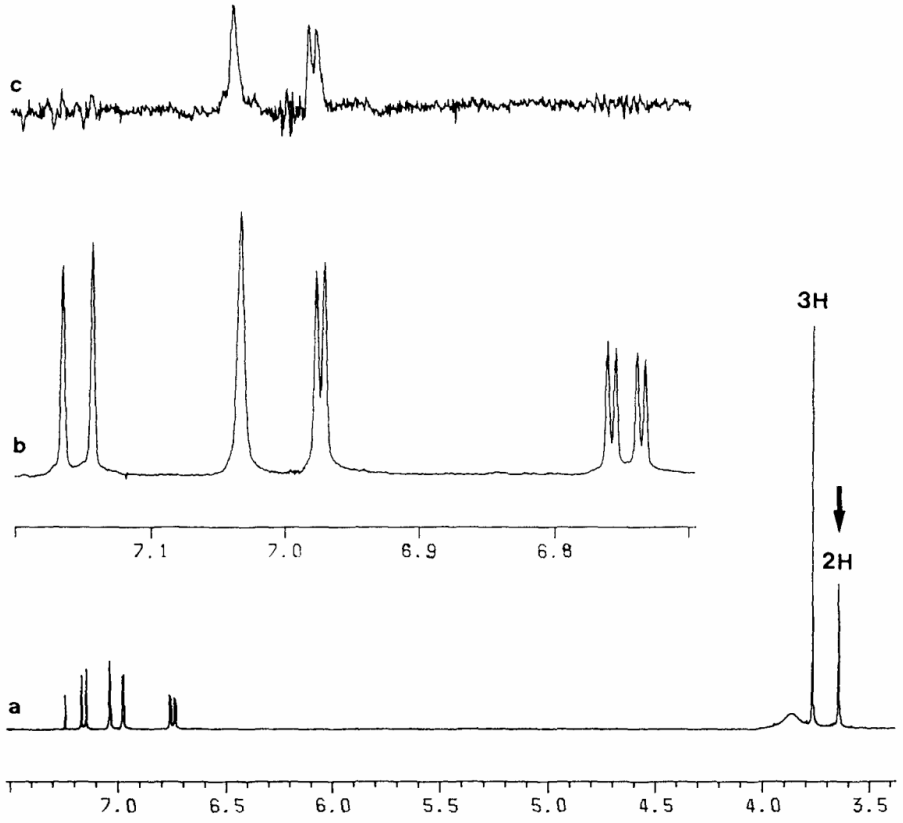
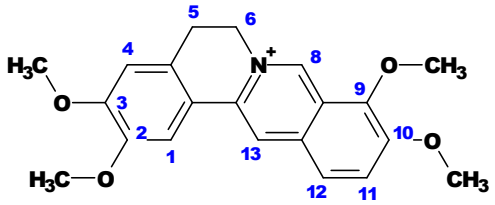
