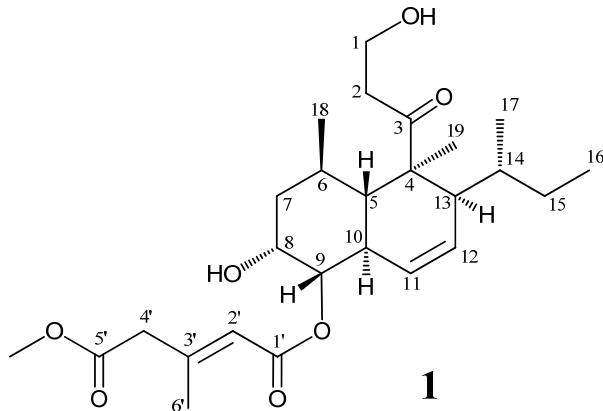


# Supplementary Information

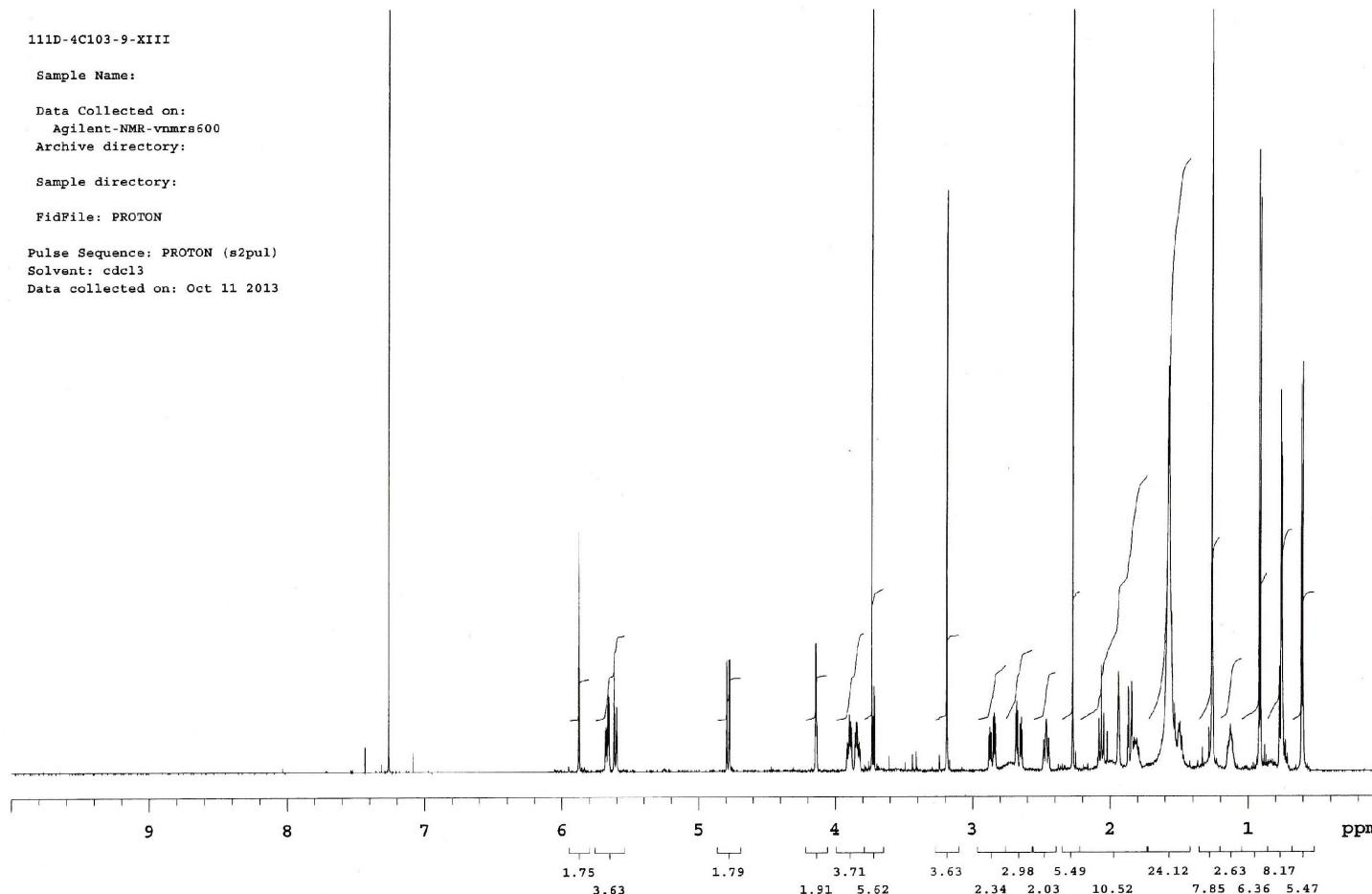


**Figure S1.** Compound 1.

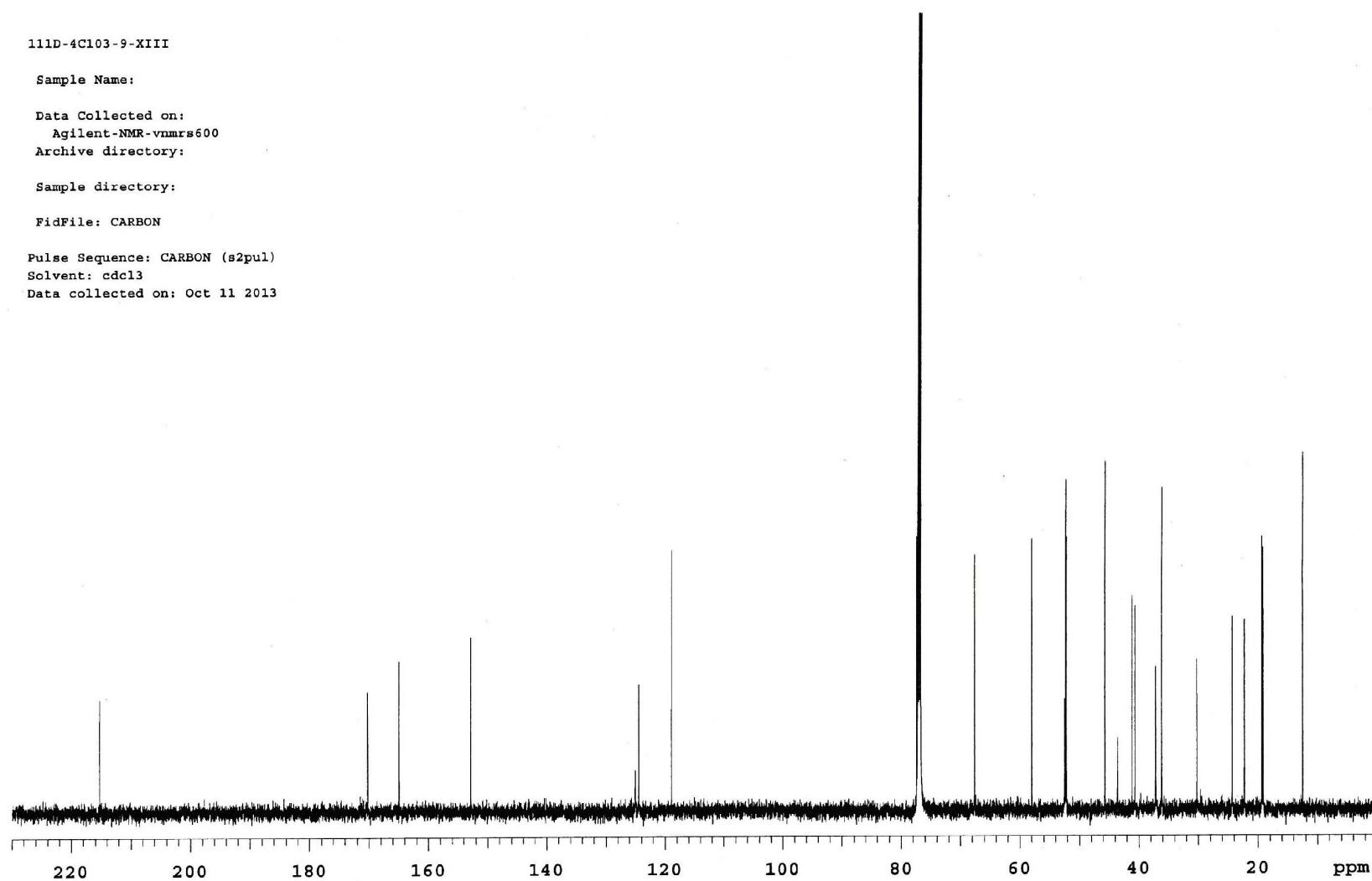
**Table S1.** Spectral data including 2D NMR data for 1.

Position	$\delta_{\text{H}}$ <sup>a</sup>		J/Hz	$^1\text{H}$ - $^1\text{H}$ COSY	NOE <sup>b</sup>	$\delta_{\text{C}}$	HMBC (C) <sup>c</sup>
1A	3.83	ddd	11.5 (2B), 6.2 (2A), 3.5 (2B)	1B, 2A, 2B		58.0 (t)	3
1B	3.89	ddd	11.5 (1B), 6.2 (2B), 3.5 (2A)	1A, 2A, 2B	14		3
2A	2.66	ddd	18.5 (2B), 6.2 (1A), 3.5 (1B)	1A, 1B, 2B	13, 14, 19	41.2 (t)	1, 3
2B	2.87	ddd	18.5 (1B), 6.2 (1B), 3.5 (1A)	1A, 1B, 2A	18, 19		1, 3
3						215.2 (s)	
4						52.5 (s)	
5	2.06	t	11.5 (6, 10)	6, 10	7 $\beta$ , 9, 15B, 16, 18	43.6 (d)	3, 4, 6, 7, 9, 10, 13
6	1.82	m		5, 7 $\beta$ , 18	10, 19	30.3 (d)	
7 $\alpha$	1.86	dt	12.3 (7 $\alpha$ ), 2.5 (6, 8)	7 $\beta$ , 8	18	40.7 (t)	5, 8, 9
7 $\beta$	1.55	td	12.3 (6, 7 $\beta$ ), 2.5 (8)	6, 7 $\alpha$ , 8	5, 9, 18		5, 6
8	4.13	q	2.5 (7 $\alpha$ , 7 $\beta$ , 9)	7 $\alpha$ , 7 $\beta$ , 9		67.6 (d)	6, 10
9	4.78	dd	11.5 (10), 2.5 (8)	8, 10	5, 7 $\beta$ , 11	77.4 (d)	1', 10
10	2.46	tdd	11.5 (5, 9), 5.2 (11), 2.5 (12)	5, 9, 11, 12	6, 19	36.2 (d)	9, 11
11	5.62	br d	10.5 (12)	10, 12, 13	9, 16	125.0 (d)	5, 9, 10, 13
12	5.74	ddd	10.5 (11), 5.2 (13), 2.5 (10)	10, 11, 13	16, 17	124.5 (d)	10, 13
13	1.94	m		11, 12, 14	2A, 17, 19	52.4 (d)	19
14	1.12	m		13, 15A, 15B, 17	1B, 2A, 16	37.2 (t)	
15A	0.74	m		14		24.4 (d)	
15B	1.50	m		14, 16	5		13, 14, 16, 17
16	0.76	t	7.3 (15)	15B	5, 11, 12, 14, 17	12.5 (q)	14, 15
17	0.92	d	6.6 (14)	14	12, 13, 16	19.1 (q)	14, 15, 16
18	0.60	d	7.0 (6)	6	2B, 5, 7 $\alpha$ , 7 $\beta$	22.3 (q)	5, 6, 7
19	1.26	s			2A, 2B, 6, 10, 13	19.4 (q)	3, 4, 5, 13
1'						164.9 (s)	
2'	5.88	s			4'	118.9 (d)	1', 3', 4', 6'
3'						152.0 (s)	
4'	3.19	s			2', 6'	45.7 (t)	2', 3', 5', 6'
5'						170.2 (s)	
6'	2.27	s			4'	19.2 (q)	2', 3', 4'
5'-OCH <sub>3</sub>	3.73	s				52.2 (q)	5'

<sup>a</sup>  $^1\text{H}$  chemical shift values ( $\delta$  ppm from SiMe<sub>4</sub>) followed by multiplicity and then the coupling constants (J/Hz). Figures in parentheses indicate the proton coupling with that position; <sup>b</sup> The correlations with geminal and vicinal protones are removed; <sup>c</sup> Long range  $^1\text{H}$ - $^{13}\text{C}$  correlations from H to C observed in the HMBC experiment.



