

**General comments:**

All reactions were carried out under air. Reactions were monitored by TLC analysis (pre-coated silica gel plates with fluorescent indicator UV254, 0.2 mm) and visualized with 254 nm UV light or iodine. Chemicals were purchased from Aldrich and were used without further purification otherwise noted. All compounds were characterized by <sup>1</sup>H NMR, <sup>13</sup>C NMR. <sup>1</sup>H spectra were recorded on Bruker AV 300 and AV 400 spectrometers. <sup>13</sup>C NMR spectra were recorded at 282 MHz. GC was performed on Agilent 6890 chromatograph with a 30 m HP5 column. All yields reported refer to isolated yields. All the products are commercially available.

**General procedure for the oxidative synthesis of primary amides from acetophenones:**

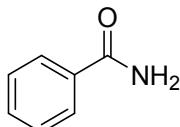
Ammonia (25 % in water; 1 mL), and acetophenones (1 mmol) were added to pressure equipped with a stirring bar. Then, TBHP (70 % in H<sub>2</sub>O; 1.0 mL; 8 equiv.), TBAI (20 mol %) were added and the final solution was kept at 100°C temperature for 16h. The mixture was cooled to room temperature and solvent was removed under vacuum. The pure product can be isolated by column.

**General procedure for the oxidative synthesis of primary amides from carbinols:**

Ammonia (25 % in water; 1 mL), and carbinols (1 mmol) were added to pressure equipped with a stirring bar. Then, TBHP (70 % in H<sub>2</sub>O; 1.0 mL; 8 equiv.), TBAI (20 mol %) were added and the final solution was kept at 100°C temperature for 16h. The mixture was cooled to room temperature and solvent was removed under vacuum. The pure product can be isolated by column.

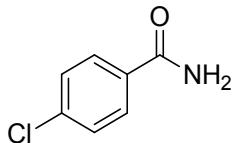
**Analytic data.**

**Benzamide**



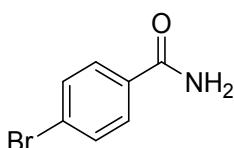
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.39-7.56 (m, 4H), 7.89-7.94 (m, 2H), 8.03 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 128.4 (2CH<sub>2</sub>), 129.1 (2CH<sub>2</sub>), 132.1 (CH), 135.2 (C), 168.8 (CO). GC-MS (EI, 70 eV): m/z (%) [M<sup>+</sup>]121 (81), 105 (100), 77 (95), 51 (40), 50 (23), 44 (10).

**4-Chlorobenzamide**



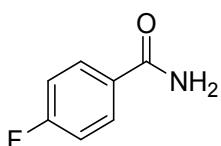
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.48-7.57 (m, 3H), 7.91-7.95 (m, 2H), 8.09 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 129.2 (2CH), 130.3 (2CH), 133.9 (C), 136.9 (C), 167.7 (CO); **GC-MS** (EI, 70 eV): m/z (%) [M<sup>+</sup>]155 (55), 141 (33), 139 (100), 111 (52), 75 (35), 50 (17); **HRMS** (EI): Calc for C<sub>7</sub>H<sub>6</sub>N<sub>1</sub>O<sub>1</sub>Cl: 155.01324; found: 155.15501291.

**4-Bromobenzamide**



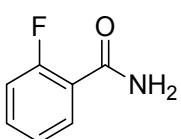
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.51'0 (s, 1H, NH<sub>2</sub>), 7.67-7.72 (m, 2H), 7.83-7.88 (m, 2H), 8.09 (s, 1H, NH<sub>2</sub>); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 125.9 (C), 130.5 (2CH), 132.1 (2CH), 134.3 (C), 167.8 (CO); **GC-MS** (EI, 70 eV): m/z (%) [M<sup>+</sup>] 199 (57), 185 (99), 183 (100), 157 (44), 155 (44), 76 (34); **HRMS** (EI): Calc for C<sub>7</sub>H<sub>6</sub>N<sub>1</sub>O<sub>1</sub>Br: 198.96273; found: 198.96294.

**4-Fluorobenzamide**



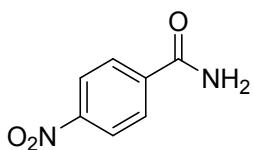
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 6.09 (s, 1H), 6.90-7.07 (m, 3H), 7.70-7.77 (m, 2H), 8.09 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 116.2 (d, J = 25.6 Hz, 2CH), 129.9 (d, J = 9.60 Hz, 2CH), 131.6 (C), 135.9 (C), 164.6 (d, J = 248.6 Hz, CF), 167.7 (CO); **GC-MS** (EI, 70 eV): m/z (%) [M<sup>+</sup>]139 (55), 123 (100), 95 (76), 75 (34), 74 (10), 50 (10); **HRMS** (EI): Calc for C<sub>7</sub>H<sub>6</sub>N<sub>1</sub>O<sub>1</sub>F: 139.04279; found: 139.04315.

**2-Fluorobenzamide**



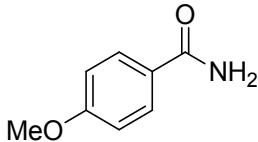
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.27-7.34 (m, 2H), 7.52 -7.57(m, 1H), 7.59-7.79 (m, 3H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 116.9 (d, *J* = 22.5 Hz, CH), 124.7 (d, *J* = 14.5 Hz, CH), 125.3 (CH), 131.1 (CH), 133.3 (CH), 160.2 (d, *J* = 248.8 Hz, CF), 166.2 (CO). GC-MS (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 139 (63), 123 (100), 95 (46), 75 (30), 74 (12), 50 (10)

#### 4-Nitrobenzamide



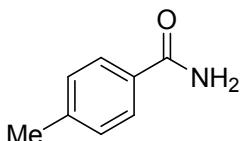
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 8.12-8.16 (m, 3H), 8.24-8.28 (m, 3H) **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 124.0, 131.1 (CH), 143.2, 149.6 (C), 168.0 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 166 (71), 150 (100), 120 (24), 104 (28), 92 (24), 50 (19); **HRMS** (EI): Calc for C<sub>7</sub>H<sub>6</sub>N<sub>2</sub>O<sub>3</sub>: 166.03729; found: 166.03782.

#### 4-Methoxybenzamide



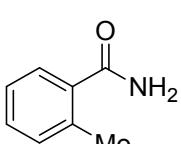
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 3.84 (s, 3H), 6.99-7.03 (m, 2H), 7.23 (s, 1H), 7.85-7.92 (m, 3H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 56.4 (OCH<sub>3</sub>), 114.5 (2CH), 127.6 (C), 130.5 (2CH), 162.7 (C), 168.6 (CO).

#### 4-Methylbenzamide



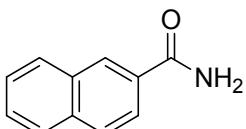
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 2.37 (s, 3H), 7.27 (d, *J* = 8.15 Hz, 2H), 7.34 (s, 1H, NH<sub>2</sub>), 7.81 (d, *J* = 8.15 Hz, 2H), 7.94 (s, 1H, NH<sub>2</sub>); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 21.8 (CH<sub>3</sub>), 128.4 (2CH), 129.6 (2CH), 132.3 (C), 141.9 (C), 168.7 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 135 (62), 119 (100), 91 (66), 89 (11), 65 (23), 63 (11); **HRMS** (EI): Calc for C<sub>8</sub>H<sub>9</sub>N<sub>1</sub>O<sub>1</sub>: 135.06787; found: 135.06777.

#### 2-Methylbenzamide



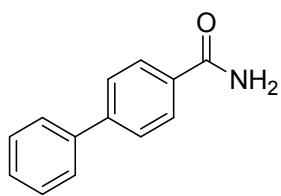
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 2.40 (s, 3H), 7.21-7.28 (m, 2H), 7.31-7.42 (m, 3H), 7.73 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 20.5 (CH<sub>3</sub>), 126.3, 127.9, 130.0, 131.3 (CH), 135.9, 137.9 (C), 171.9 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 135 (85), 119 (92), 91 (100), 90 (27), 89 (23).

#### 2-Naphthamide



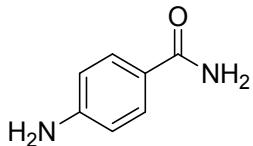
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.55 (s, 1H), 7.59-7.66 (m, 2H), 7.99-8.06 (m, 4H), 8.21 (s, 1H), 8.55 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 125.3, 127.5, 128.5, 128.7, 129.8, 132.5, 133.1, 135.1, 168.9 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 171 (80), 156 (12), 155 (99), 127 (100), 75 (10).

#### [1,1'-Niphenyl]-4-carboxamide



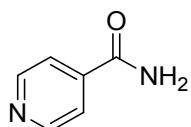
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.24-7.55 (m, 4H), 7.74-7.81 (m, 4H), 8.00-8.0 (m, 2H), 8.09 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 127.3 (2CH), 127.7 (2CH), 128.9 (CH), 129.1 (2CH), 129.9 (2CH), 134.0 (C), 140.1 (C), 143.6 (C), 168.5 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 197 (71), 181 (100), 153 (34), 152 (66), 151 (20), 76 (14).

#### 4-Aminobenzamide



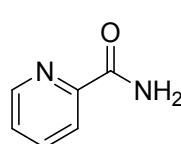
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 5.63 (s, 2H, NH<sub>2</sub>), 6.52-6.58 (m, 2H), 6.87 (s, 1H), 7.54-7.64 (m, 3H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 113.3 (2CH), 121.8 (C), 130.0 (2CH), 152.5 (C), 168.9 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 136 (69), 120 (100), 92 (38), 65 (33); **HRMS** (EI) [M]<sup>+</sup>: Calc for C<sub>7</sub>H<sub>8</sub>N<sub>2</sub>O: 136.06311; found: 136.06345.

#### Isonicotinamide



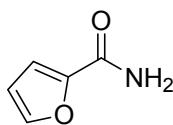
**<sup>1</sup>H NMR** (δ = 7.74-7.82 (m, 3H), 8.28 (s, 1H), 8.73-8.77 (m, 2H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 122.4 (2CH), 142.2 (C), 151.2 (2CH), 167.3 (CO).

#### Picolinamide



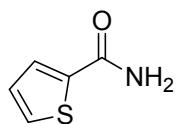
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.11-7.14 (m, 1H), 7.34 (s, 1H, NH<sub>2</sub>), 7.72-7.75 (m, 3H), 7.96 (s, 1H, NH<sub>2</sub>); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 125.4 (CH), 128.8 (CH), 129.5 (CH), 131.8 (CH), 141.2 (C), 163.7 (CO).

