

Synthesis of Polysubstituted Cyclopentene and Cyclopenta[*b*]carbazole Analogues from Unsymmetrical 4-Arylidene-3,6-diarylhex-2-en-5-ynal and Indole Derivatives via an Iodine Mediated Electrocyclization Reaction

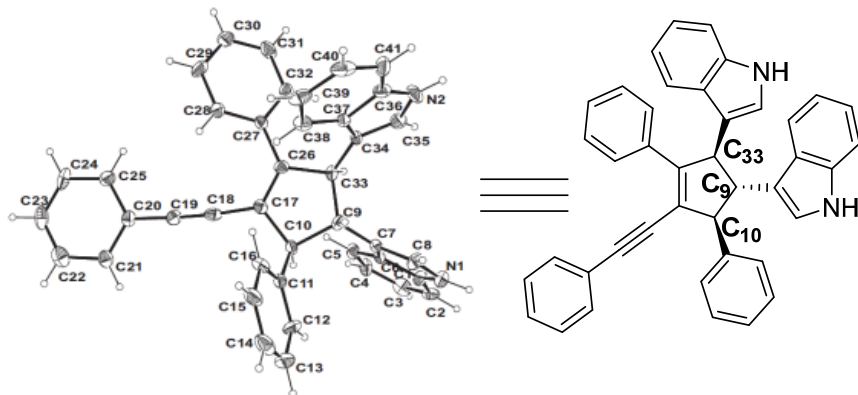
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Table S1. Crystal data and structure refinement for **3a**.



Identification code	d20433a	
Empirical formula	C ₄₁ H ₃₀ N ₂	
Formula weight	550.67	
Temperature	200(2) K	
Wavelength	0.71073 Å	
Crystal system	Monoclinic	
Space group	P 21/c	
Unit cell dimensions	a = 9.1373(13) Å	α = 90°.
	b = 18.993(3) Å	β = 100.289(4)°.
	c = 18.606(2) Å	γ = 90°.
Volume	3177.0(8) Å ³	
Z	4	
Density (calculated)	1.151 Mg/m ³	
Absorption coefficient	0.067 mm ⁻¹	
F(000)	1160	
Crystal size	0.10 x 0.08 x 0.02 mm ³	
Theta range for data collection	2.27 to 25.06°.	
Index ranges	-10 ≤ h ≤ 10, -22 ≤ k ≤ 22, -22 ≤ l ≤ 22	
Reflections collected	43966	
Independent reflections	5621 [R(int) = 0.1105]	
Completeness to theta = 25.06°	99.8 % Absorption correction	multi-scan
Max. and min. transmission	0.9987 and 0.9934	
Refinement method	Full-matrix least-squares on F ²	
Data / restraints / parameters	5621 / 0 / 388	
Goodness-of-fit on F ²	1.017	
Final R indices [I > 2σ(I)]	R1 = 0.0601, wR2 = 0.1310	

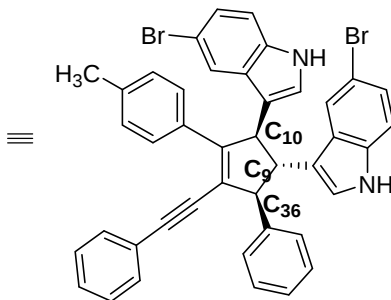
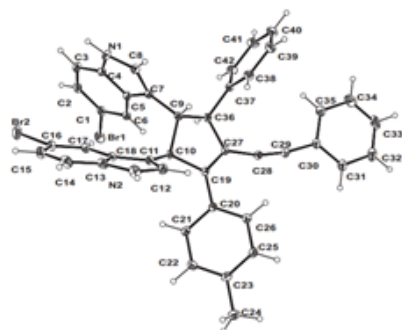
R indices (all data)

R1 = 0.1142, wR2 = 0.1498

Largest diff. peak and hole

0.159 and -0.174 e.Å⁻³

Table S2. Crystal data and structure refinement for **3o** (ORTEP diagram with ellipsoid contour 30% probability).



Identification code

d20250

Empirical formula

C₄₂ H₃₀ Br₂ N₂

Formula weight

722.50

Temperature

200(2) K

Wavelength

0.71073 Å

Crystal system

Triclinic

Space group

P -1

Unit cell dimensions

a = 9.4052(4) Å

α = 80.4220(10)°.

b = 12.5506(5) Å

β = 80.6500(10)°.

c = 16.5291(7) Å

γ = 89.9930(10)°.

Volume

1897.63(14) Å³

Z

2

Density (calculated)

1.264 Mg/m³

Absorption coefficient

2.165 mm⁻¹

F(000)

732

Crystal size

0.34 x 0.04 x 0.02 mm³

Theta range for data collection

2.24 to 25.04°.

Index ranges

-11 ≤ h ≤ 11, -14 ≤ k ≤ 14, -19 ≤ l ≤ 19

Reflections collected

67740

Independent reflections

6691 [R(int) = 0.0677]

Completeness to theta = 25.04°

99.8 %

Absorption correction

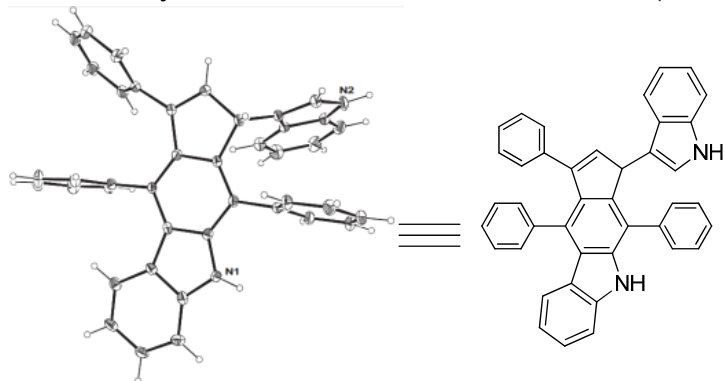
multi-scan

Max. and min. transmission

0.9580 and 0.5264

Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	6691 / 0 / 416
Goodness-of-fit on F ²	1.082
Final R indices [I>2sigma(I)]	R1 = 0.0374, wR2 = 0.0906
R indices (all data)	R1 = 0.0526, wR2 = 0.0971
Largest diff. peak and hole	0.222 and -0.562 e.Å ⁻³

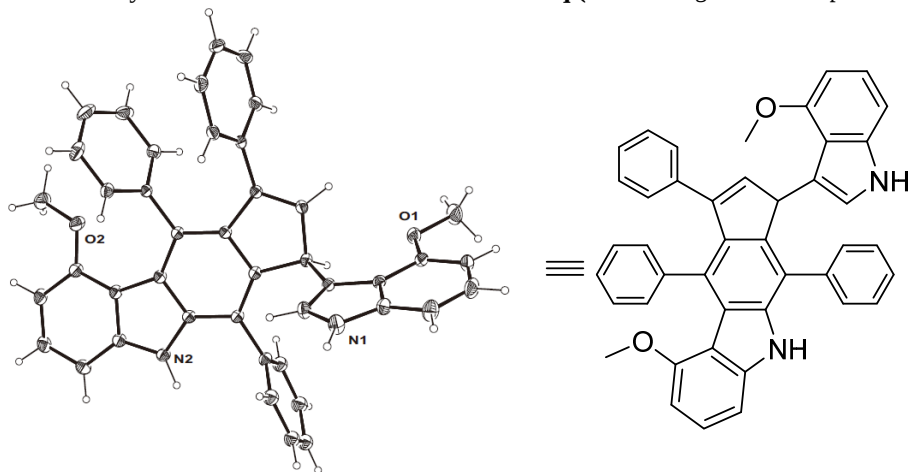
Table S3. Crystal data and structure refinement for **4a** (ORTEP diagram with ellipsoid contour 30% probability)



Identification code	a16363
Empirical formula	C ₄₅ H ₃₆ N ₂ O ₂ (C ₄₁ H ₂₈ N ₂ +C ₄ H ₈ O ₂ (Ethylacetate))
Formula weight	636.76
Temperature	200(2) K
Wavelength	0.71073 Å
Crystal system	Triclinic
Space group	P -1
Unit cell dimensions	a = 11.134(2) Å α = 83.142(7)°. b = 12.176(3) Å β = 77.790(7)°. c = 13.759(3) Å γ = 67.611(8)°.
Volume	1684.1(7) Å ³
Z	2
Density (calculated)	1.256 Mg/m ³
Absorption coefficient	0.076 mm ⁻¹
F(000)	672
Crystal size	0.13 x 0.11 x 0.05 mm ³
Theta range for data collection	2.01 to 25.10°.
Index ranges	-13<=h<=13, -11<=k<=14, -16<=l<=16
Reflections collected	13220
Independent reflections	5954 [R(int) = 0.0957]
Completeness to theta	= 25.10° 99.2 %

Absorption correction	multi-scan
Max. and min. transmission	0.9962 and 0.9901
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	5954 / 0 / 442
Goodness-of-fit on F ²	0.947
Final R indices [I > 2σ(I)]	R1 = 0.0681, wR2 = 0.1581
R indices (all data)	R1 = 0.1791, wR2 = 0.2229
Largest diff. peak and hole	0.765 and -0.337 e.Å ⁻³

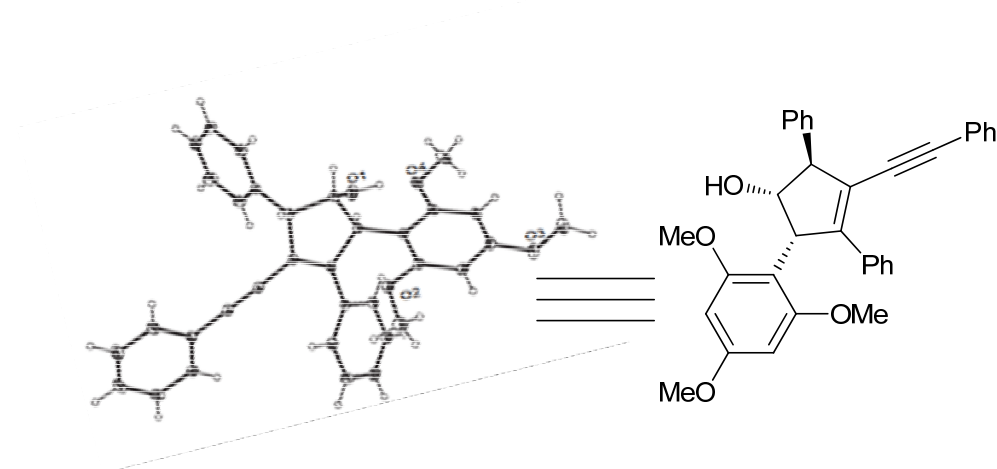
Table S4. Crystal data and structure refinement for **4q** (ORTEP diagram with ellipsoid contour 30% probability)



Identification code	d19307
Empirical formula	C ₄₃ H ₃₂ N ₂ O ₂
Formula weight	608.71
Temperature	200(2) K
Wavelength	0.71073 Å
Crystal system	Triclinic
Space group	P -1
Unit cell dimensions	a = 7.0575(5) Å α = 88.791(3)° b = 10.2773(8) Å β = 88.610(3)° c = 22.134(2) Å γ = 81.318(2)°
Volume	1586.3(2) Å ³
Z	2
Density (calculated)	1.274 Mg/m ³
Absorption coefficient	0.078 mm ⁻¹
F(000)	640
Crystal size	0.24 x 0.12 x 0.02 mm ³
Theta range for data collection	2.19 to 25.06°

Index ranges	-8<=h<=8, -12<=k<=12, -26<=l<=26
Reflections collected	49319
Independent reflections	5617 [R(int) = 0.0733]
Completeness to theta	= 25.06° 99.6 %
Absorption correction	multi-scan
Max. and min. transmission	0.9984 and 0.9815
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	5617 / 4 / 420
Goodness-of-fit on F ²	1.052
Final R indices [I>2sigma(I)]	R1 = 0.0540, wR2 = 0.1117
R indices (all data)	R1 = 0.0836, wR2 = 0.1279
Extinction coefficient	0.0141(14)
Largest diff. peak and hole	0.352 and -0.306 e.Å ⁻³

Table S5. Crystal data and structure refinement for **5a** (ORTEP diagram with ellipsoid contour 30% probability)

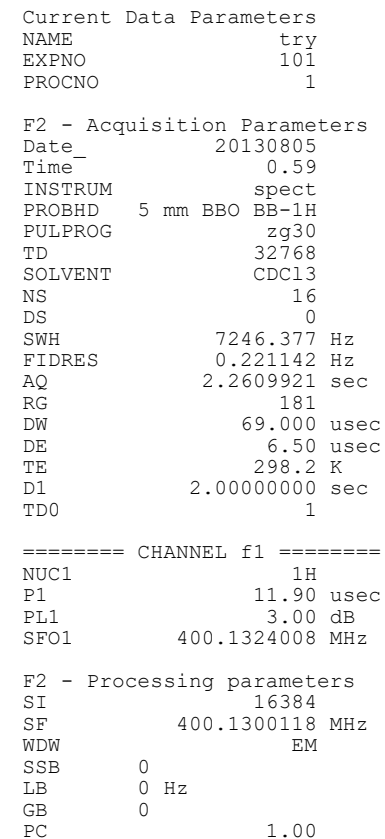


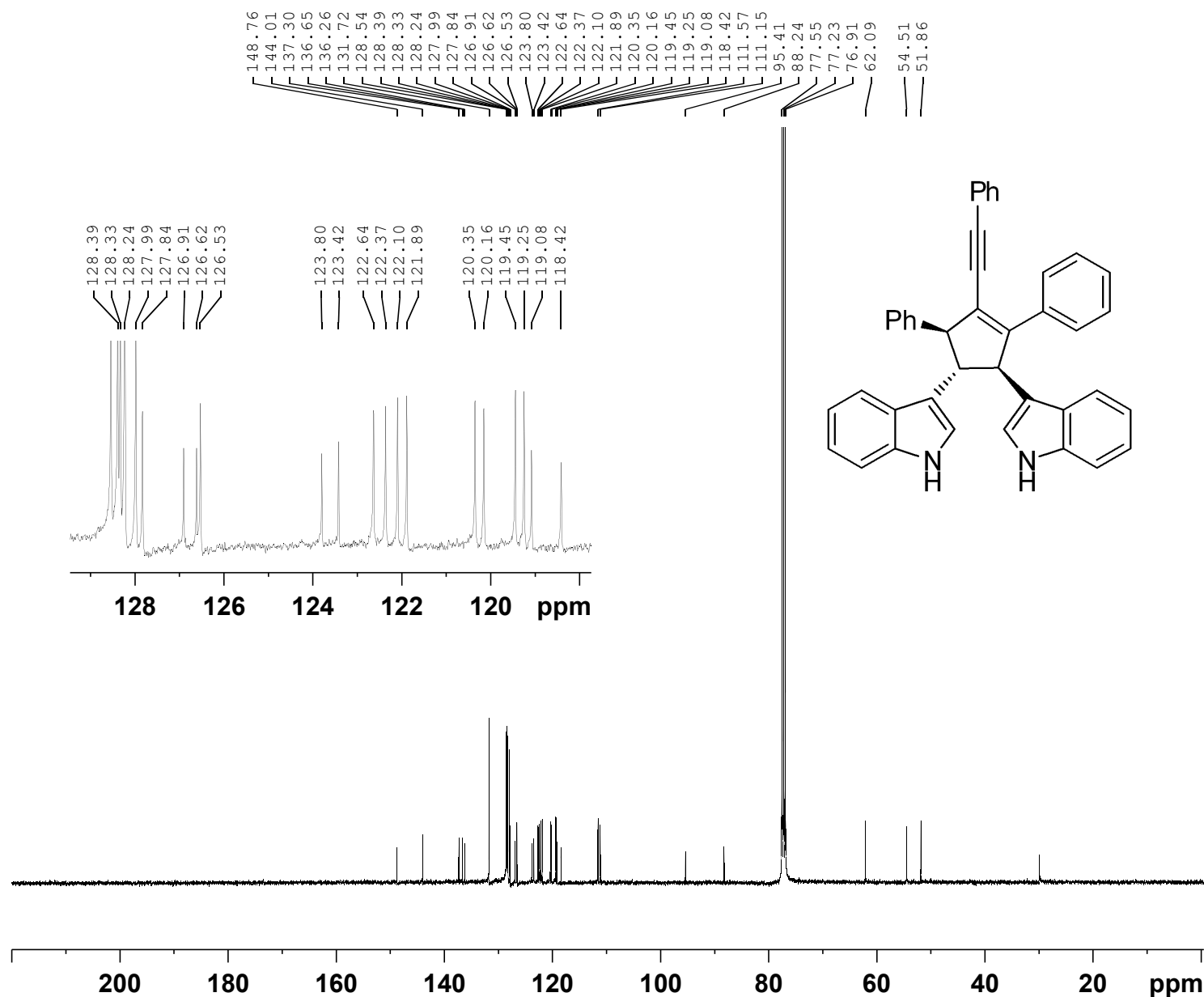
Identification code	d20430	
Empirical formula	C ₃₄ H ₃₀ O ₄	
Formula weight	502.58	
Temperature	238(2) K	
Wavelength	0.71073 Å	
Crystal system	Triclinic	
Space group	P -1	
Unit cell dimensions	a = 10.044(3) Å	α = 81.343(7)°.
	b = 12.223(3) Å	β = 71.583(7)°.
	c = 12.450(3) Å	γ = 66.010(7)°.
Volume	1324.6(6) Å ³	

Z	2
Density (calculated)	1.260 Mg/m ³
Absorption coefficient	0.081 mm ⁻¹
F(000)	532
Crystal size	0.32 x 0.26 x 0.07 mm ³
Theta range for data collection	2.31 to 25.25°.
Index ranges	-11<=h<=11, -14<=k<=14, -14<=l<=14
Reflections collected	43775
Independent reflections	4710 [R(int) = 0.0684]
Completeness to theta = 25.25°	98.5 %
Absorption correction	multi-scan
Max. and min. transmission	0.9943 and 0.9744
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	4710 / 0 / 347
Goodness-of-fit on F ²	1.070
Final R indices [I>2sigma(I)]	R1 = 0.0476, wR2 = 0.1053
R indices (all data)	R1 = 0.0656, wR2 = 0.1165
Extinction coefficient	0.018(3)
Largest diff. peak and hole	0.432 and -0.436 e.Å ⁻³

¹H NMR and ¹³C NMR Spectra copies

58





```
Current Data Parameters
NAME                      try
EXPNO                     102
PROCNO                    1
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```

F2 - Acquisition Parameters
Date_                20130803
Time_                23.33
INSTRUM              spect
PROBHD      5 mm BBO BB-1H
PULPROG              zgpg30
TD                   32768
SOLVENT              CDCl3
NS                   14722
DS                     0
SWH                  24038.461 Hz
FIDRES              0.733596 Hz
AQ                  0.6815744 sec
RG                   8192
DW                   20.800 usec
DE                   6.50 usec
TE                   297.7 K
D1                   2.00000000 sec
D11                  0.03000000 sec
TD0                   1

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===== CHANNEL f1 =====
NUC1                13C
P1                  9.40 usec
PL1                 7.00 dB
SFO1               100.6233325 MHz
```

```

===== CHANNEL f2 =====
CPDPRG[2          waltz16
NUC2              1H
PCPD2             90.00 usec
PL2               3.00 dB
PL12              20.70 dB
PL13              23.70 dB
SFO2              400.1316005 MHz

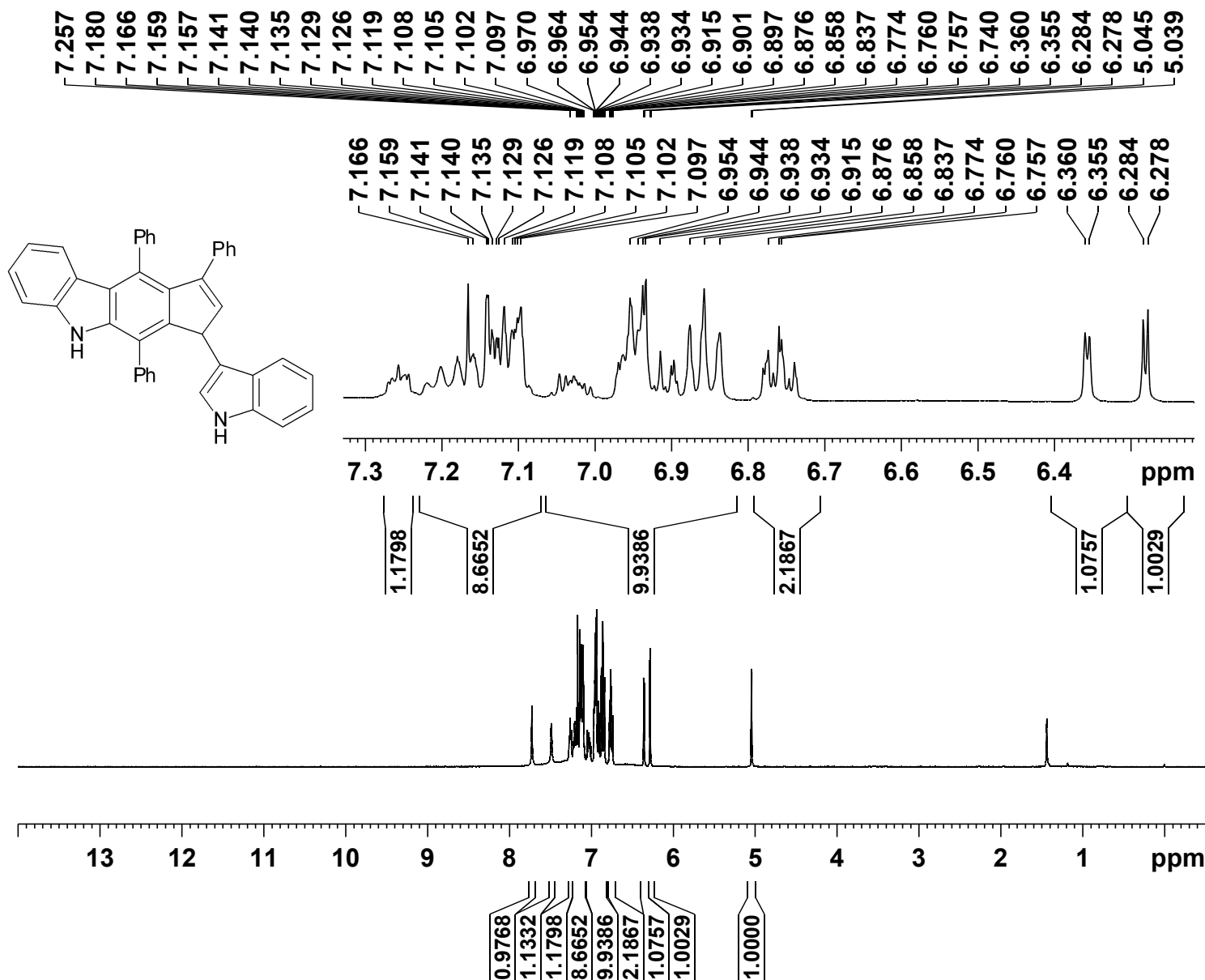
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F2 - Processing parameters
SI                      32768
SF                      100.6127484 MHz
WDW                      EM
SSB                      0
LB                      1.00 Hz
GB                      0
PC                      1.00

```

3-(1H-indol-3-yl)-1,4,10-triphenyl-3,5-dihydrocyclopenta[b]carbazole (4a)



Current Data Parameters
NAME try
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20140716
Time_ 22.24
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 100
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 158.74
DW 69.333 usec
DE 10.52 usec
TE 298.4 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 400.1324008 MHz
NUC1 1H
P1 12.80 usec
PLW1 15.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1300470 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3-(1H-indol-3-yl)-1,4,10-triphenyl-3,5-dihydrocyclopenta[b]carbazole (4a)



C
NAME try
EXPNO 2
PROCNO 1

F2 - Acquisition Parameters

Date_ 20140716
Time 22.32
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 14836
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.5 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

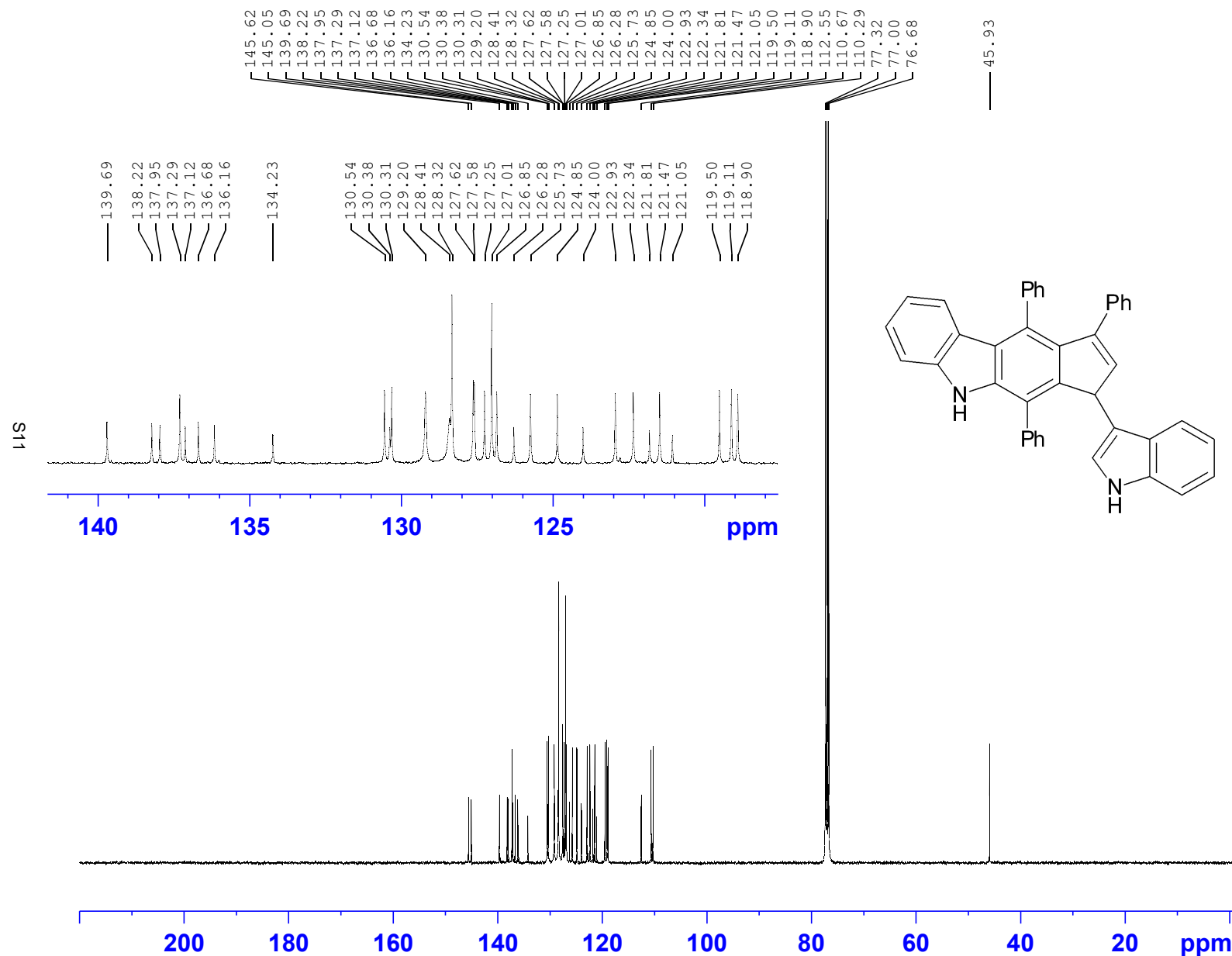
SFO1 100.623319 MHz
NUC1 13C
P1 10.00 usec
PLW1 46.00000000 W