**Figure S2.**  $^1\text{H}$  NMR spectrum of 1.



**Figure S3.**  $^{13}\text{C}$  NMR spectrum of 1.

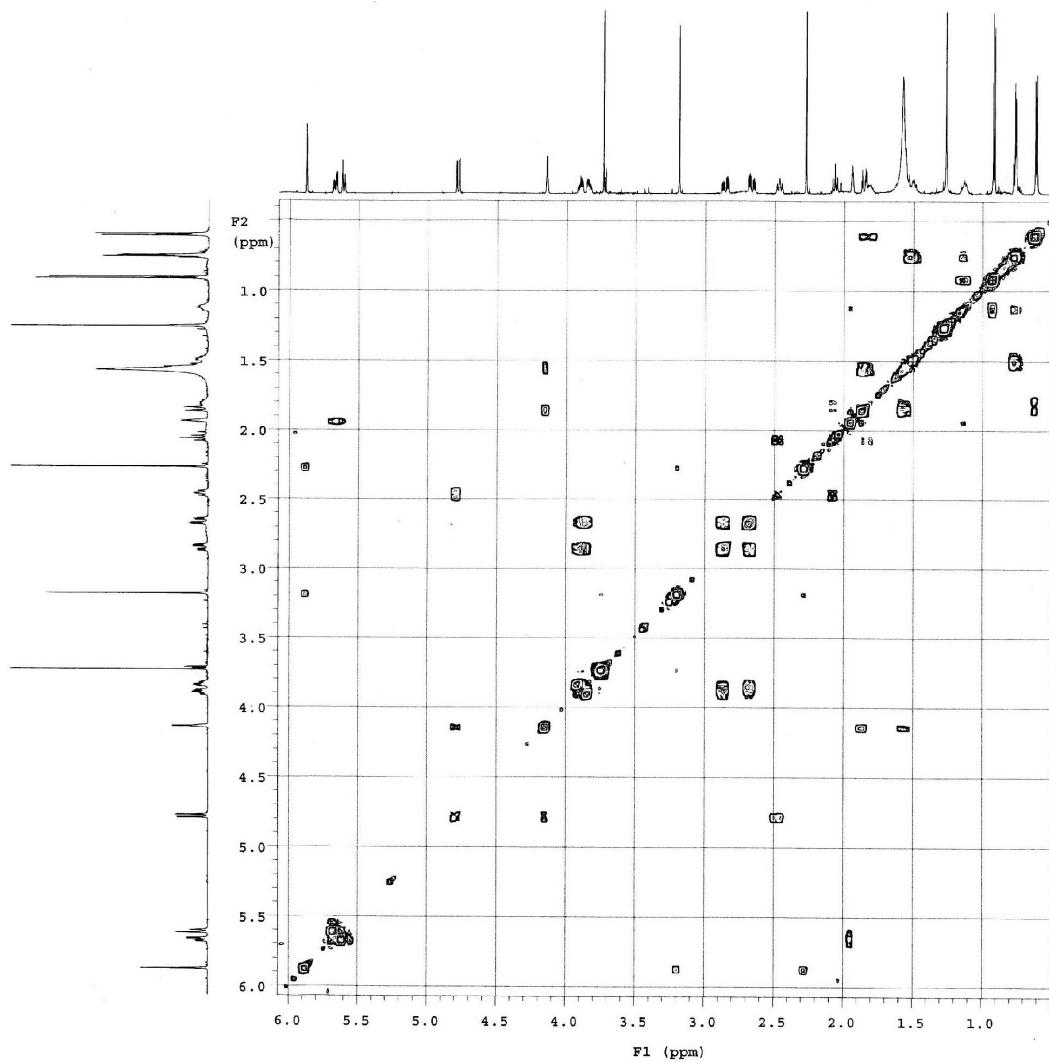
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111D-4C103-9-XXXI
exp33 gCOSY

SAMPLE          FLAGS
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solvent edels3 sspl     y
sample hsgv1      6540

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sw    4699.2 temp  not used
at    0.150 gain   48
np    1410 spin   0
tb    4000 P2 PROCESSING
ss    32 ab    -0.075
d1    1.000 abs   not used
nt    16 tn    2048
2D ACQUISITION  F1 PROCESSING
sw1   4699.2 ab1  -0.027
ni    256 abel  not used
d2    0 proc1 lp
PRESATURATION  f1l   2048
satmode n      DISPLAY
wst    n sp    217.2
TRANSMITTER wp    3423.5
tn    H1 sp1  221.8
sfrq  599.898 wpl  3423.5
tof   -792.3 rfl   140.7
tpwr  58 rfp   0
pw    9.100 rfl1  140.7
GRADIENTS rfp1  0
gzlvR  5458 PLOT
gtE   0.001000 wc   200.0
EDratio 1.000 sc   0
gstab  0.000500 wc2  200.0
DECOUPLER sc2   0
dn    C13 vs   583
dm    nnn th   2
ai    cdc av

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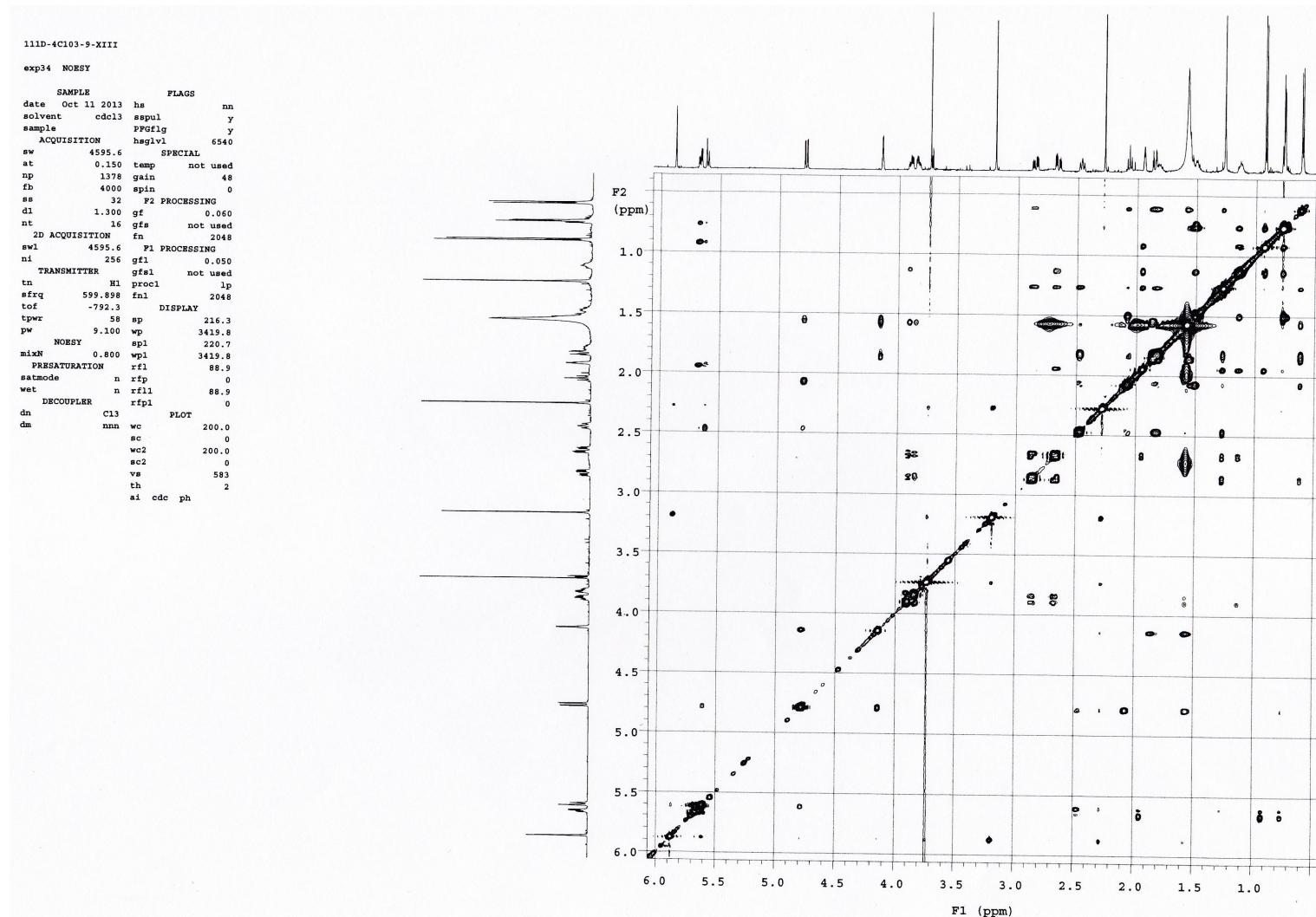


Figure S5. NOESY of 1.

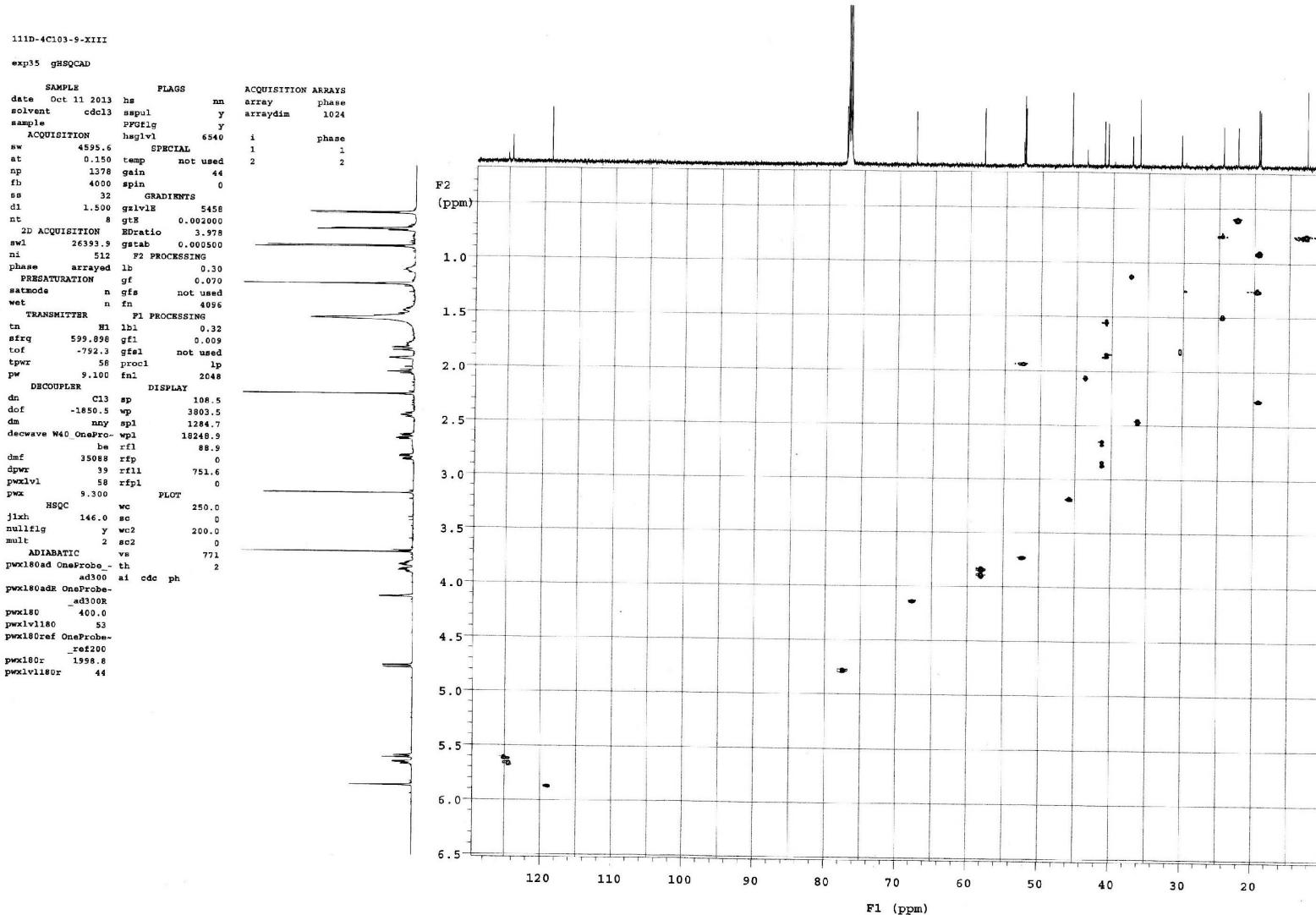


Figure S6. HSQC of 1.