#### Furan-2-carboxamide



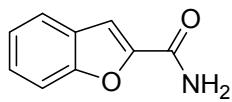
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 6.62–6.64 (m, 1H), 7.12 (d, *J* = 4.80 Hz, 1H), 7.40 (1, 1H), 7.79 (s, 1H), 7.83-7.84 (m, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 112.6 (CH), 114.5 (CH), 145.9 (CH), 148.9 (C), 160.3 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 111 (100), 95 (96), 44 (18), 39 (40).

#### Thiophene-2-carboxamide.



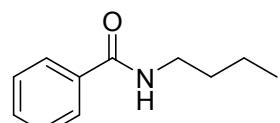
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.12–7.19 (m, 1H), 7.41 (s, 1H), 7.75–7.79 (m, 2H), 7.99 (s, 1H); **<sup>13</sup>C NMR** (DMSO-d<sub>6</sub>): δ = 128.8 (CH), 129.6 (CH), 131.9 (CH), 141.2 (C), 163.8 (CO). **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 127 (67), 111 (100), 39 (20).

### Benzofuran-2-carboxamide



**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.33–7.39 (m, 1H), 7.46–7.49 (m, 1H), 7.52 (s, 1H), 7.57–7.59 (s, 1H), 7.65 (s, 1H), 7.69–7.82 (m, 1H), 8.16 (s, 1H); **<sup>13</sup>C NMR** (DMSO-d<sub>6</sub>): δ = 110.4 (CH), 112.7 (CH), 123.6 (CH), 124.5 (CH), 127.6 (CH), 128.1 (C), 150.2 (C), 155.1 (C), 160.7 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 161 (98), 145 (100), 90 (11), 89 (58), 63 (23).

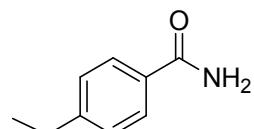
### N-Butylbenzamide



177.11482; found: 177.11443.

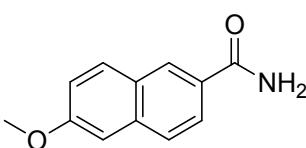
**<sup>1</sup>H NMR** (300 MHz, CDCl<sub>3</sub>): δ = 0.93 (t, *J* = 7.28 Hz, 3H), 1.35 (sext., 2H), 1.53 (pent., 2H), 3.30 (q, *J* = 7.13 Hz, 2H), 7.45–7.54 (m, 3H), 7.84–7.87 (m, 2H), 8.45 (s, 1H, NH<sub>2</sub>); **<sup>13</sup>C NMR** (CDCl<sub>3</sub>): δ = 14.7 (CH<sub>3</sub>), 20.6 (CH<sub>2</sub>), 32.2 (CH<sub>2</sub>), 39.7 (CH<sub>2</sub>), 128.1 (2CH), 129.2 (2CH), 131.9 (CH), 135.7 (C), 167.0 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 177 (9), 135 (19), 134 (19), 105 (100), 77 (40), 51 (11); **HRMS** (EI) [M]<sup>+</sup> : Calc for C<sub>11</sub>H<sub>15</sub>N<sub>1</sub>O<sub>1</sub>: 177.11482; found: 177.11443.

### 4-Ethylbenzamide



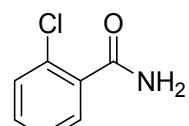
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 1.22 (t, *J* = 7.50 Hz, 3H), 2.67 (q, *J* = 15.3 Hz, 2H), 7.28–7.34 (m, 3H), 7.81–7.86 (m, 2H), 7.94 (s, 1H, NH<sub>2</sub>); **<sup>13</sup>C NMR** (DMSO-d<sub>6</sub>): δ = 16.5 (CH<sub>3</sub>), 28.9 (CH<sub>2</sub>), 128.4 (2CH), 128.5 (2CH), 132.6 (C), 148.1 (C), 168.7 (CO); **GC-MS** (EI, 70 eV): *m/z* (%) [M<sup>+</sup>] 149 (50), 134 (11), 133 (100), 106 (14), 79 (21), 77 (26)

### 6-Methoxy-2-naphthamide



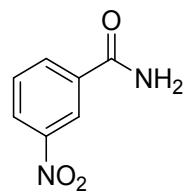
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 3.93 (s, OCH<sub>3</sub>), 7.23–7.29 (m, 1H), 7.38–7.44 (m, 2H), 7.76–7.99 (m, 3H), 8.09 (s, 1H), 8.45 (s, 1H); **<sup>13</sup>C NMR** (DMSO-d<sub>6</sub>): δ = 56.2 (OCH<sub>3</sub>), 106.7 (CH), 120.2 (CH), 125.8 (CH), 127.4 (CH), 128.4 (CH), 130.2 (CH), 131.3 (C), 136.7 (C), 159.4 (C), 168.9 (CO).

### 2-Chlorobenzamide



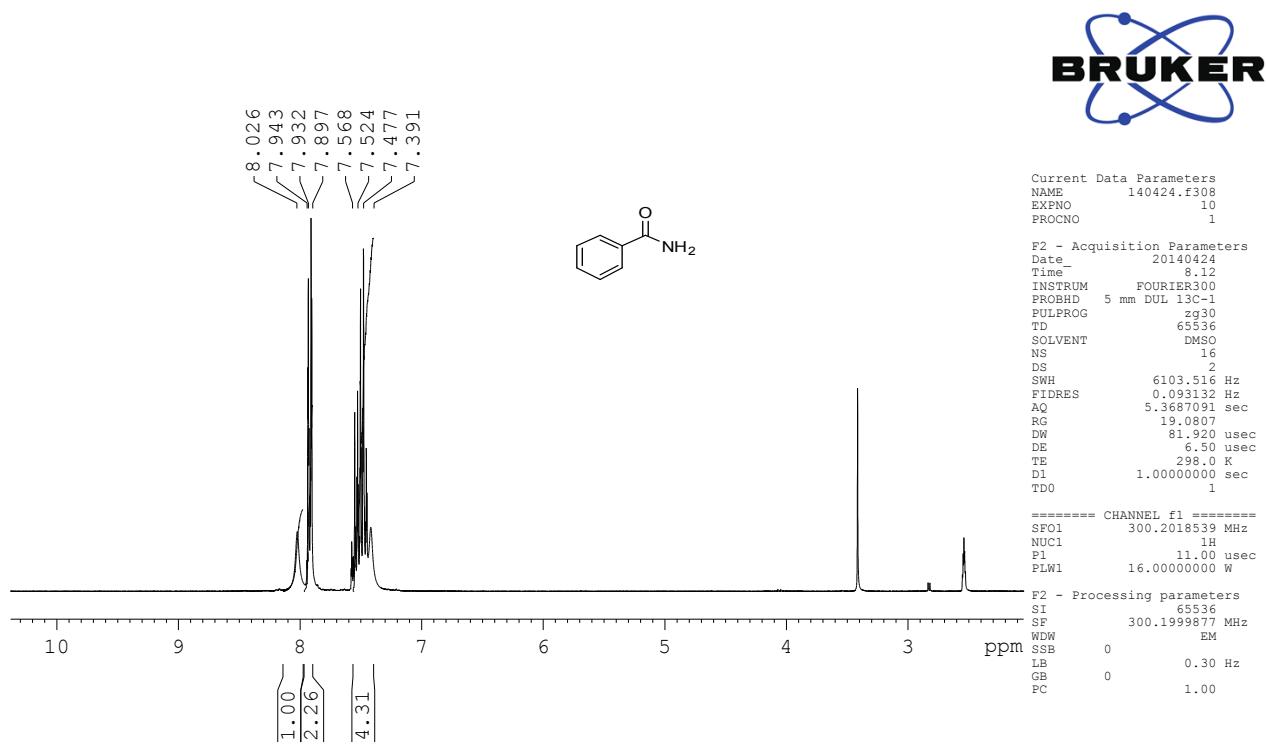
**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.39–7.54 (m, 4H), 7.62 (s, 1H), 7.90 (s, 1H); **<sup>13</sup>C NMR** (DMSO-d<sub>6</sub>): δ = 127.9 (CH), 129.5 (CH), 130.5 (CH), 131.4 (CH), 138.0 (C), 169.0 (CO).

### 3-Nitrobenzamide

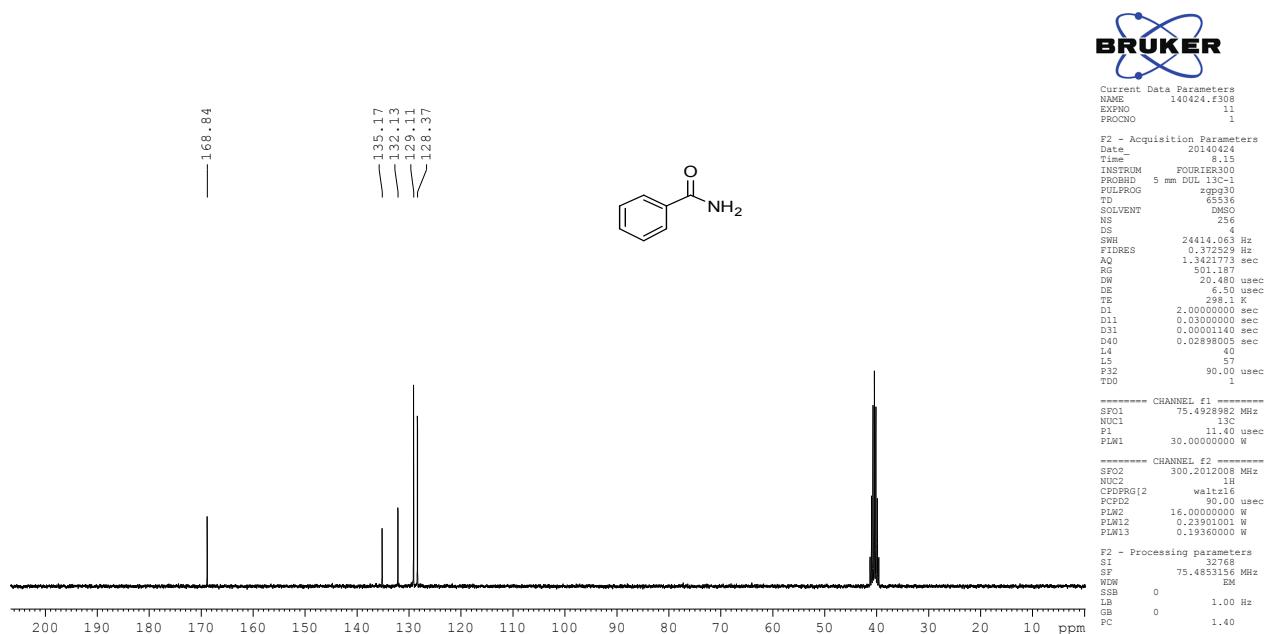


**<sup>1</sup>H NMR** (300 MHz, DMSO-d<sub>6</sub>): δ = 7.73-7.82 (m, 2H), 8.33-8.41 (m, 3H), 8.73 (s, 1H); **<sup>13</sup>CNMR** (DMSO-d<sub>6</sub>): δ = 123.2 (CH), 126.7 (CH), 130.9 (CH), 134.7 (CH), 136.7 (CH), 148.7 (C), 166.6 (CO).