===== CHANNEL f2 =====

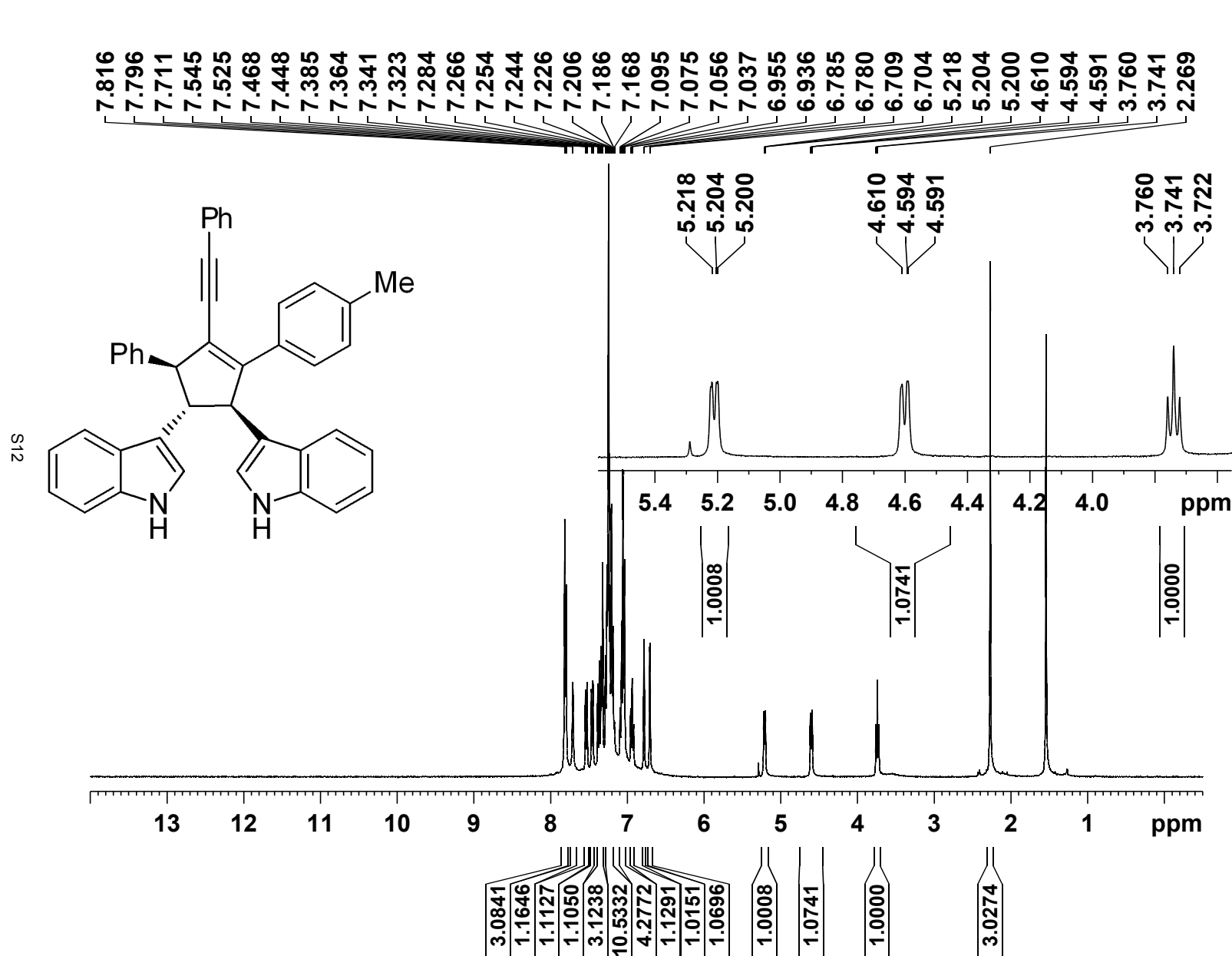
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.34252000 W
PLW13 0.27744001 W

F2 - Processing parameters

SI 32768
SF 100.6127711 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00



3,3'-(5-phenyl-4-(phenylethynyl)-3-(p-tolyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3b)



Current Data Parameters

NAME test
EXPNO 11
PROCNO 1

F2 - Acquisition Parameters

Date_ 20170710
Time 17.25
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 32
DS 0
SWH 7246.377 Hz
FIDRES 0.221142 Hz
AQ 2.2609921 sec
RG 181
DW 69.000 usec
DE 6.50 usec
TE 300.1 K
D1 2.00000000 sec
TD0 1

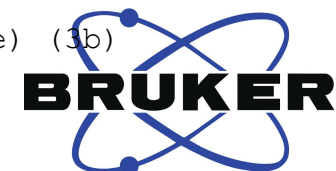
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NUC1 1H
P1 16.80 usec
PL1 1.80 dB
SFO1 400.1324008 MHz

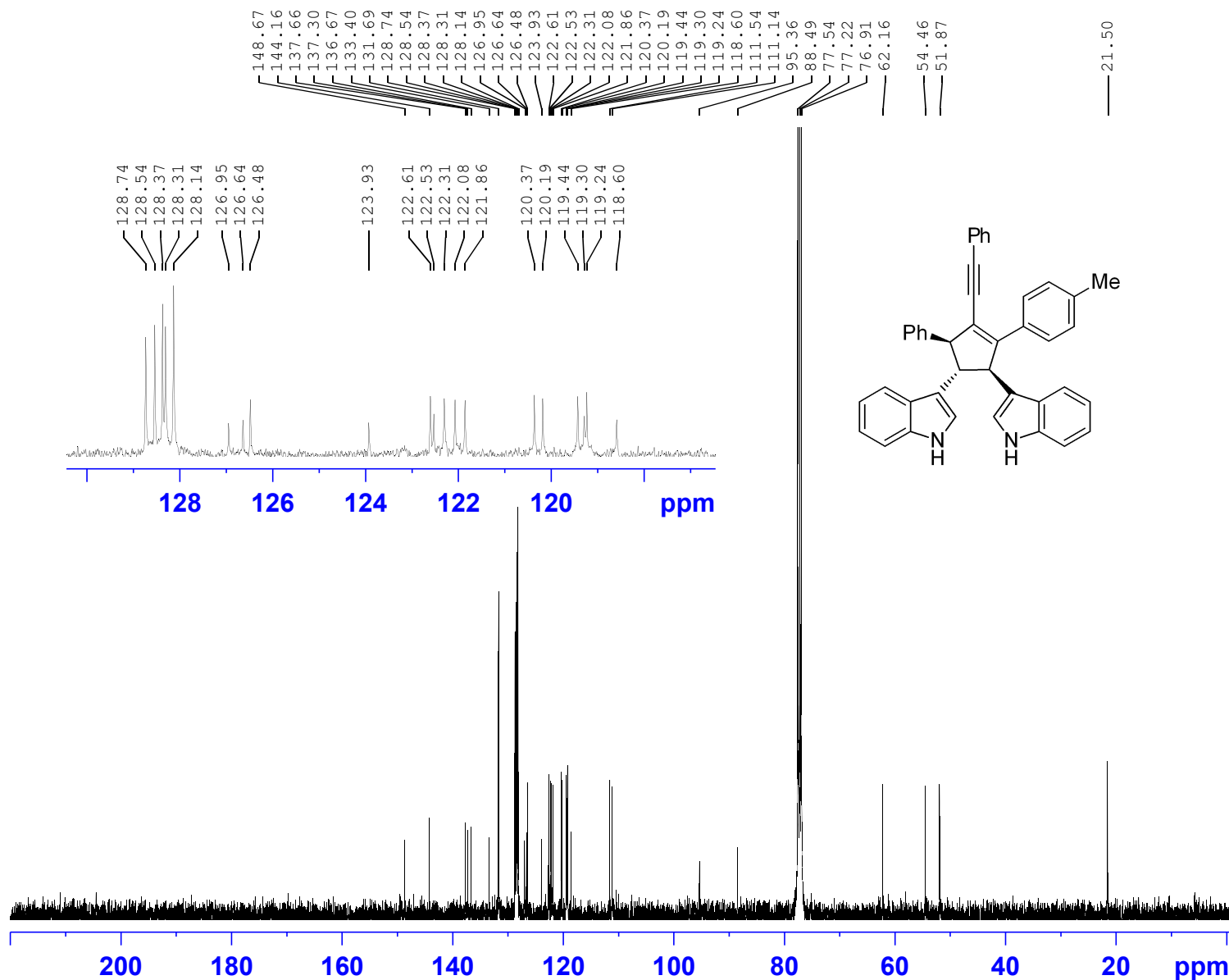
F2 - Processing parameters

SI 16384
SF 400.1300118 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(5-phenyl-4-(phenylethynyl)-3-(p-tolyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3b)



SIS



Current Data Parameters
 NAME test
 EXPNO 12
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20170710
 Time 17.28
 INSTRUM spect
 PROBHD 5 mm BBO BB-1H
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 1000
 DS 0
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 5792.6
 DW 20.800 usec
 DE 6.50 usec
 TE 300.2 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

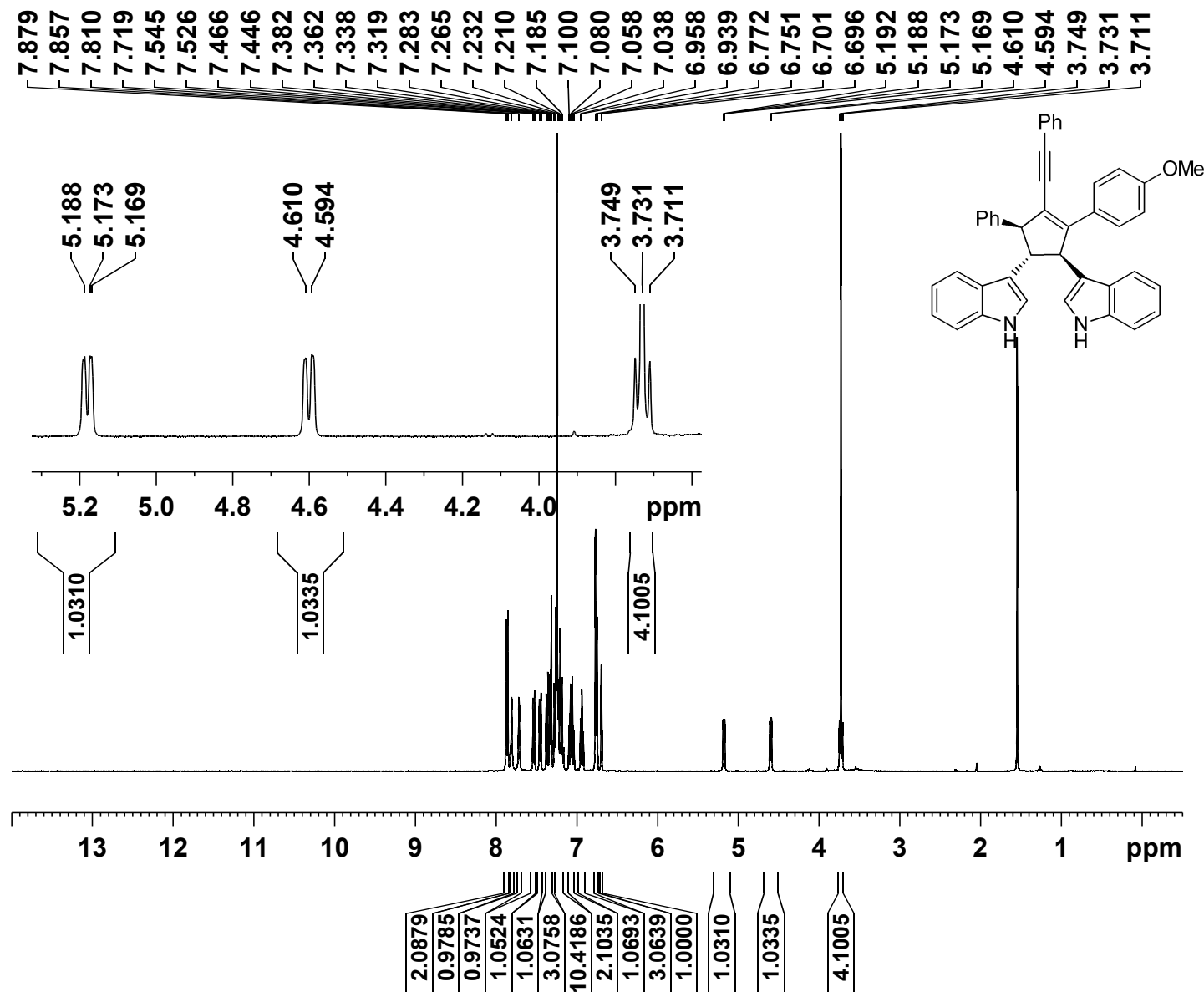
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 P1 10.45 usec
 PL1 7.00 dB
 SFO1 100.623325 MHz

===== CHANNEL f2 =====
 CPDPRG[2] waltz16
 NUC2 1H
 PCPD2 90.00 usec
 PL2 1.80 dB
 PL12 15.68 dB
 PL13 20.00 dB
 SFO2 400.1316005 MHz

F2 - Processing parameters
 SI 32768
 SF 100.6127481 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.00

3,3'-(3-(4-methoxyphenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3c)

S14



Current Data Parameters

NAME test
EXPNO 13
PROCNO 1

F2 - Acquisition Parameters

Date_ 20170711
Time_ 15.12
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 32
DS 0
SWH 7246.377 Hz
FIDRES 0.221142 Hz
AQ 2.2609921 sec
RG 161.3
DW 69.000 usec
DE 6.50 usec
TE 298.1 K
D1 2.00000000 sec
TD0 1

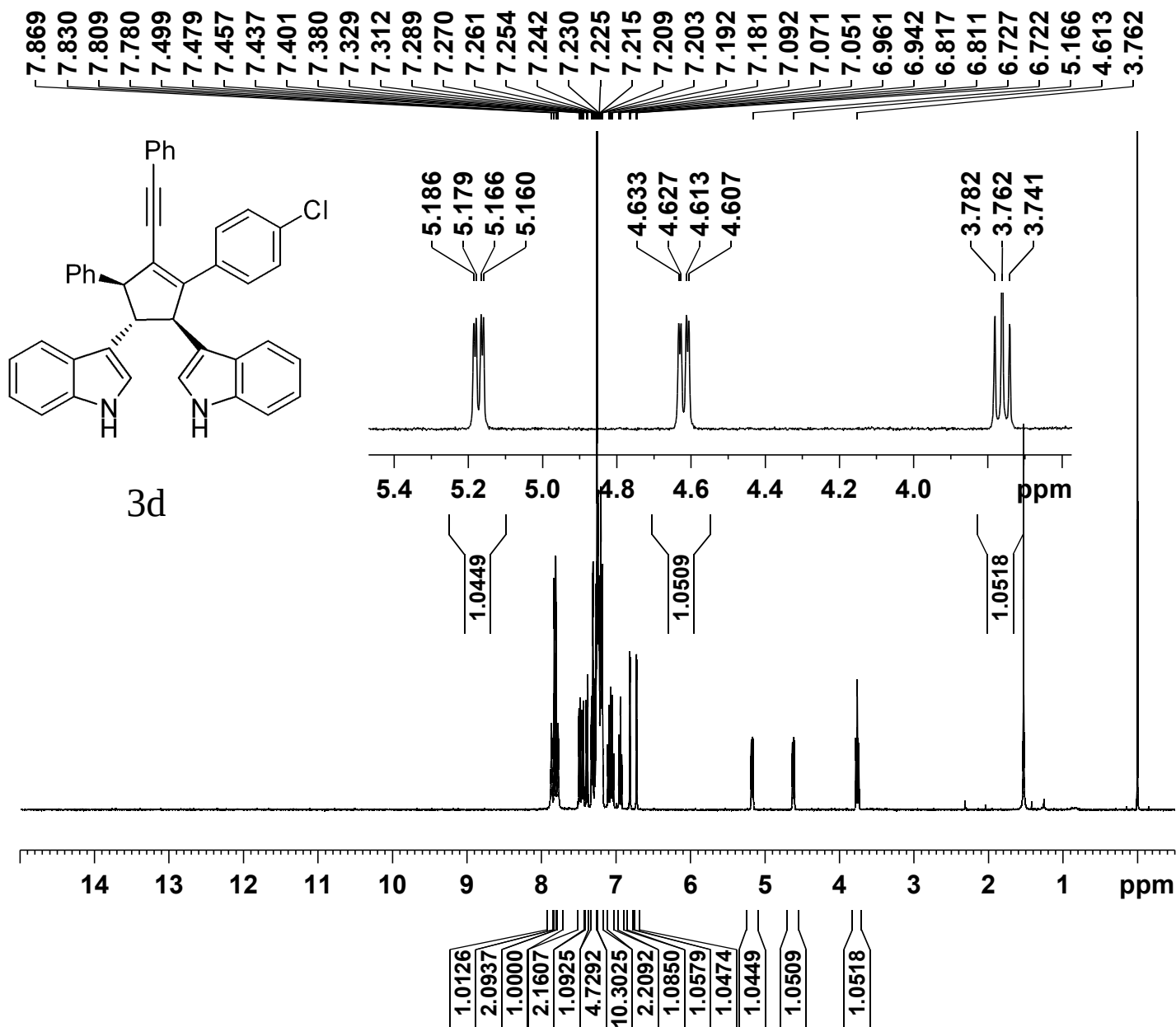
===== CHANNEL f1 =====

NUC1 1H
P1 16.80 usec
PL1 1.80 dB
SFO1 400.1324008 MHz

F2 - Processing parameters

SI 16384
SF 400.1300118 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-(4-chlorophenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3d)



Current Data Parameters

NAME try
EXPNO 88
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180718
Time_ 19.09
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 33
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 299.5 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

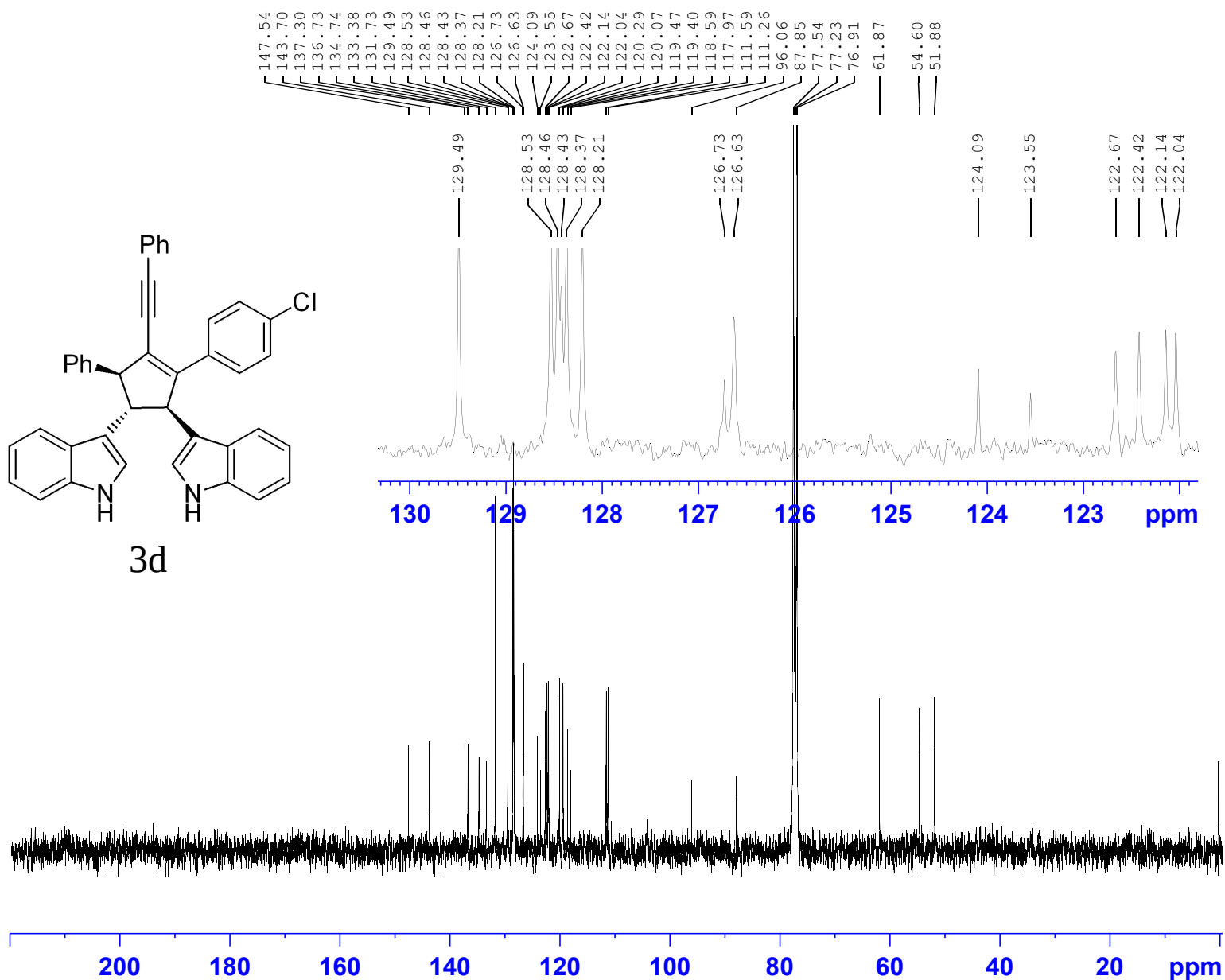
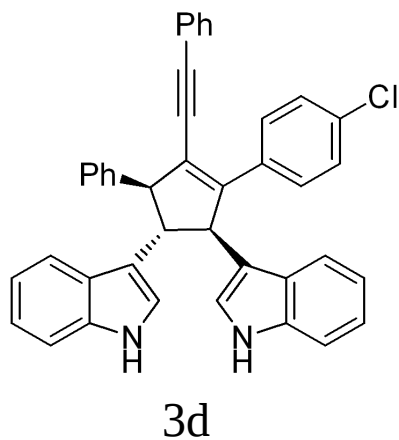
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300116 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-(4-chlorophenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3d)

S17



Current Data Parameters
NAME try
EXPNO 91
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180718
Time_ 21.59
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 2772
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 299.5 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

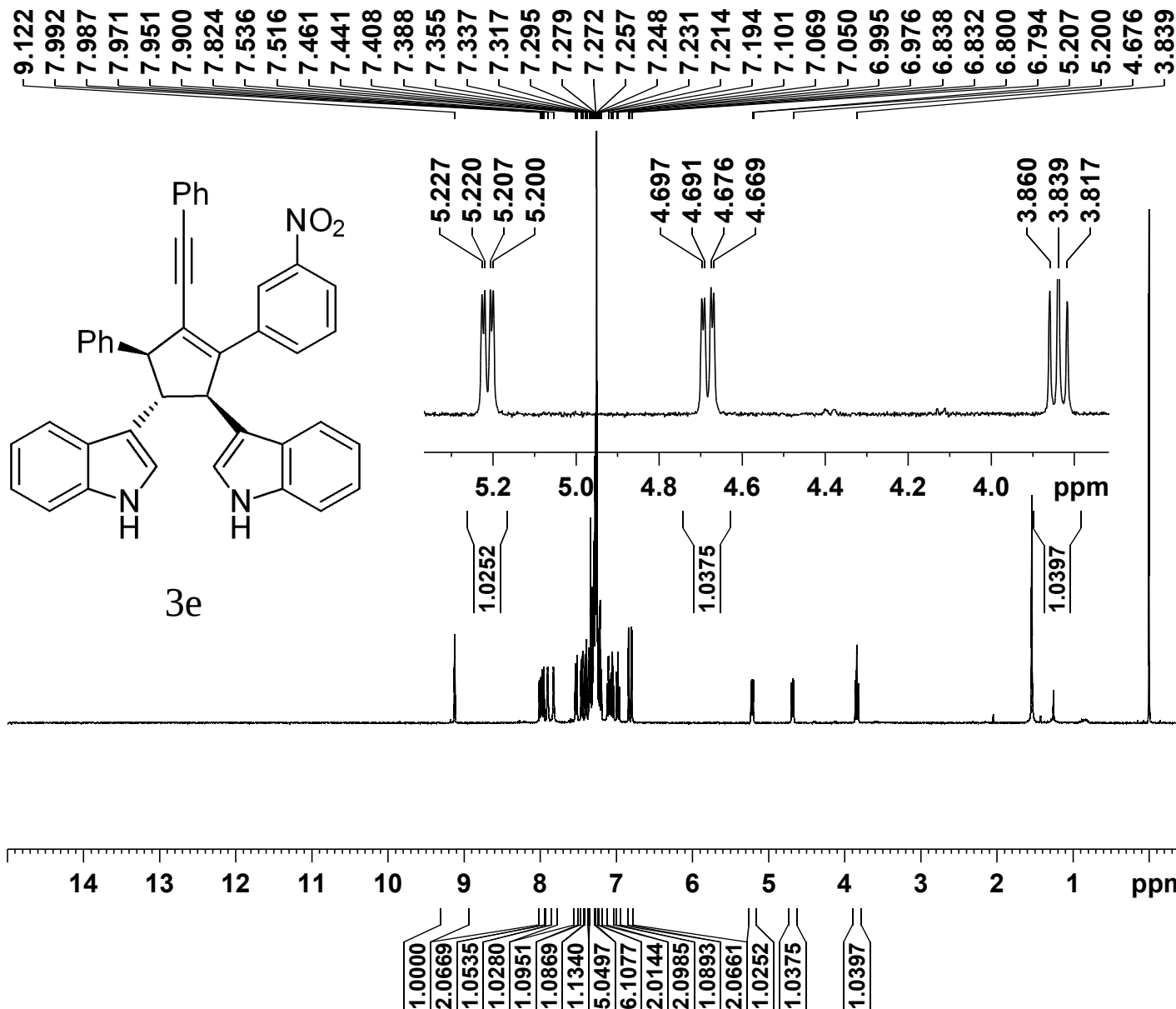
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NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====
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NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters
SI 32768
SF 100.6127468 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00

3,3'-(3-(3-nitrophenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3e)

81S



Current Data Parameters
NAME try
EXPNO 101
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180807
Time_ 14.43
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 32
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 297.6 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1300107 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-(3-nitrophenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3e)