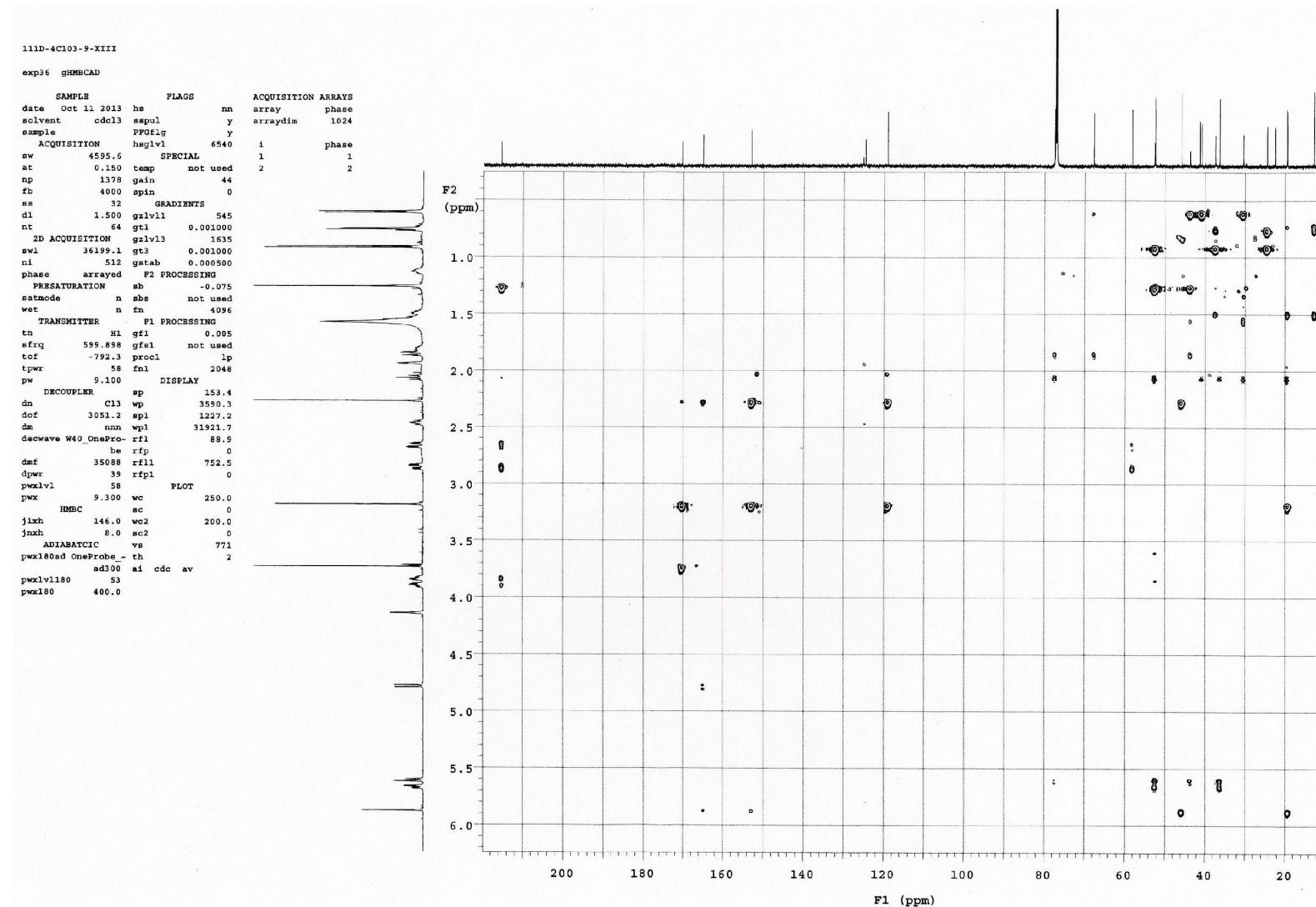
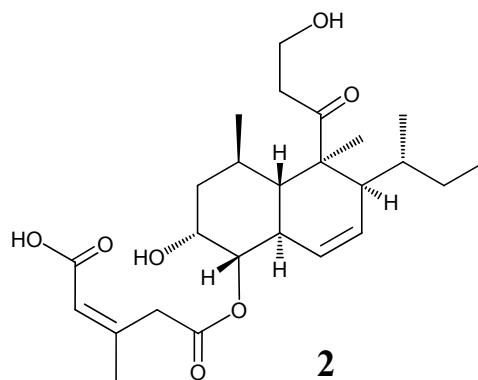
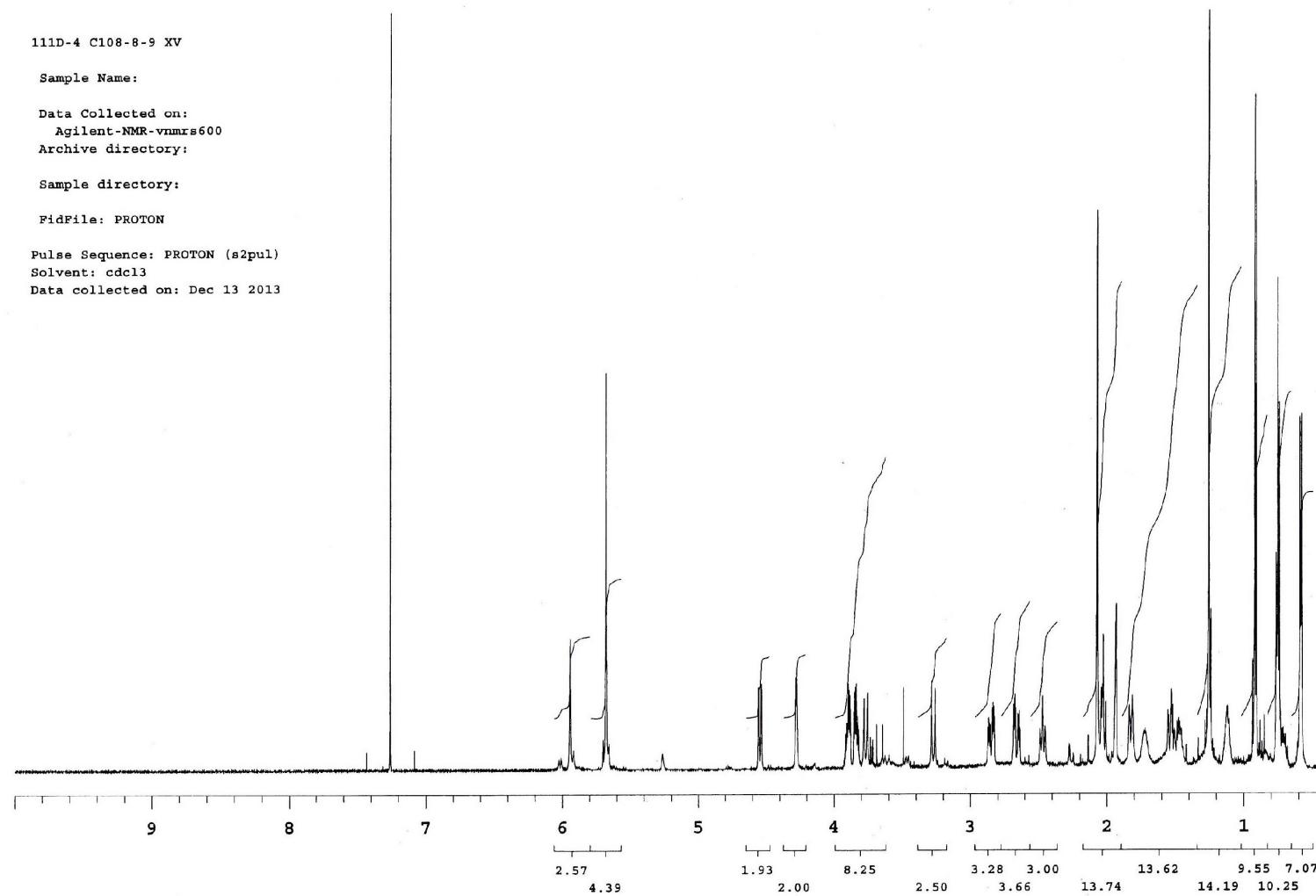


Figure S7. HMBC of 1.

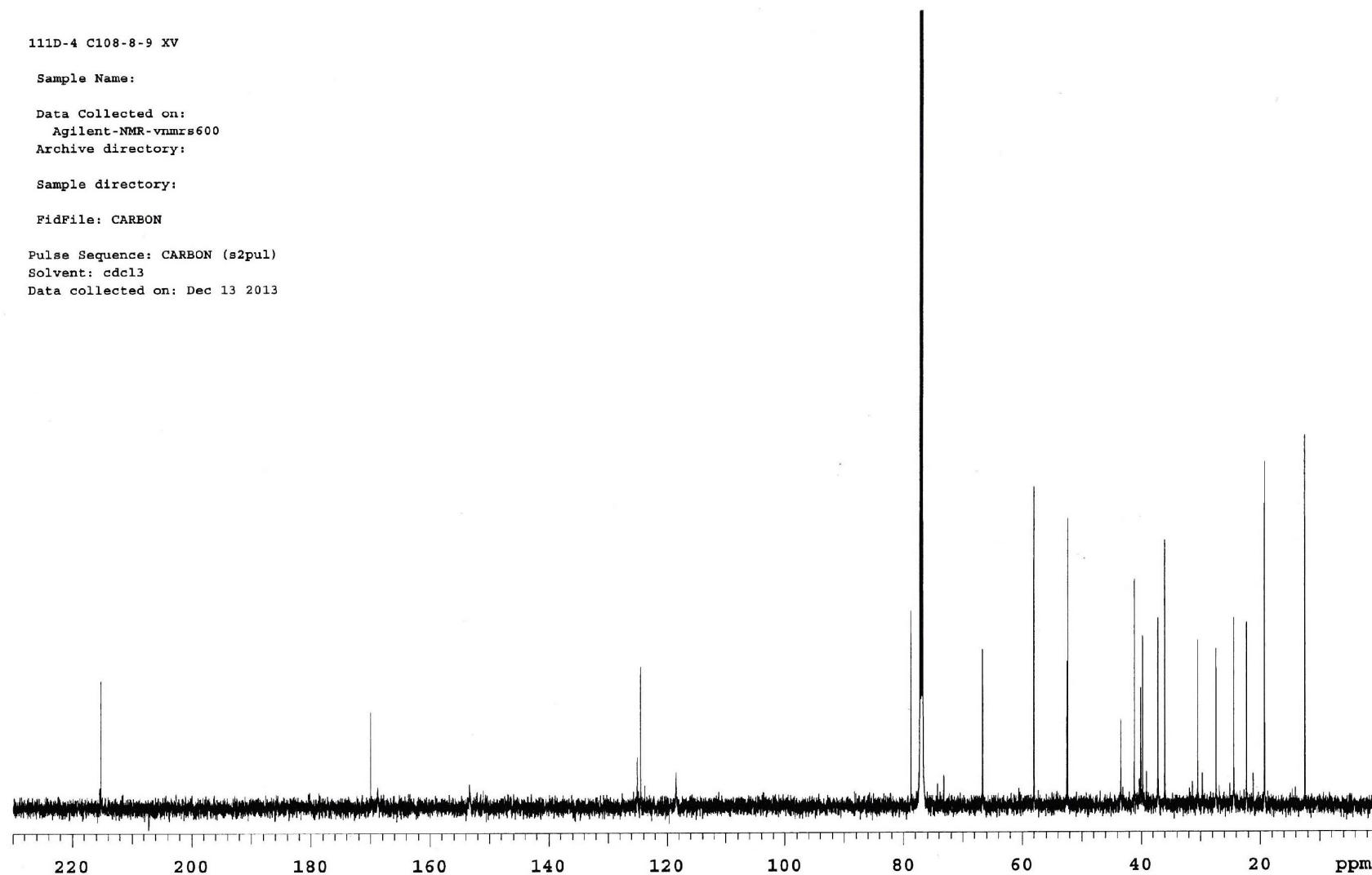
**Figure S8.** Compound 2.**Table S2.** Spectral data including 2D NMR data for 2.