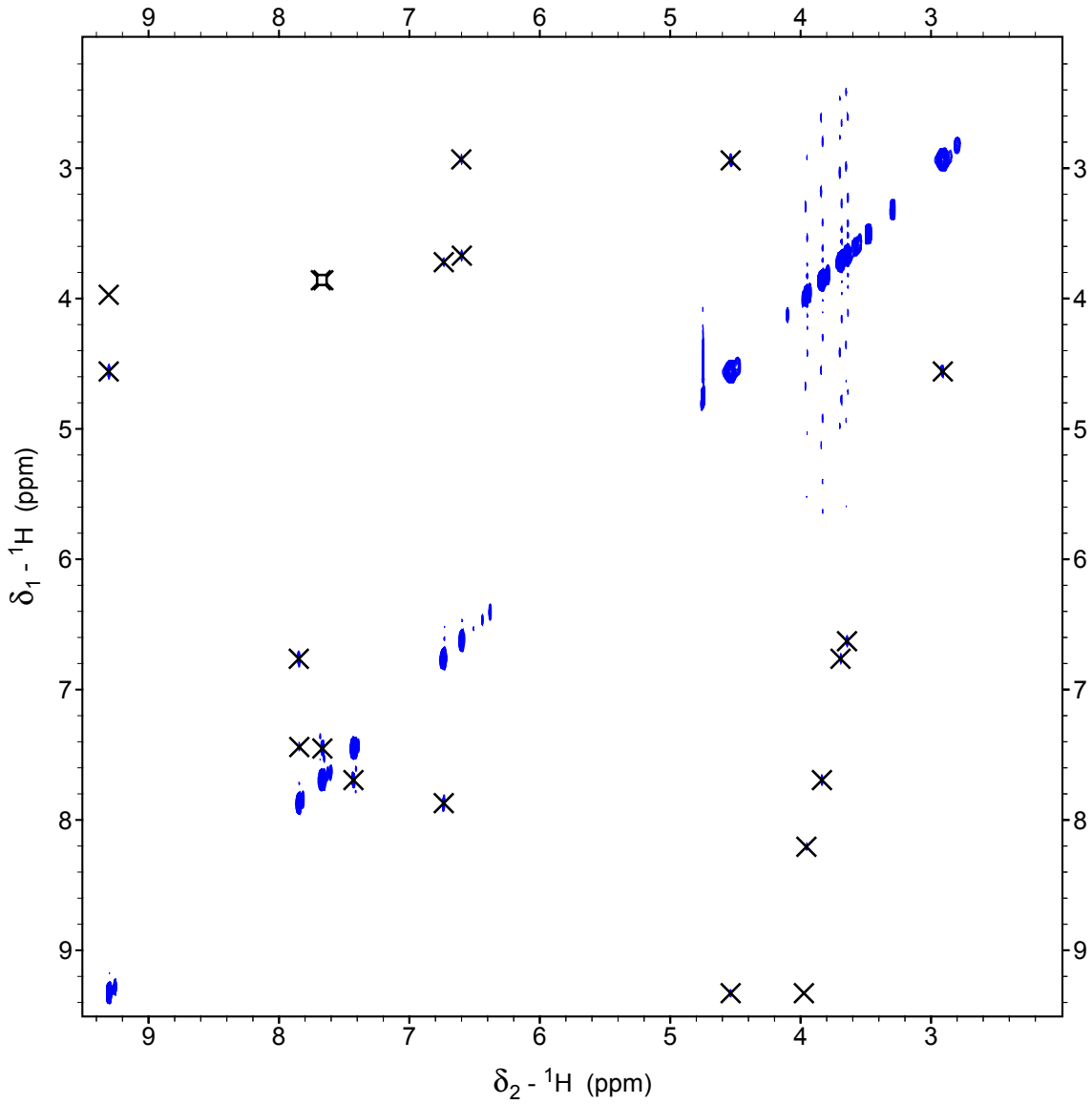