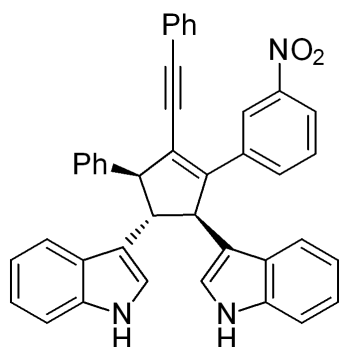
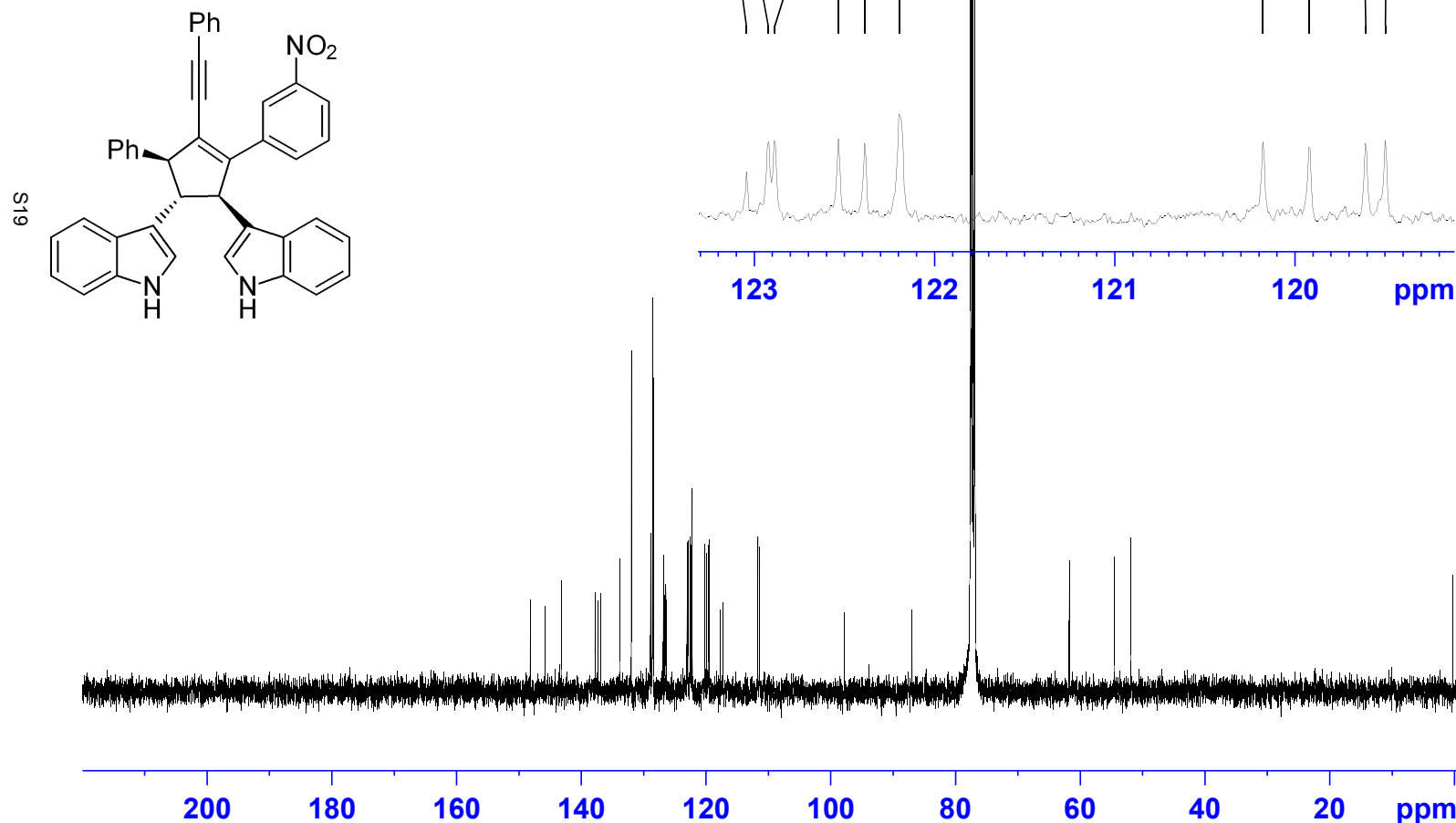
Current Data Parameters
NAME test
EXPNO 57
PROCNO 1

F2 - Acquisition Parameters
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Time 22.37
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PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 10000
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5792.6
DW 20.800 usec
DE 6.50 usec
TE 298.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

F2 - Processing parameters
SI 32768
SF 100.6127474 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00



3,3'-(3-(naphthalen-1-yl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3f)

Current Data Parameters

NAME try
EXPNO 892
PROCNO 1

F2 - Acquisition Parameters

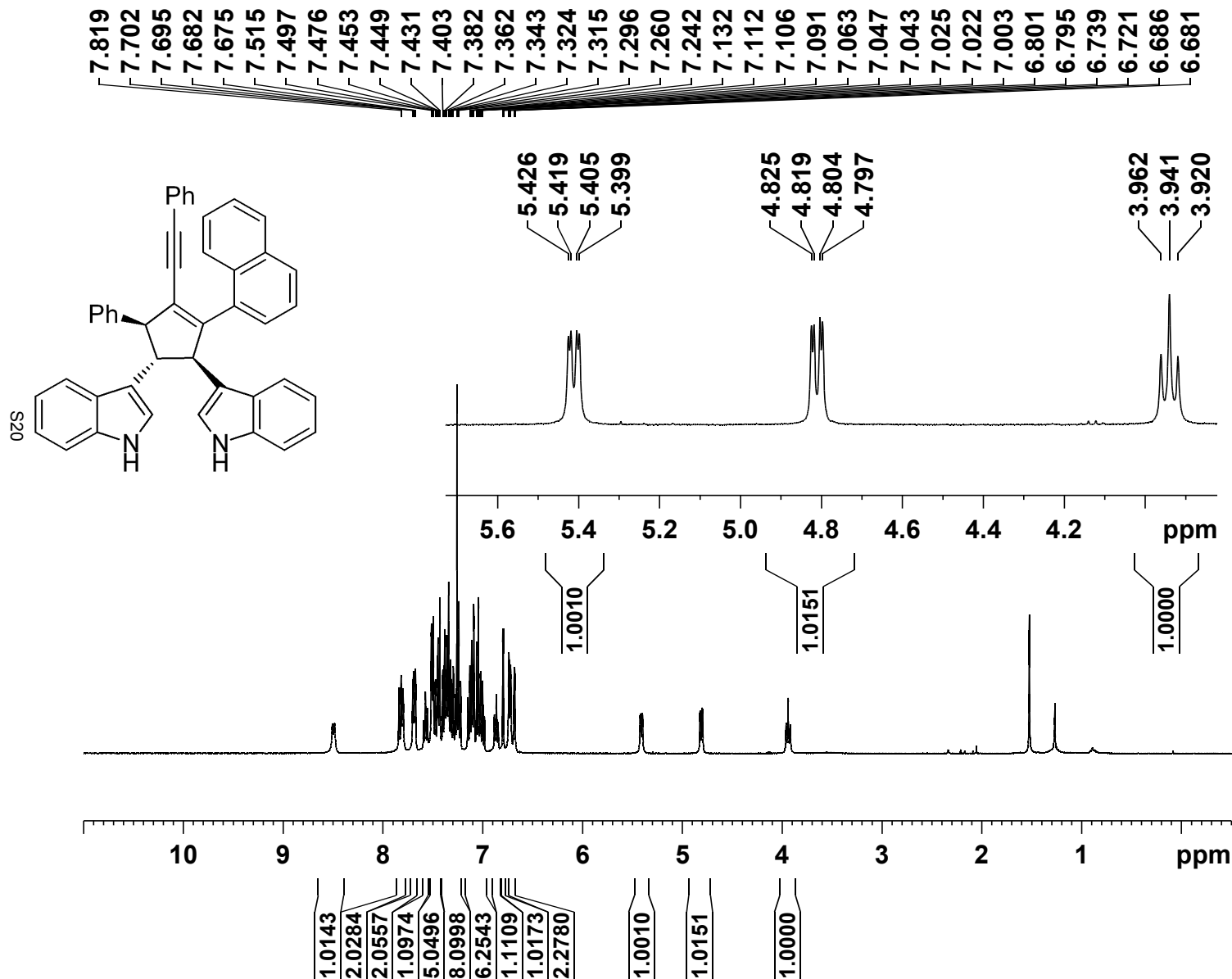
Date_ 20170810
Time_ 22.45
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 64
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 128.9
DW 69.333 usec
DE 10.50 usec
TE 298.6 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

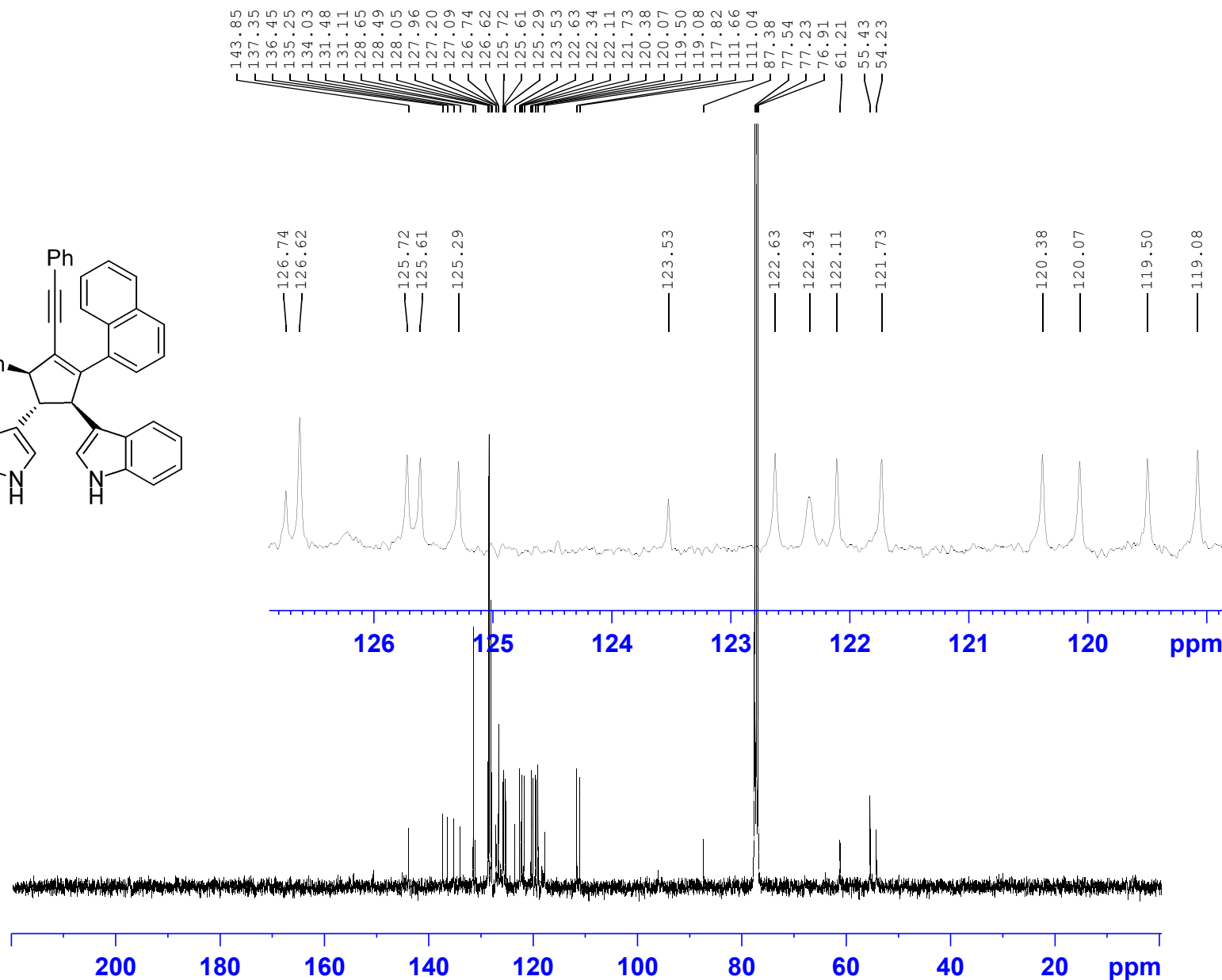
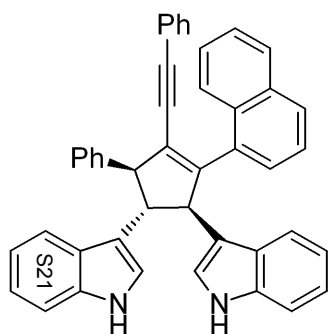
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300100 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00



3,3'-(3-(naphthalen-1-yl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3f)



Current Data Parameters
NAME try
EXPNO 893
PROCNO 1

F2 - Acquisition Parameters
Date_ 20170810
Time_ 22.50
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 1500
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.7 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters
SI 32768
SF 100.6127479 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00

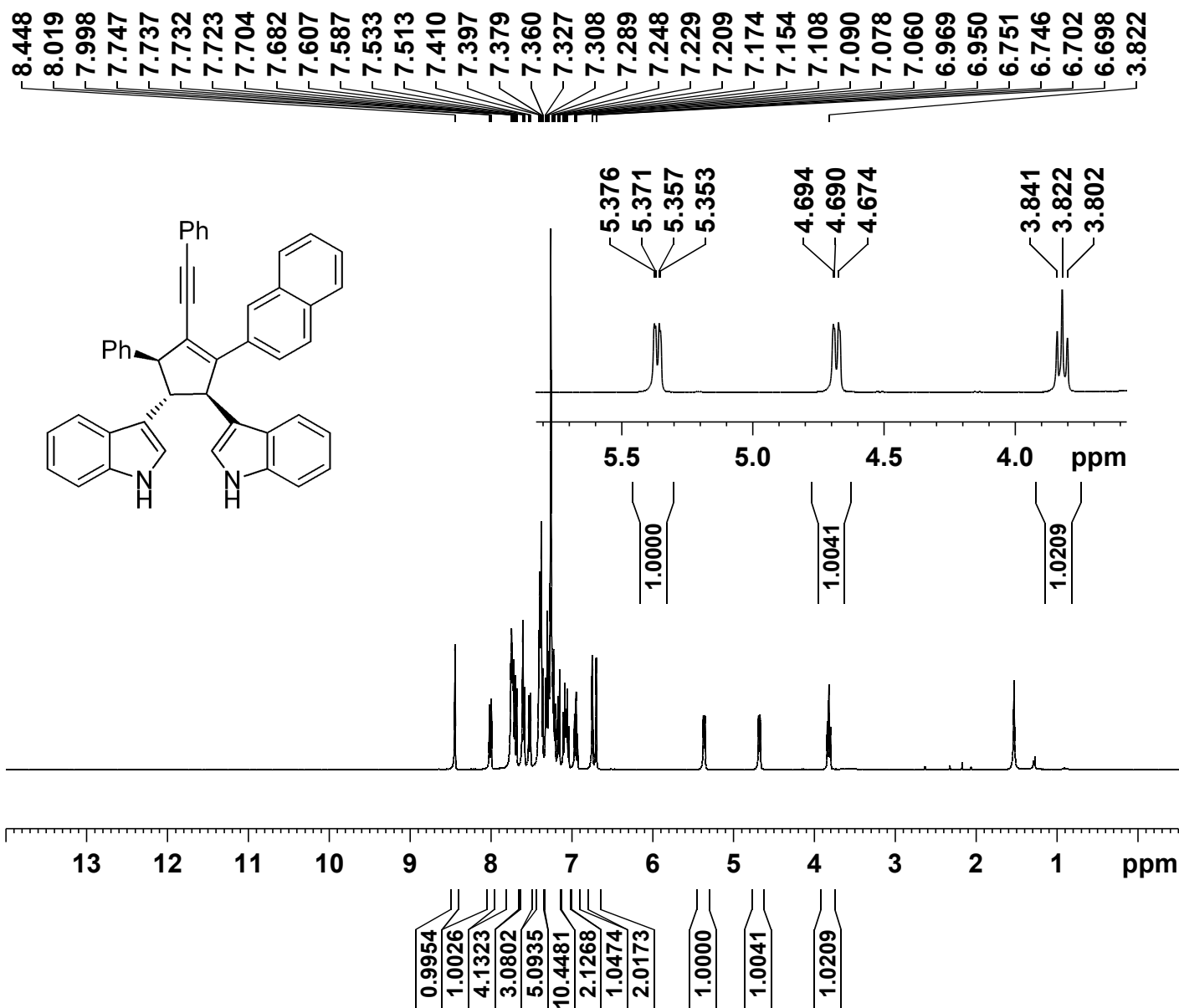
3,3'-(3-(naphthalen-2-yl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3g)

Current Data Parameters
NAME test
EXPNO 36
PROCNO 1

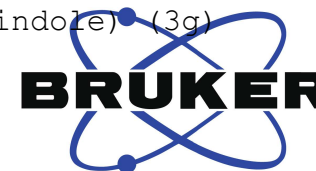
F2 - Acquisition Parameters
Date_ 20170901
Time_ 16.30
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 64
DS 0
SWH 7246.377 Hz
FIDRES 0.221142 Hz
AQ 2.260921 sec
RG 114
DW 69.000 usec
DE 6.50 usec
TE 297.5 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====
NUC1 1H
P1 15.00 usec
PL1 0 dB
SFO1 400.1324008 MHz

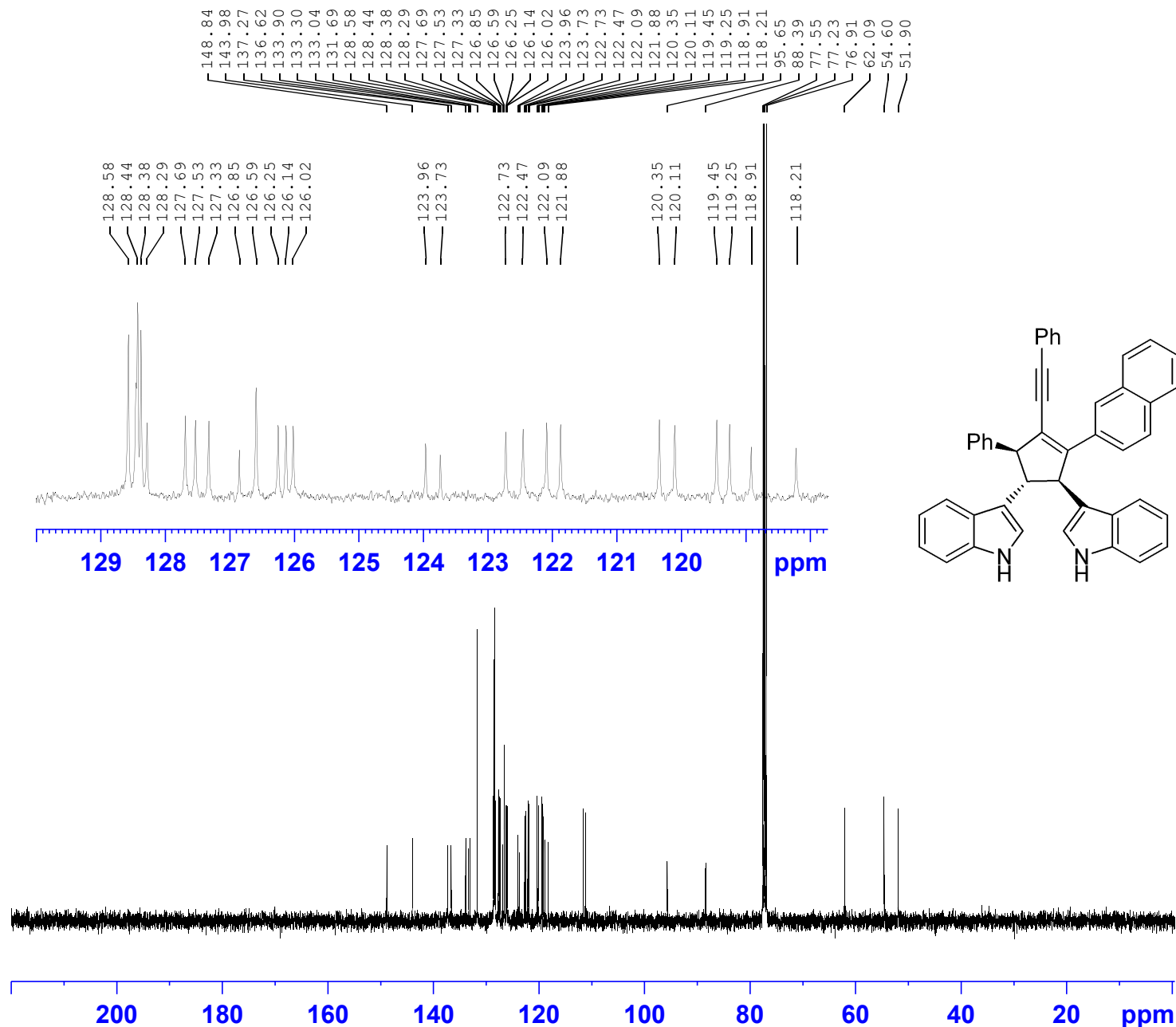
F2 - Processing parameters
SI 16384
SF 400.1300118 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00



3,3'-(3-(naphthalen-2-yl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3g)



S23



Current Data Parameters

NAME test
EXPNO 37
PROCNO 1

F2 - Acquisition Parameters

Date_ 20170901
Time_ 16.34
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 372
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5792.6
DW 20.800 usec
DE 6.50 usec
TE 297.6 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====

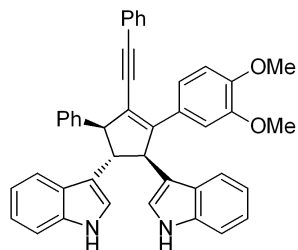
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

F2 - Processing parameters

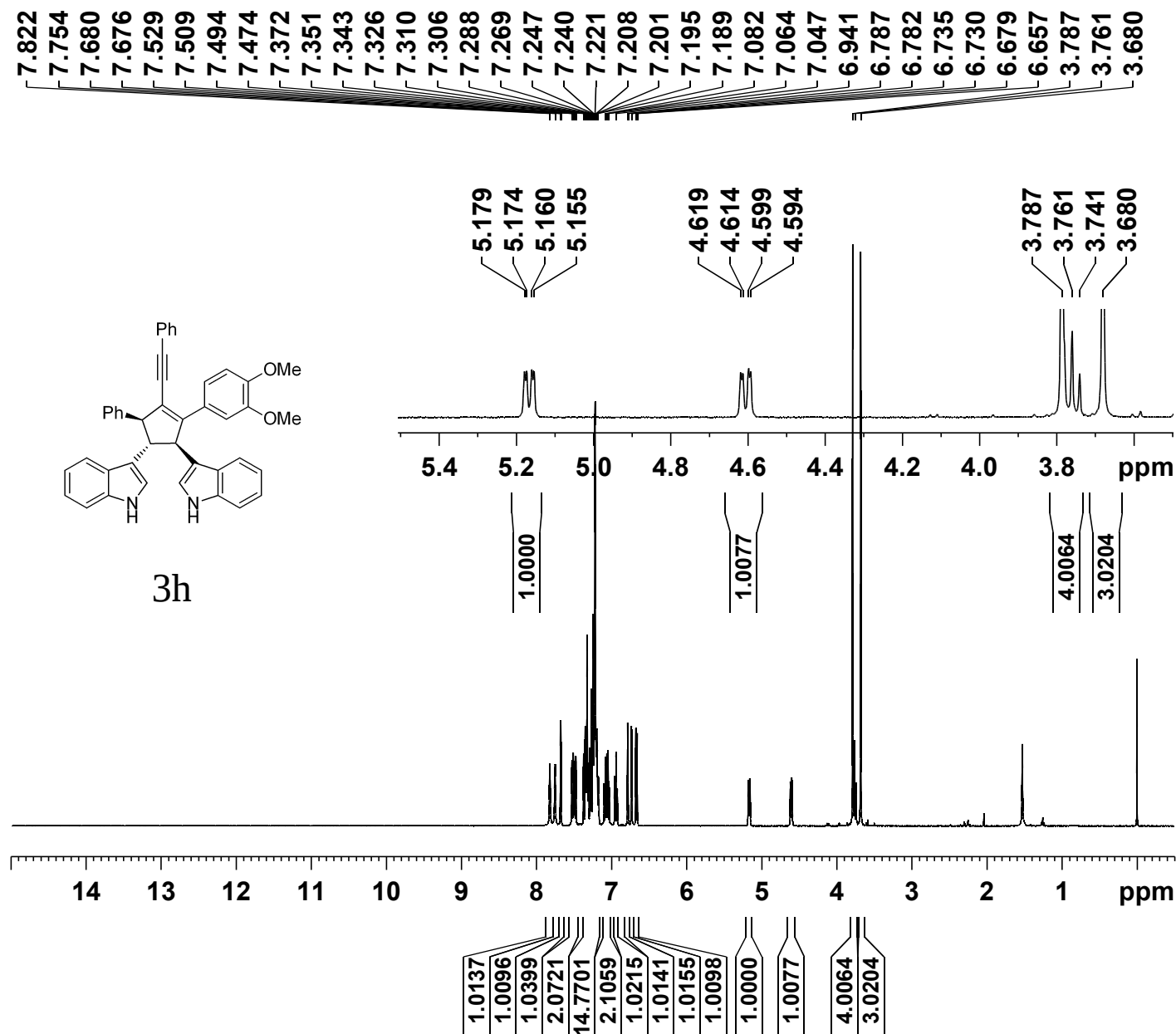
SI 32768
SF 100.6127521 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

3,3'-(3-(3,4-dimethoxyphenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3h)

S24



3h



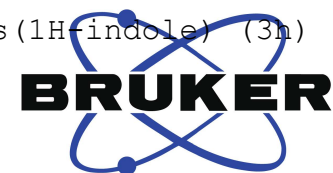
Current Data Parameters
NAME try
EXPNO 85
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180717
Time_ 19.40
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 32
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 177.16
DW 69.333 usec
DE 10.50 usec
TE 298.9 K
D1 2.00000000 sec
TD0 1

