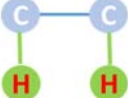
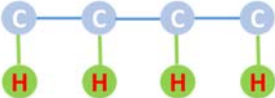

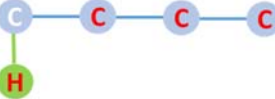
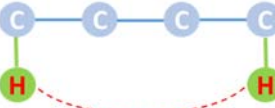


## Basic 1-D NMR Experiments

Method	Information	Parameter Set	Pulse Sequence
<b><sup>1</sup>H</b>	$\delta_{1H}, J_{HH}$	<b>PROTON</b>	zg30
	<i>homo-nuclear decoupling</i>	<b>1d_1H_homoDec</b>	zghd.2
<b><sup>1</sup>H + Solvent Suppression</b>	<i>solvent suppression using excitation sculpting</i>	<b>ZGESGP</b>	zgesgp
	<i>solvent suppression using pre-saturation with complex pulses</i>	<b>ZGCPPR</b>	zgcppr
	<i>multi-solvent suppression using pre-saturation</i>	<b>LC1D12GP</b>	lc1gppnf2
<b><sup>13</sup>C + <sup>1</sup>H decoupling</b>	$\delta_{13C}, ^1H$ decoupled (NOE effect, integration not possible)	<b>C13CPD</b>	zpgp30
	$\delta_{13C}, ^1H$ decoupled (fast relaxation, short $d_1$ needed)	<b>1d_13C_UDEFT</b>	udeft
	$\delta_{13C}, ^1H$ decoupled (no NOE effect, integration possible)	<b>C13IG</b>	zgig30
<b><sup>13</sup>C + <math>CH_n</math> test</b>	$\delta_{13C}$ , select CH, CH <sub>2</sub> , CH <sub>3</sub> signals only (all up)	<b>C13DEPT45</b>	dept45
	$\delta_{13C}$ , select CH signals only	<b>C13DEPT90</b>	dept90
	$\delta_{13C}$ , select CH, CH <sub>2</sub> , CH <sub>3</sub> signals only (CH <sub>2</sub> down)	<b>C13DEPT135</b>	dept135
	$\delta_{13C}$ , all <sup>13</sup> C signals (CH, CH <sub>3</sub> up; C, CH <sub>2</sub> down)	<b>1d_13C_DEPTq</b>	deptqgpsp

## Basic 2-D NMR Experiments

Correlation	Information	Method	Parameter Set	Pulse Sequence
	$^1\text{H}/^1\text{H}$ nearest neighbor, via $^3J_{\text{HH}}$	COSY	COSYGPSW	cosygpppqf
	$^1\text{H}/^1\text{H}$ nearest neighbor, via $^3J_{\text{HH}}$ (with $J$ information)	DQF-COSY	COSYGPDPHSW	cosygpmpfphpp
	$^1\text{H}/^1\text{H}$ total spin system, via $^nJ_{\text{HH}}$ ( $n=2,3,4,\dots$ )	TOCSY	MLEVPHSW	mlevphpp
	$^1\text{H}/^{13}\text{C}$ one-bond correlation, via $^1J_{\text{CH}}$	HSQC	HSQCGP	hsqcetgpsi2
	lower resolution in $^{13}\text{C}$ dimension	HMQC	HMQCPH	hmqcph
	$^1\text{H}/^{13}\text{C}$ long-range correlation, via $^nJ_{\text{CH}}$ ( $n=2,3,4$ )	HMBC	HMBCLPND	hmbclpndqf
	$^1\text{H}/^1\text{H}$ through space, via $D_{\text{HH}}$ (small & large molecules)	NOESY	NOESYPHSW	noesyghphpp
	$^1\text{H}/^1\text{H}$ through space, via $D$ -coupling (medium molecules)	ROESY	ROESYPHSW	roesyphpp.2
N/A	Diffusion coefficients vs. molecule sizes, using $^1\text{H}$ NMR	DOSY	2d_DOSY	stebpgplls

## NMR Acronyms

<b>COSY</b>	<b>C</b> ORrelation <b>S</b> pectroscop <b>Y</b>
<b>CPD</b>	<b>C</b> omposite <b>P</b> ulse <b>D</b> ecoupling
<b>DEPT</b>	<b>D</b> istortionless <b>E</b> nhancement by <b>P</b> olarization <b>T</b> ransfer
<b>DEPTQ</b>	<b>DEPT</b> with retention of <b>Q</b> uaternaries
<b>DOSY</b>	<b>D</b> iffusion <b>O</b> rdered <b>S</b> pectroscop <b>Y</b>
<b>DQF</b>	<b>D</b> ouble <b>Q</b> uantum <b>F</b> ilter
<b>HMBC</b>	<b>H</b> eteronuclear <b>M</b> ultiple- <b>B</b> ond <b>C</b> orrelation
<b>HMQC</b>	<b>H</b> eteronuclear <b>M</b> ultiple- <b>Q</b> uantum <b>C</b> orrelation
<b>HSQC</b>	<b>H</b> eteronuclear <b>S</b> ingle- <b>Q</b> uantum <b>C</b> orrelation
<b>NOESY</b>	<b>N</b> uclear <b>O</b> verhauser <b>E</b> ffect <b>S</b> pectroscop <b>Y</b>
<b>ROESY</b>	<b>R</b> otating-frame <b>NOE</b> <b>S</b> pectroscop <b>Y</b>
<b>TOCSY</b>	<b>T</b> Otal <b>C</b> orrelation <b>S</b> pectroscop <b>Y</b>
<b>UDEFT</b>	<b>U</b> niform <b>D</b> riven <b>E</b> quilibrium <b>F</b> ourier <b>T</b> ransform