Position	$\delta_{\text{H}}$ <sup>a</sup>	J/Hz	<sup>1</sup> H- <sup>1</sup> H COSY	NOE <sup>b</sup>	$\delta_{\text{C}}$	HMBC (C) <sup>c</sup>
1A	3.84	ddd	10.8 (2B), 6.5 (2A), 3.5 (2B)	1B, 2A, 2B	58.0	(t) 2, 3
1B	3.90	ddd	10.8 (1B), 7.5 (2B), 3.5 (2A)	1A, 2A, 2B	14	2, 3
2A	2.67	ddd	18.5 (2B), 6.5 (1A), 3.5 (1B)	1A, 1B, 2B	13, 14, 19	41.1 (t) 1, 3
2B	2.85	ddd	18.5 (1B), 7.5 (1B), 3.5 (1A)	1A, 1B, 2A	18, 19	1, 3
3					215.2	(s)
4					52.5	(s)
5	2.03	t	10.8 (6, 10)	6, 10	7 $\beta$ , 9, 15B, 18	43.4 (d) 3, 4, 6, 7, 9, 10, 13, 19
6	1.73	m		5, 7 $\beta$ , 18	10, 18, 19	30.5 (d)
7 $\alpha$	1.83	dt	12.0 (7 $\alpha$ ), 3.0 (6, 8)	7 $\beta$ , 8	18	40.1 (t) 5, 8, 9
7 $\beta$	1.53	td	12.0 (6, 7 $\beta$ ), 3.0 (8)	6, 7 $\alpha$ , 8	5, 9, 18	6
8	4.28	q	3.0 (7 $\alpha$ , 7 $\beta$ , 9)	7 $\alpha$ , 7 $\beta$ , 9	66.7	(d)
9	4.55	dd	12.5 (10), 3.0 (8)	8, 10	5, 7 $\beta$ , 11	78.8 (d) 1'
10	2.47	br t	12.5 (5, 9)	5, 9	6, 19	36.0 (d) 5, 6, 9
11	5.69	dr d	10.8 (12)		9	125.1 (d) 5, 9, 10, 13
12	5.67	dd	10.8 (11), 4.2 (13)	13	16, 17	124.5 (d) 4, 10, 13
13	1.94	m		12, 14	2A, 17, 19	52.3 (d) 4, 5, 14, 15, 17, 19
14	1.12	m		13, 15A, 17	1B, 2A, 16	37.2 (t)
15A	0.72	m		14, 15B, 17	17	24.5 (d) 14, 16, 17
15B	1.47	m		15A, 16	5	14, 16, 17
16	0.75	t	7.1 (15)	15B	12, 14, 17	12.5 (q) 14, 15
17	0.92	d	7.1 (14)	14	12, 13, 15A, 16	19.2 (q) 14, 15, 16
18	0.59	d	6.0 (6)	6	2B, 5, 7 $\alpha$ , 7 $\beta$	22.3 (q) 5, 6, 7
19	1.26	s			2A, 2B, 6, 10, 13	19.3 (q) 3, 4, 5, 13
1'					170.0	(s)
2'A	3.29	d	15.2 (2'B)	2'B	6'	39.7 (t) 1', 3', 4', 6'
2'B	3.77	d	15.2 (2'A)	2'A		1', 3', 4', 6'
3'					153.4	(s)
4'	5.94	s			6'	118.5 (d) 2', 5', 6'
5'					168.9	(s)
6'	2.08	s		2'A, 4'	27.4	(q) 2', 3', 4'

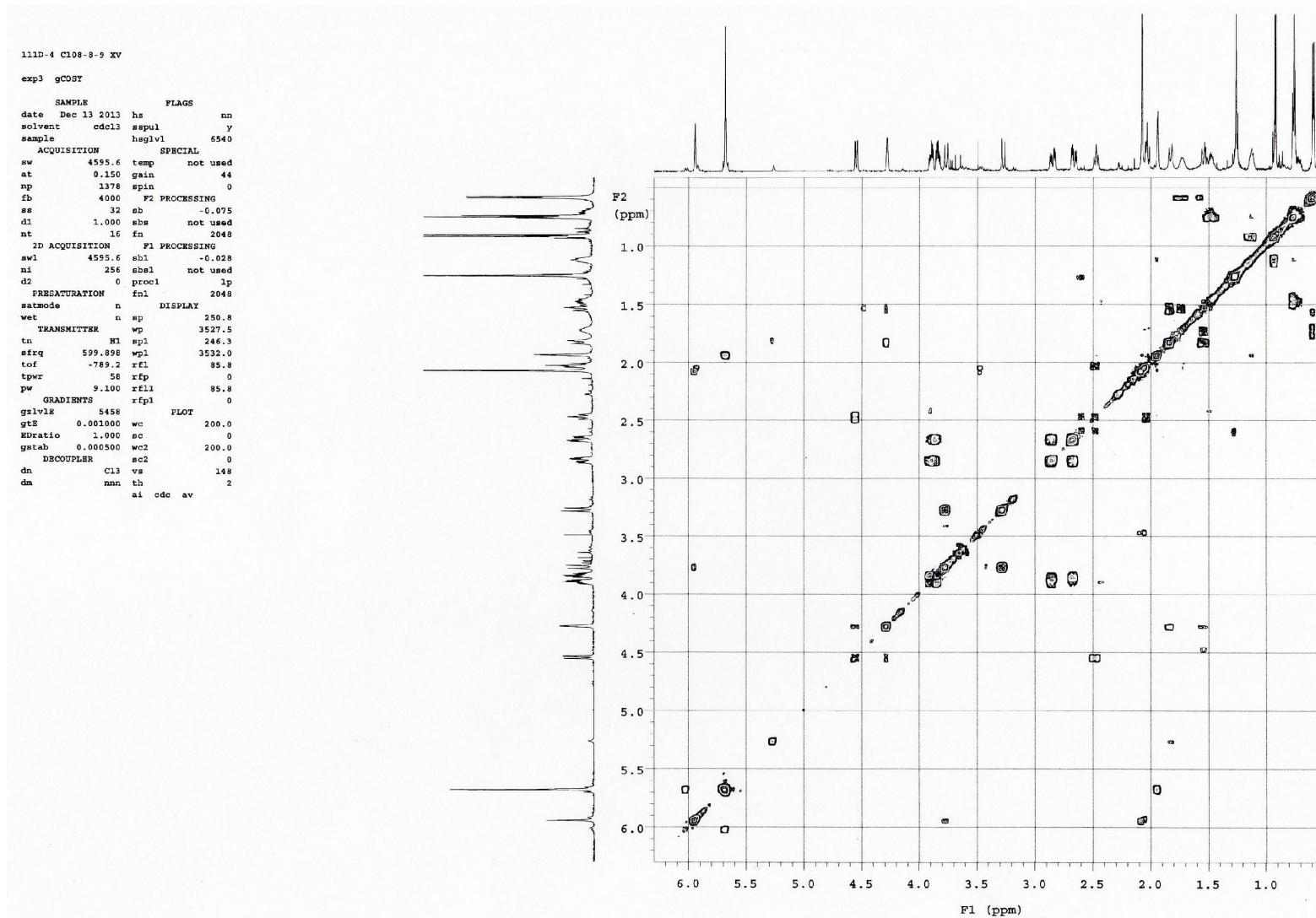
<sup>a</sup> <sup>1</sup>H chemical shift values ( $\delta$  ppm from SiMe<sub>4</sub>) followed by multiplicity and then the coupling constants (J/Hz). Figures in parentheses indicate the proton coupling with that position; <sup>b</sup> The correlations with geminal and vicinal protones are removed; <sup>c</sup> Long range <sup>1</sup>H-<sup>13</sup>C correlations from H to C observed in the HMBC experiment.



**Figure S9.**  $^1\text{H}$  NMR spectrum of 2.



**Figure S10.**  $^{13}\text{C}$  NMR spectrum of 2.

Figure S11.  $^1\text{H}$ - $^1\text{H}$  COSY of 2.

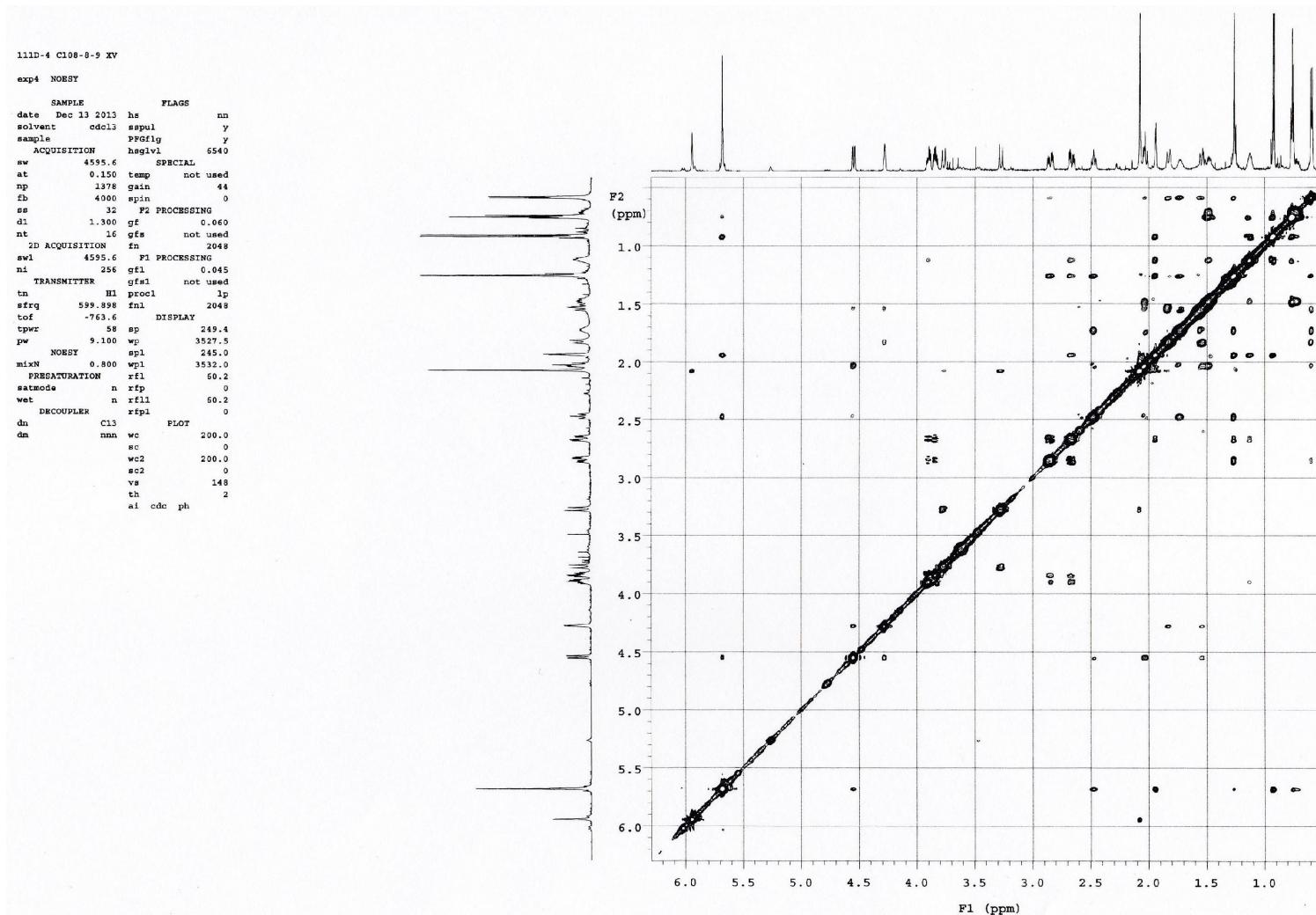


Figure S12. NOESY of 2.