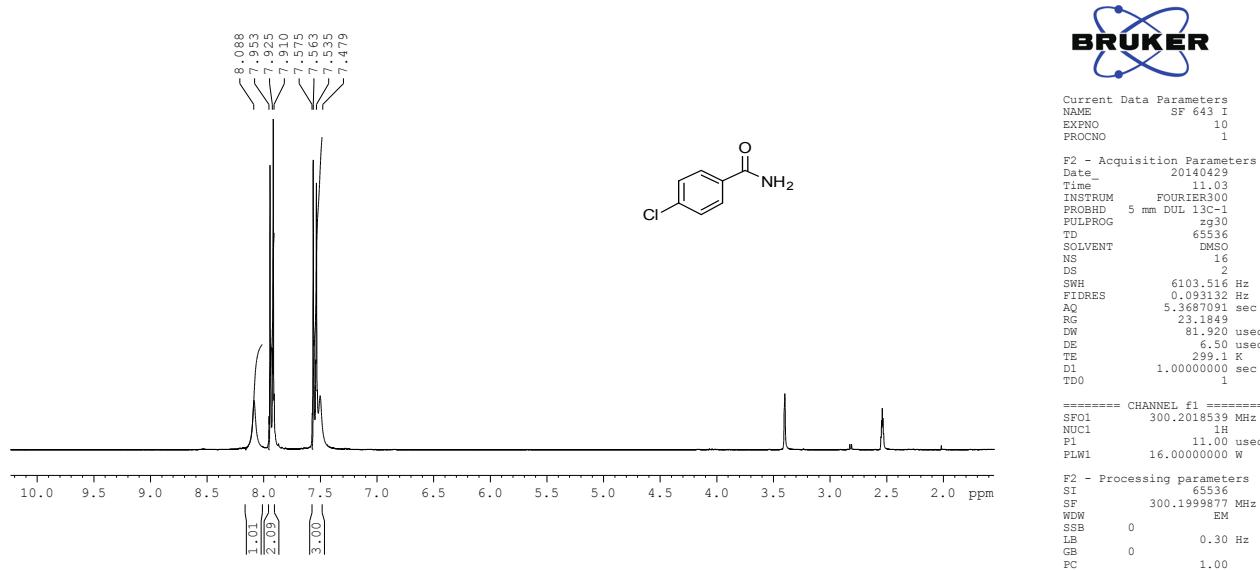
<sup>1</sup>H NMR



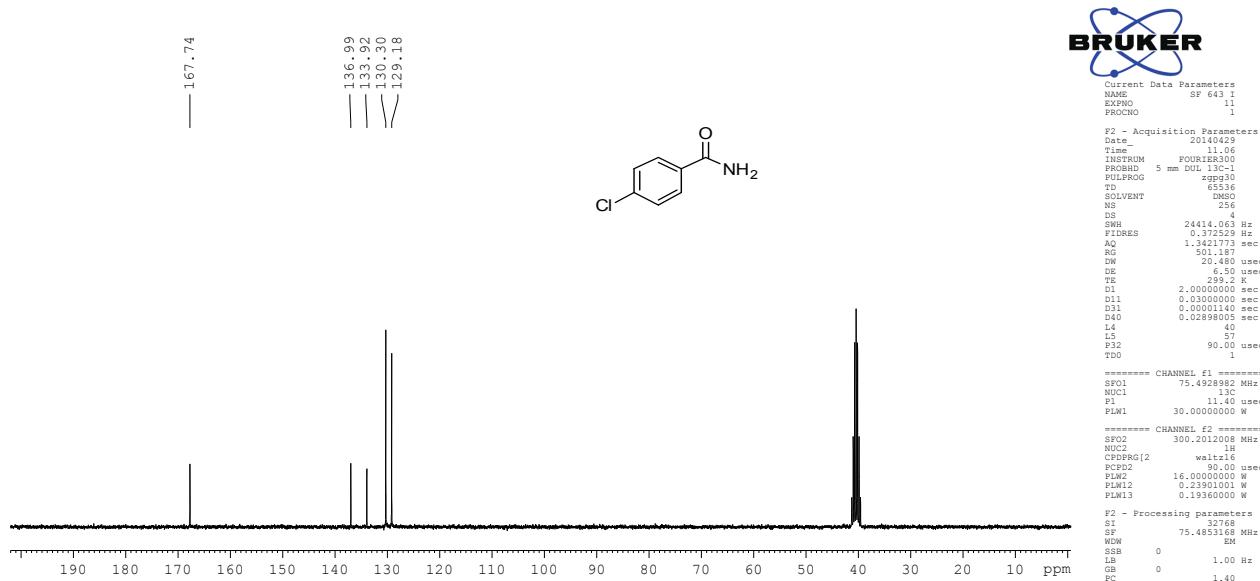
<sup>13</sup>C NMR



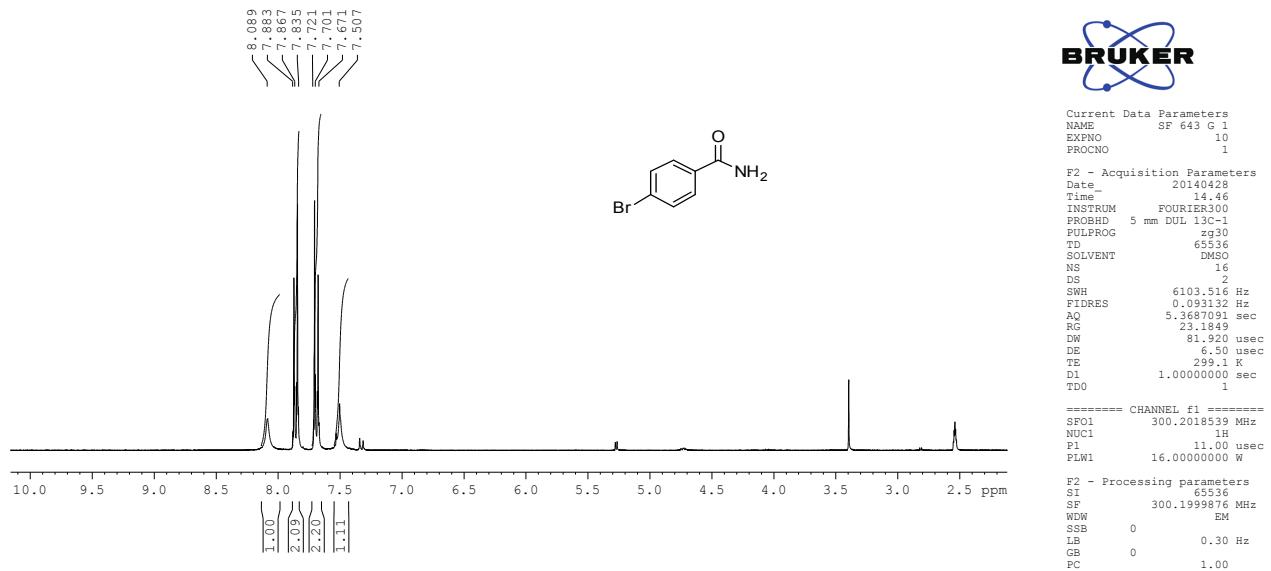
<sup>1</sup>H NMR



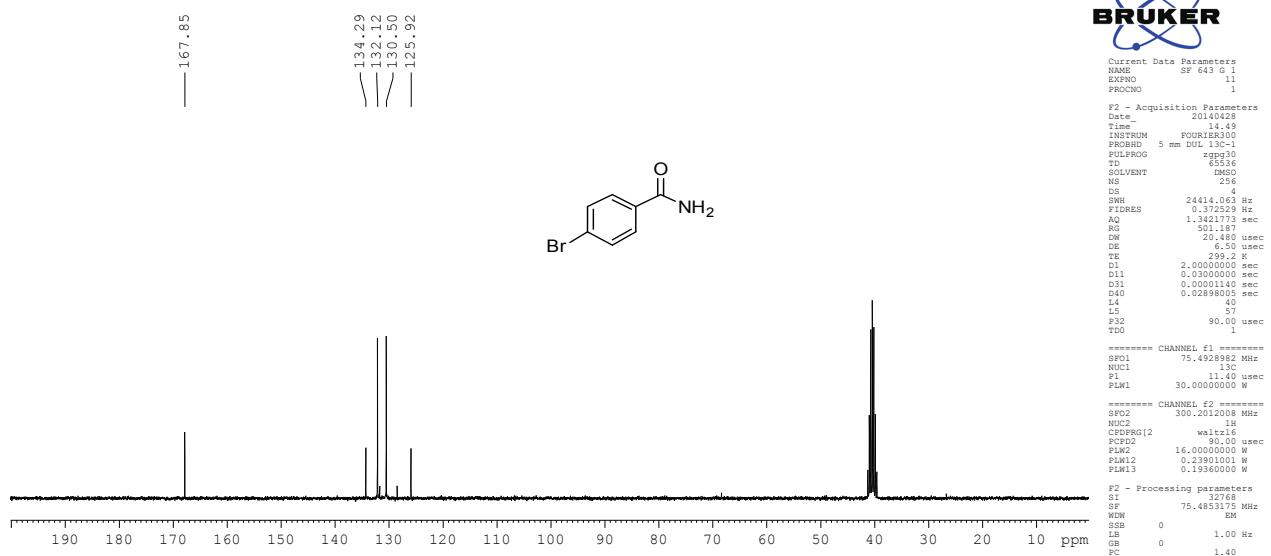
<sup>13</sup>C NMR