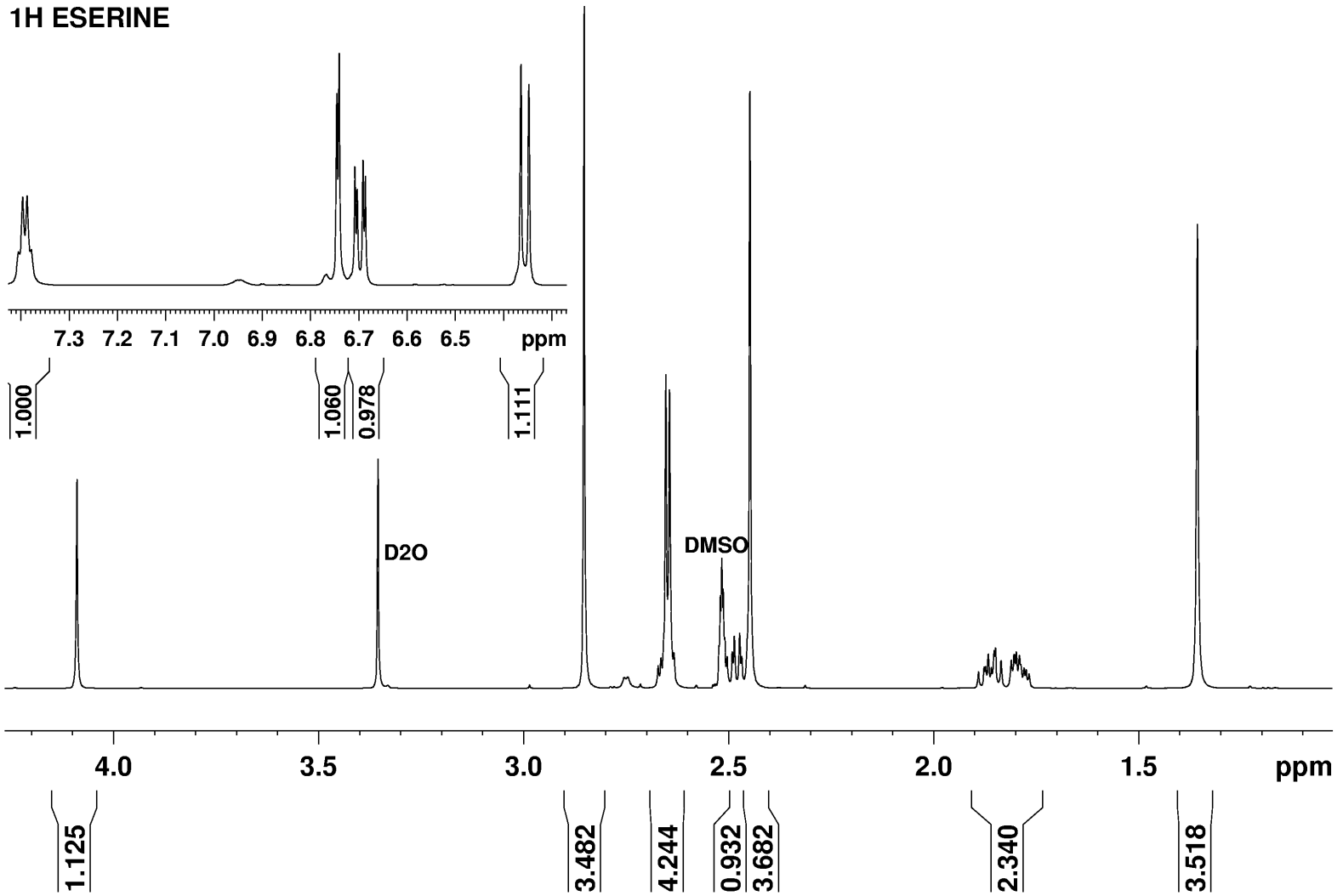
Fig. 4.1. 400 MHz ^{1}H NMR spectrum of **13** in a mixture of $CDCl_3$ and CD_3OD. **a** Full spectrum; **b** expanded section of the aromatic proton signals; **c** ^{1}H NOE difference spectrum, same section as in **b**, irradiation position at $\delta = 3.64$.