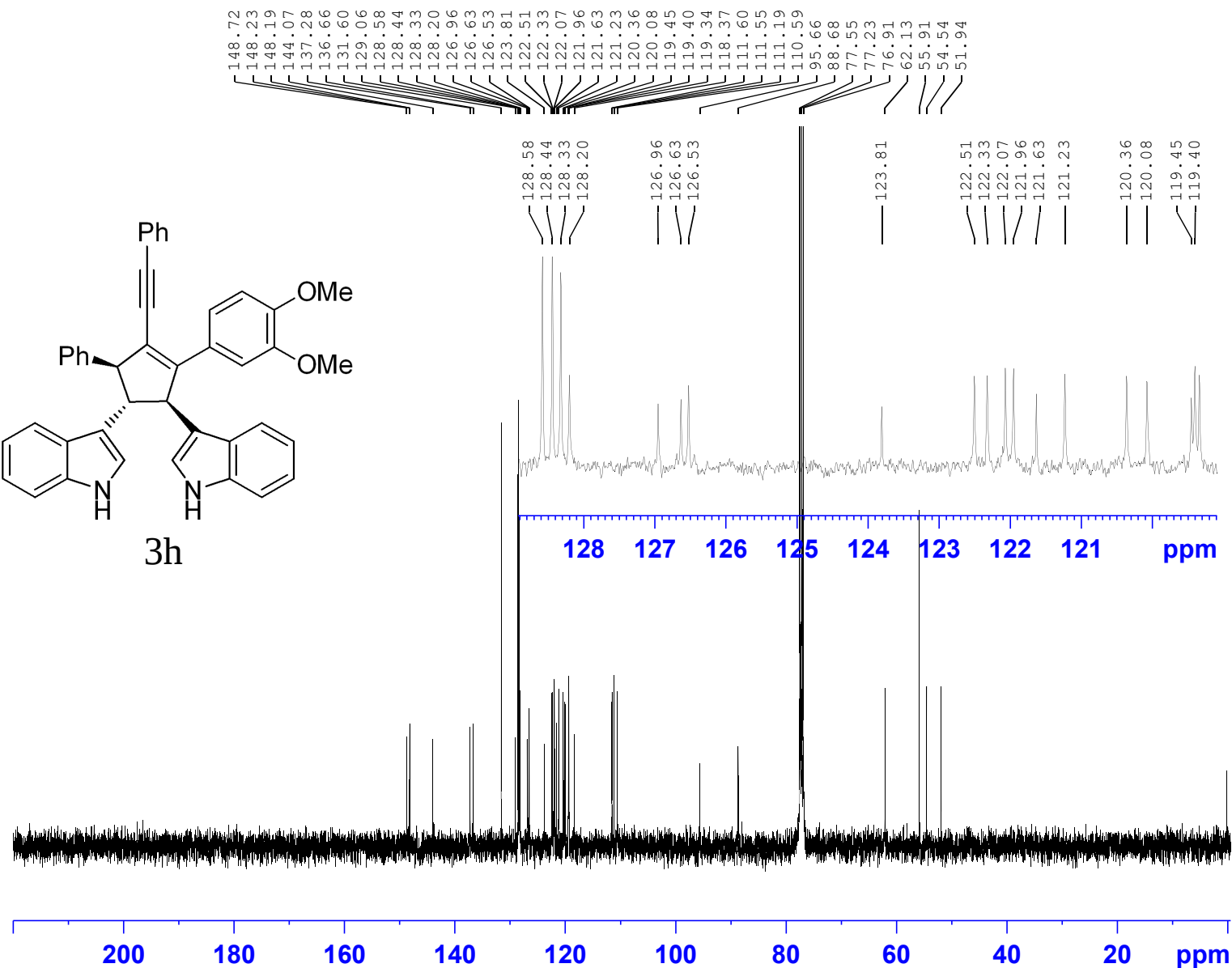
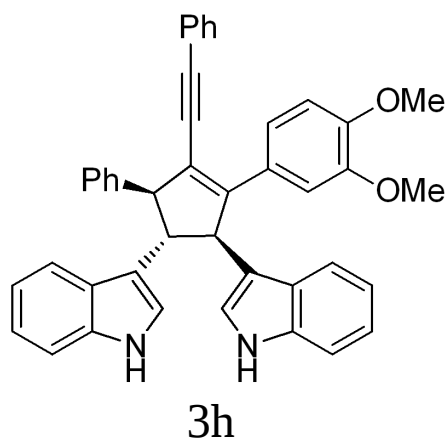
===== CHANNEL f1 =====
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1300147 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-(3,4-dimethoxyphenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3h)



926



Current Data Parameters
NAME test
EXPNO 69
PROCNO 1

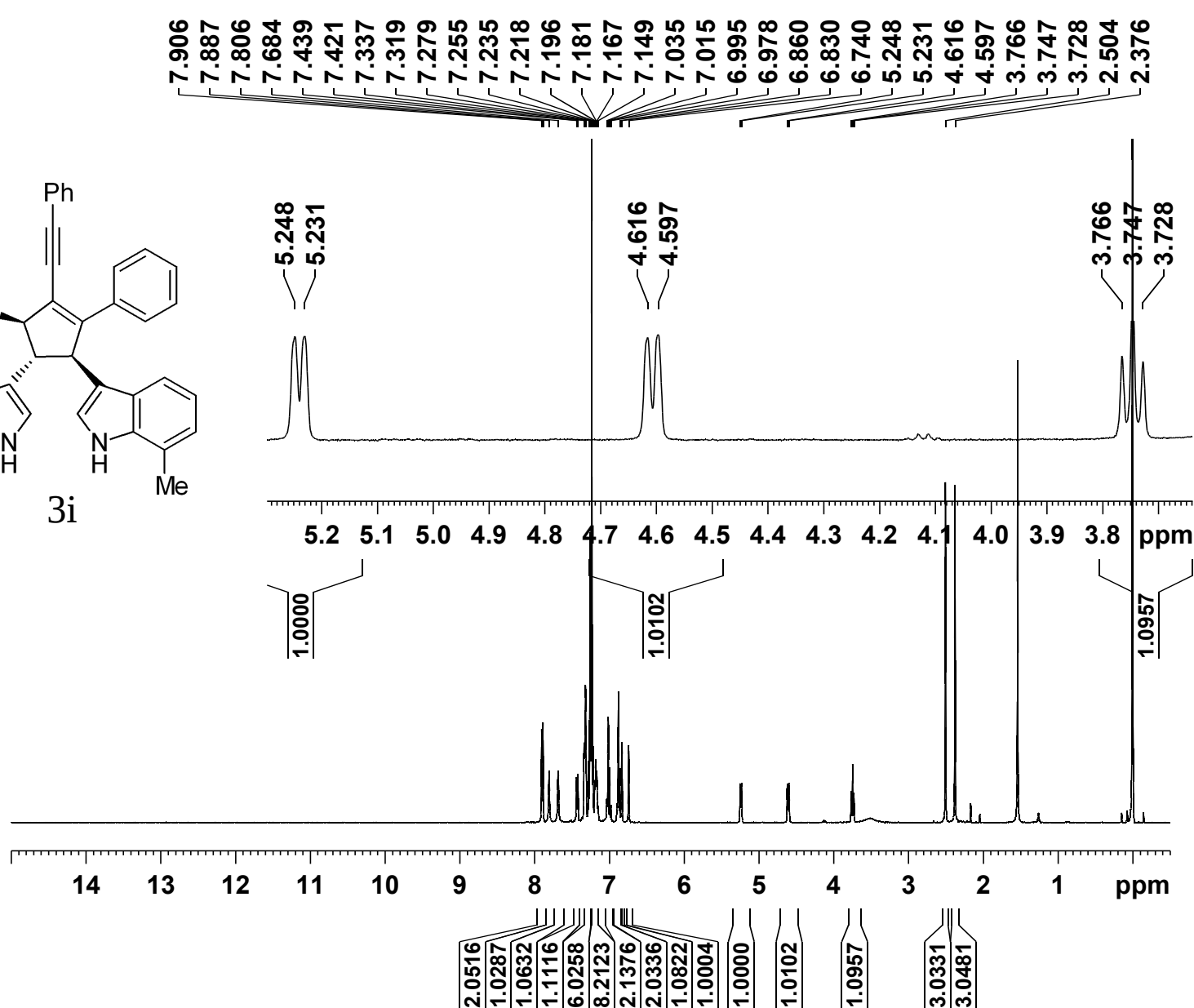
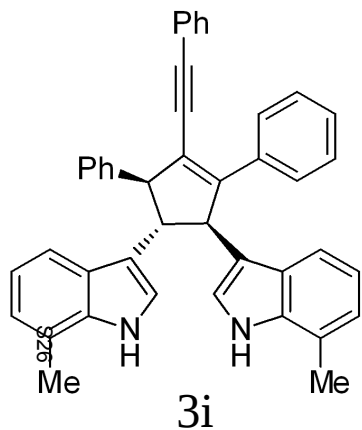
F2 - Acquisition Parameters
Date_ 20180102
Time 23.09
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 965
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5792.6
DW 20.800 usec
DE 6.50 usec
TE 298.6 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

F2 - Processing parameters
SI 32768
SF 100.6127484 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3i)



Current Data Parameters

NAME test
EXPNO 118
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180606
Time_ 15.18
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 64
DS 0
SWH 7246.377 Hz
FIDRES 0.221142 Hz
AQ 2.2609921 sec
RG 181
DW 69.000 usec
DE 6.50 usec
TE 300.1 K
D1 2.00000000 sec
TD0 1

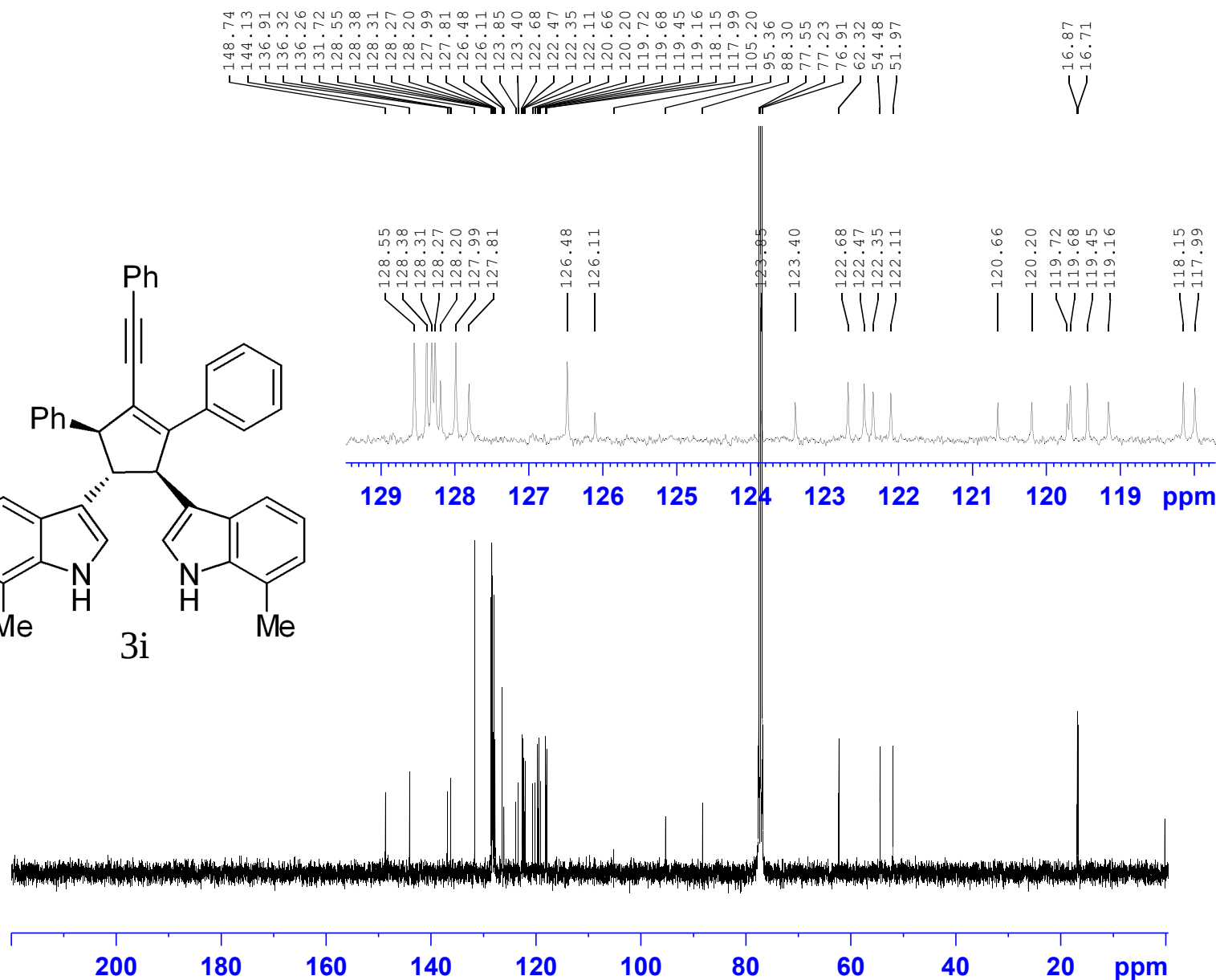
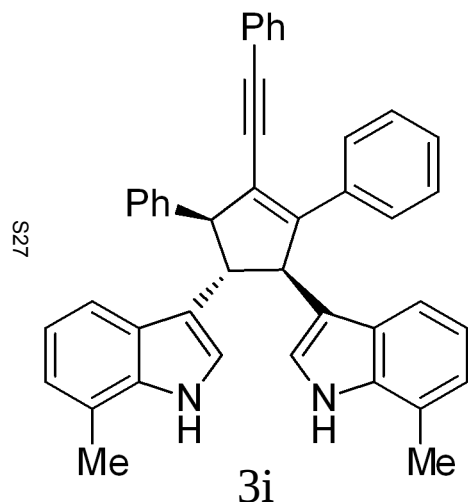
===== CHANNEL f1 =====

NUC1 1H
P1 15.00 usec
PL1 0 dB
SFO1 400.1324008 MHz

F2 - Processing parameters

SI 16384
SF 400.1300118 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3i)



Current Data Parameters
NAME test
EXPNO 67
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180102
Time 21.57
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 1319
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5792.6
DW 20.800 usec
DE 6.50 usec
TE 298.5 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

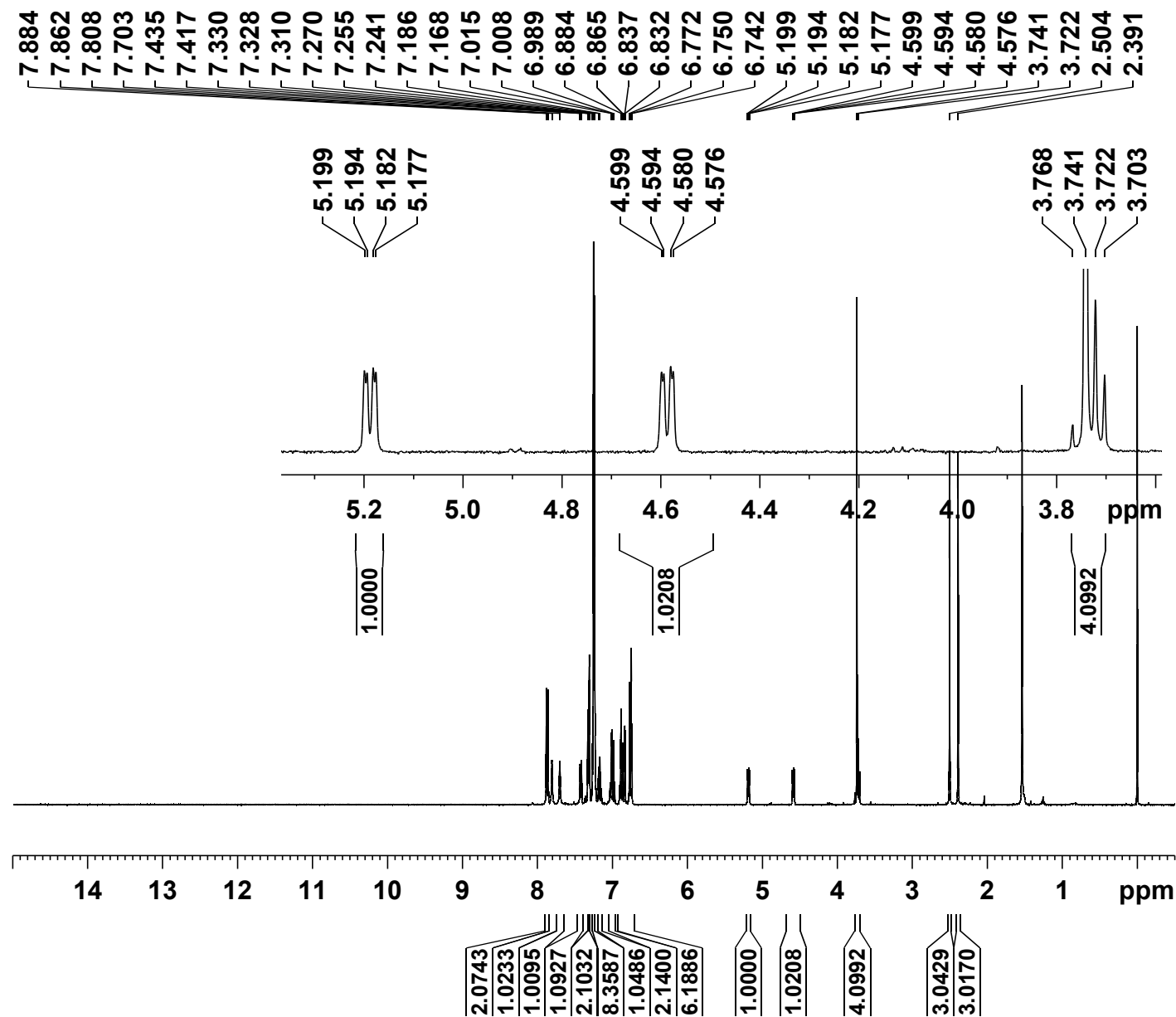
===== CHANNEL f1 =====
NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

F2 - Processing parameters
SI 32768
SF 100.6127477 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

3,3'-(3-(4-methoxyphenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3j)

823



Current Data Parameters

NAME try
EXPNO 84
PROCNO 1

F2 - Acquisition Parameters

Date 20180717
Time 19.35
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 30
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 299.0 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

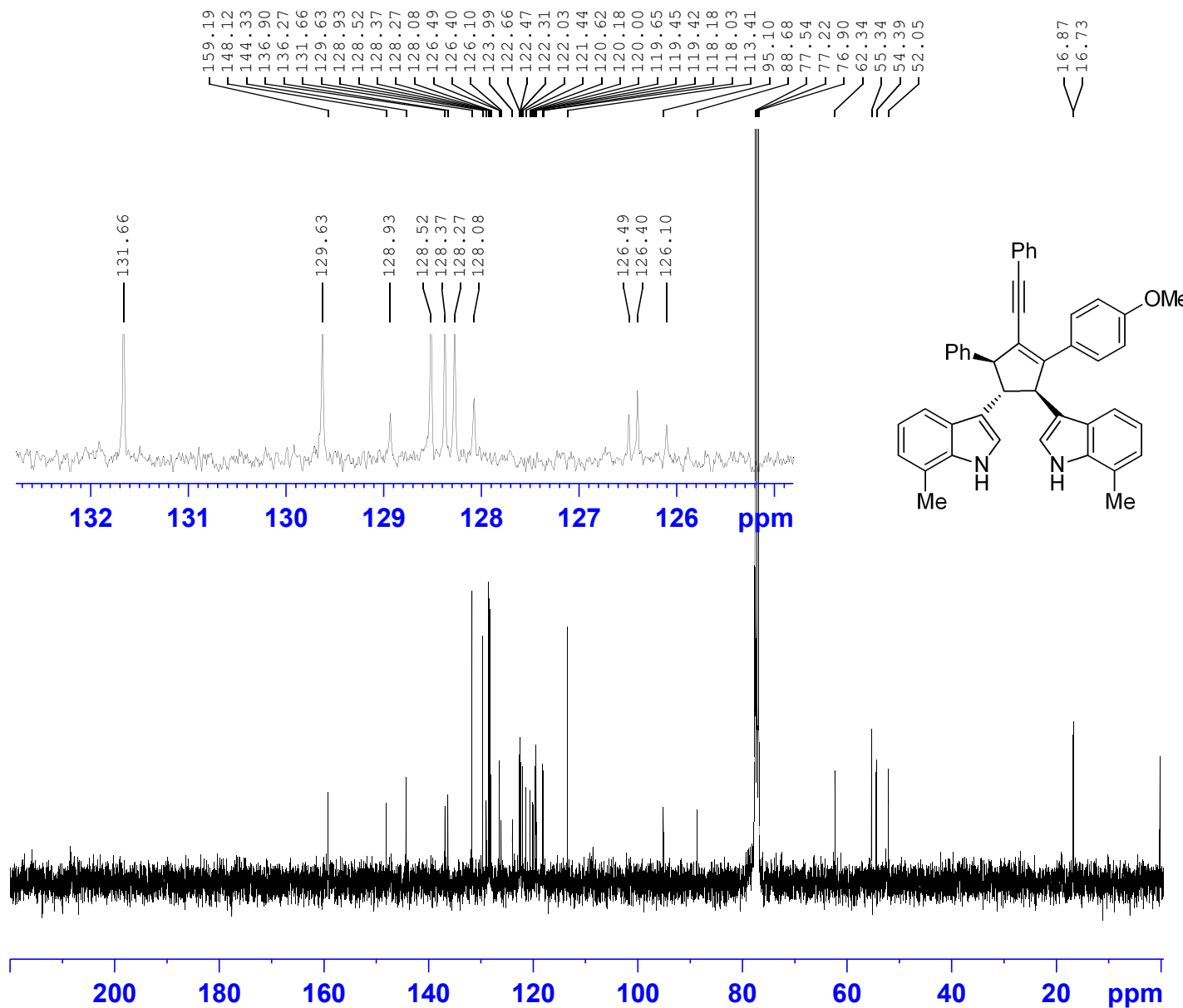
F2 - Processing parameters

SI 16384
SF 400.1300112 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-(4-methoxyphenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3j)



62S



Current Data Parameters

NAME test
EXPNO 80
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180114
Time_ 20.25
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 2000
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 8192
DW 20.800 usec
DE 6.50 usec
TE 298.6 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====

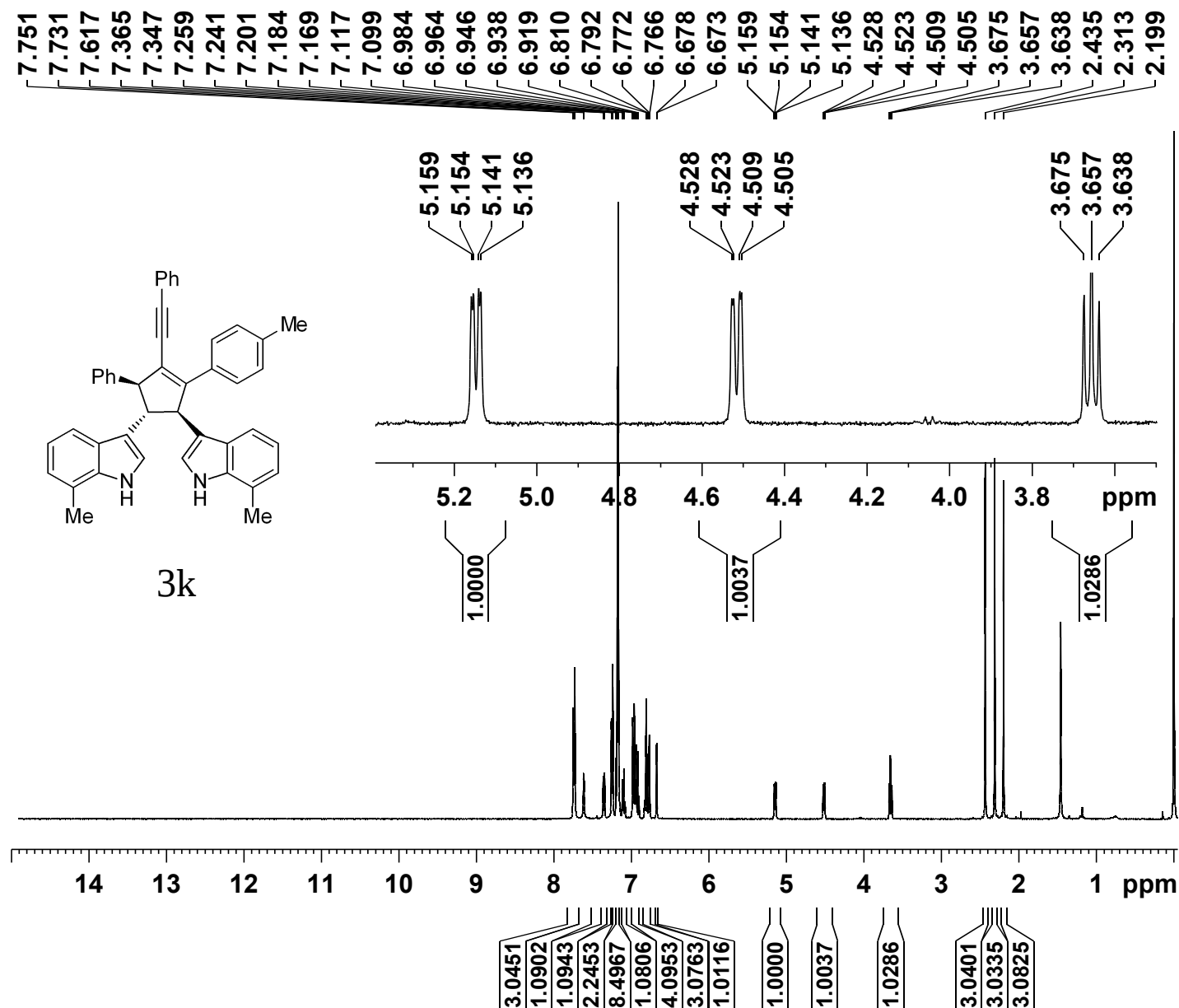
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

F2 - Processing parameters

SI 32768
SF 100.6127477 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

3,3'-(5-phenyl-4-(phenylethynyl)-3-(p-tolyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3k)

083



Current Data Parameters

NAME try
EXPNO 129
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180815
Time_ 22.12
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 33
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 298.0 K
D1 2.00000000 sec
TD0 1