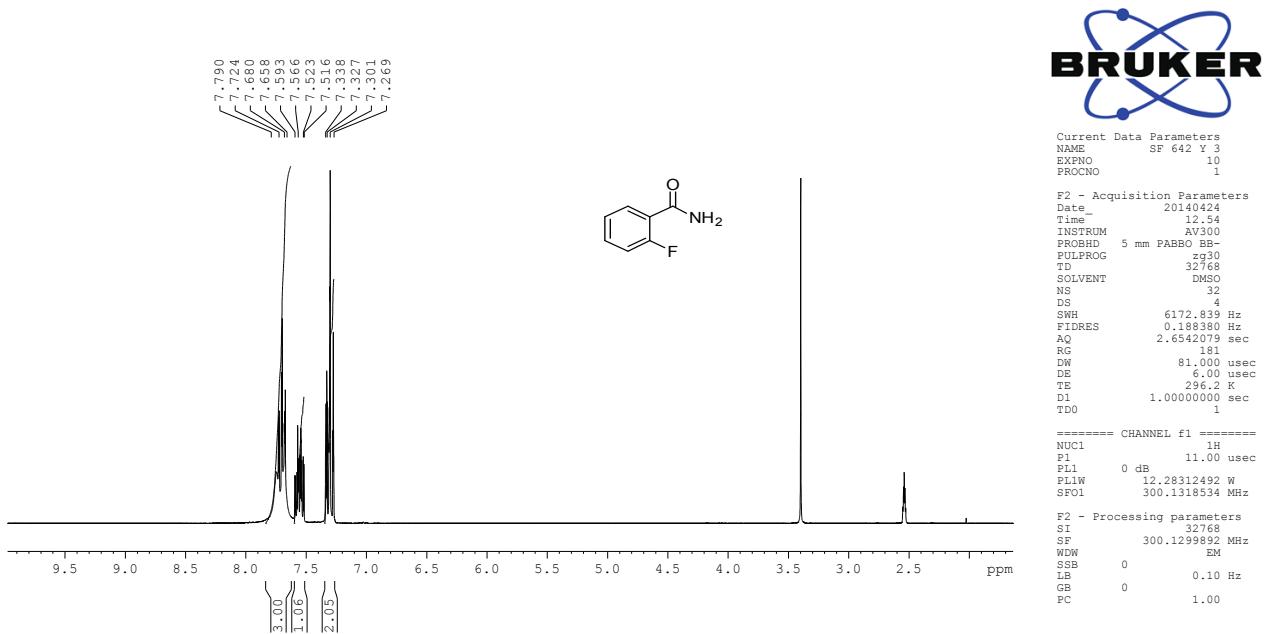
**<sup>1</sup>HNMR**



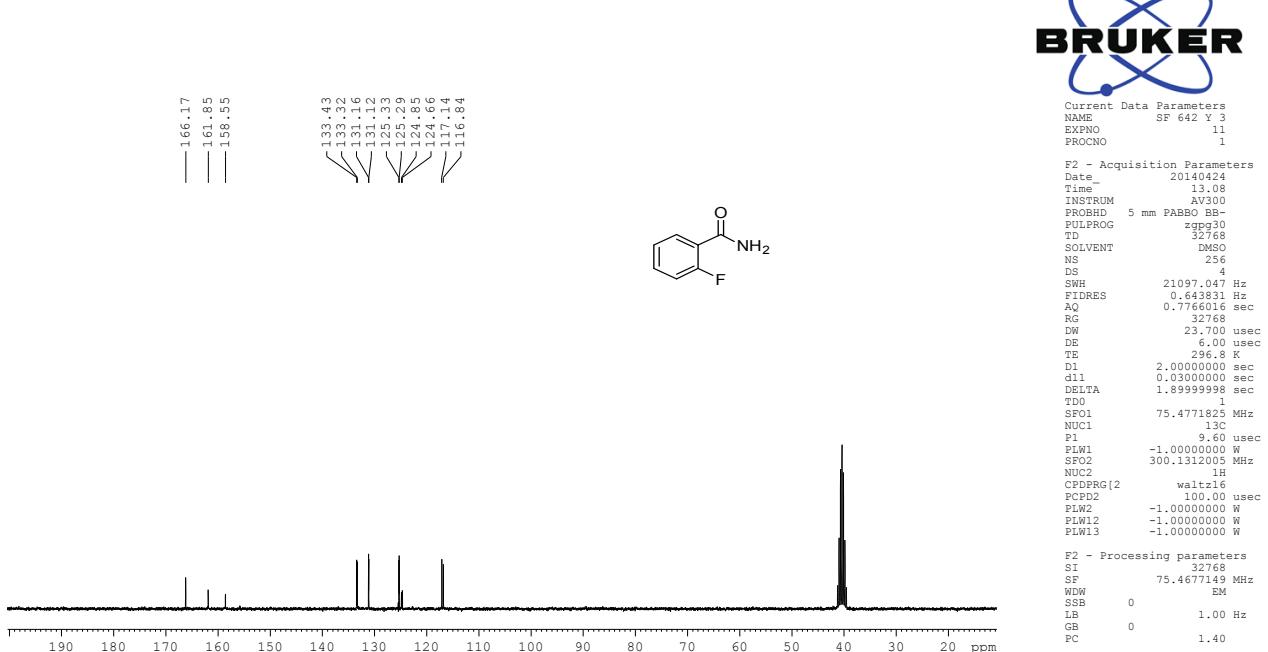
**<sup>13</sup>CNMR**



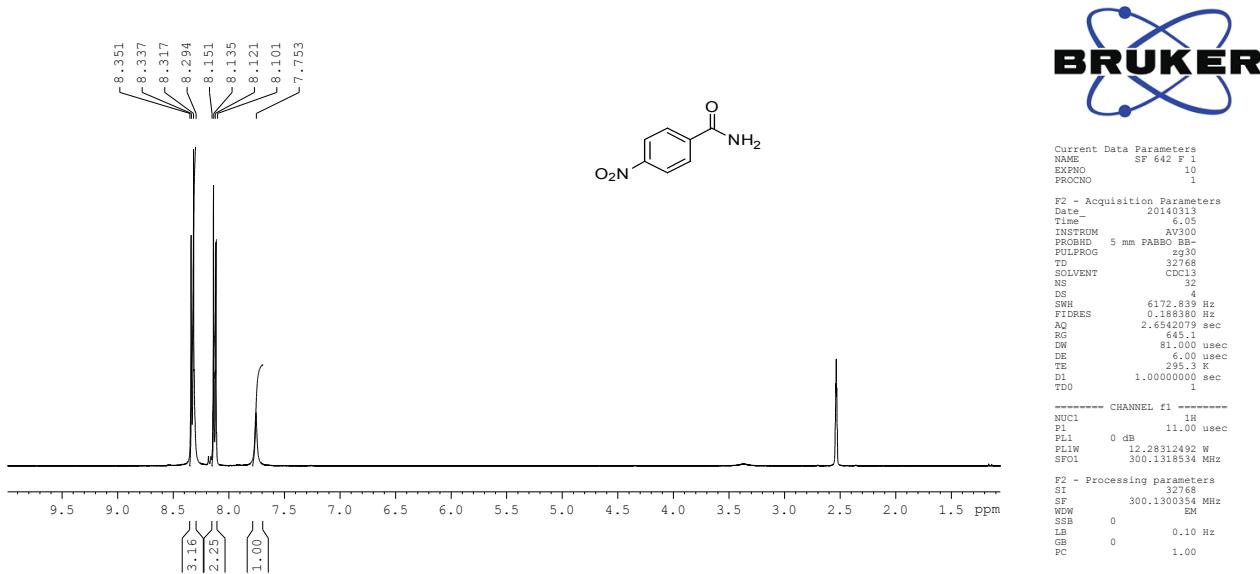
**<sup>1</sup>H NMR**



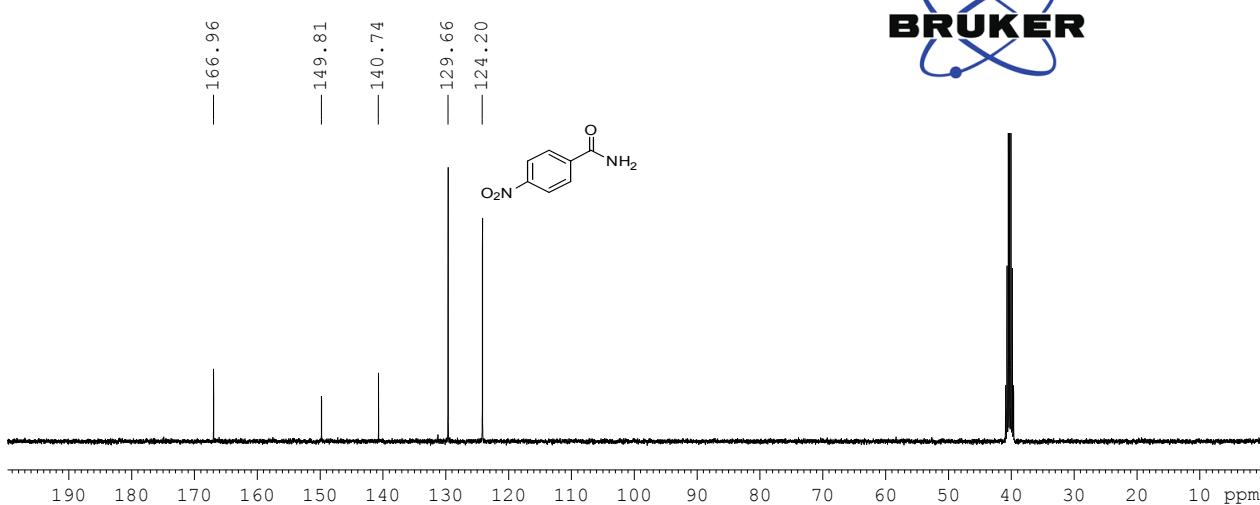