NOESY - Palmatine

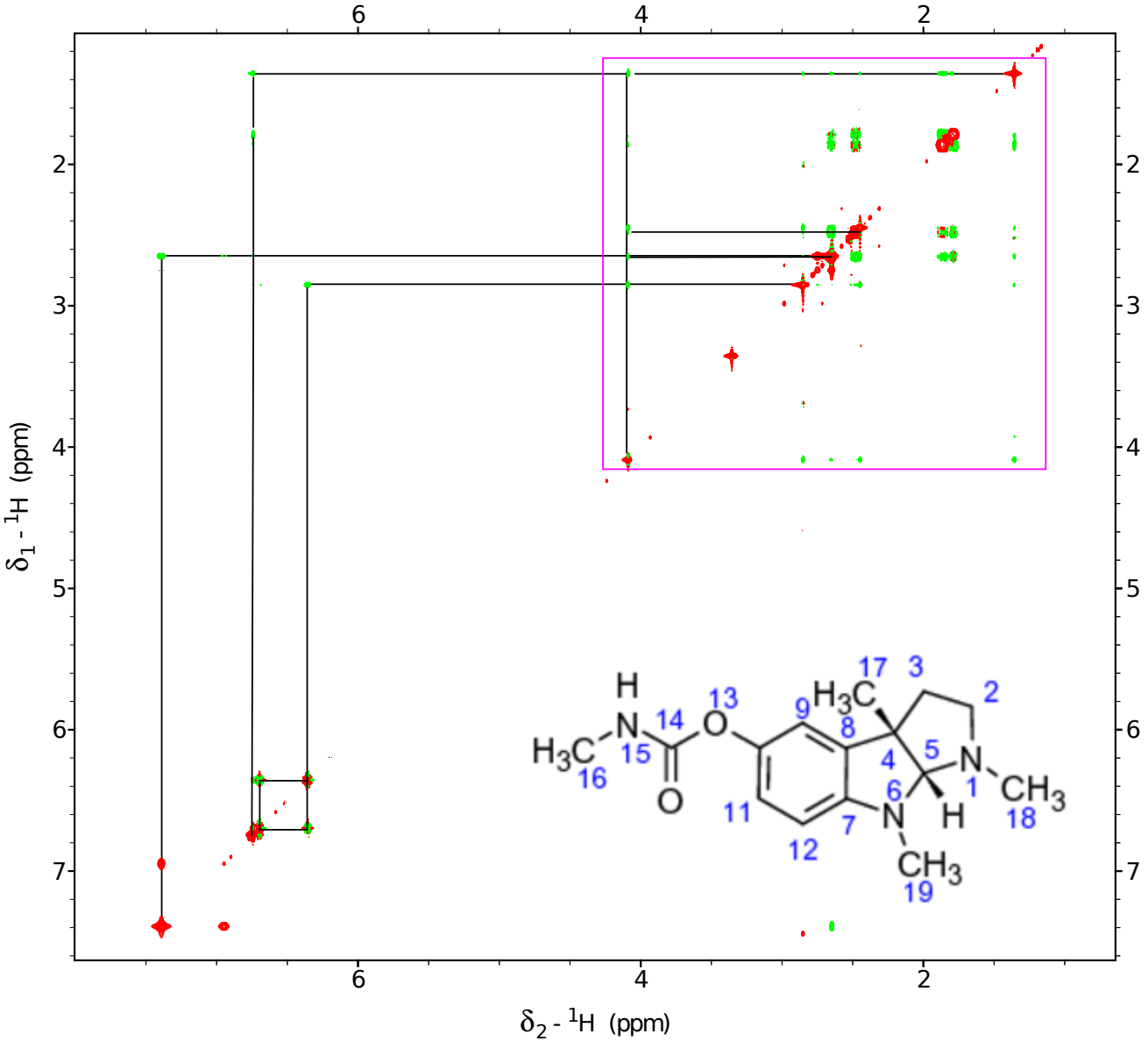




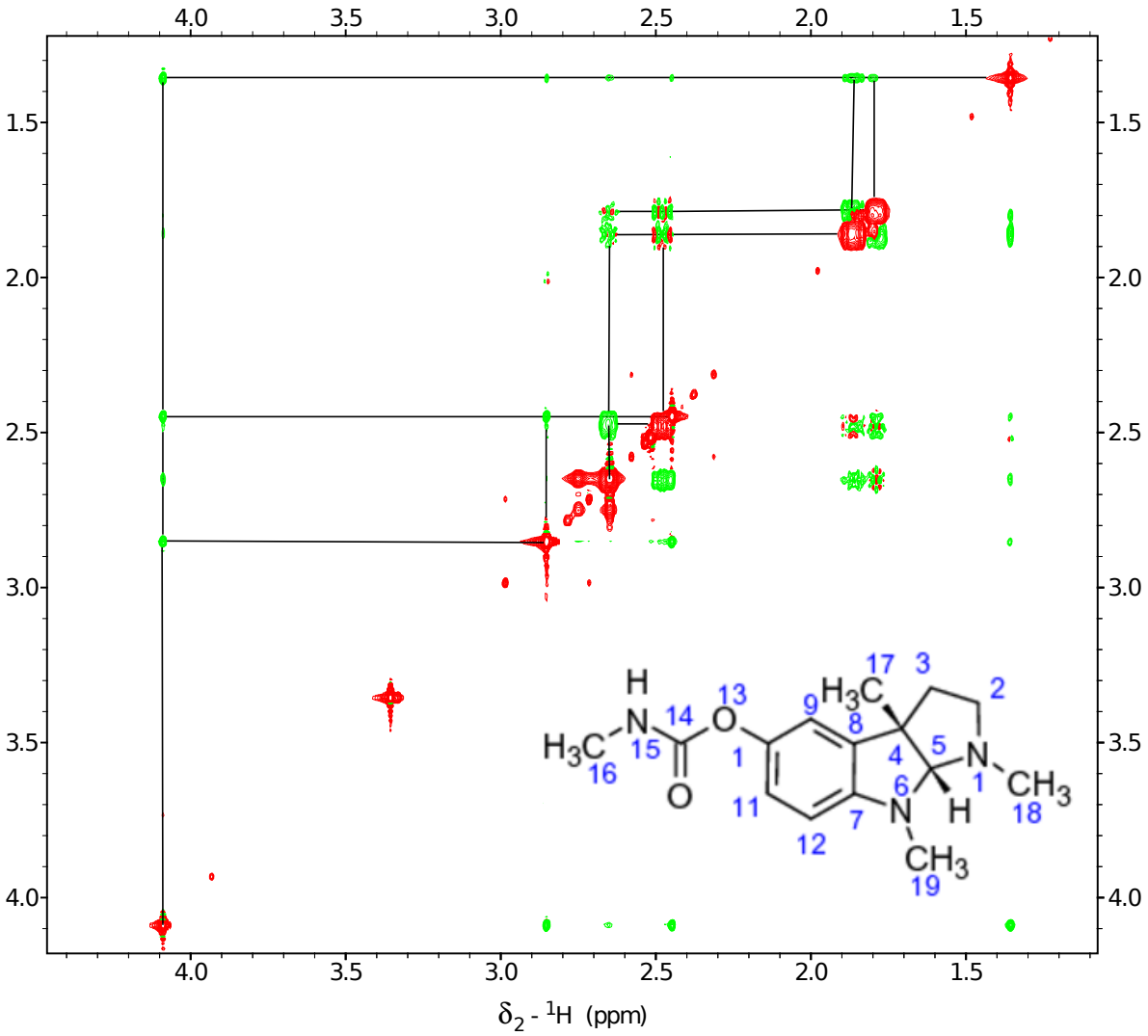
Eserine ^{1}H

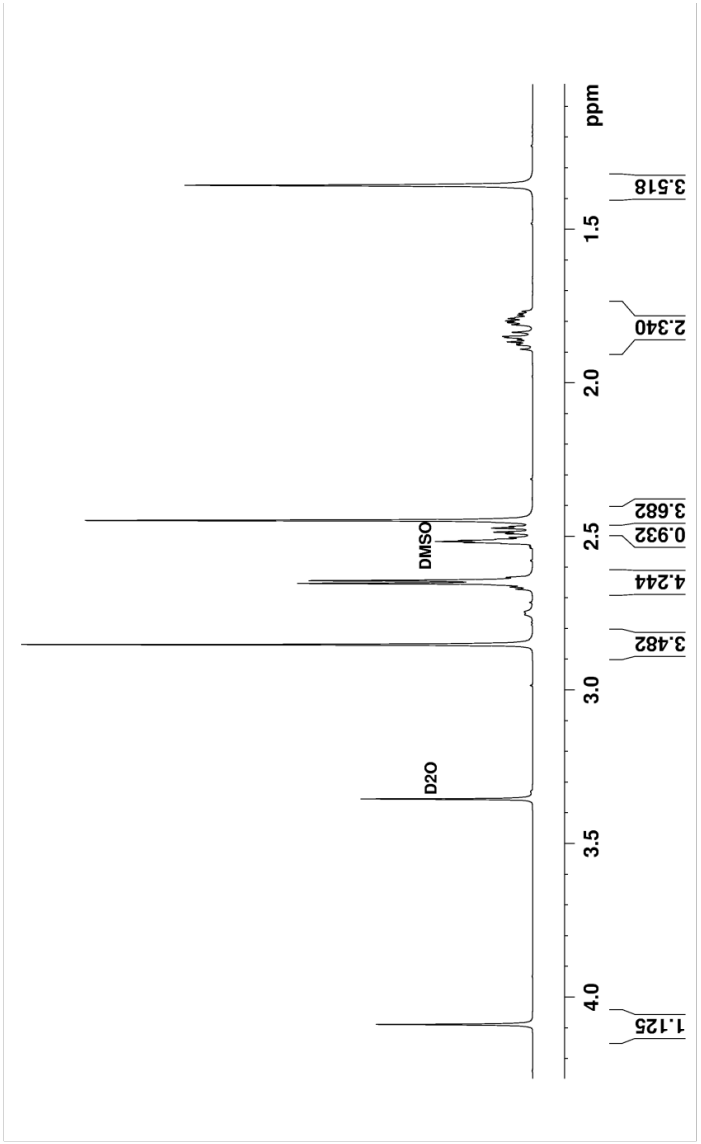

1H ESERINE
7.3 7.2 7.1 7.0 6.9 6.8 6.7 6.6 6.5 ppm
1.000
1.060
0.978
1.111
D2O
DMSO
4.0 3.5 3.0 2.5 2.0 1.5 ppm
1.125
3.482
4.244
0.932
3.682
2.340
3.518