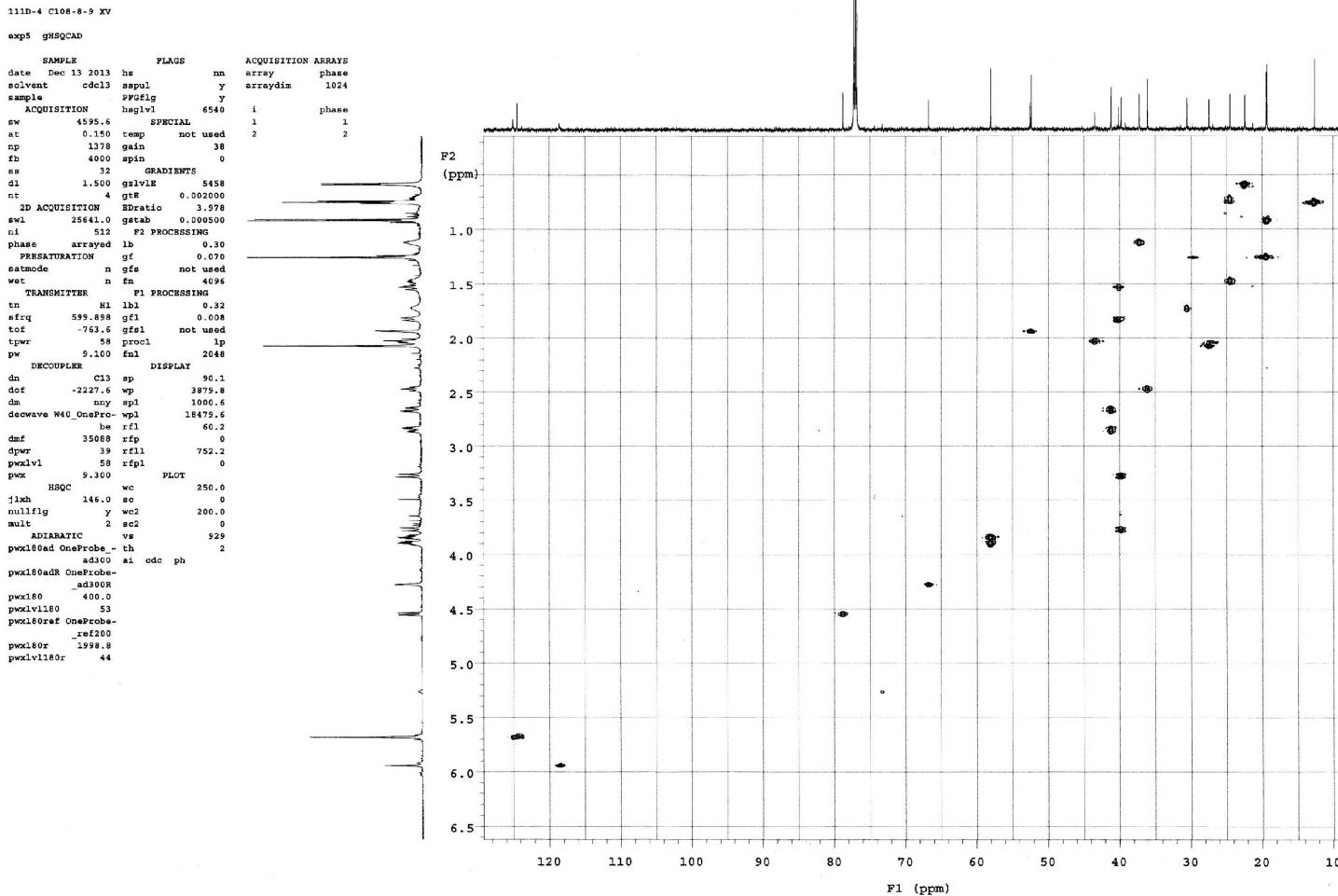


Figure S13. HSQC of 2.

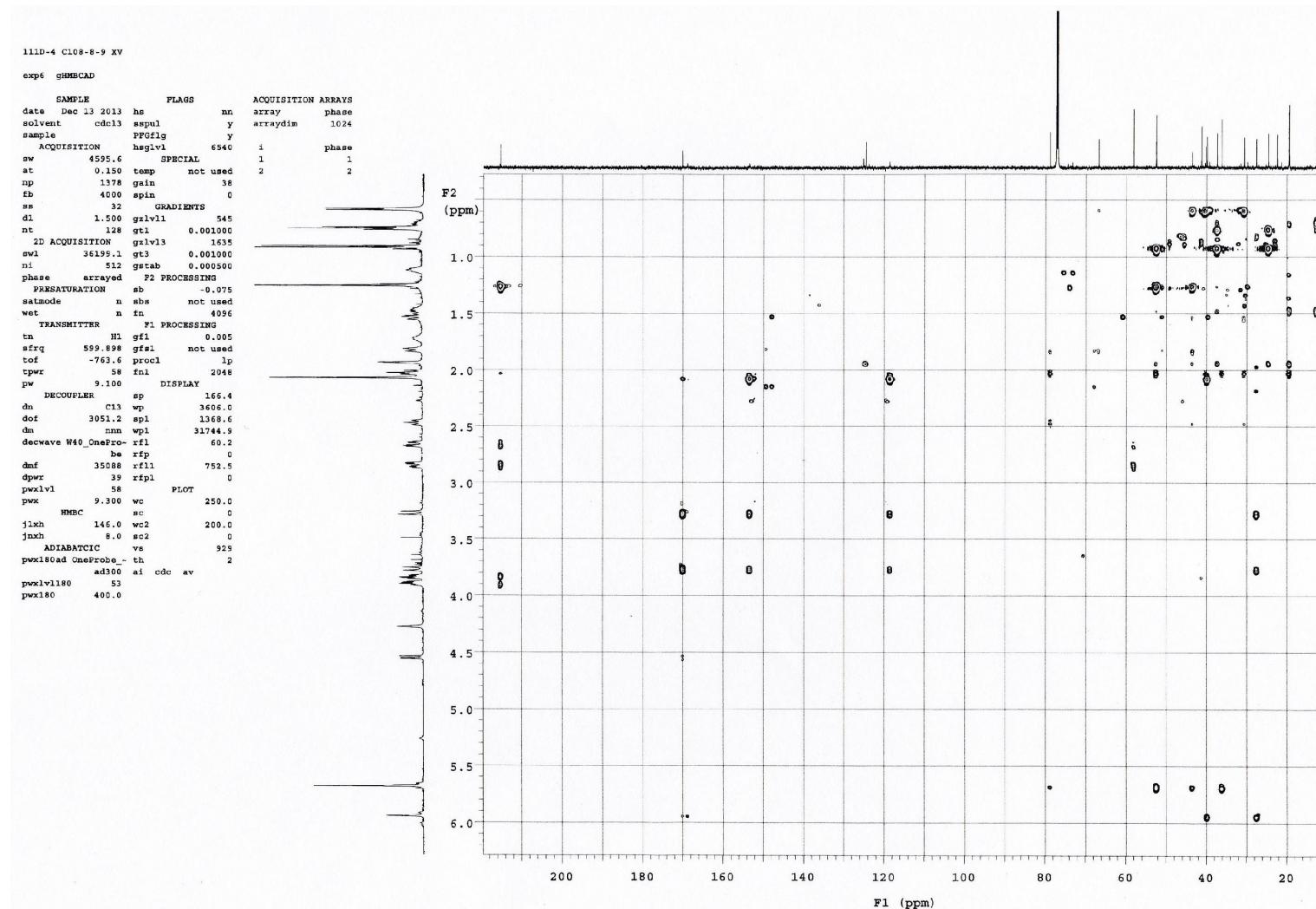
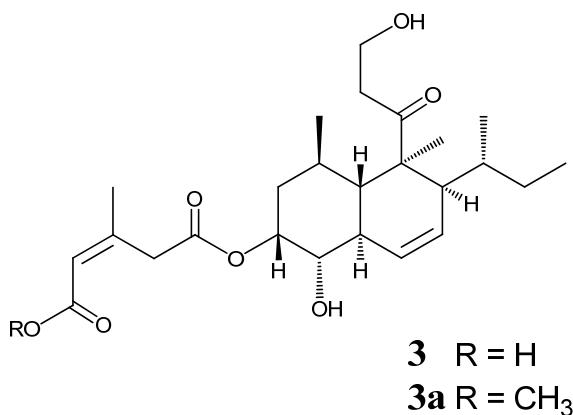
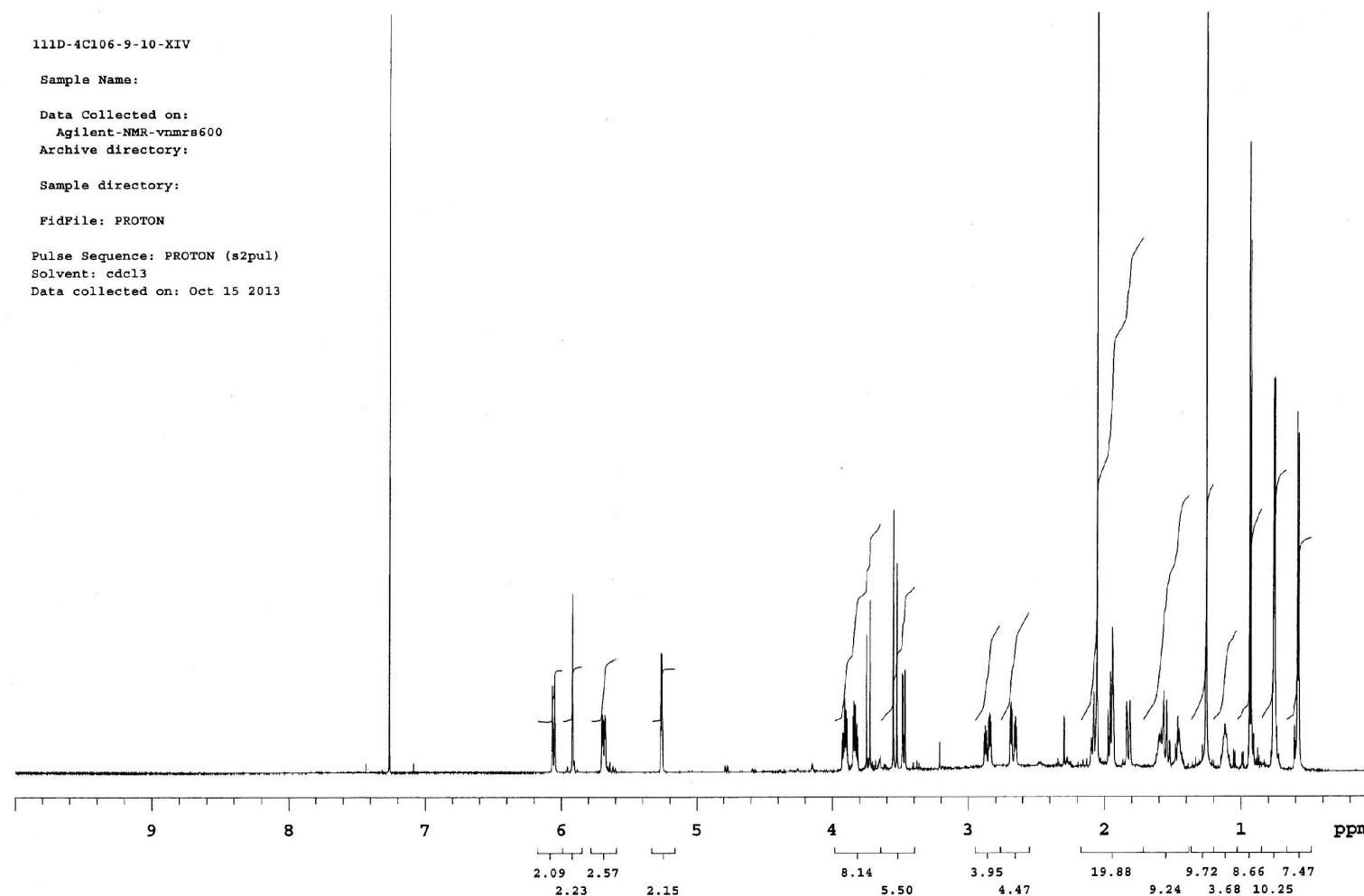


Figure S14. HMBC of 2.