**<sup>13</sup>C NMR**



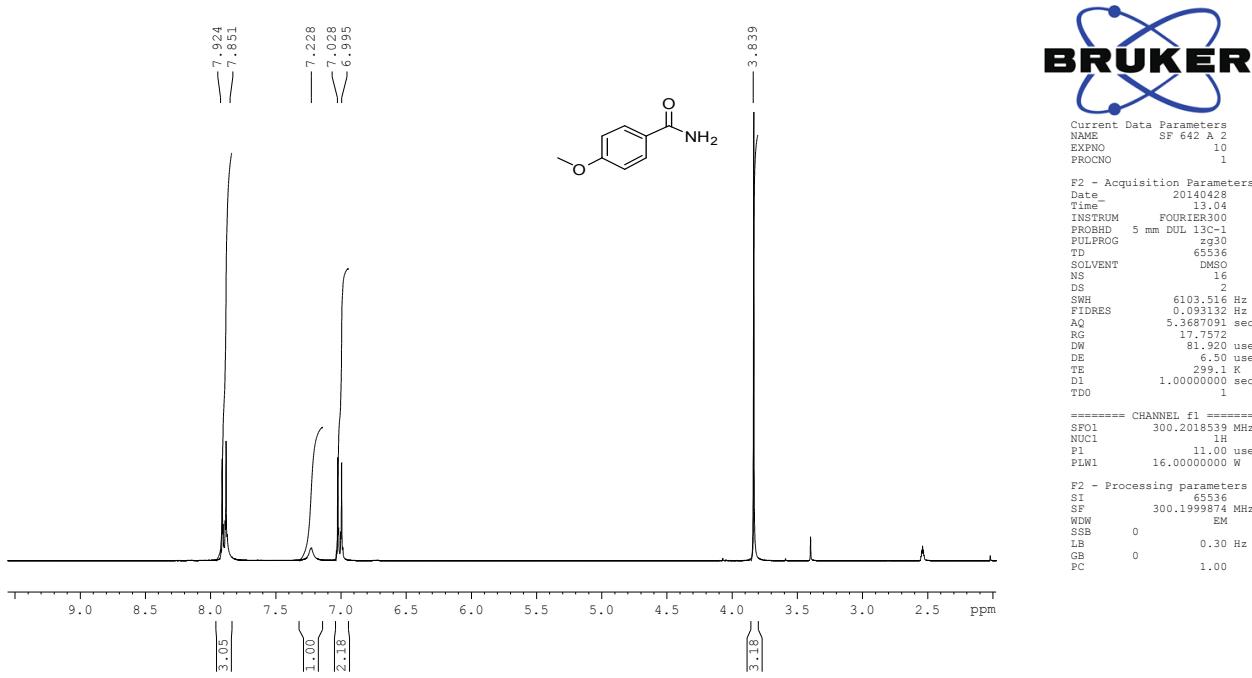
**<sup>1</sup>H NMR**



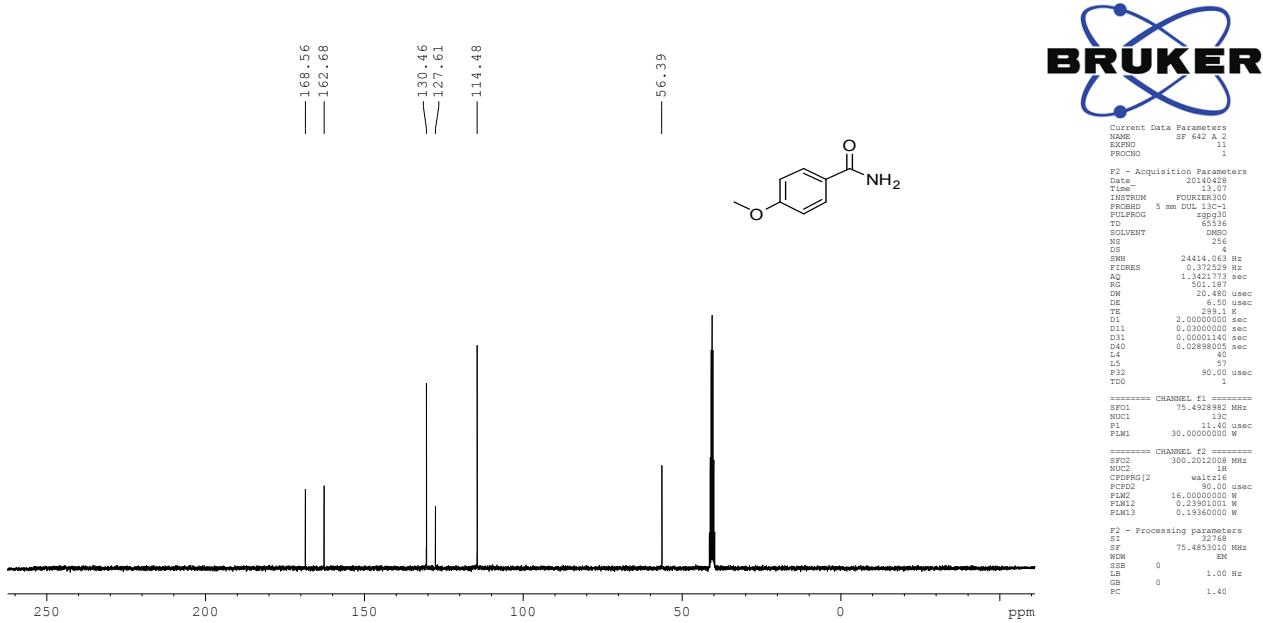
**<sup>13</sup>C NMR**



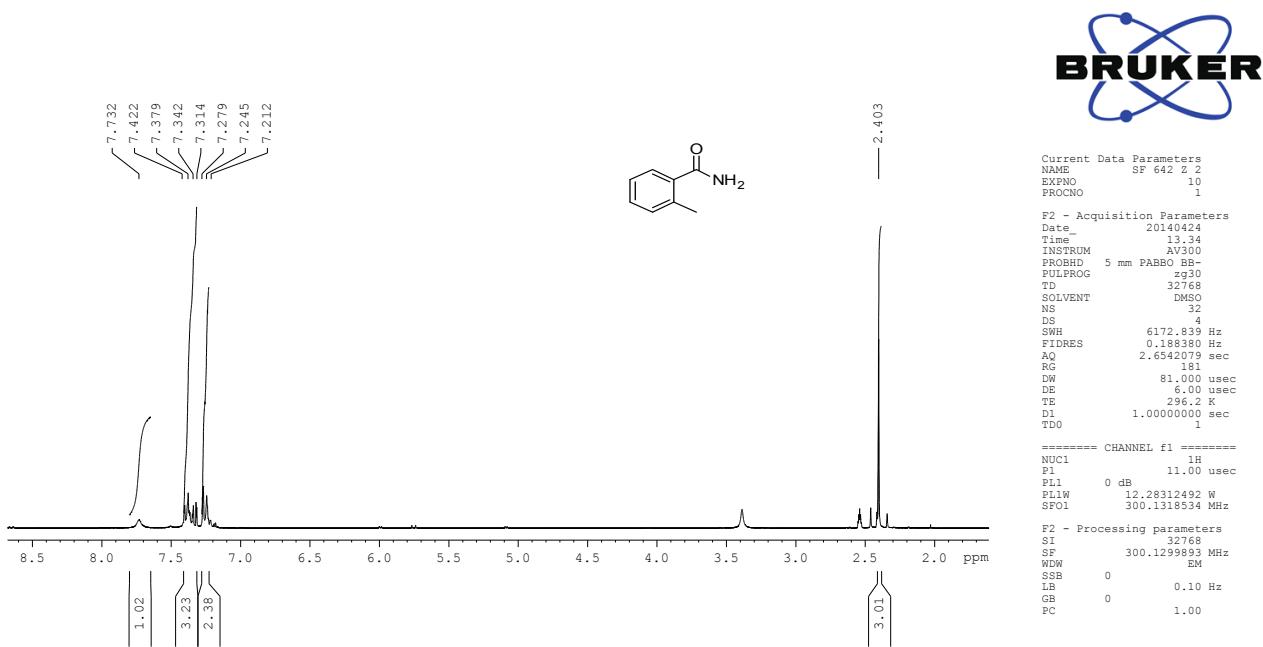
<sup>1</sup>H NMR



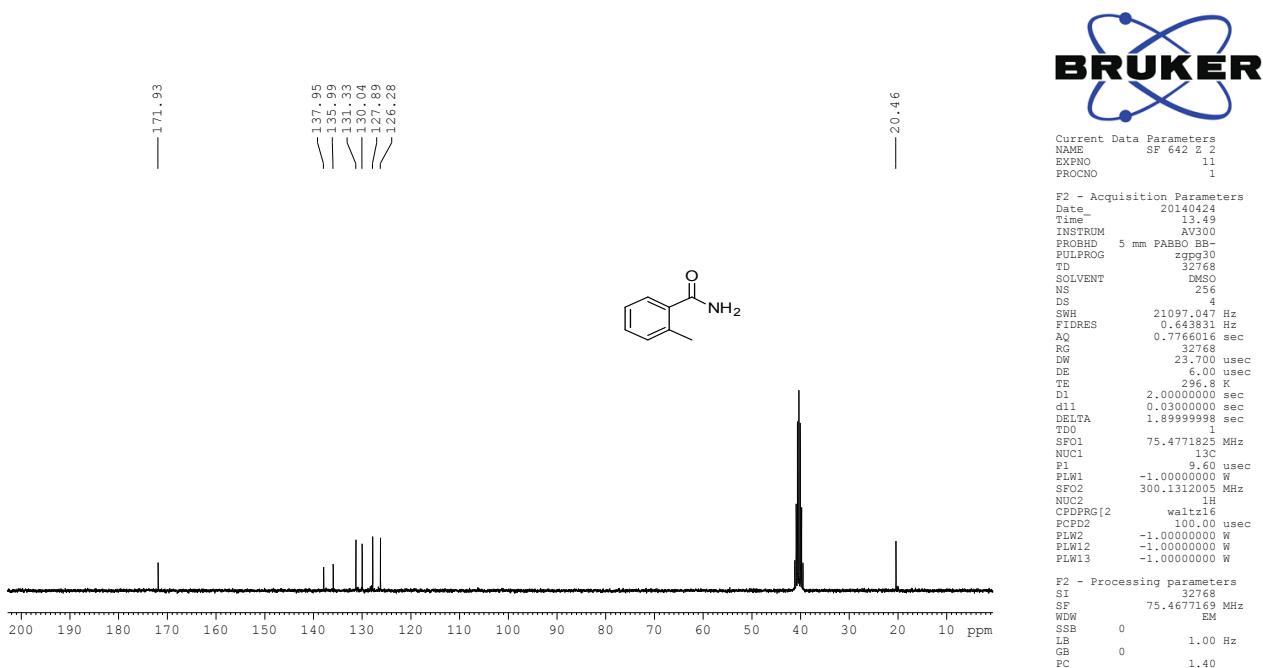