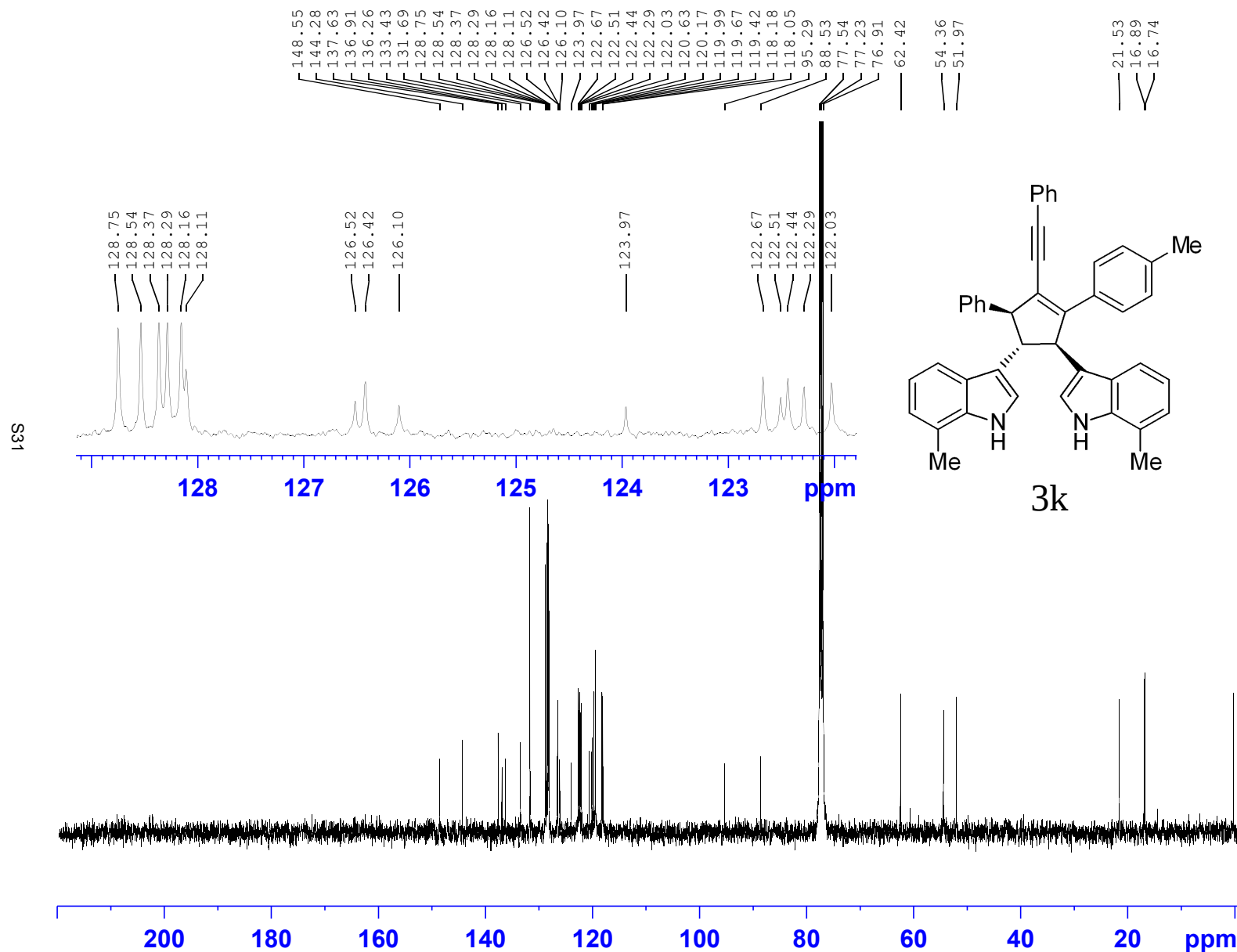
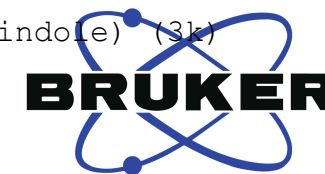
===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300398 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(5-phenyl-4-(phenylethynyl)-3-(p-tolyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3k)



Current Data Parameters
NAME try
EXPNO 12
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180121
Time 7.33
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 4784
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters
SI 32768
SF 100.6127473 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00

3,3'-(3-(4-chlorophenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (31)

Current Data Parameters

NAME try
EXPNO 132
PROCNO 1

F2 - Acquisition Parameters

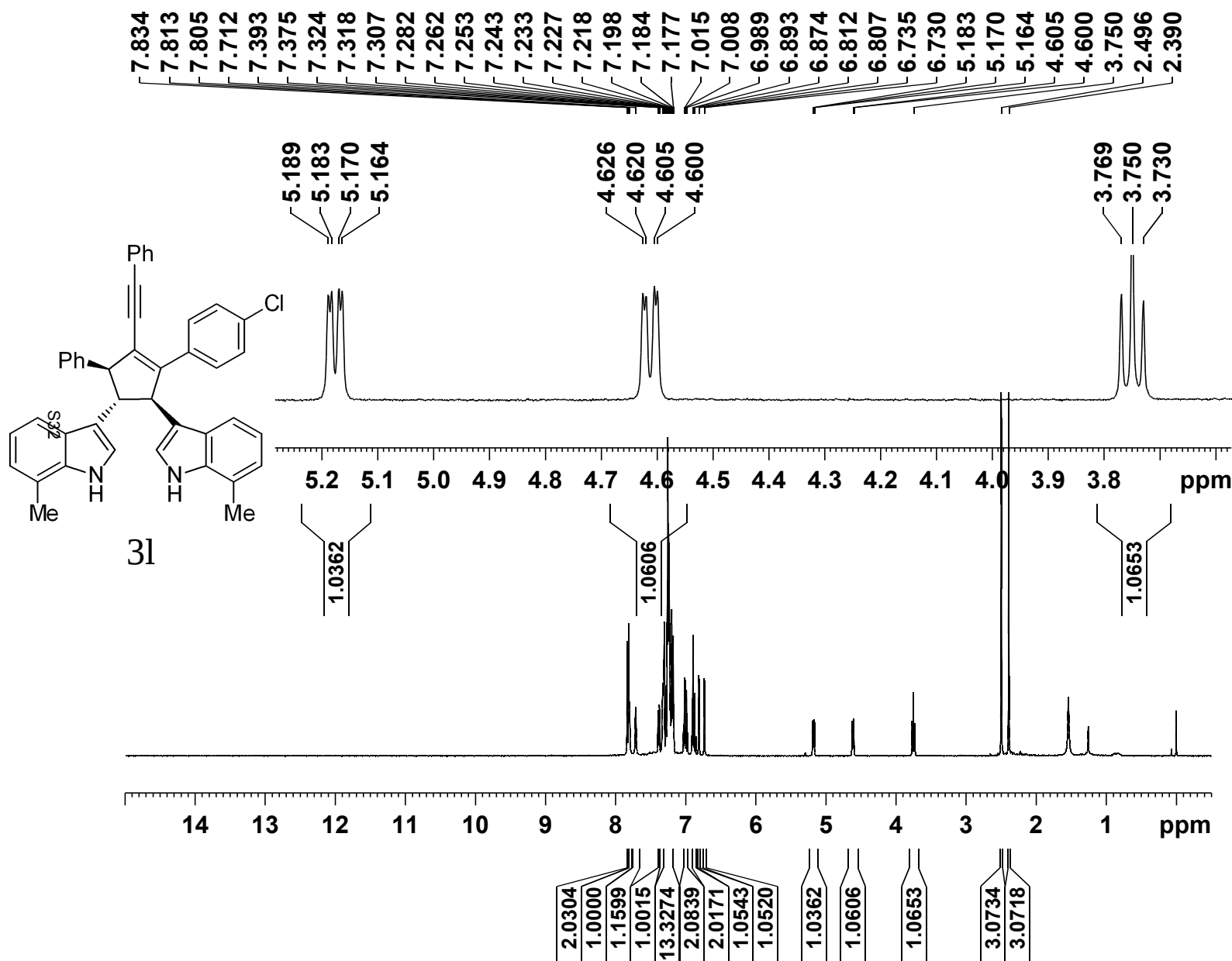
Date_ 20180817
Time_ 20.57
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 64
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 297.9 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300121 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00



3,3'-(3-(4-chlorophenyl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (31)



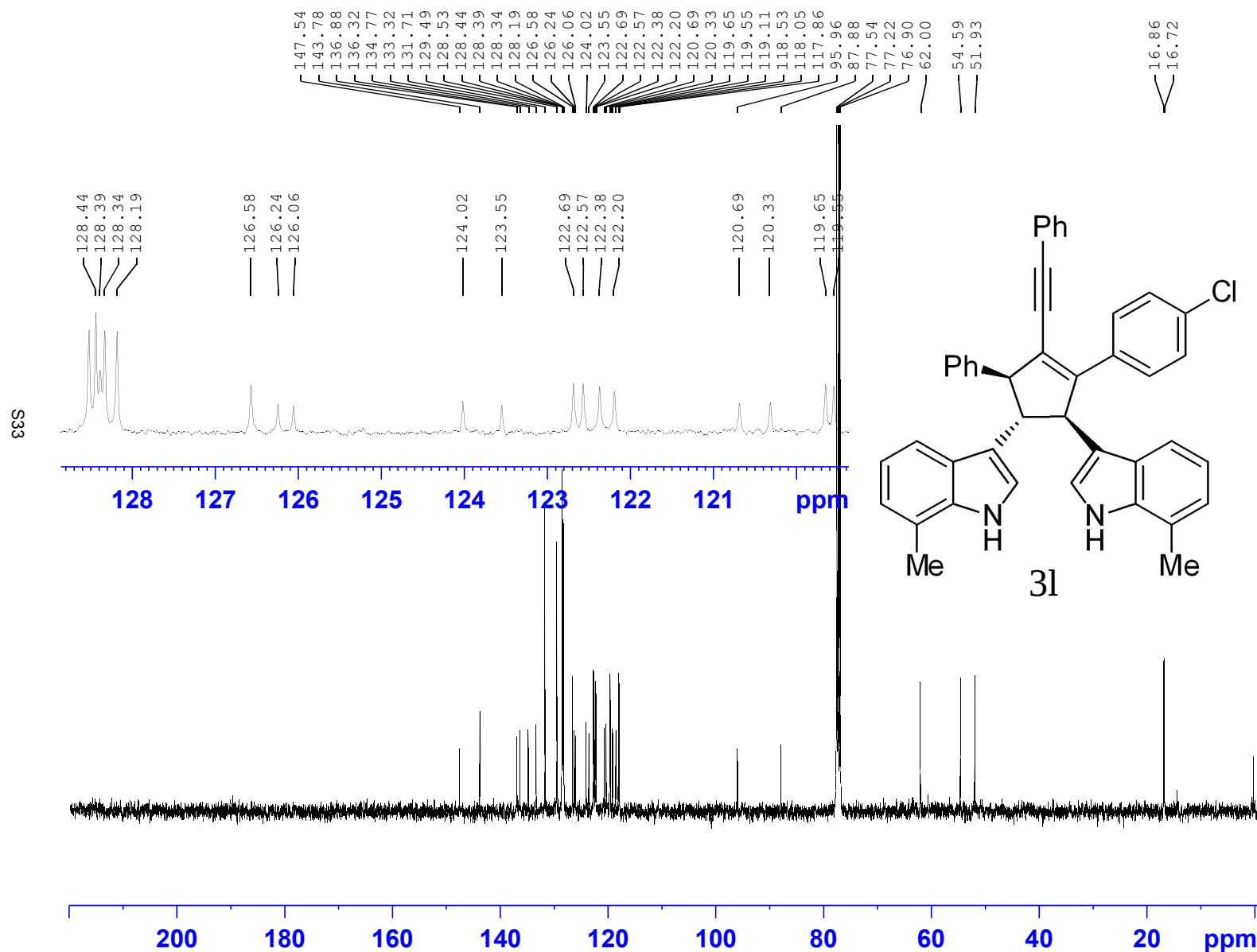
Current Data Parameters
NAME try
EXPNO 49
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180224
Time 21.47
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 1779
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 297.9 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

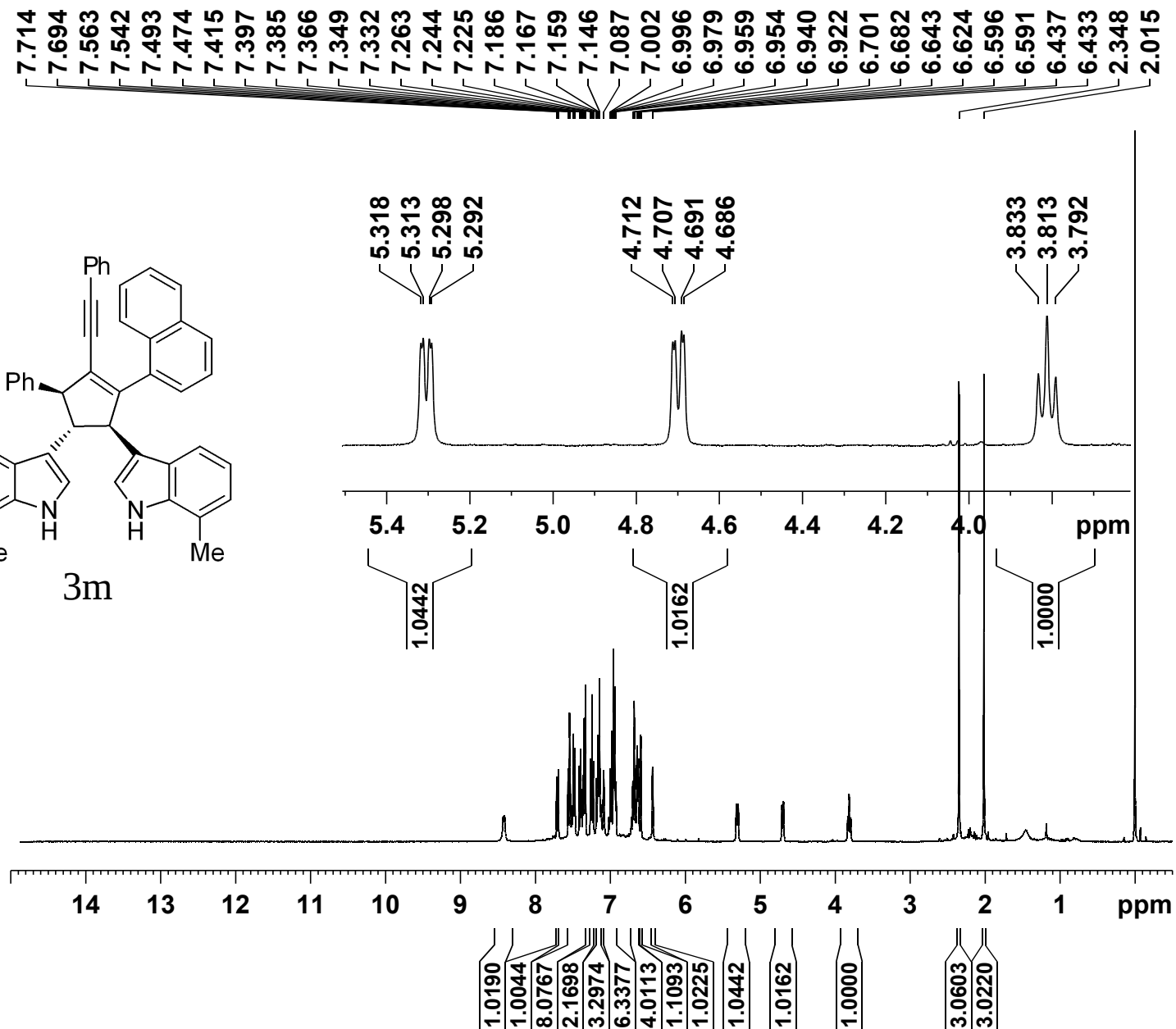
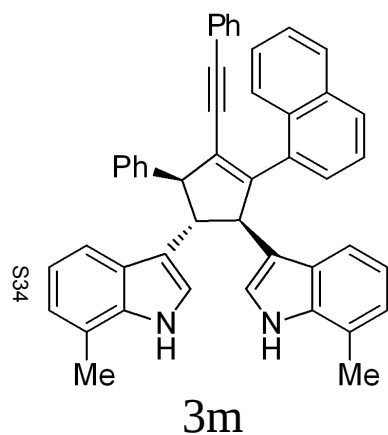
===== CHANNEL f1 =====
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters
SI 32768
SF 100.6127494 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00



3,3'-(3-(naphthalen-1-yl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3m)



Current Data Parameters

NAME try
EXPNO 122
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180812
Time_ 20.52
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 64
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 63.58
DW 69.333 usec
DE 10.50 usec
TE 297.8 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

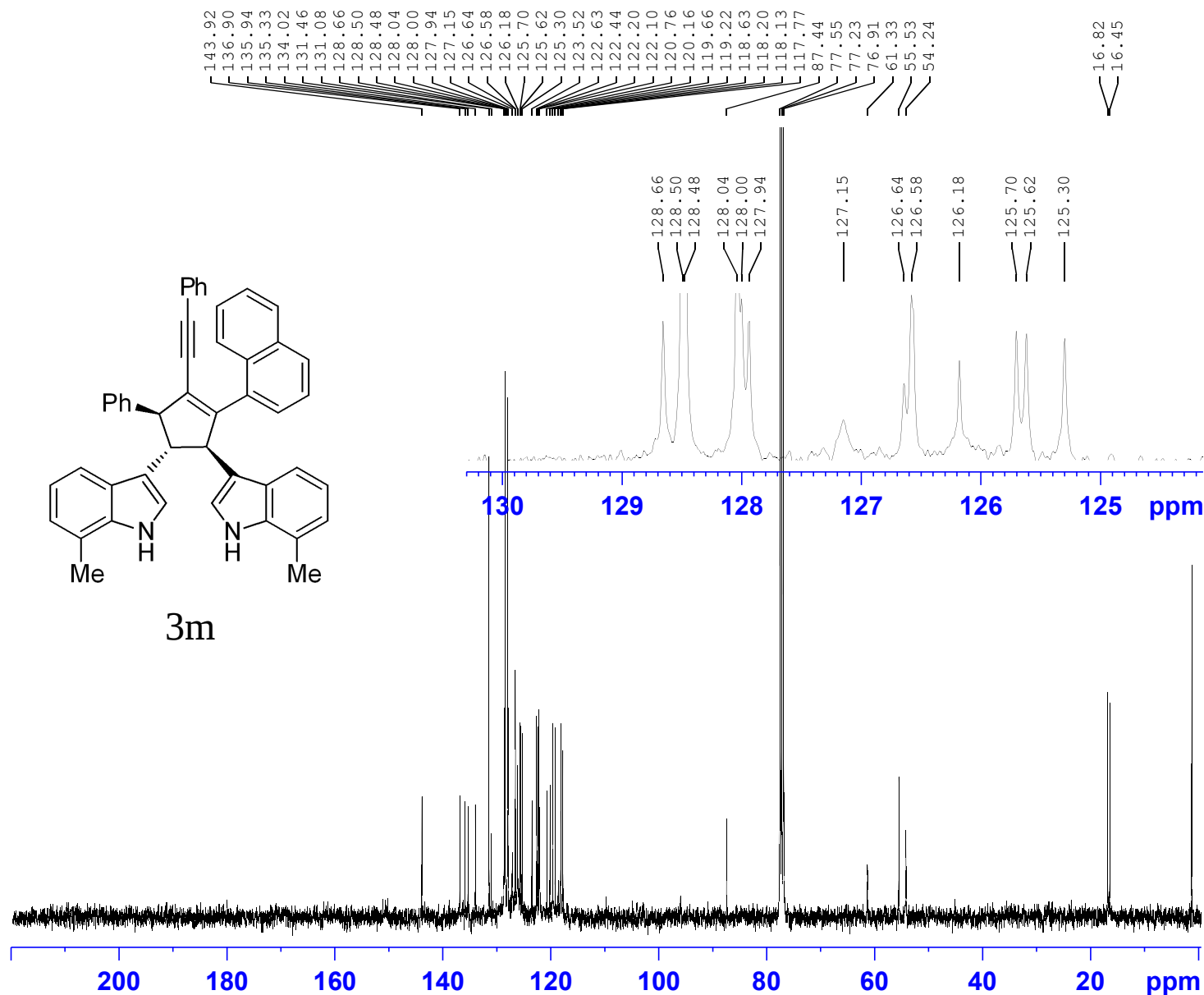
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300551 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-(naphthalen-1-yl)-5-phenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(7-methyl-1H-indole) (3m)

938



Current Data Parameters

NAME try
EXPNO 123
PROCNO 1

F2 - Acquisition Parameters

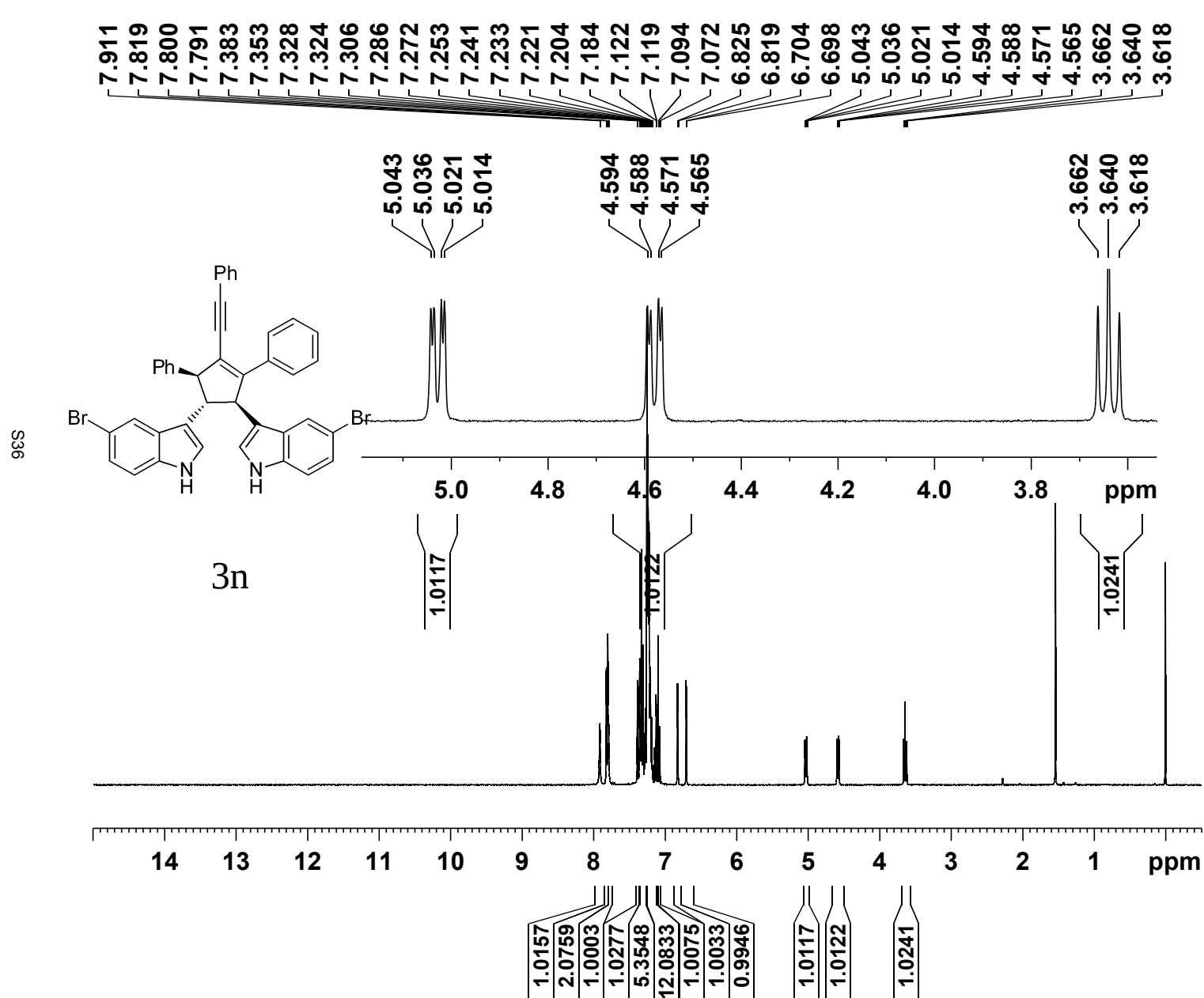
Date_ 20180812
Time_ 20.58
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 536
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.0 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters
SI 32768
SF 100.6127516 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(5-bromo-1H-indole) (3n)



Current Data Parameters
NAME try
EXPNO 92
PROCNO 1

F2 - Acquisition Parameters

Date 20180719
Time 18.14
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 34
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 297.5 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300121 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(5-bromo-1H-indole) (3n)



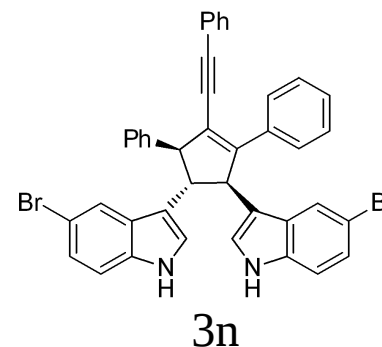
148.64
143.29
135.96
135.69
135.12
131.75
128.51
128.47
128.41
128.34
128.19
128.03
127.97
126.83
125.09
124.84
123.94
123.71
123.59
123.54
122.67
122.51
118.19
116.99
113.00
112.75
112.71
112.62
95.65
87.90
77.54
77.23
76.91
60.92
55.15
51.67

128.51
128.47
128.41
128.34
128.19
128.03
127.97

126.83

125.09
124.84

123.94
123.71
123.59
123.54



Current Data Parameters

NAME try
EXPNO 72
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180418
Time_ 8.00
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 428
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