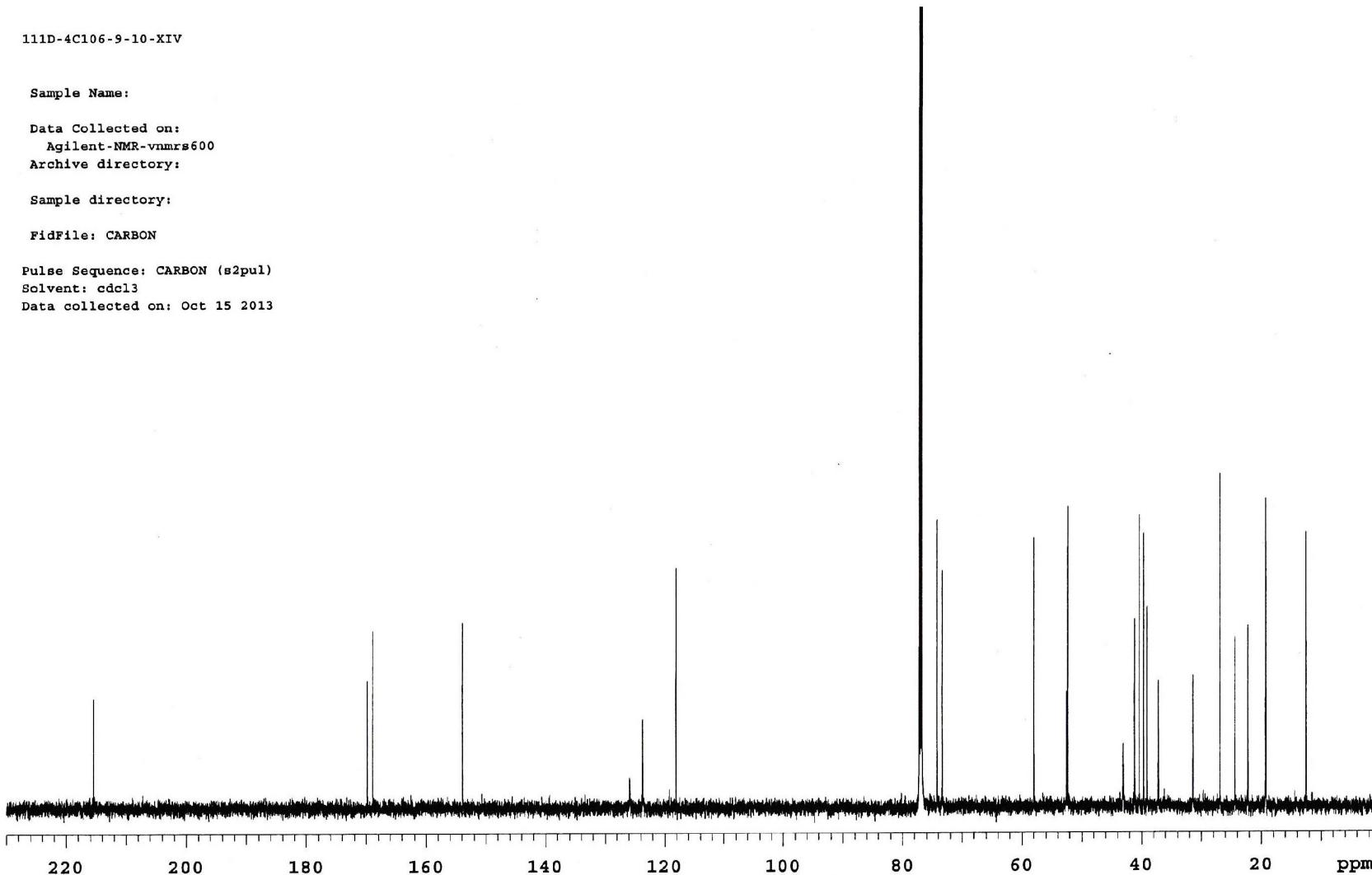
**Figure S15.** Compound 3.**Table S3.** Spectral data including 2D NMR data for 3.

Position	$\delta_{\text{H}}$ <sup>a</sup>	J/Hz	<sup>1</sup> H- <sup>1</sup> H COSY	NOE <sup>b</sup>	$\delta_{\text{C}}$	HMBC (C) <sup>c</sup>
1A	3.83	ddd	11.8 (2B), 6.5 (2A), 3.5 (2B)	1B, 2A, 2B	58.0	(t) 3
1B	3.91	ddd	11.8 (1B), 6.5 (2B), 3.5 (2A)	1A, 2A, 2B	14	3
2A	2.67	ddd	18.5 (2B), 6.5 (1A), 3.5 (1B)	1A, 1B, 2B	13, 14, 19	41.2 (t) 1, 3
2B	2.86	ddd	18.5 (1B), 6.5 (1B), 3.5 (1A)	1A, 1B, 2A	18, 19	1, 3
3					215.2	(s)
4					52.5	(s)
5	1.96	t	10.8 (6, 10)	6, 10	7 $\beta$ , 9, 15B, 18	43.1 (d) 3, 4, 6, 7, 9, 10, 13, 19
6	1.59	m		5, 7 $\beta$ , 18	10, 19	31.4 (d) 18
7 $\alpha$	1.83	dt	12.0 (7 $\alpha$ ), 3.0 (6, 8)	7 $\beta$ , 8	18	39.1 (t) 5, 8, 9
7 $\beta$	1.55	td	12.0 (6, 7 $\beta$ ), 3.0 (8)	6, 7 $\alpha$ , 8	5, 9, 18	6
8	5.26	q	3.0 (7 $\alpha$ , 7 $\beta$ , 9)	7 $\alpha$ , 7 $\beta$ , 9	73.3	(d)
9	3.48	dd	10.8 (10), 3.0 (8)	8, 10	5, 7 $\beta$ , 11	74.2 (d) 8, 10
10	2.08	tdd	10.8 (5, 9), 4.2 (11), 2.4 (12)	5	6, 19	40.4 (d) 9
11	6.06	dt	10.8 (12), 4.2 (11, 13)	12, 13	9	125.9 (d) 5, 9, 10, 13
12	5.69	ddd	10.8 (11), 4.2 (13), 2.4 (10)	11, 13	16, 17	123.7 (d) 10, 13, 14
13	1.94	m		11, 12, 14	2A, 17, 19	52.4 (d) 4, 12, 19
14	1.12	m		13, 15A, 17	1B, 2A, 16	37.2 (t)
15A	0.74	m		14	17	24.4 (d) 16
15B	1.47	m		16	5	14, 16, 17
16	0.76	t	7.3 (15)	15B	12, 14, 17	12.5 (q) 14, 15
17	0.93	d	7.2 (14)	14	12, 13, 15A, 16	19.2 (q) 14, 15, 16
18	0.59	d	6.0 (6)	6	2B, 5, 7 $\alpha$ , 7 $\beta$	22.3 (q) 5, 6, 7
19	1.25	s			2A, 2B, 6, 10, 13	19.3 (q) 3, 4, 5, 13
1'					169.8	(s)
2'A	3.54	d	15.2 (2'B)		6'	39.6 (t) 1', 3', 4', 6'
2'B	3.73	d	15.2 (2'A)		6'	1', 3', 4', 6'
3'					153.9	(s)
4	5.92	s			6'	118.1 (d) 2', 5', 6'
5'					168.9	(s)
6'	2.05	s		2'A, 2'B, 4'	26.9	(q) 2', 3', 4'

<sup>a</sup> <sup>1</sup>H chemical shift values ( $\delta$  ppm from SiMe<sub>4</sub>) followed by multiplicity and then the coupling constants (J/Hz). Figures in parentheses indicate the proton coupling with that position; <sup>b</sup> The correlations with geminal and vicinal protones are removed; <sup>c</sup> Long range <sup>1</sup>H-<sup>13</sup>C correlations from H to C observed in the HMBC experiment.



**Figure S16.**  $^1\text{H}$  NMR spectrum of 3.



**Figure S17.**  $^{13}\text{C}$  NMR spectrum of 3.

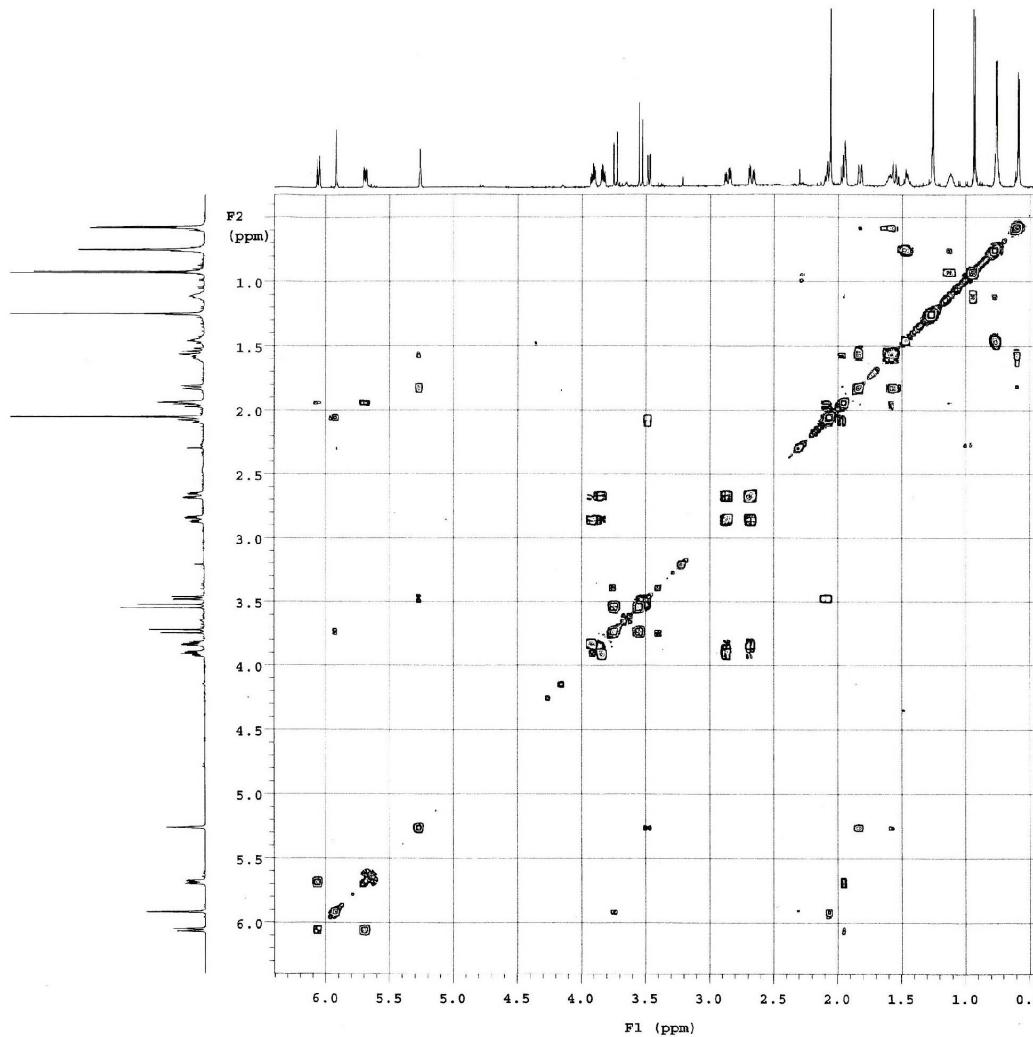
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111D-4C106-9-10-XIV
exp33 gCOSY

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solvent cdc13 spul    y
sample hgv1v1  6540

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at     0.150 gain     42
np     1368 swin    0
fb     4000 F2 PROCESSING
se     32 ab     -0.075
dl     1.000 abs    not used
nt     8 fn     2048
      2D ACQUISITION  F1 PROCESSING
sw1    4562.0 sb1    -0.028
ni     256 sbel    not used
d2     0 proc1    1p
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satmode n      DISPLAY
wt     n      sp    197.0
      TRANSMITTER        wP    3639.8
tn     H1 spl    192.6
sfrq   599.898 wpl    3648.7
tof    -751.0 rf1    70.3
tpwr   58 rfp    0
pw     9.100 rf11   70.3
      GRADIENTS         rf11  0
g1v1R  5456 PLOT
gtk    0.0001000 wc    200.0
EDratio 1.000 sc    0
getab  0.0005000 wc2   200.0
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dn     C13 vs    316
dm     mn1 th    2
ai     cdc av

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**Figure S18.**  $^1\text{H}$ - $^1\text{H}$  COSY of 3.