<sup>13</sup>C NMR



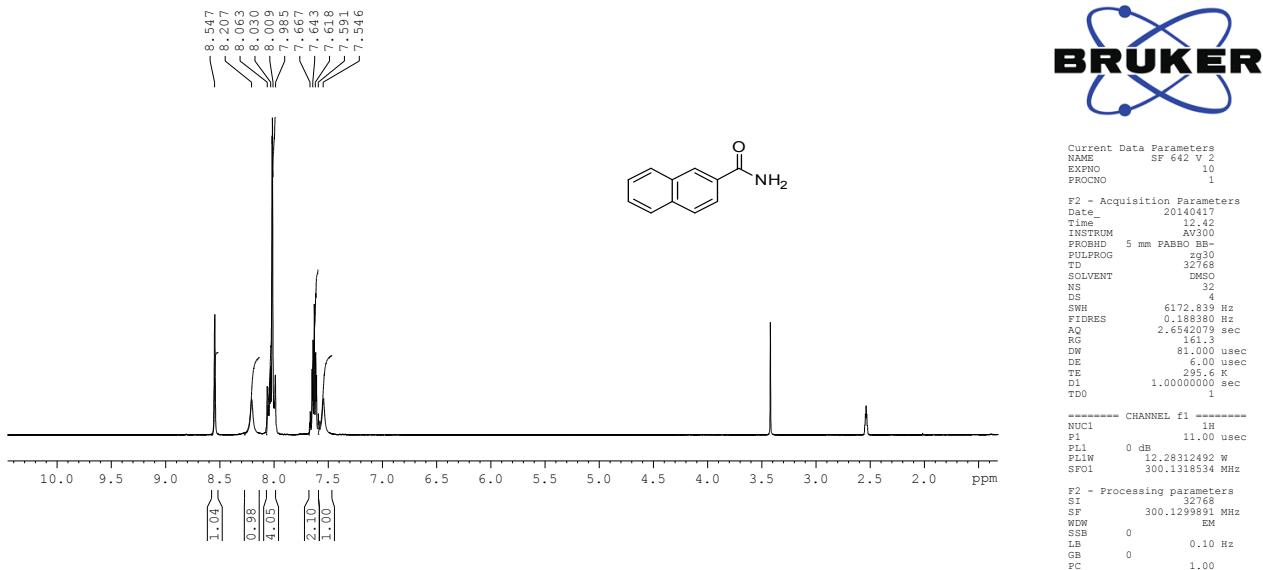
**<sup>1</sup>H NMR**



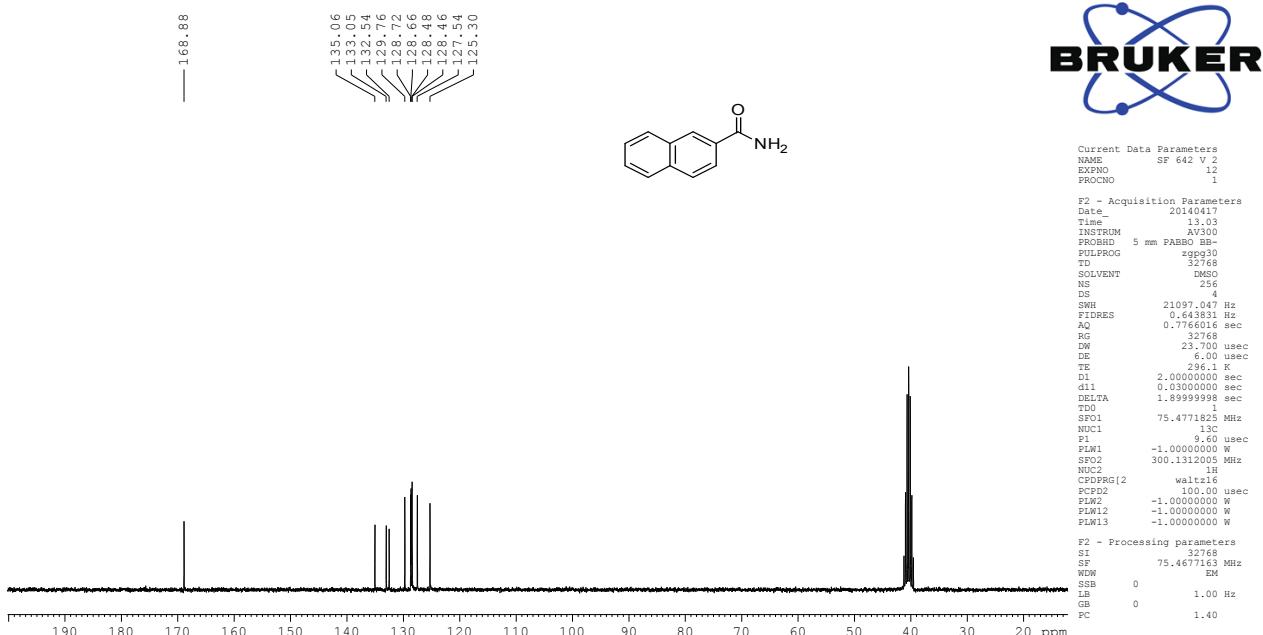
**<sup>13</sup>C NMR**



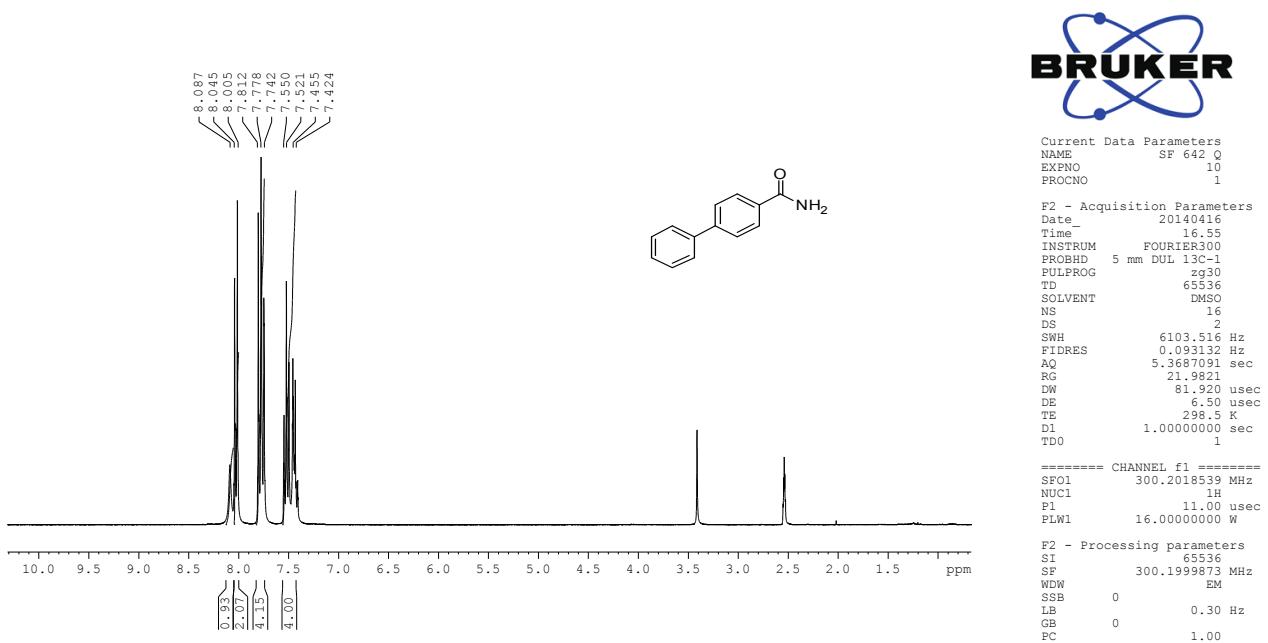
<sup>1</sup>H NMR



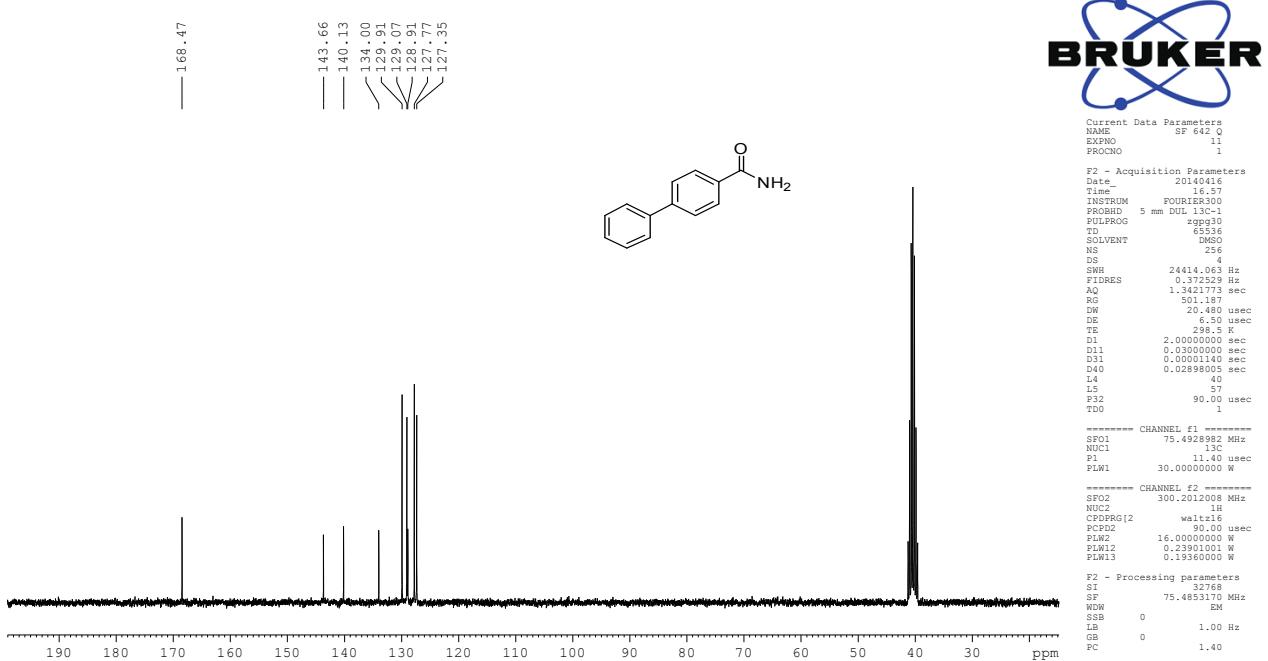