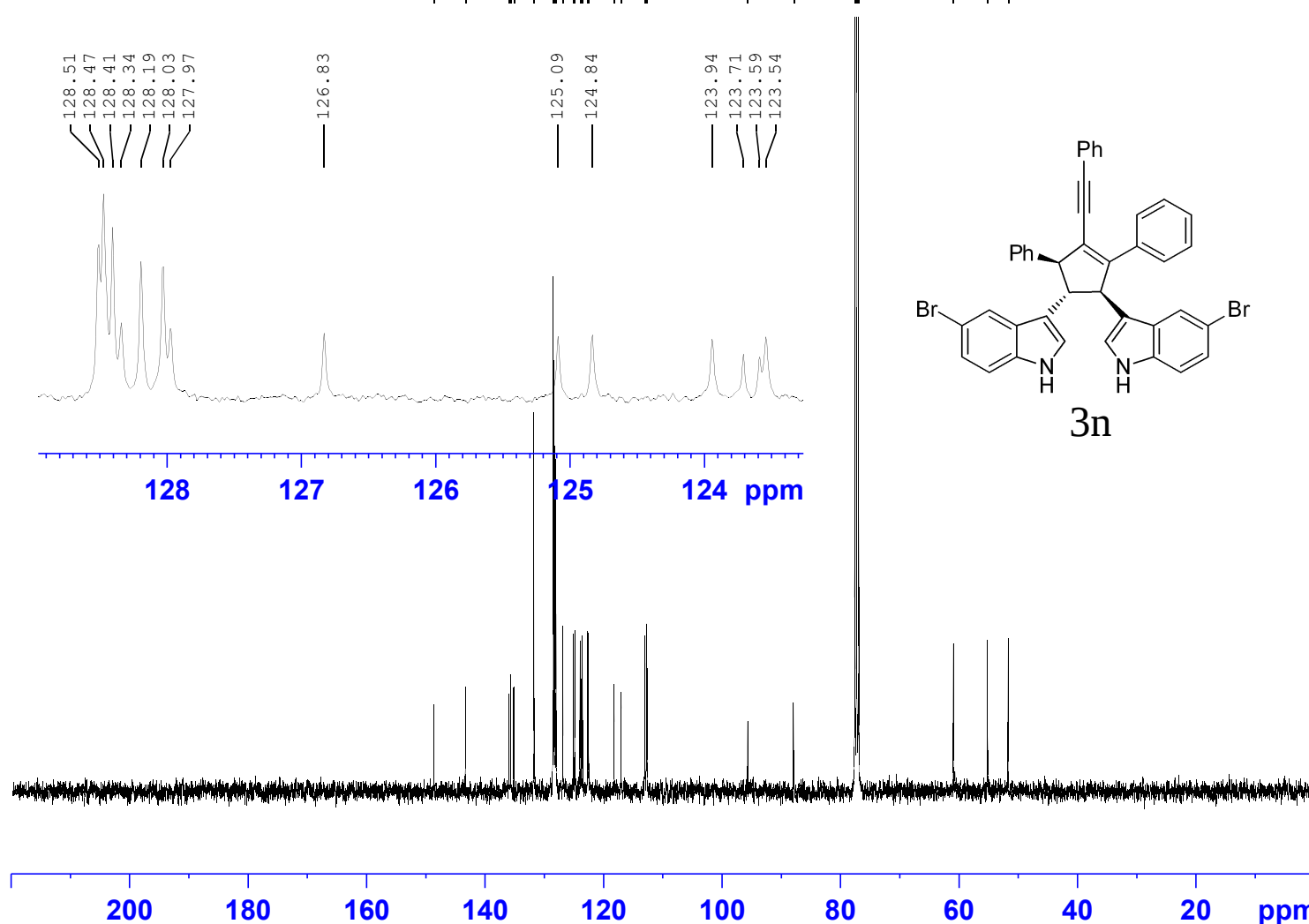
===== CHANNEL f2 =====

SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

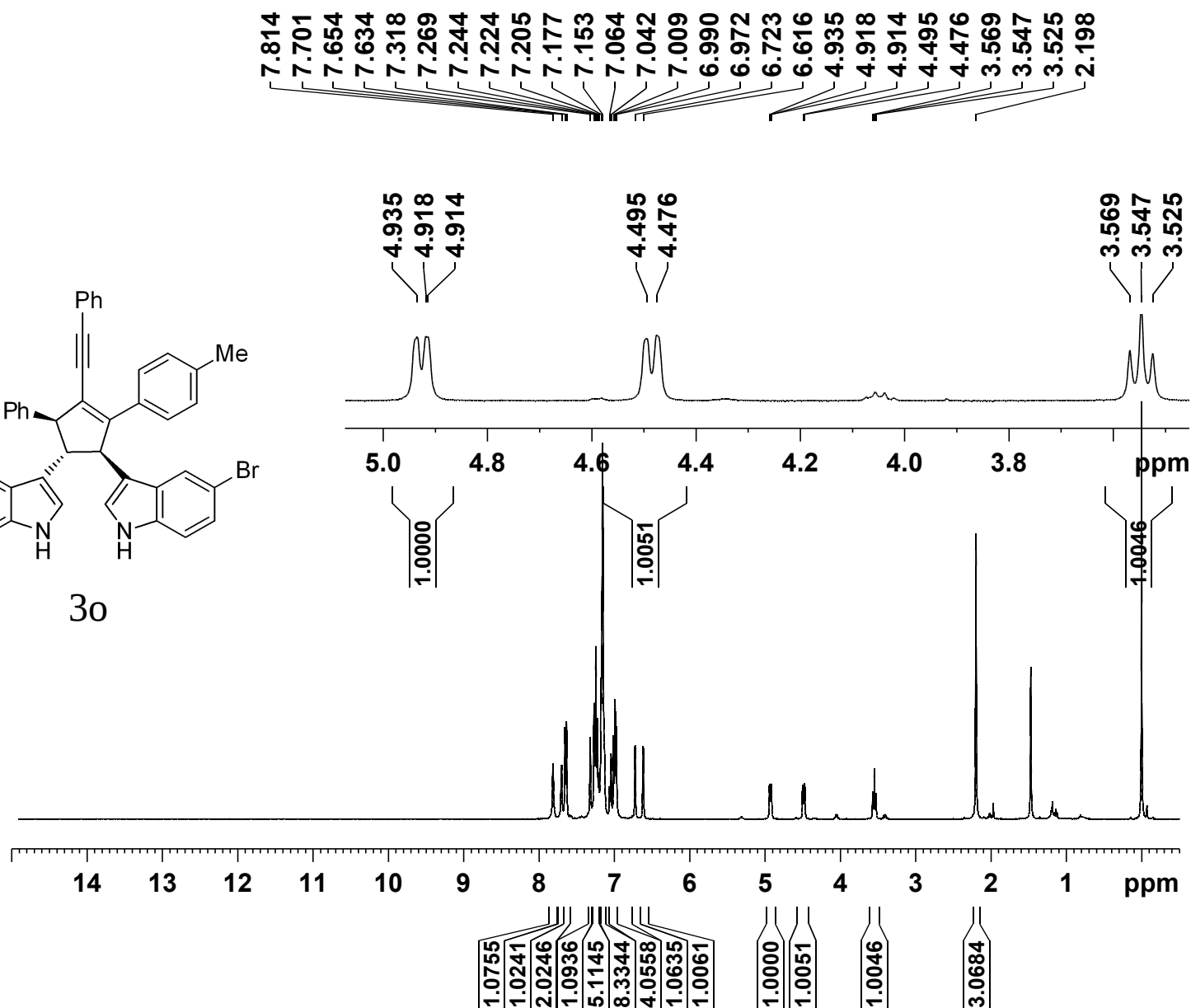
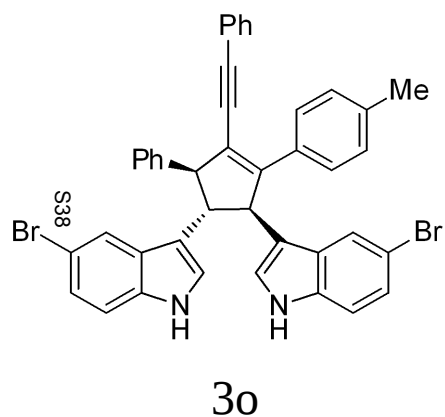
F2 - Processing parameters

SI 32768
SF 100.6127498 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00

128



3,3'-(5-phenyl-4-(phenylethynyl)-3-(p-tolyl)cyclopent-3-ene-1,2-diyl)bis(5-bromo-1H-indole) (3o)



Current Data Parameters

NAME try
EXPNO 128
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180815
Time_ 22.05
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 54
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 113.31
DW 69.333 usec
DE 10.50 usec
TE 298.1 K
D1 2.00000000 sec
TD0 1

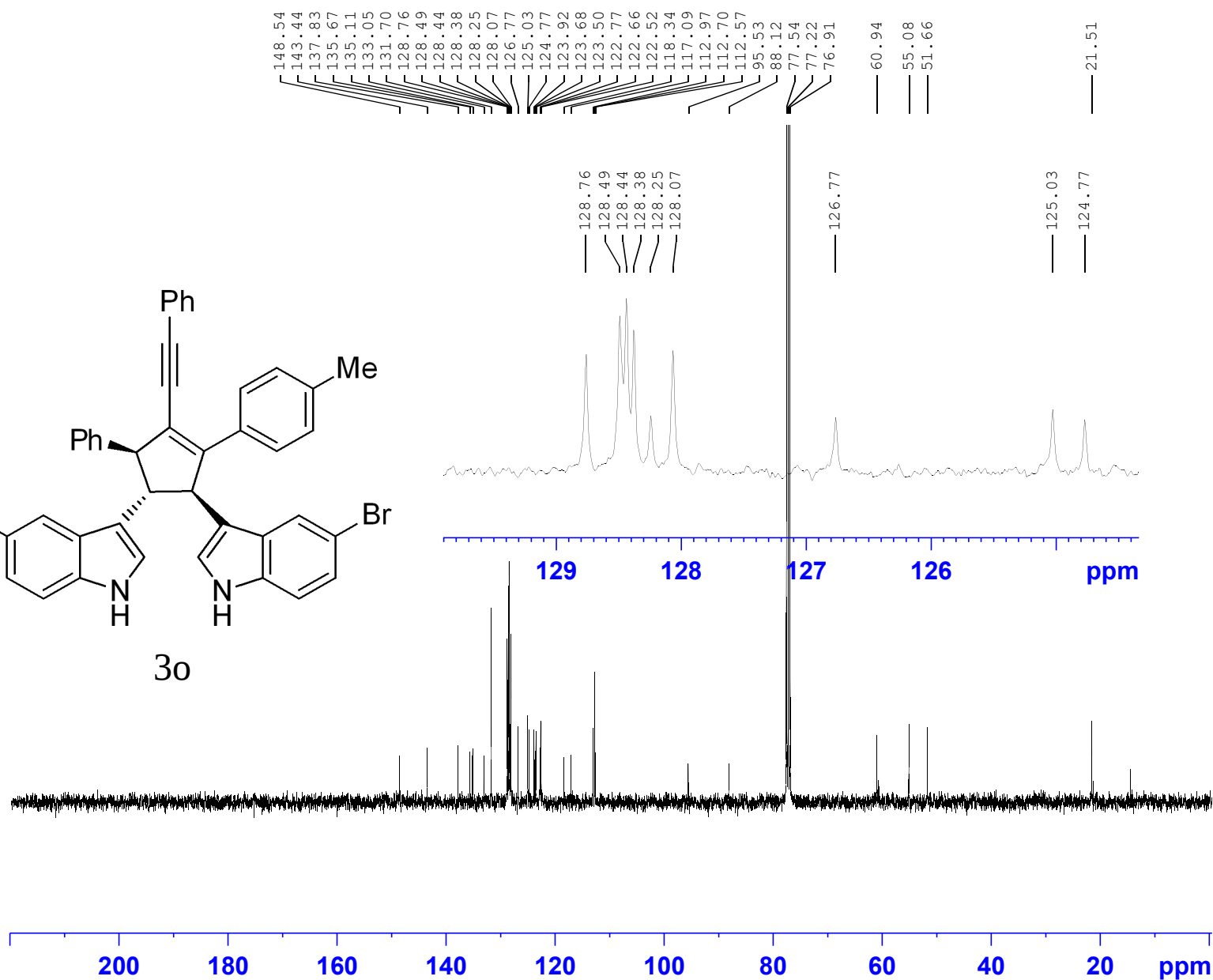
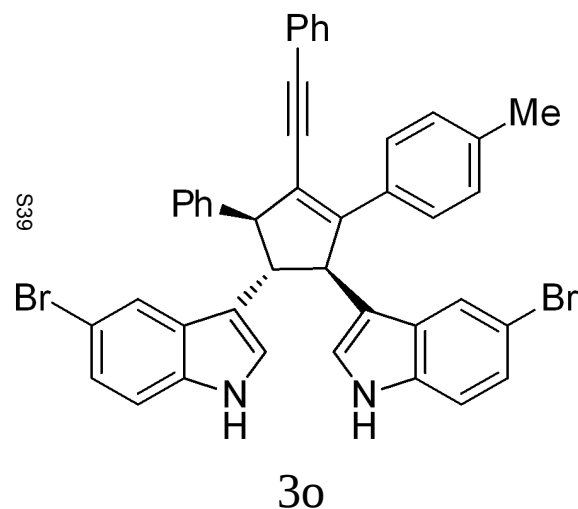
===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300426 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(5-phenyl-4-(phenylethynyl)-3-(p-tolyl)cyclopent-3-ene-1,2-diyl)bis(5-bromo-1H-indole) (3o)



Current Data Parameters
 NAME try
 EXPNO 67
 PROCNO 1

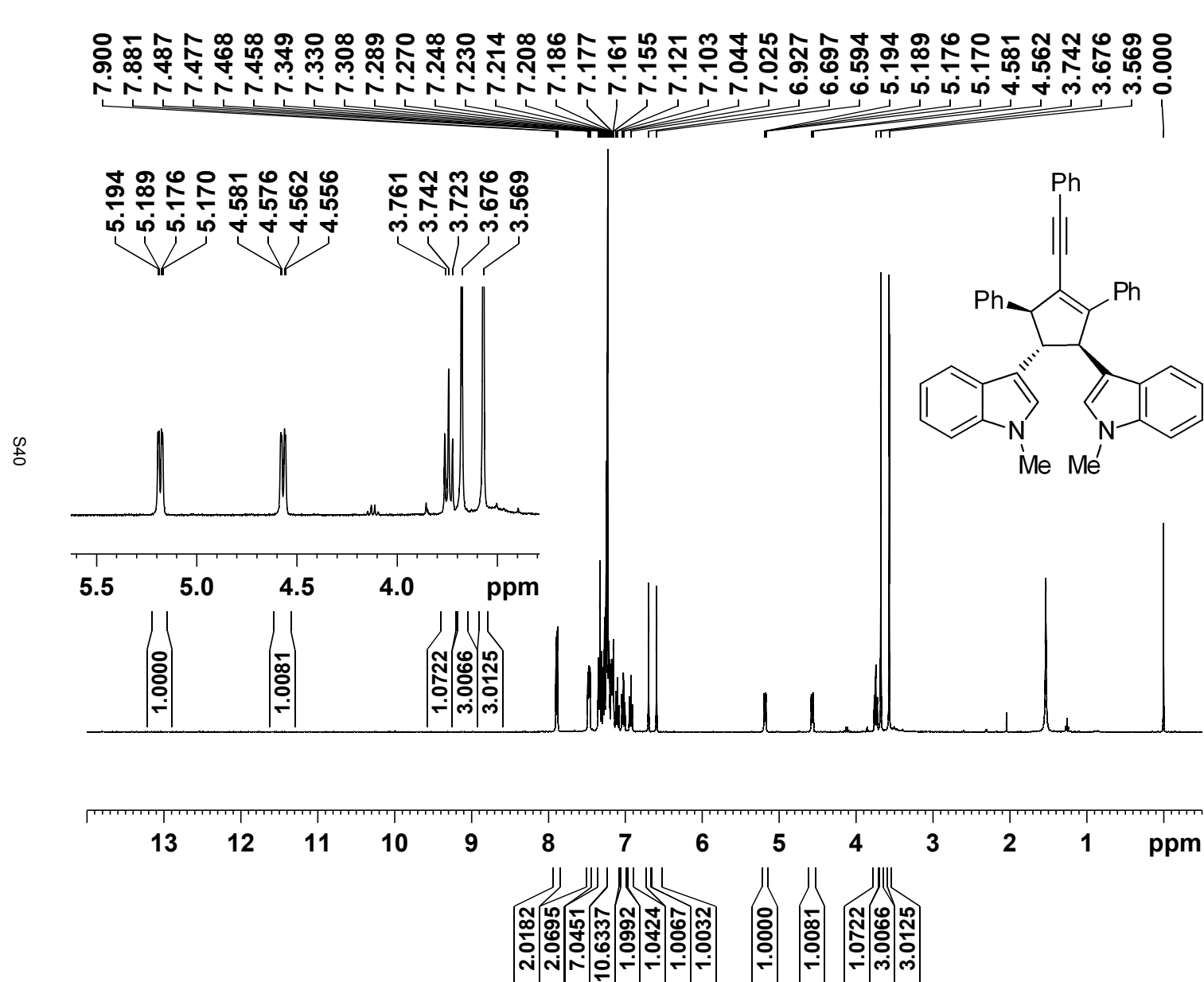
F2 - Acquisition Parameters
 Date_ 20180416
 Time_ 11.31
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 134
 DS 0
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 198.09
 DW 20.800 usec
 DE 6.50 usec
 TE 297.2 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====
 SFO1 100.6228298 MHz
 NUC1 13C
 P1 10.00 usec
 PLW1 47.50000000 W

===== CHANNEL f2 =====
 SFO2 400.1316005 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 15.00000000 W
 PLW12 0.33750001 W
 PLW13 0.27338001 W

F2 - Processing parameters
 SI 32768
 SF 100.6127516 MHz
 WDW EM
 SSB 0
 LB 2.00 Hz
 GB 0
 PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1-methyl-1H-indole) (3t)



Current Data Parameters

NAME test
EXPNO 60
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180103
Time_ 0
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 32
DS 0
SWH 7246.377 Hz
FIDRES 0.221142 Hz
AQ 2.2609921 sec
RG 181
DW 69.000 usec
DE 6.50 usec
TE 298.7 K
D1 2.00000000 sec
TD0 1

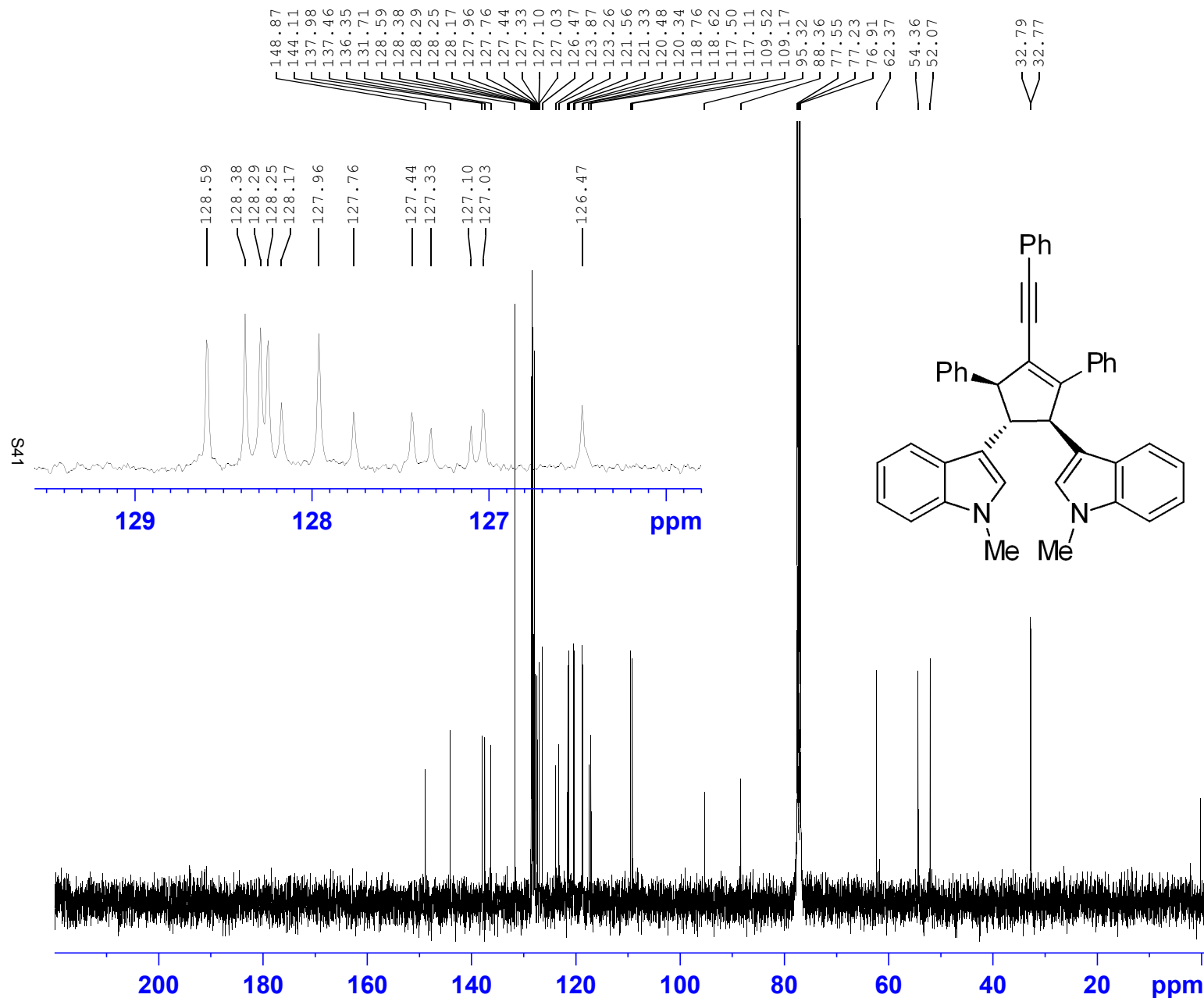
===== CHANNEL f1 =====

NUC1 1H
P1 15.00 usec
PL1 0 dB
SFO1 400.1324008 MHz

F2 - Processing parameters

SI 16384
SF 400.1300141 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1-methyl-1H-indole) (3t)



Current Data Parameters
NAME test
EXPNO 61
PROCNO 1

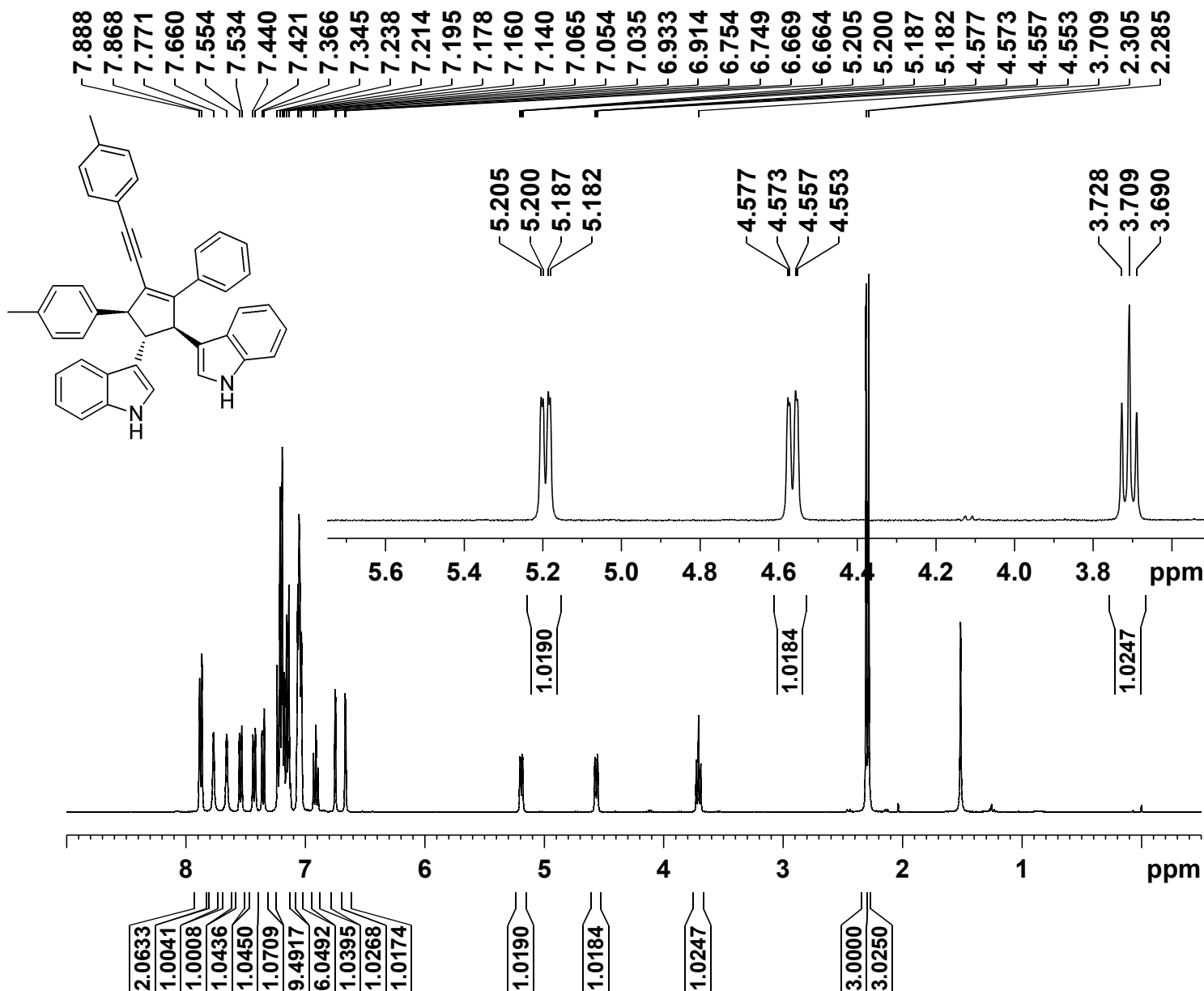
F2 - Acquisition Parameters
Date_ 20180103
Time_ 0.03
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 1239
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5792.6
DW 20.800 usec
DE 6.50 usec
TE 298.7 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

F2 - Processing parameters
SI 32768
SF 100.6127477 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

3,3'-(3-phenyl-5-(p-tolyl)-4-(p-tolyneethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3u)



Current Data Parameters

NAME try
EXPNO 163
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180927
Time_ 12.55
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 27
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 144.49
DW 69.333 usec
DE 10.50 usec
TE 297.8 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300182 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3-phenyl-5-(p-tolyl)-4-(p-tolylethynyl)cyclopent-3-ene-1,2-diyl)bis(1H-indole) (3u)



Cur
NAME try
EXPNO 164
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180927
Time 12.59
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 1359
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

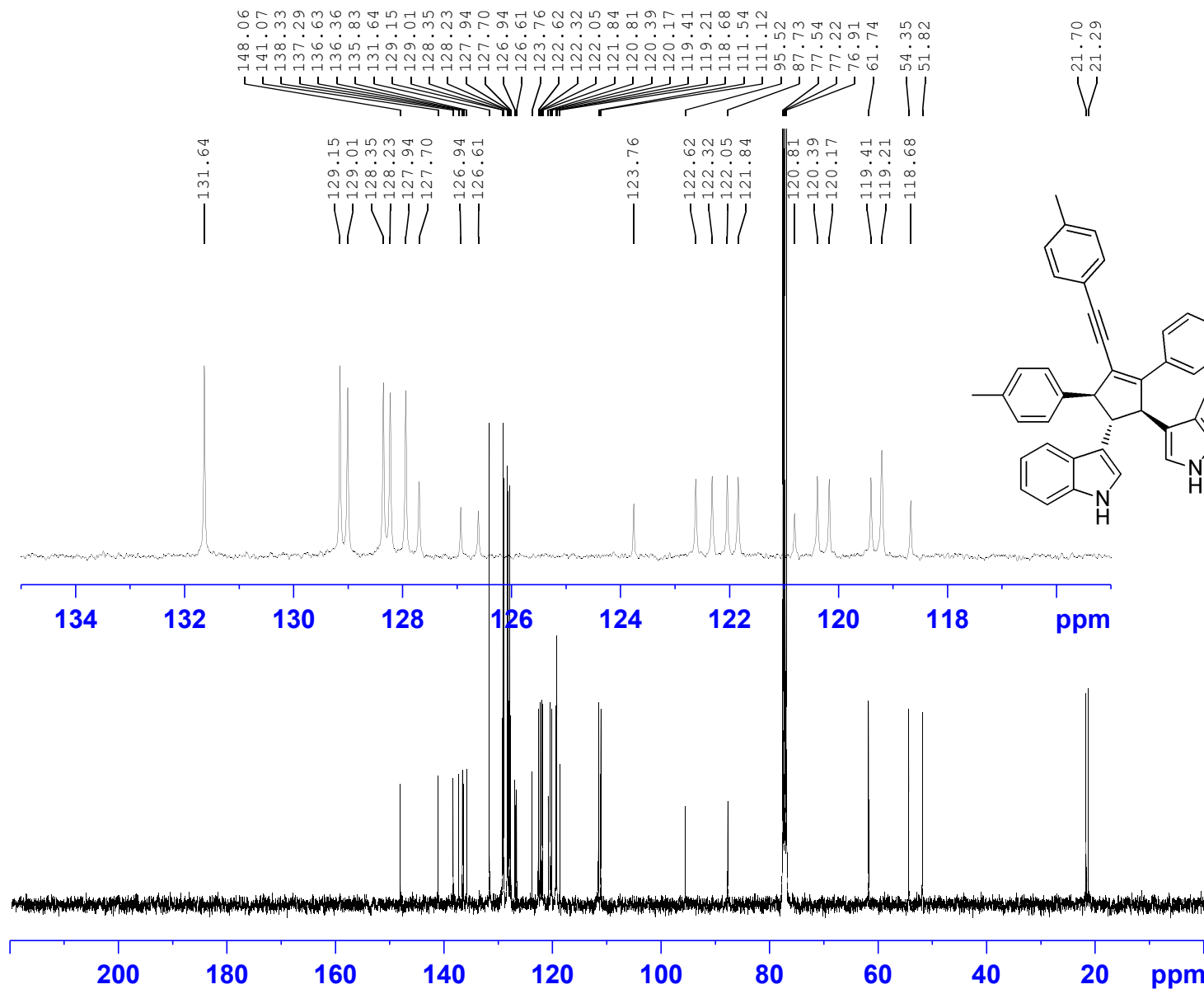
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====