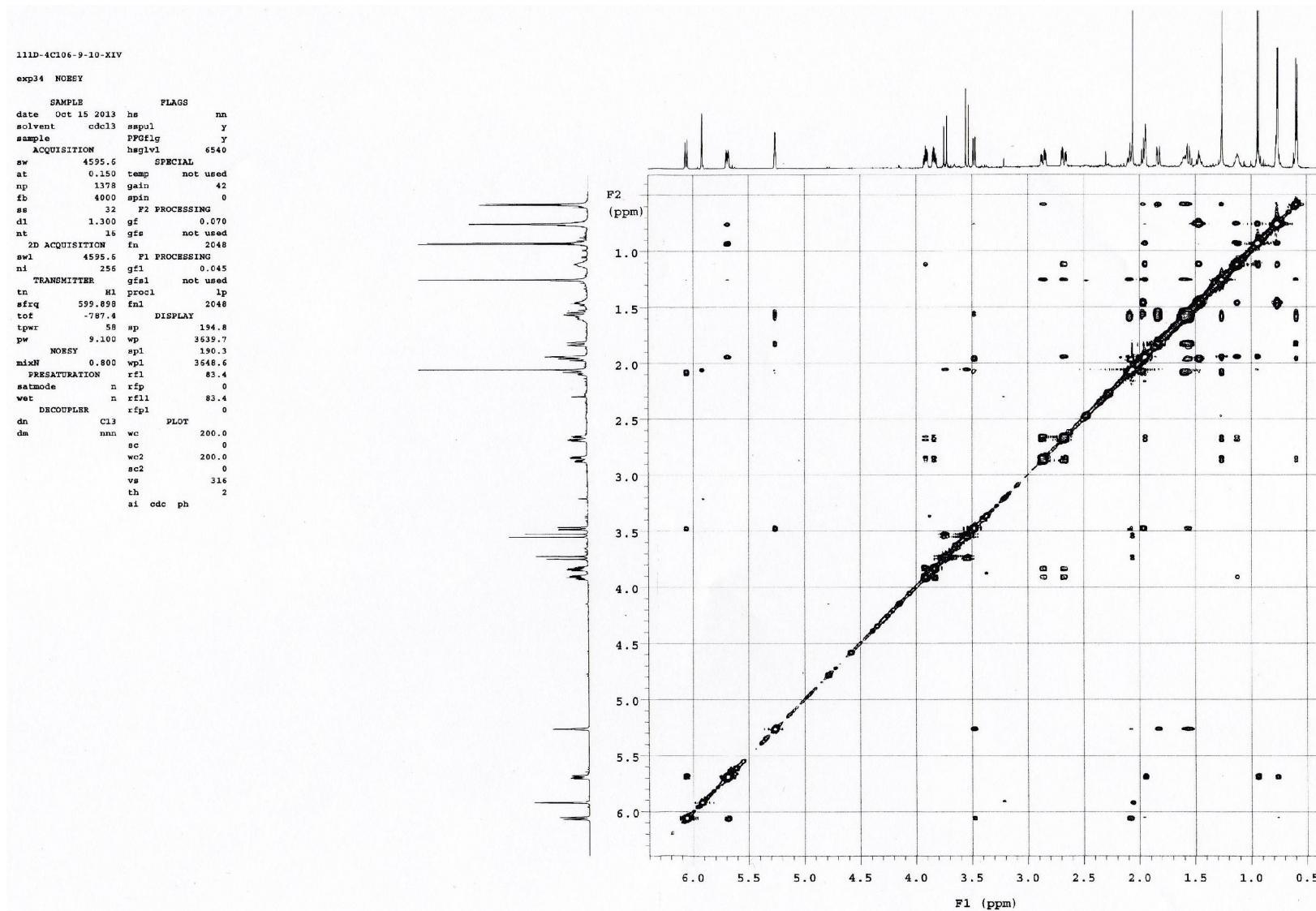


Figure S19. NOESY of 3.

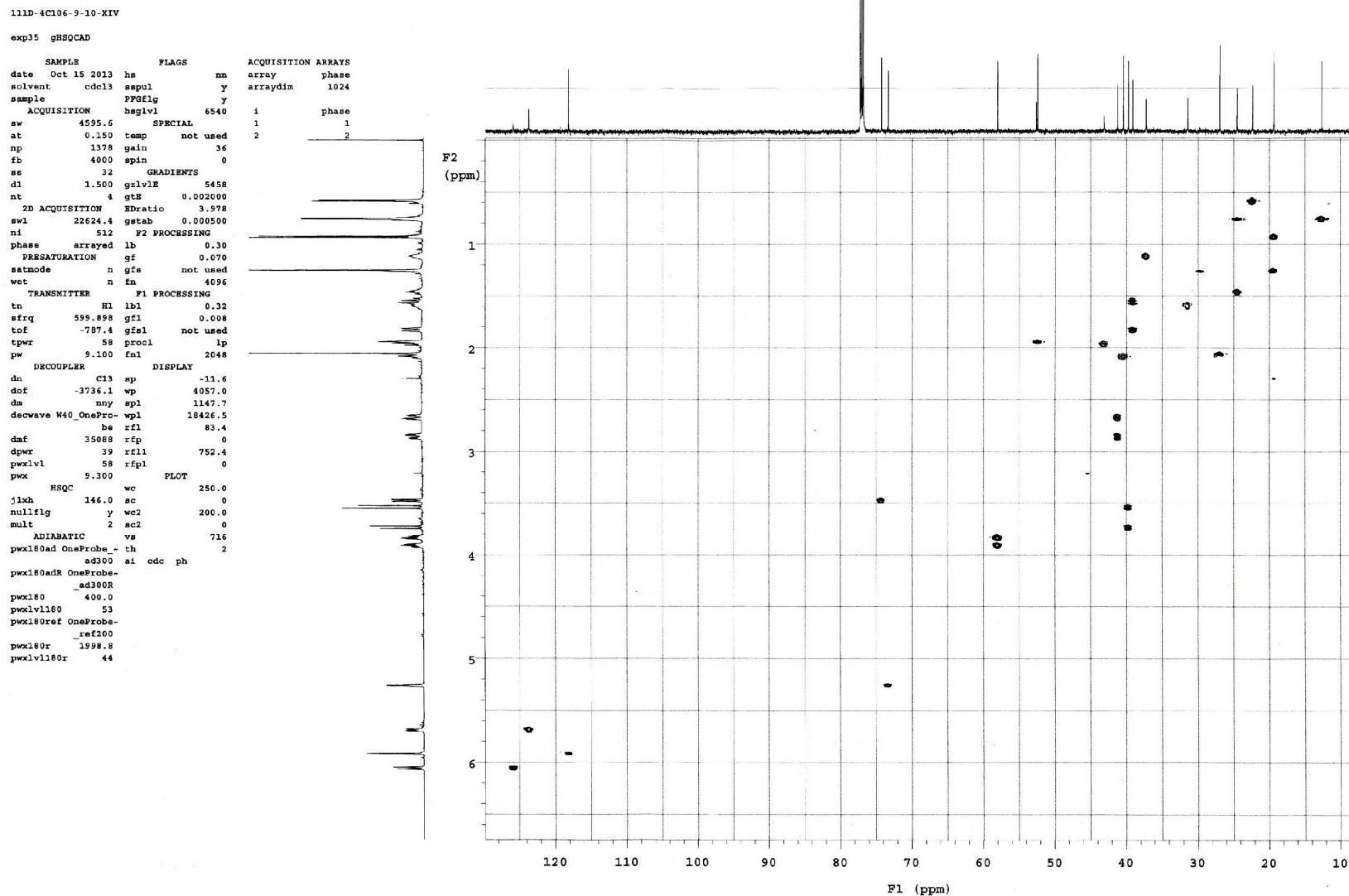


Figure S20. HSQC of 3.