<sup>13</sup>C NMR



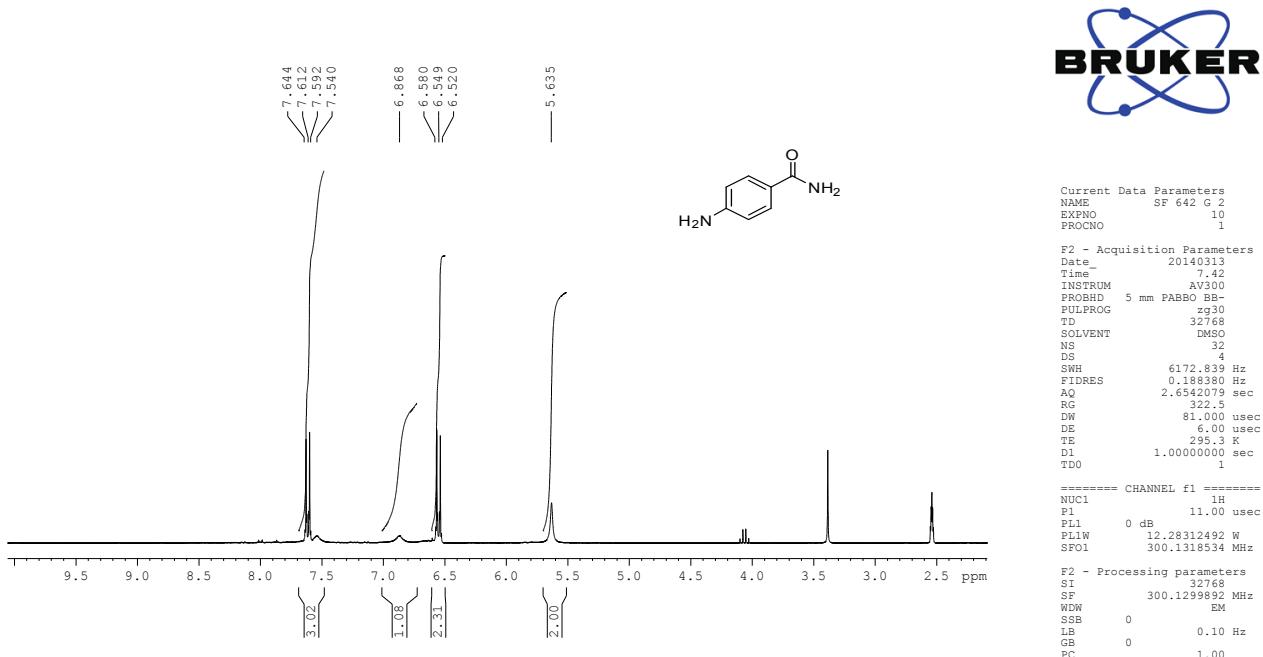
<sup>1</sup>H NMR



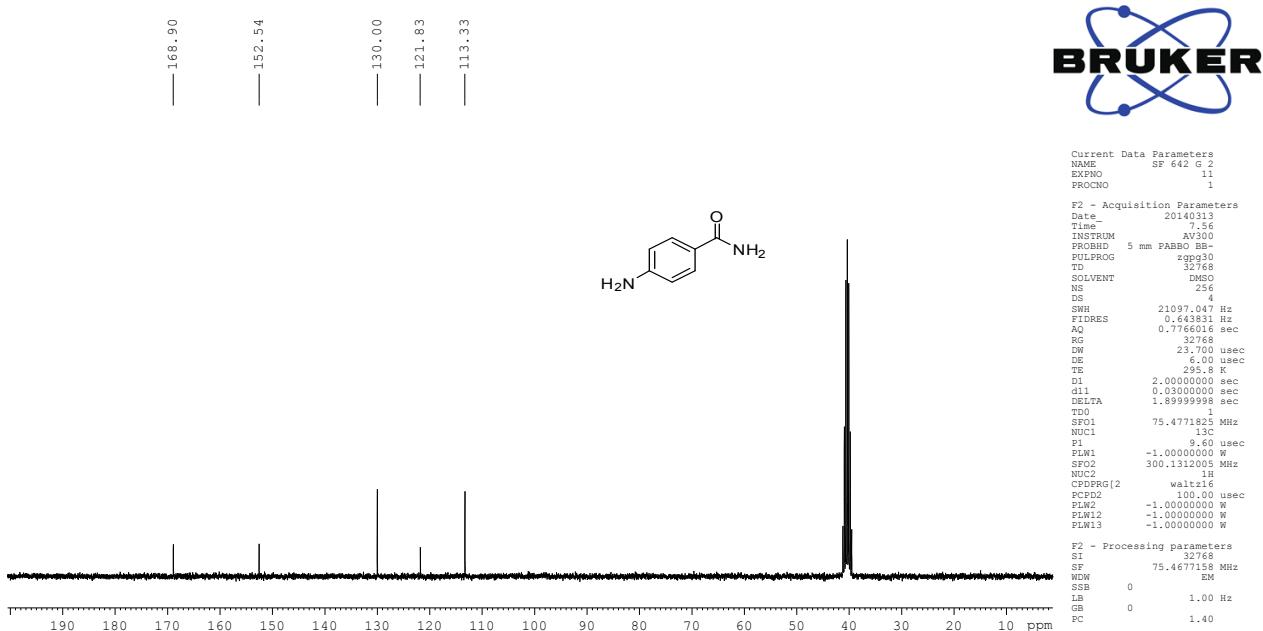
<sup>13</sup>C NMR



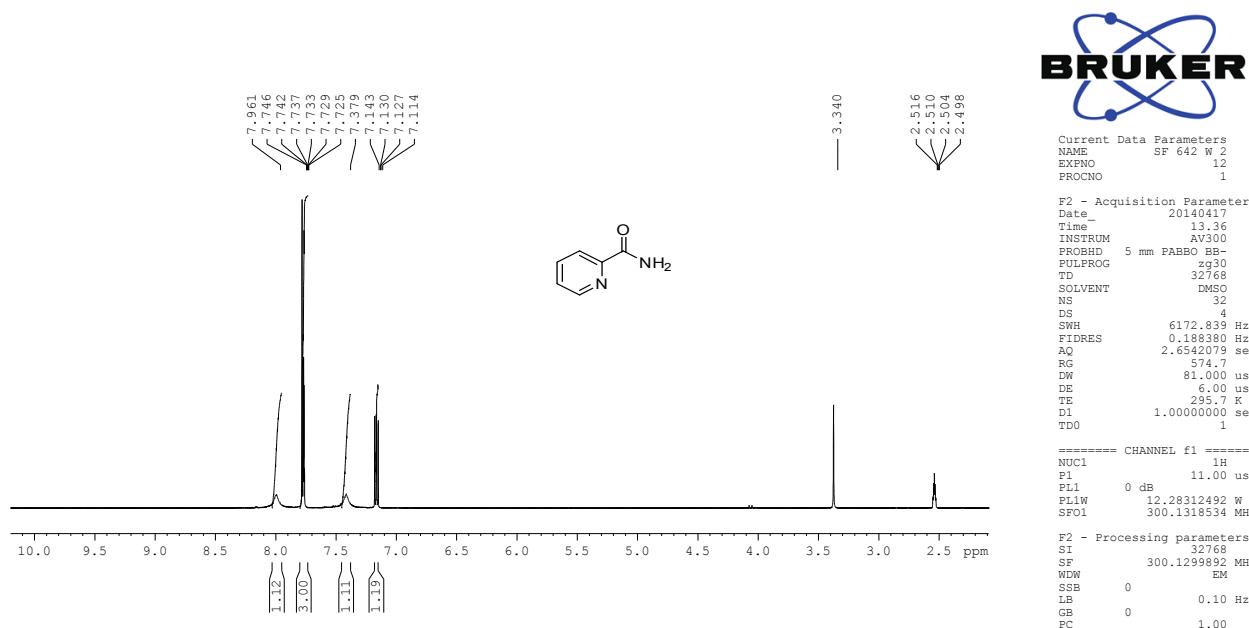
<sup>1</sup>H NMR



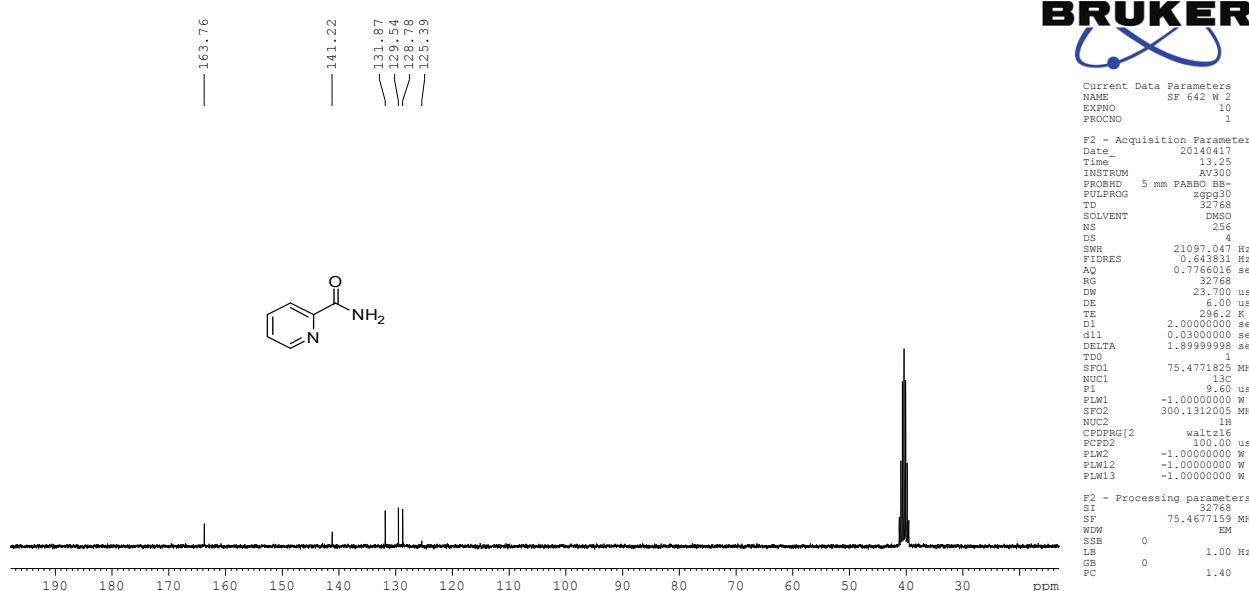