NOESY - Eserine in DMSO

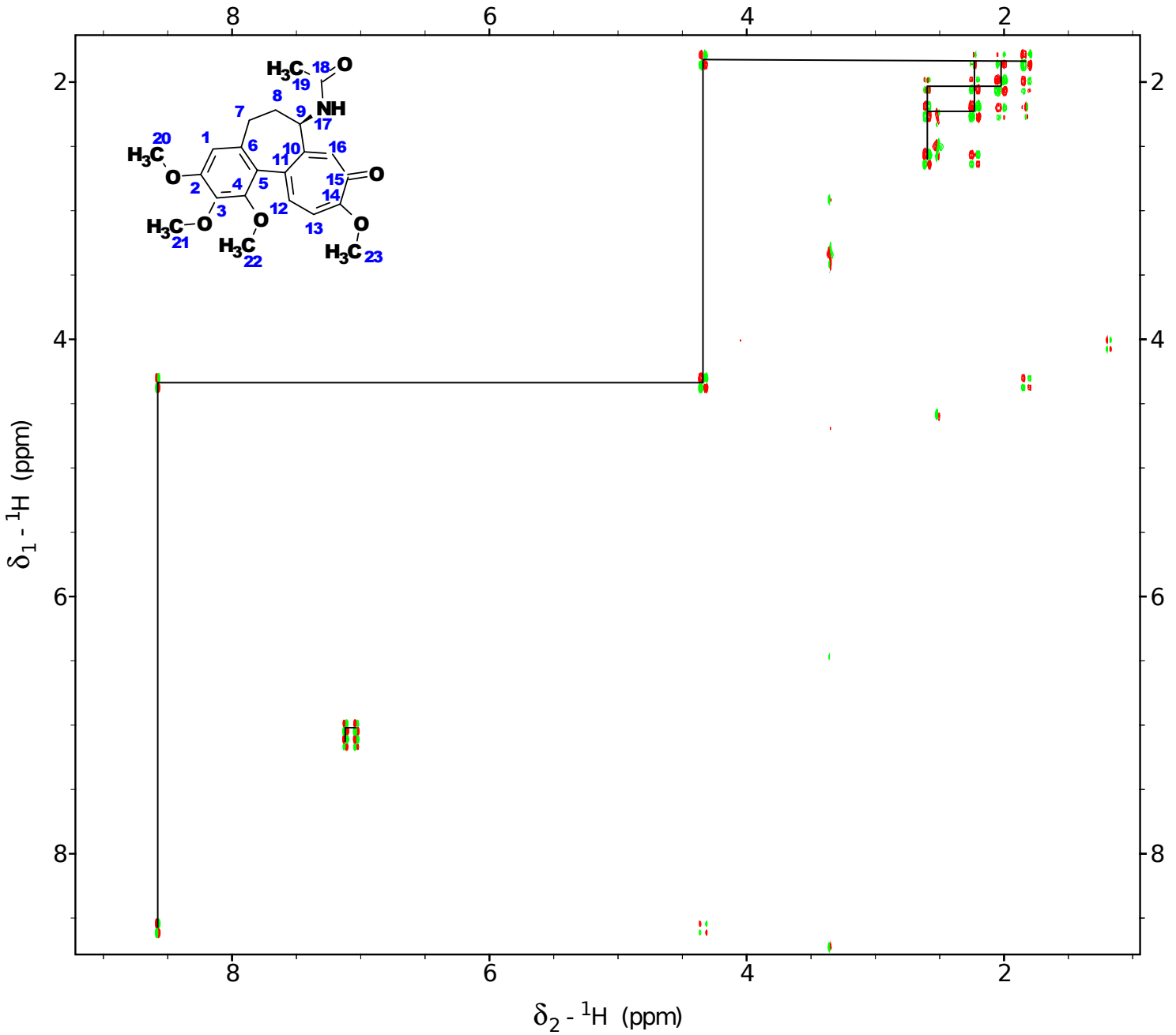


NOESY - Eserine





Colchicine - DQF-COSY



Colchicine - NOESY