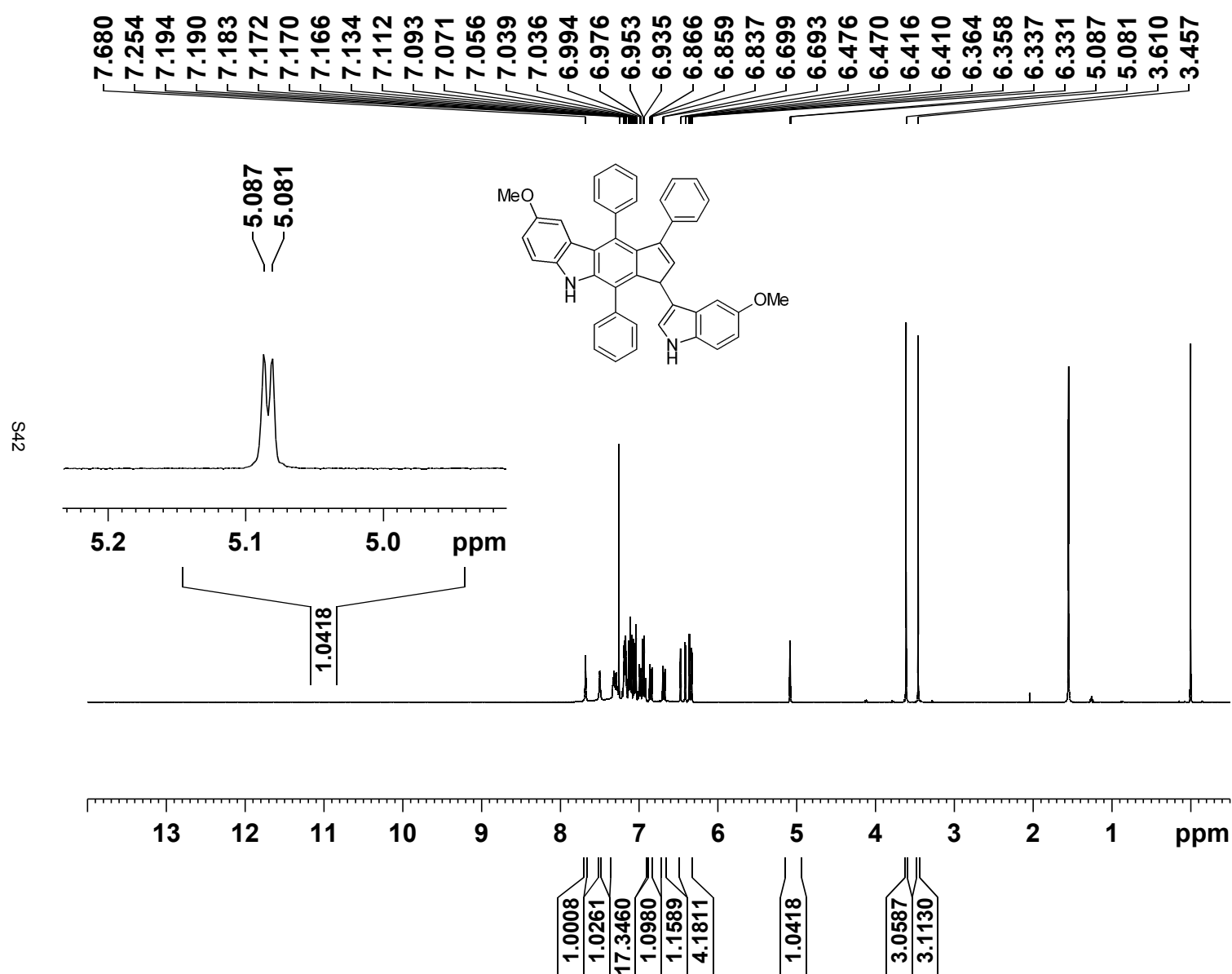
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters

SI 32768
SF 100.6127494 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00



8-methoxy-3-(5-methoxy-1H-indol-3-yl)-1,4,10-triphenyl-3,5-dihydrocyclopenta[b]carbazole (4p)



Current Data Parameters

NAME try
EXPNO 13
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180121
Time_ 12.39
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 33
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 297.5 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

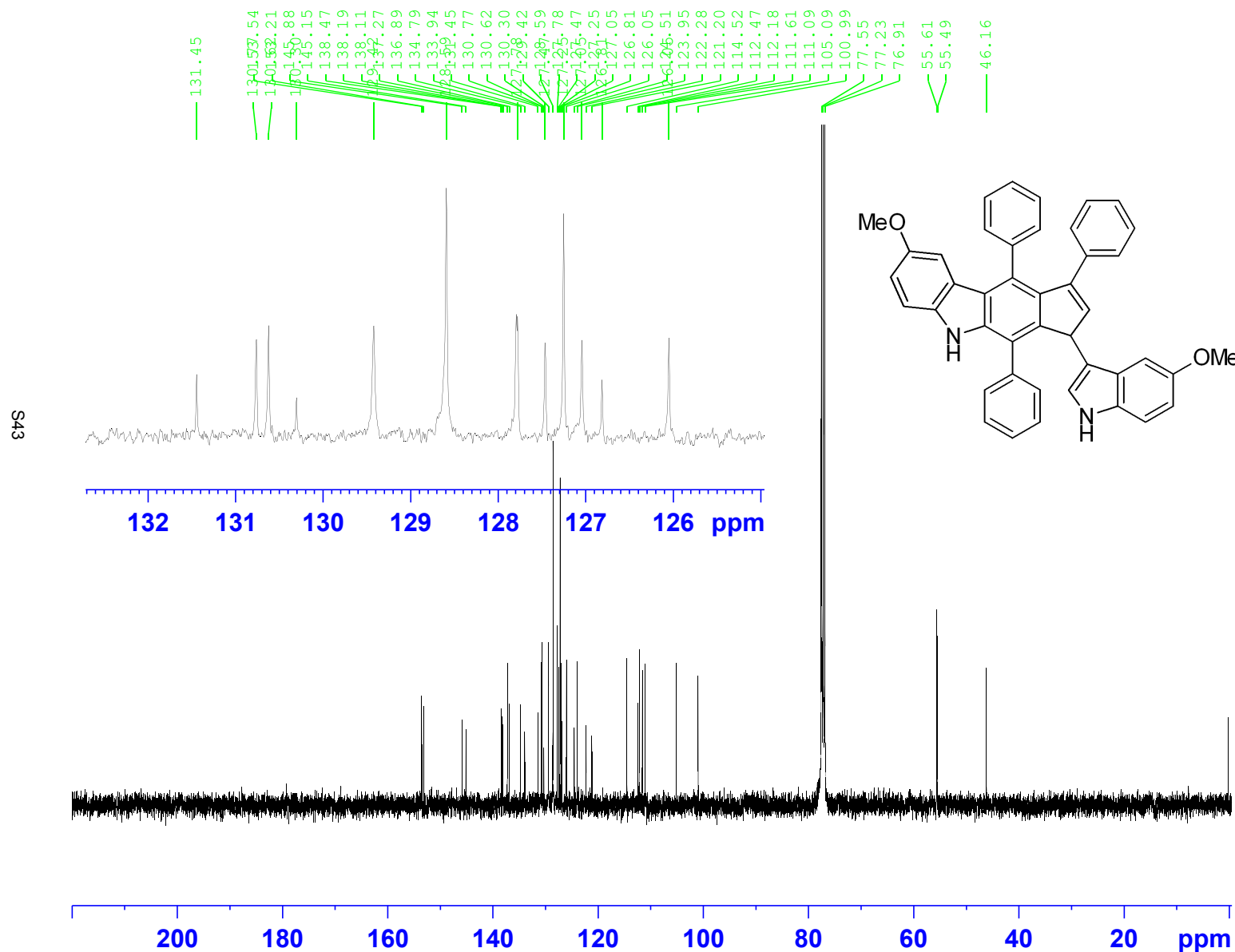
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300120 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

bazole (4p)

BRUKER



```
Current Data Parameters
NAME                test
EXPNO                86
PROCNO              1
```

```

F2 - Acquisition Parameters
Date_                20180122
Time_                22.11
INSTRUM              spect
PROBHD      5 mm BBO BB-1H
PULPROG              zgpg30
TD                   32768
SOLVENT              CDC13
NS                   3069
DS                   0
SWH                 24038.461 Hz
FIDRES              0.733596 Hz
AQ                 0.6815744 sec
RG                 5792.6
DW                 20.800 usec
DE                 6.50 usec
TE                 298.1 K
D1                 2.00000000 sec
D11                0.03000000 sec
TD0                 1

```

```
===== CHANNEL f1 =====
NUC1                13C
P1                  10.45 usec
PL1                 7.00 dB
SFO1                100.6233325 MHz
```

```

===== CHANNEL f2 =====
CPDPRG[2          waltz16
NUC2              1H
PCPD2            90.00 usec
PL2              0 dB
PL12             15.00 dB
PL13             20.00 dB
SFO2            400.1316005 MHz

```

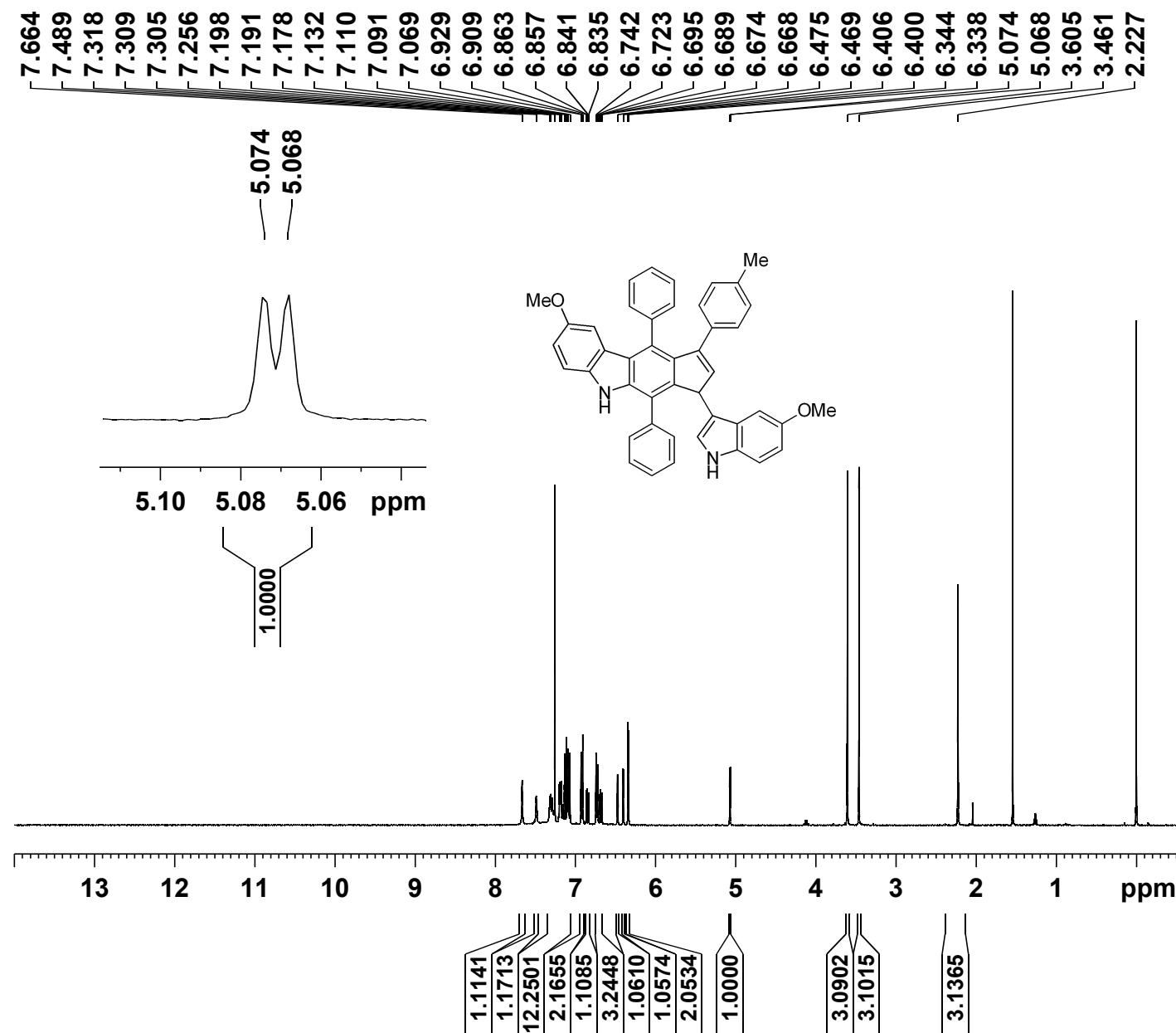
```

F2 - Processing parameters
SI                32768
SF              100.6127470 MHz
WDW                EM
SSB                0
LB                1.00 Hz
GB                0
PC                1.00

```

8-methoxy-3-(5-methoxy-1H-indol-3-yl)-4,10-diphenyl-1-(p-tolyl)-3,5-dihydrocyclopenta[b]carbazole (4q)

S44



Current Data Parameters

NAME try
EXPNO 22
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180124
Time_ 21.38
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 16
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 297.7 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300113 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

8-methoxy-3-(5-methoxy-1H-indol-3-yl)-4,10-diphenyl-1-(p-tolyl)-3,5-dihydrocyclopenta[b]carbazole (4q)



Current Data Parameters
 NAME test
 EXPNO 96
 PROCNO 1

F2 - Acquisition Parameters

Date_ 20180127
 Time 23.15
 INSTRUM spect
 PROBHD 5 mm BBO BB-1H
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 2977
 DS 0
 SWH 24038.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 5792.6
 DW 20.800 usec
 DE 6.50 usec
 TE 296.5 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TD0 1

===== CHANNEL f1 =====

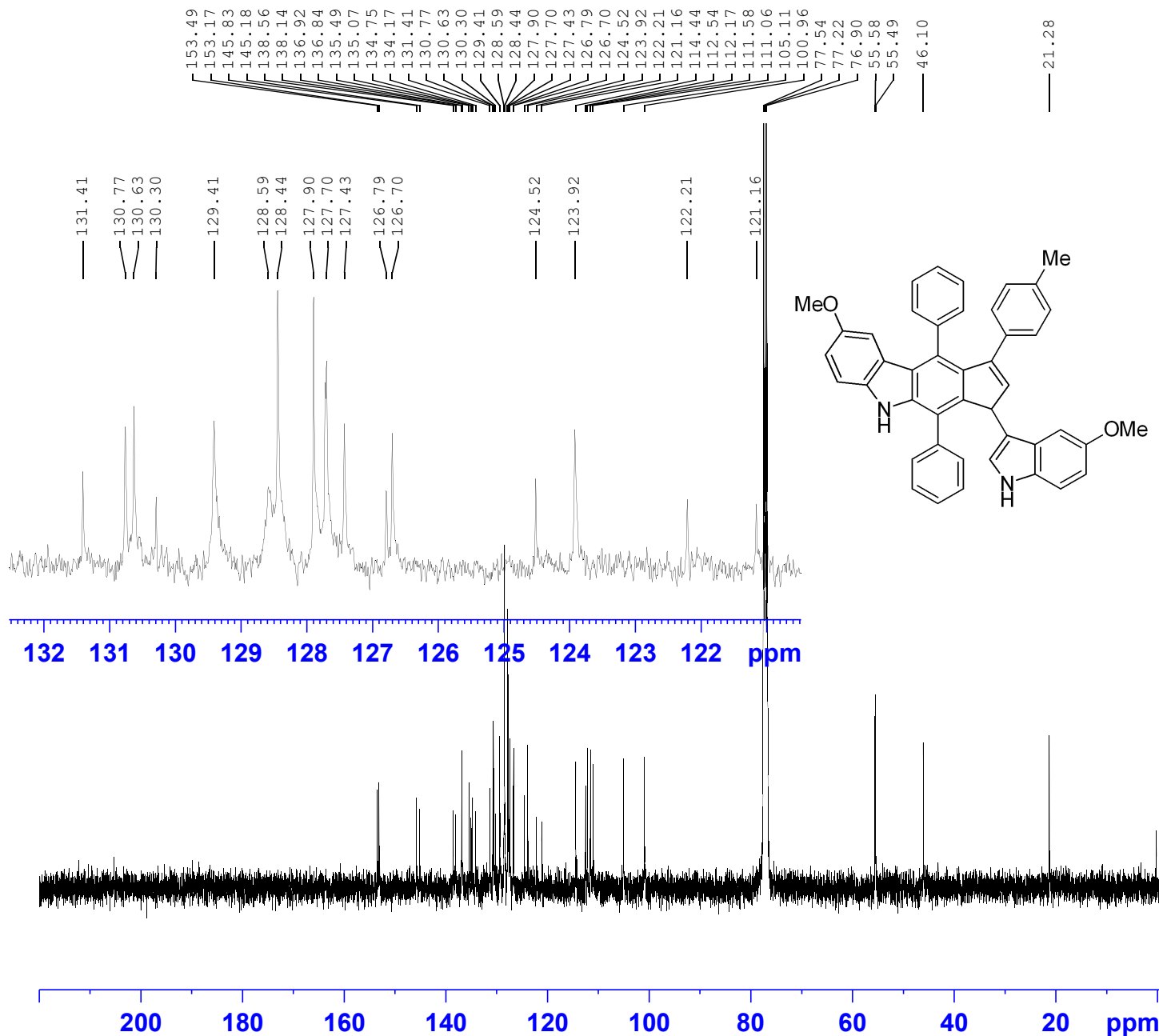
NUC1 13C
 P1 10.45 usec
 PL1 7.00 dB
 SFO1 100.6233325 MHz

===== CHANNEL f2 =====

CPDPRG[2] waltz16
 NUC2 1H
 PCPD2 90.00 usec
 PL2 0 dB
 PL12 15.00 dB
 PL13 20.00 dB
 SFO2 400.1316005 MHz

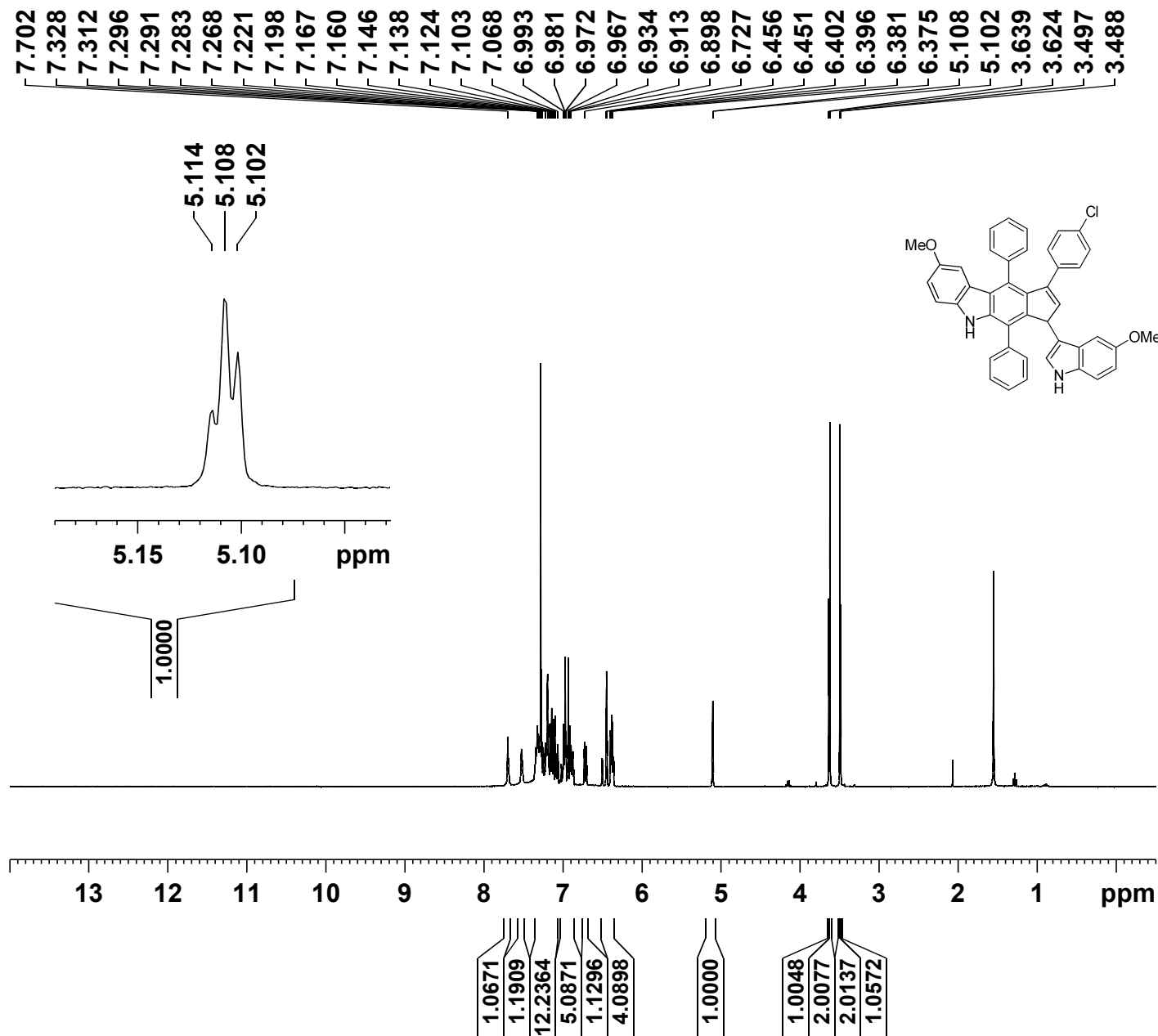
F2 - Processing parameters

SI 32768
 SF 100.6127484 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.00



1-(4-chlorophenyl)-8-methoxy-3-(5-methoxy-1H-indol-3-yl)-4,10-diphenyl-3,5-dihydrocyclopenta[b]carbazole (4r)

94S



Current Data Parameters

NAME try
EXPNO 37
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180221
Time_ 15.29
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 40
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 302.4 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300000 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

1-(4-chlorophenyl)-8-methoxy-3-(5-methoxy-1H-indol-3-yl)-4,10-diphenyl-3,5-dihydrocyclopenta[b]carbazole (4



Current Data Parameters

NAME try
EXPNO 38
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180221
Time_ 15.31
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 5469
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 302.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

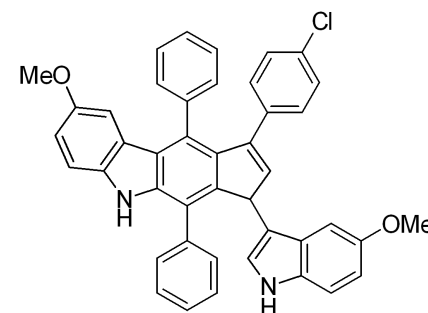
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====

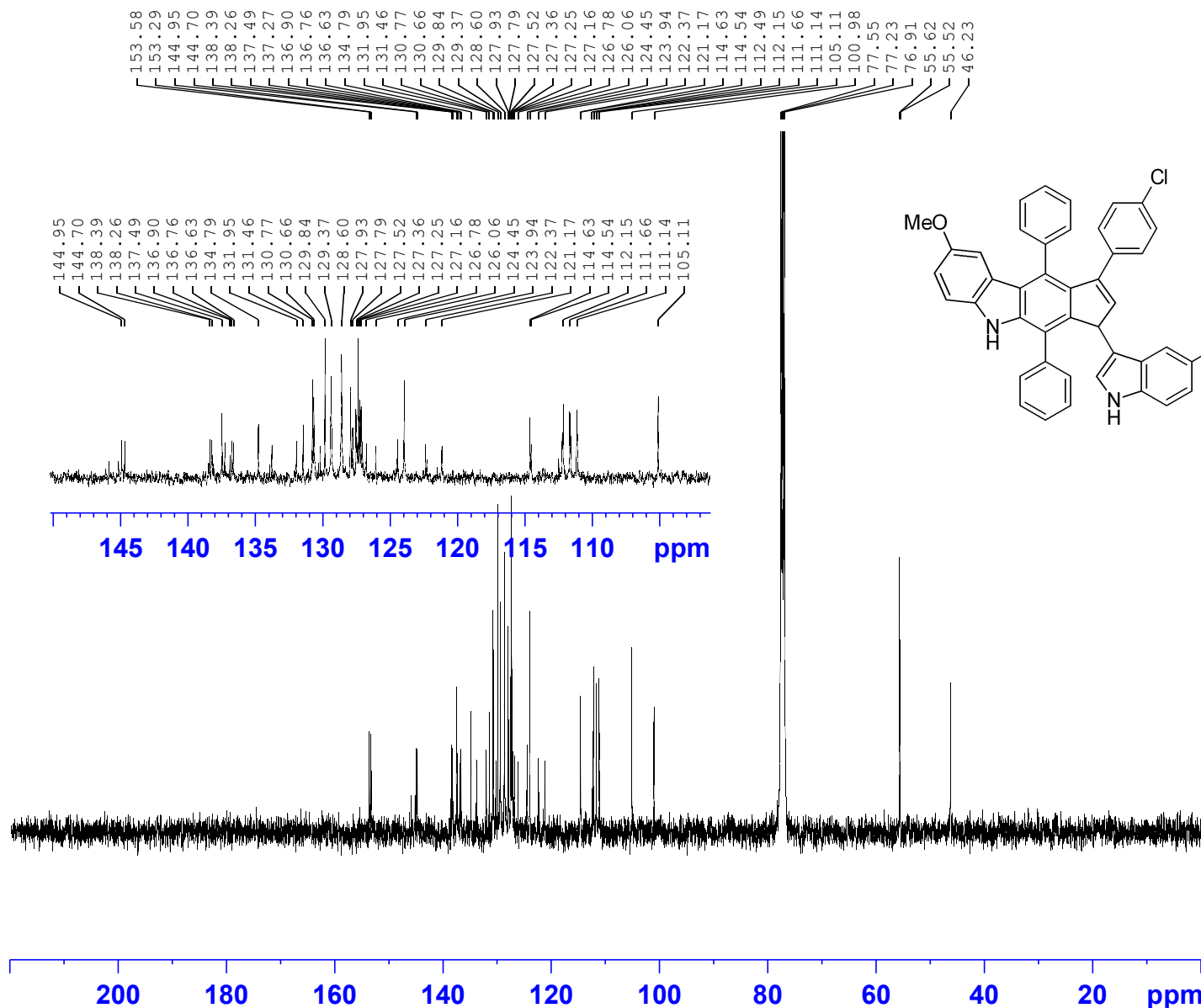
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters