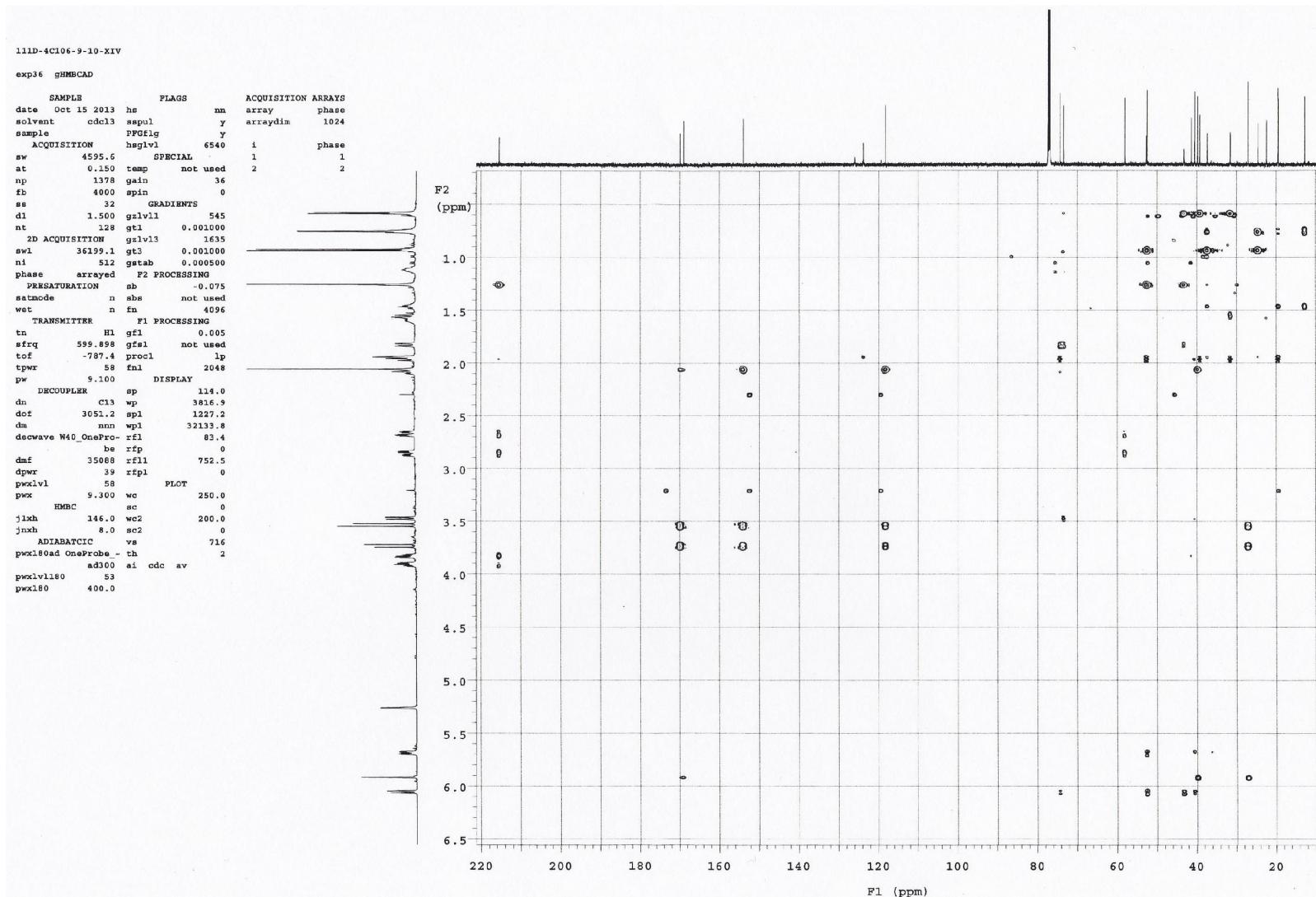
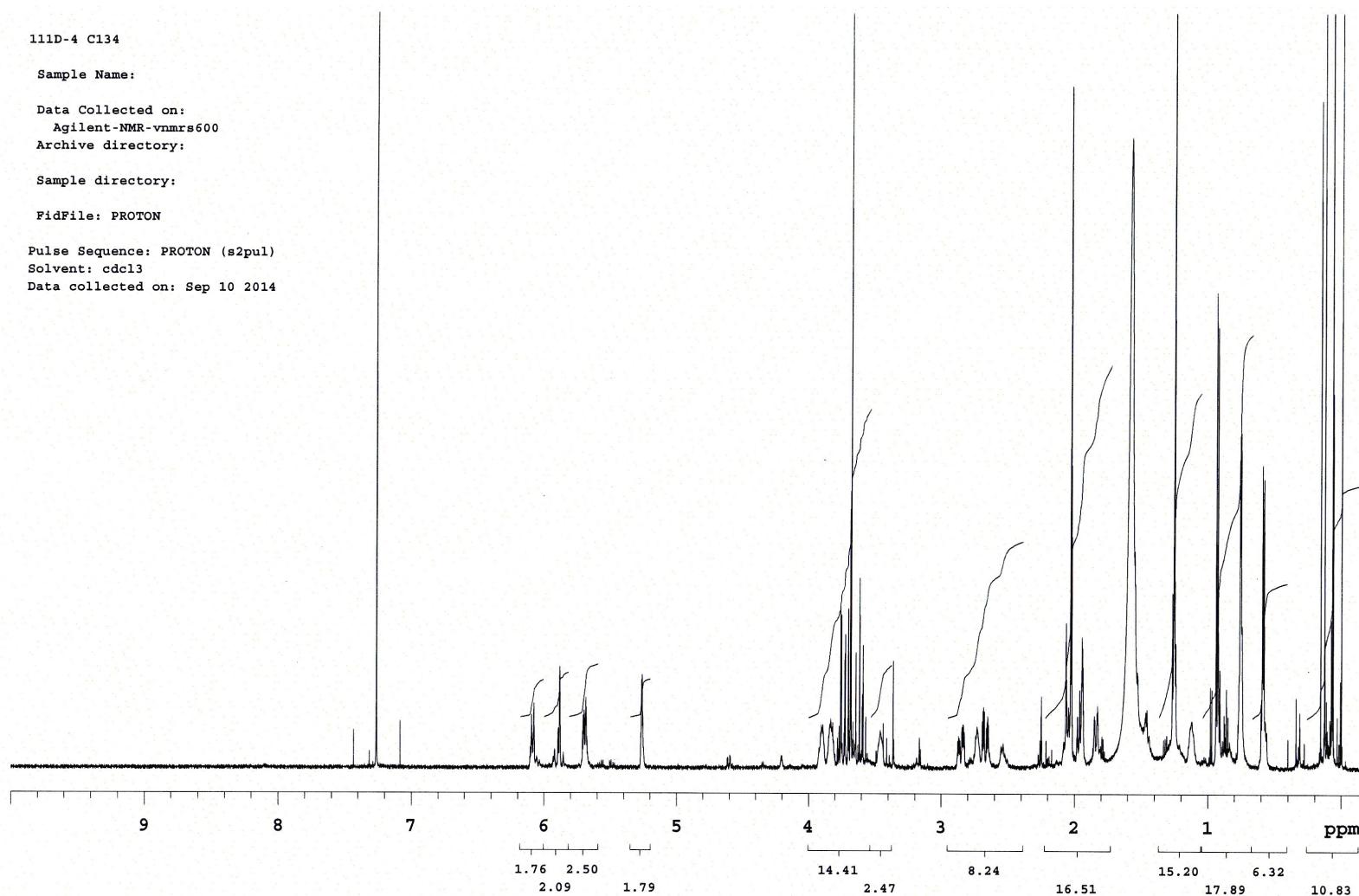


Figure S21. HMBC of 3.

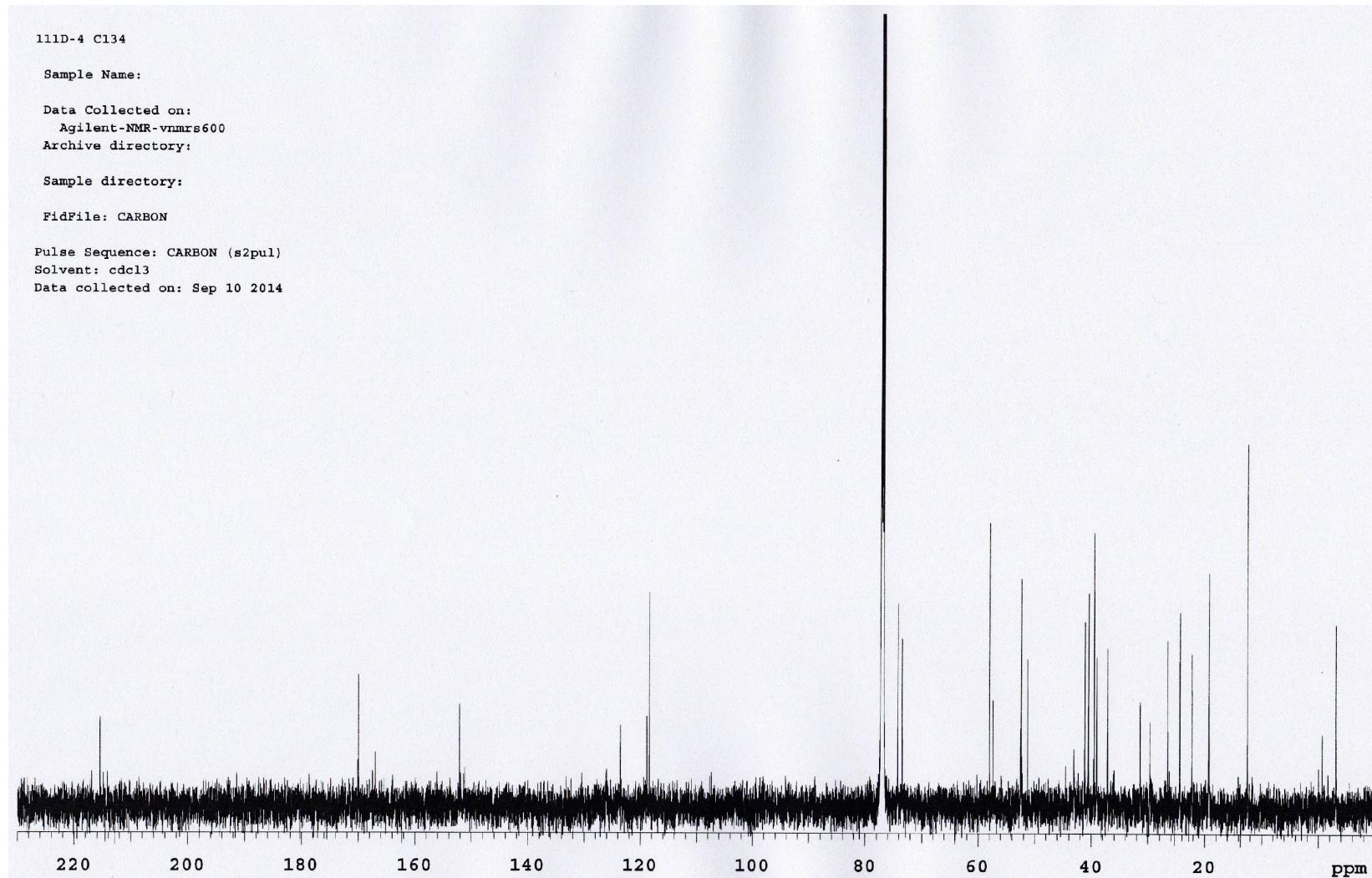
**Table S4.** Spectral data including 2D NMR data for 3a.

Position	$\delta_{\text{H}}$ <sup>a</sup>	J/Hz	$^1\text{H}$ - $^1\text{H}$ COSY	NOE <sup>b</sup>	$\delta_{\text{C}}$	HMBC (C) <sup>c</sup>
1A	3.84	m		1B, 2A, 2B	58.0	(t)
1B	3.90	m		1A, 2A, 2B		
2A	2.67	ddd	18.5 (2B), 6.0 (1A), 3.0 (1B)	1A, 1B, 2B	13, 14, 19	41.2 (t)
2B	2.85	ddd	18.5 (1B), 6.0 (1B), 3.0 (1A)	1A, 1B, 2A	18, 19	3, 1, 3
3					215.5	(s)
4					52.6	(s)
5	1.96	t	10.2 (6, 10)	6, 10	7 $\beta$ , 9, 15B, 18	43.1 (d)
6	1.60	m		5, 7 $\alpha$ , 18	31.4	(d)
7 $\alpha$	1.84	dt	12.0 (7 $\alpha$ ), 3.0 (6, 8)	6, 7 $\beta$ , 8	18	39.1 (t)
7 $\beta$	1.58	m		7 $\alpha$ , 8	5, 9, 18	6
8	5.26	q	3.0 (7 $\alpha$ , 7 $\beta$ , 9)	7 $\alpha$ , 7 $\beta$ , 9	73.5	(d)
9	3.46	m		8, 10	5, 7 $\beta$ , 11	74.2 (d)
10	2.06	m		5	19	40.5 (d)
11	6.09	m		12	9	126.0 (d)
12	5.69	m		11, 13	16, 17	123.6 (d)
13	1.94	m		12	2A, 17, 19	52.4 (d)
14	1.12	m		15A, 17	2A, 16	37.2 (t)
15A	0.76	m		14		24.4 (d)
15B	1.46	m		16	5	16, 17
16	0.76	t	7.3 (15)	15B	12, 14, 17	12.5 (q)
17	0.93	d	7.2 (14)	14	12, 13, 16	19.2 (q)
18	0.59	d	6.0 (6)	6	2B, 5, 7 $\alpha$ , 7 $\beta$	22.3 (q)
19	1.26	s			2A, 2B, 10, 13	5, 6, 7
1'					19.3 (q)	3, 4, 5, 13
2'A	3.60	d	15.2 (2'B)	2'B	6'	170.0 (s)
2'B	3.72	d	15.2 (2'A)	2'A	6'	14, 15
3'						1', 3', 4', 6'
4	5.88	s			6'	1', 3', 4', 6'
5'						2', 5', 6'
6'	2.03	s			2'A, 2'B, 4'	167.0 (s)
5'-OCH <sub>3</sub>	3.68		s		26.6 (q)	2', 3', 4'
					51.3 (q), 5'	

<sup>a</sup>  $^1\text{H}$  chemical shift values ( $\delta$  ppm from SiMe<sub>4</sub>) followed by multiplicity and then the coupling constants (J/Hz). Figures in parentheses indicate the proton coupling with that position; <sup>b</sup> The correlations with geminal and vicinal protones are removed; <sup>c</sup> Long range  $^1\text{H}$ - $^{13}\text{C}$  correlations from H to C observed in the HMBC experiment.



**Figure S22.**  $^1\text{H}$  NMR spectrum of 3a.



**Figure S23.**  $^{13}\text{C}$  NMR spectrum of 3a.