### C8953 NMR structural analysis - seminar

Few Basic Concepts & Vector model

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# Assign correct value of chemical shift to labelled NMR active atoms<sup>1</sup>:

¹http://www.chem.wisc.edu/areas/reich/chem605/ □ ▶ ◀ 🖶 ▶ ∢ 🛢 ▶ 🧸 💆 🛷 🤉 🕞

# Diastereotopicity<sup>1</sup> Determine the equivalency of geminal protons

$$\begin{array}{c} C_6H_6 \\ Si \\ O \\ H_6C_6 \end{array}$$

$$\begin{array}{c} C_6H_6 \\ O \\ H_3C \end{array}$$

$$\begin{array}{c} C_6H_6 \\ O \\ H_3C \end{array}$$

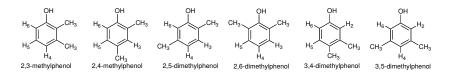
$$\begin{array}{c} CI \\ H_{MM} \\ H \\ Br \end{array}$$

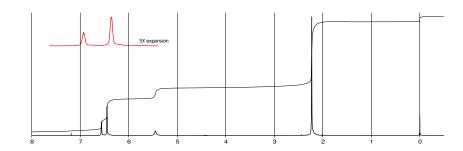
¹http://www.chem.wisc.edu/areas/reich/chem605/< □ > < ⊕ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > < ₹ > <

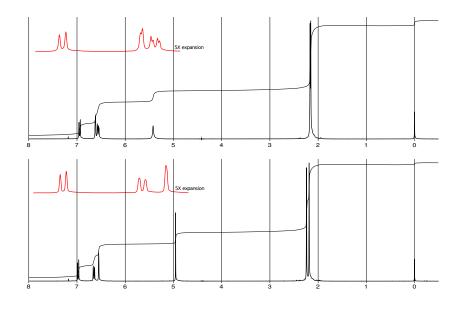
### Determination of regioisomers - note number of signals, splitting

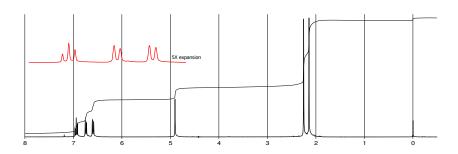
#### and $\delta$

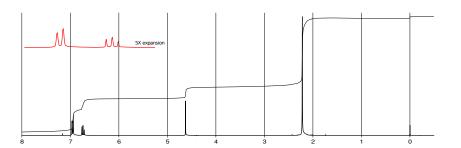
#### 300 MHz <sup>1</sup>H NMR spectra in CDCl<sub>3</sub>









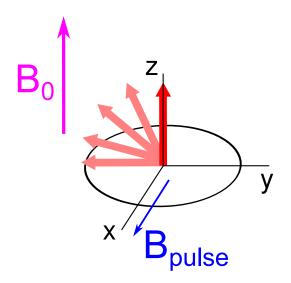


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# Draw the estimate of <sup>13</sup>C NMR spectrum (with and without <sup>1</sup>H decoupling)

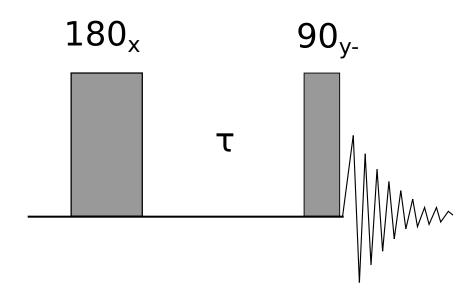
# Analysis of simple pulse sequences using vector model

- simple model based on rotation of the vector of bulk magnetization in the plane perpendicular to the vector of magnetic field, direction is determined by the "right-hand rule"
- NMR signal is detectable only as coherent magnetization oscillating in xy plane
- the free precession  $\omega$  (due to the B<sub>0</sub>) of magnetization vector is eliminated by introducing rotating frame  $\omega_0 \Rightarrow$  magnetic field of excitation pulses (B<sub>1</sub>) is motionless and the individual resonance frequencies differs in so called offset  $\Omega_i = \omega_i \omega_0$
- applicability of vector model is rather limited to simple single-quantum experiments without transfer of polarisation



#### $T_1$ relaxation

Apply following sequence (inversion recovery) to isolated spin characterized by **a)**  $\tau = 2 * T_1$  and **b)**  $\tau = 0.2 * T_1$ . Draw semi-quantitatively resulting spectrum.



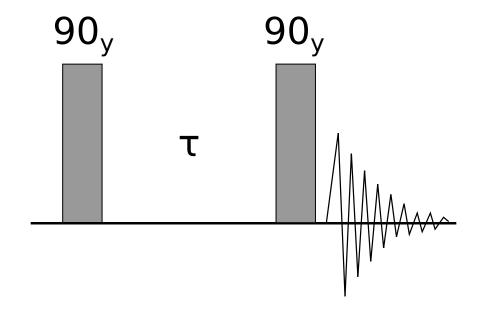
### 1-1 sequence

Draw the evolution of macroscopic magnetization through the sequence:

90(y) - 
$$\tau$$
 - 90(y) - aq

Consider the evolution of an isolated spin due to the chemical shift.

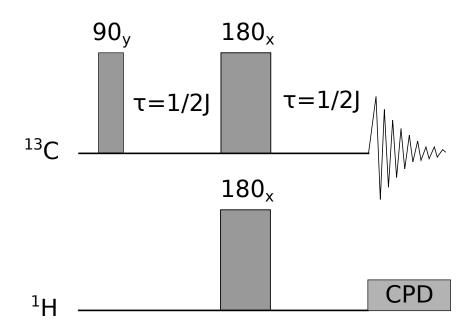
- 1. How does the result differ for the following offsets:  $\Omega \tau = 0, \pi/2, \pi$ .
- 2. Draw lineshapes of resulting signal assuming the a) y + b) x + corresponds to zero phase of receiver (prior phase correction).



#### Heteronuclear spin echo

By using vector diagrams determine the result of attached pulse sequence.

- 1. **Ignore 180 pulse** in hydrogen channel for isolated spin systems **a)** <sup>13</sup>C-<sup>1</sup>H and **b)** <sup>13</sup>C-<sup>1</sup>H<sub>2</sub>. Explain the role of CPD block.
- 2. Lets consider **the complete sequence** and isolated spin systems **a)** <sup>13</sup>C-<sup>1</sup>H and **b)** <sup>13</sup>C-<sup>1</sup>H<sub>2</sub>.



### Next topic

edited 1D 13C spectra, 2D NMR - homonuclear experiments