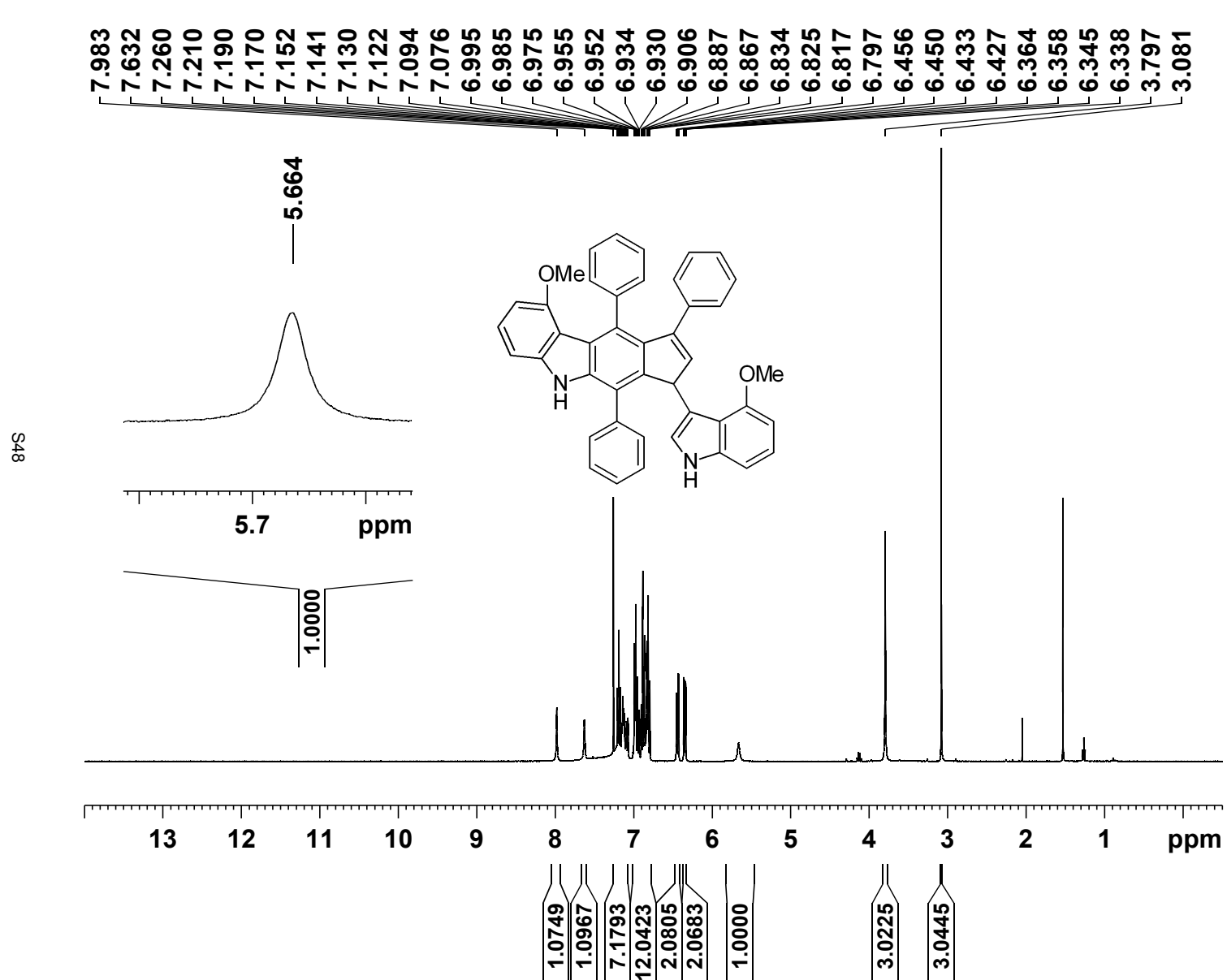
SI 32768
SF 100.6127464 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00



S47



9-methoxy-3-(4-methoxy-1H-indol-3-yl)-1,4,10-triphenyl-3,5-dihydrocyclopenta[b]carbazole (4s)



Current Data Parameters
NAME try
EXPNO 888
PROCNO 1

F2 - Acquisition Parameters
Date_ 20170810
Time_ 13.21
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 64
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 198.09
DW 69.333 usec
DE 10.50 usec
TE 298.2 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1300101 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

9-methoxy-3-(4-methoxy-1H-indol-3-yl)-1,4,10-triphenyl-3,5-dihydrocyclopenta[b]carbazole (4s)



Current Data Parameters
NAME try
EXPNO 889
PROCNO 1

F2 - Acquisition Parameters

Date_ 20170810
Time_ 13.27
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDC13
NS 904
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 298.3 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====

SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

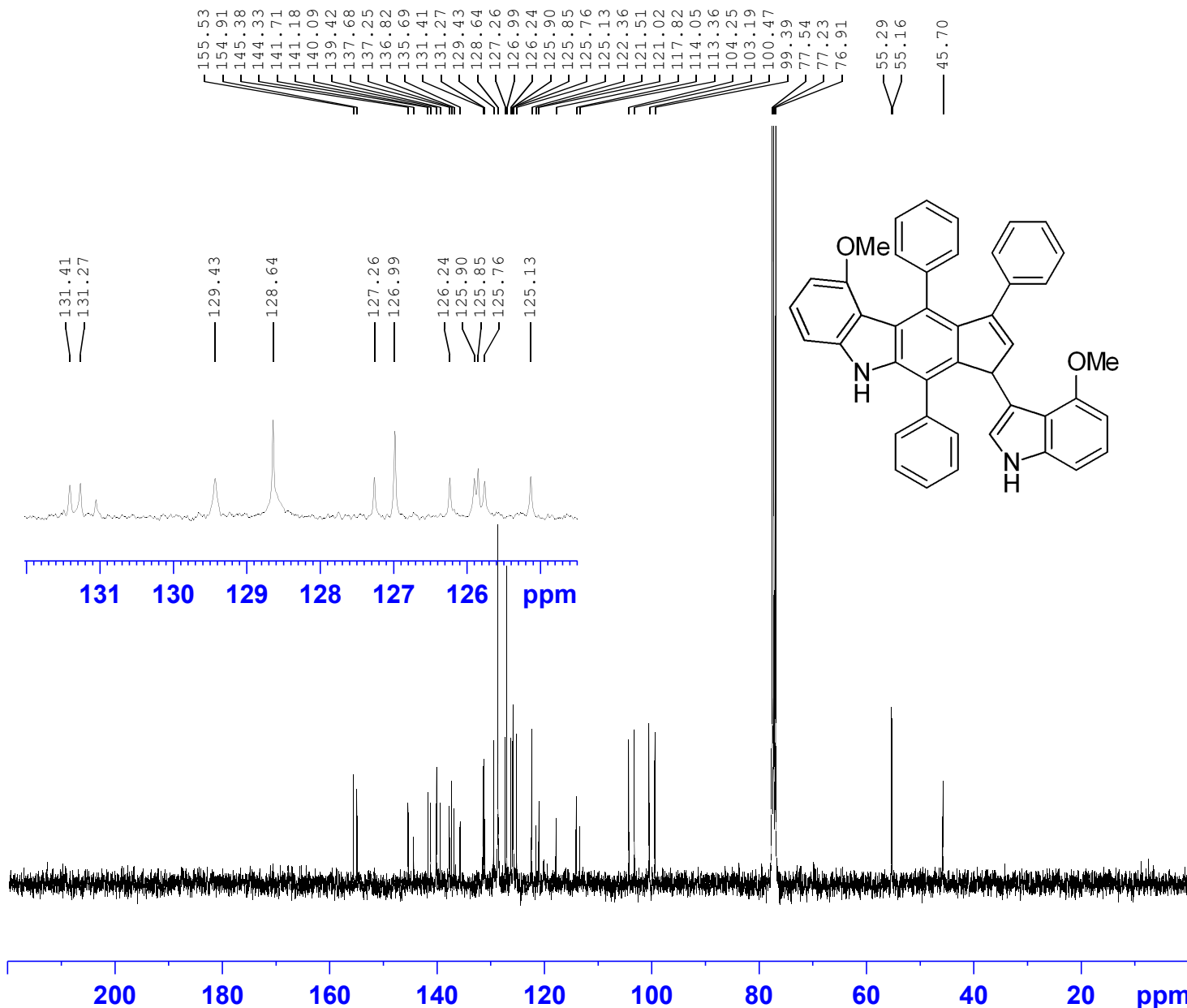
===== CHANNEL f2 =====

SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

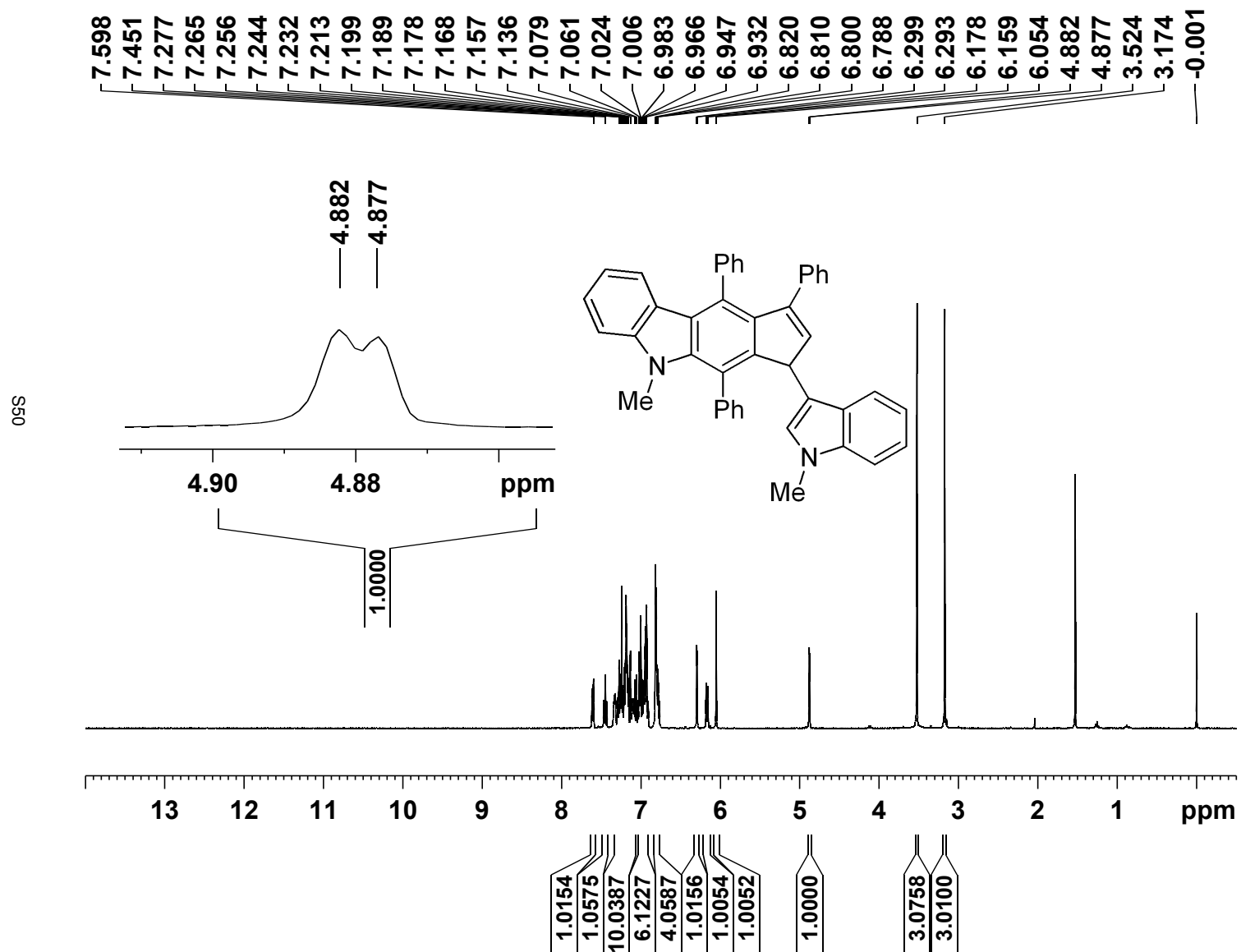
F2 - Processing parameters

SI 32768
SF 100.6127479 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00

64s



5-methyl-3-(1-methyl-1H-indol-3-yl)-1,4,10-triphenyl-3,5-dihydrocyclopenta[b]carbazole (4t)



Current Data Parameters

NAME test
EXPNO 54
PROCNO 1

F2 - Acquisition Parameters

Date_ 20170928
Time_ 21.23
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 50
DS 0
SWH 7246.377 Hz
FIDRES 0.221142 Hz
AQ 2.2609921 sec
RG 181
DW 69.000 usec
DE 6.50 usec
TE 298.1 K
D1 2.00000000 sec
TD0 1

===== CHANNEL f1 =====

NUC1 1H
P1 15.00 usec
PL1 0 dB
SFO1 400.1324008 MHz

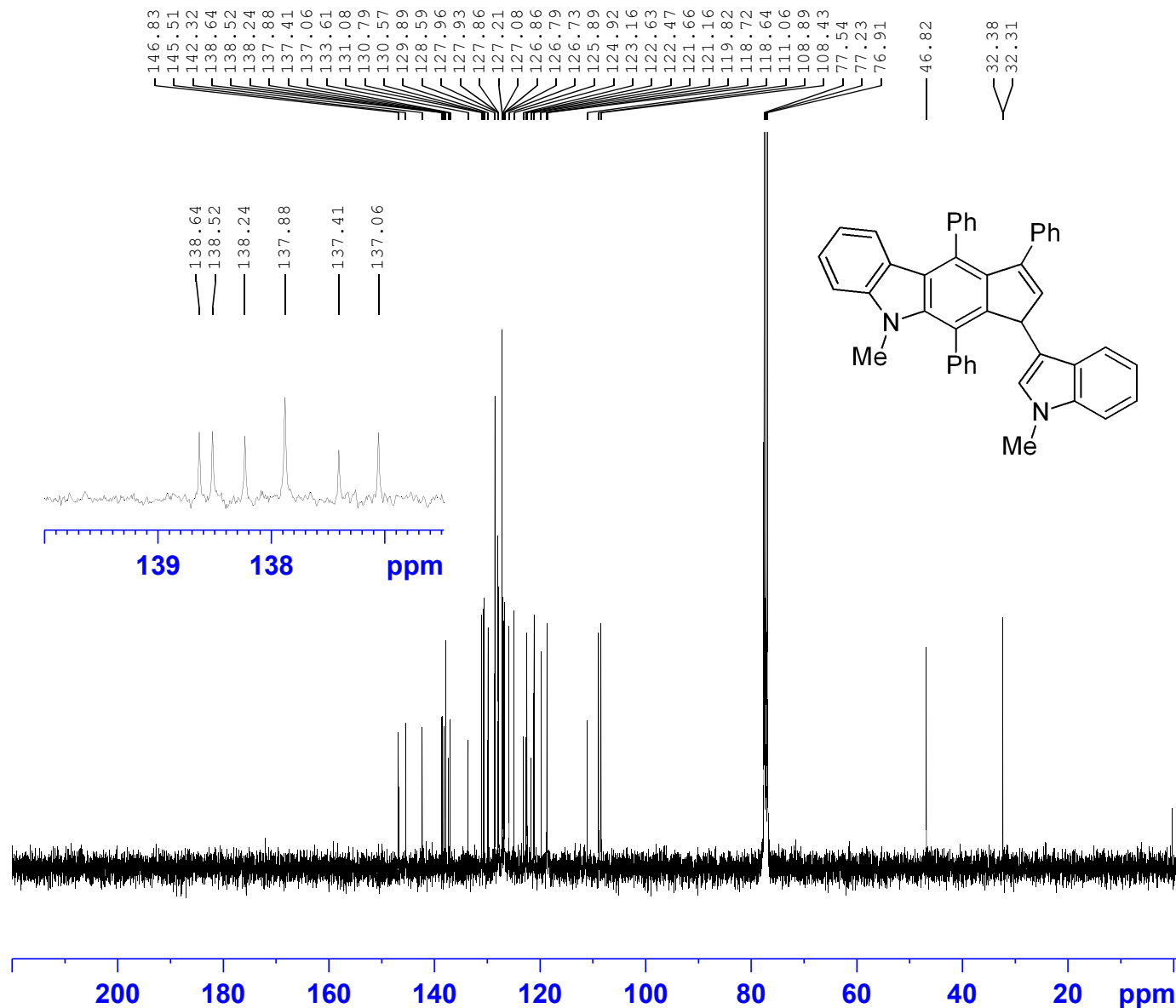
F2 - Processing parameters

SI 16384
SF 400.1300158 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

3,3'-(3,5-diphenyl-4-(phenylethynyl)cyclopent-3-ene-1,2-diyl)bis(1-methyl-1H-indole) (4)



155



Current Data Parameters
NAME test
EXPNO 55
PROCNO 1

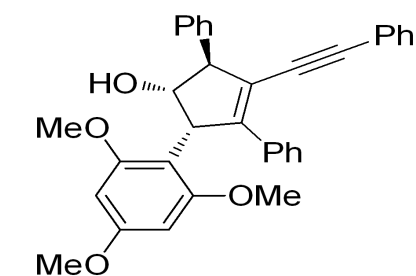
F2 - Acquisition Parameters
Date_ 20170928
Time_ 21.28
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 1149
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5792.6
DW 20.800 usec
DE 6.50 usec
TE 298.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
NUC1 13C
P1 10.45 usec
PL1 7.00 dB
SFO1 100.6233325 MHz

===== CHANNEL f2 =====
CPDPRG[2] waltz16
NUC2 1H
PCPD2 90.00 usec
PL2 0 dB
PL12 15.00 dB
PL13 20.00 dB
SFO2 400.1316005 MHz

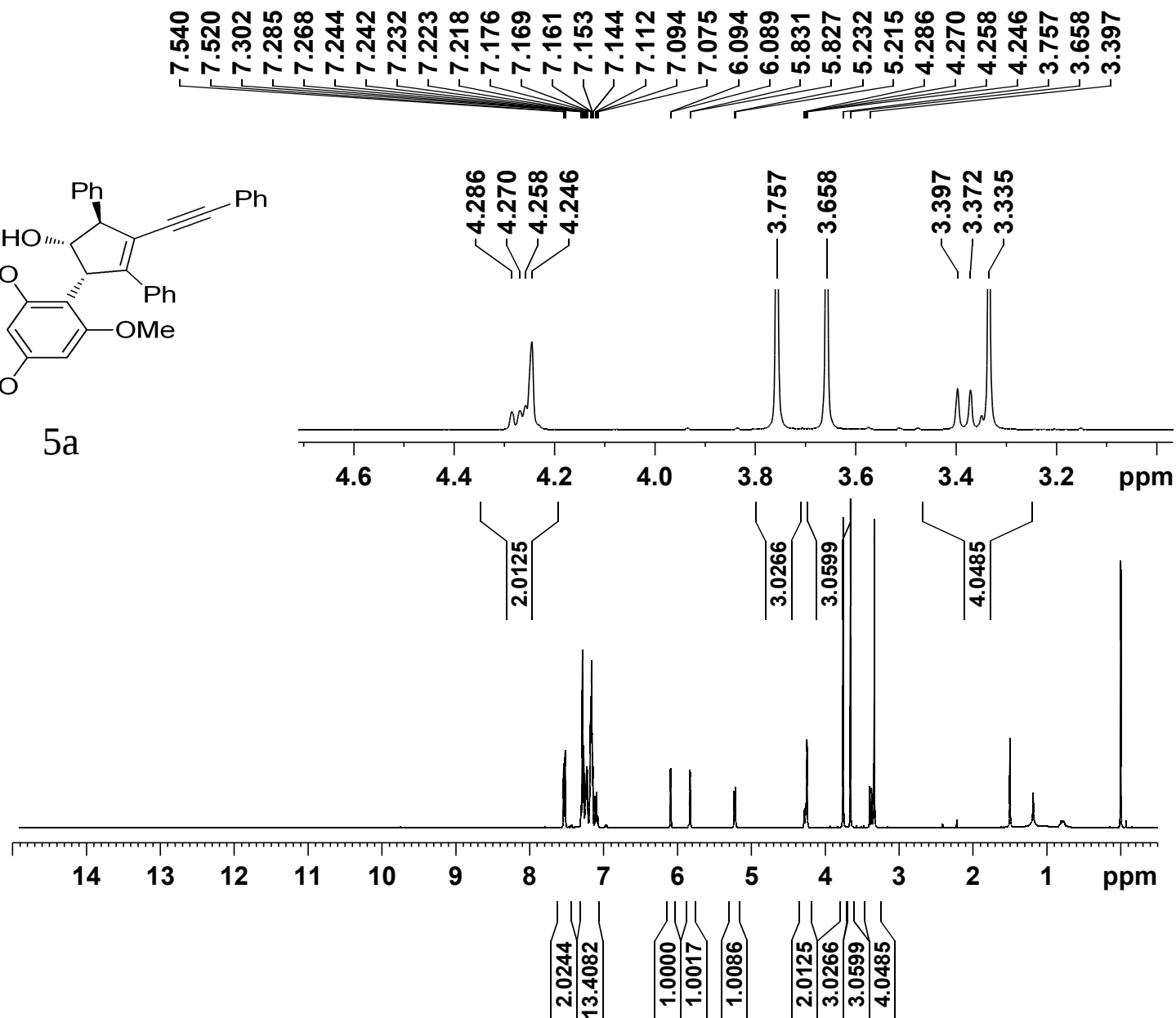
F2 - Processing parameters
SI 32768
SF 100.6127484 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

2,4-diphenyl-3-(phenylethynyl)-5-(2,4,6-trimethoxyphenyl)cyclopent-3-en-1-ol (5a)



5a

S62



Current Data Parameters
NAME try
EXPNO 117
PROCNO 1

F2 - Acquisition Parameters

Date_ 20180811
Time_ 19.17
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 56
DS 0
SWH 7211.539 Hz
FIDRES 0.220079 Hz
AQ 2.2719147 sec
RG 89.08
DW 69.333 usec
DE 10.50 usec
TE 297.5 K
D1 2.00000000 sec
TD0 1

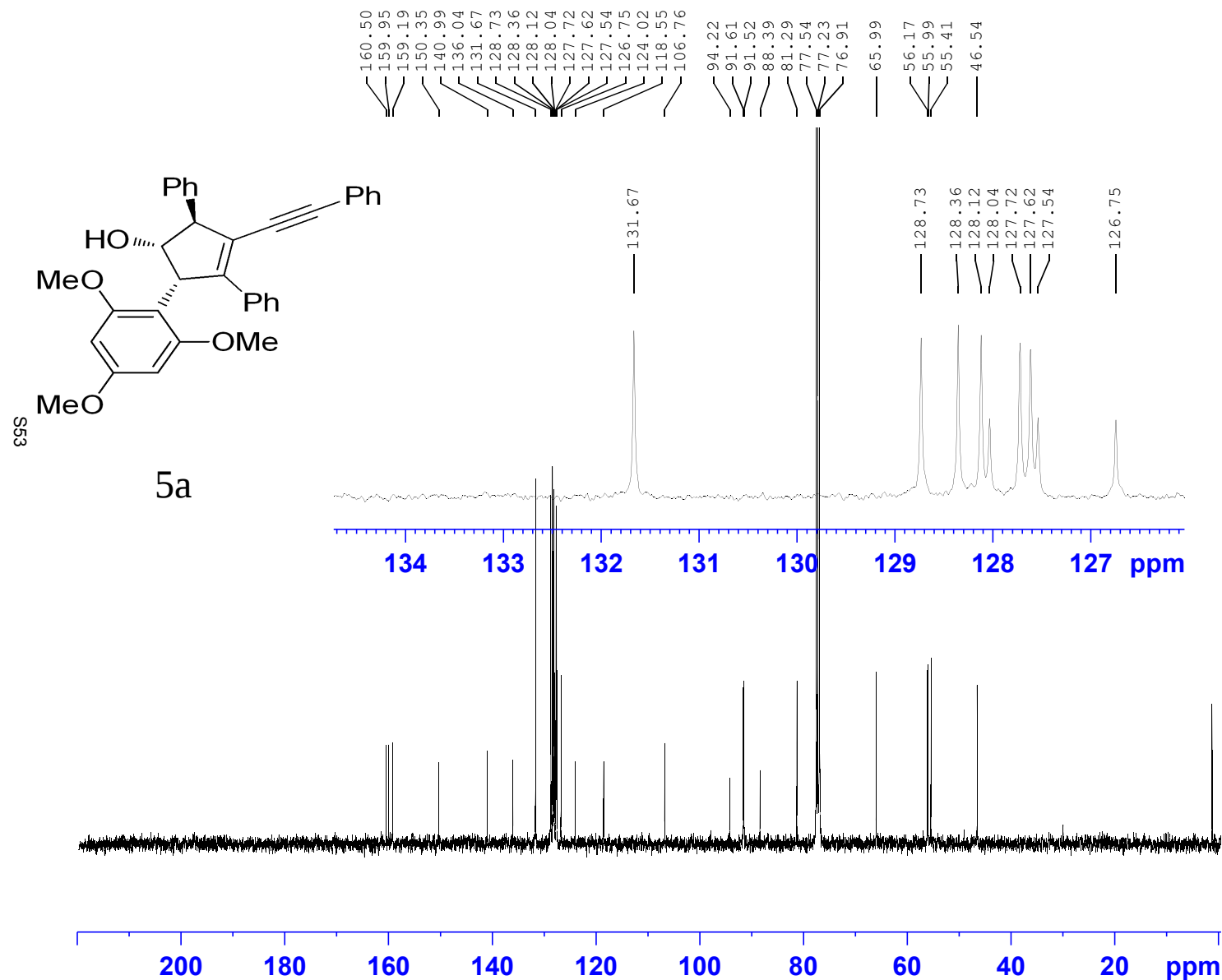
===== CHANNEL f1 =====

SFO1 400.1324008 MHz
NUC1 1H
P1 12.90 usec
PLW1 15.00000000 W

F2 - Processing parameters

SI 16384
SF 400.1300429 MHz
WDW EM
SSB 0
LB 0 Hz
GB 0
PC 1.00

2,4-diphenyl-3-(phenylethynyl)-5-(2,4,6-trimethoxyphenyl)cyclopent-3-en-1-ol (5a)



Current Data Parameters
NAME try
EXPNO 119
PROCNO 1

F2 - Acquisition Parameters
Date_ 20180811
Time_ 20.47
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 500
DS 0
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 198.09
DW 20.800 usec
DE 6.50 usec
TE 299.0 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

===== CHANNEL f1 =====
SFO1 100.6228298 MHz
NUC1 13C
P1 10.00 usec
PLW1 47.50000000 W

===== CHANNEL f2 =====
SFO2 400.1316005 MHz
NUC2 1H
CPDPRG[2] waltz16
PCPD2 90.00 usec
PLW2 15.00000000 W
PLW12 0.33750001 W
PLW13 0.27338001 W

F2 - Processing parameters
SI 32768
SF 100.6127490 MHz
WDW EM
SSB 0
LB 2.00 Hz
GB 0
PC 1.00