<sup>13</sup>C NMR



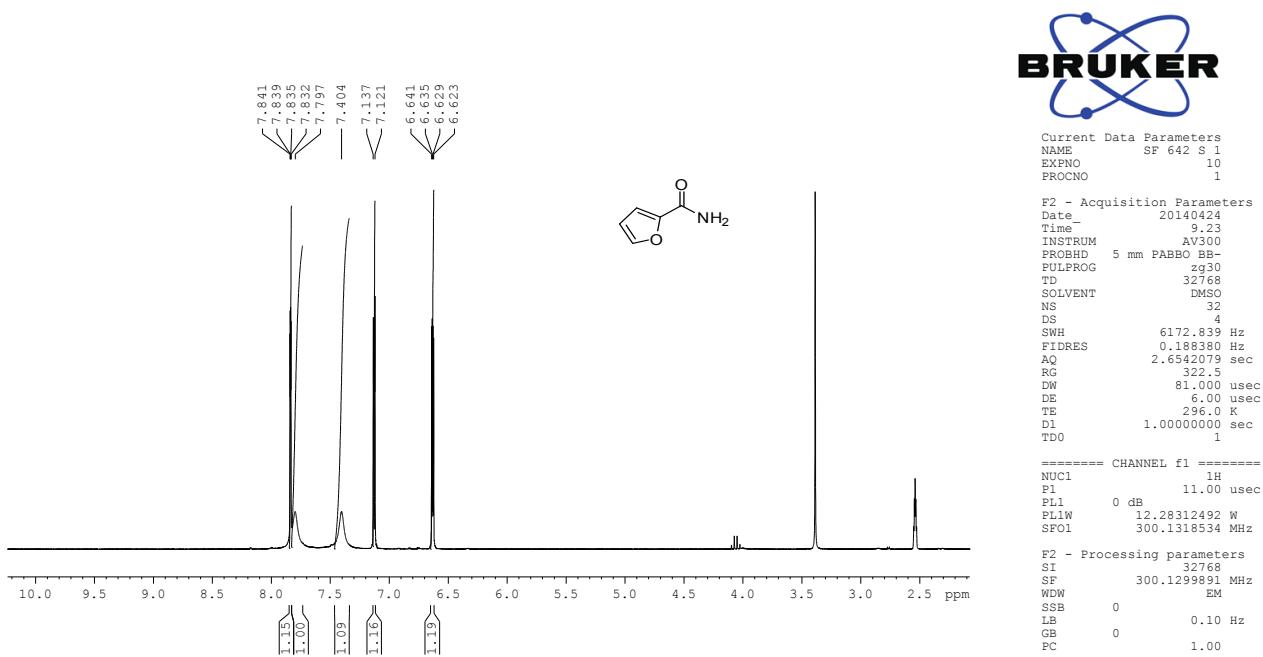
<sup>1</sup>H NMR



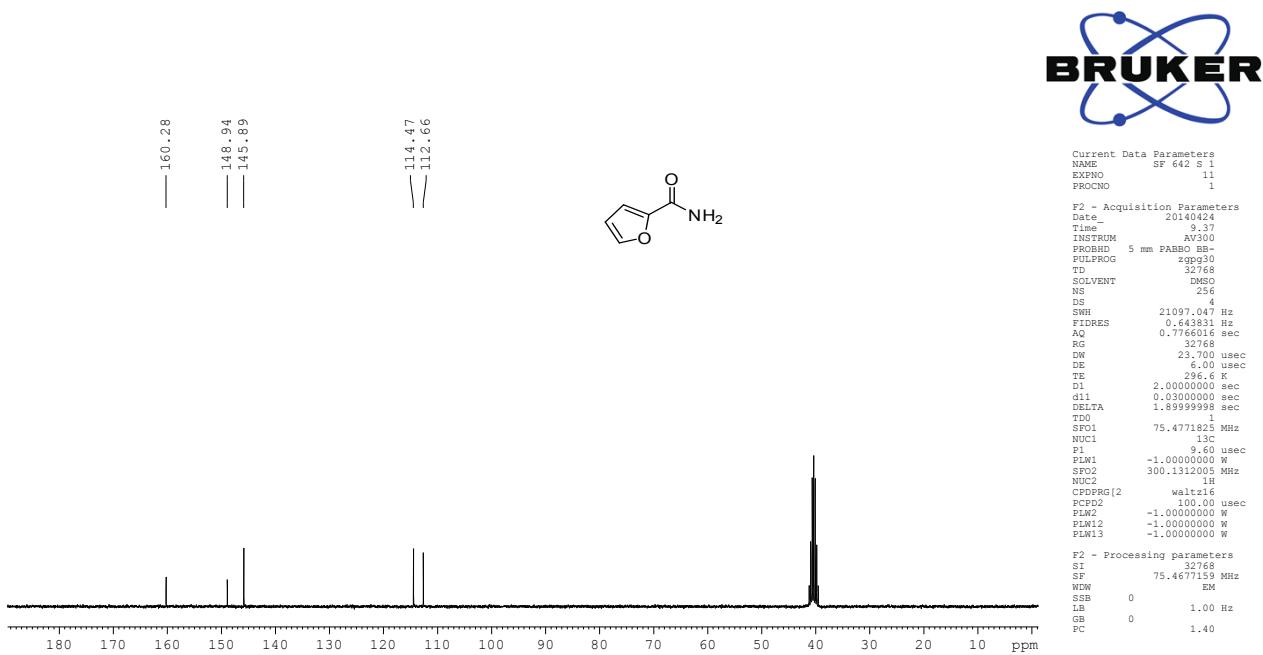
<sup>13</sup>C NMR



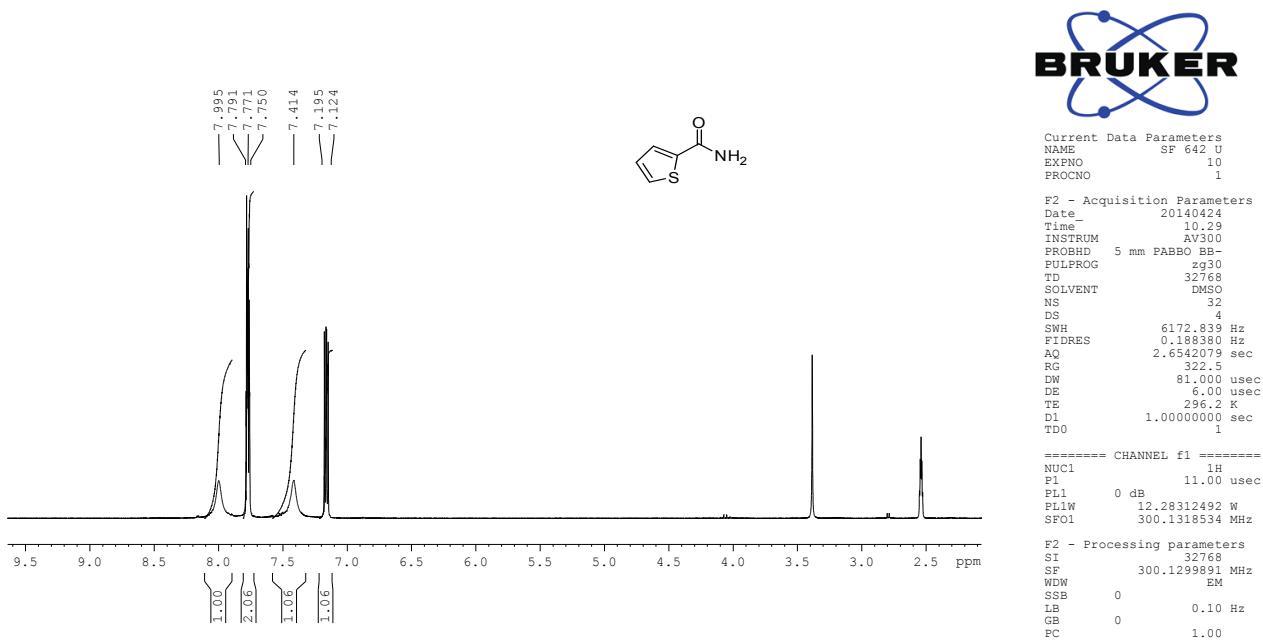
<sup>1</sup>H NMR



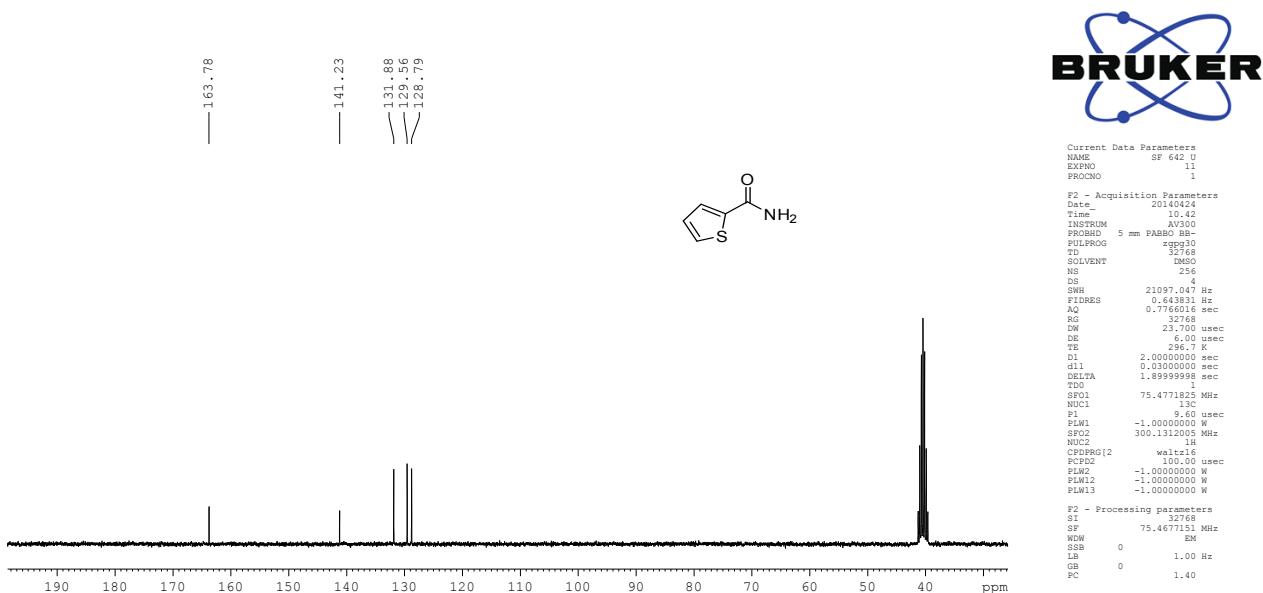
<sup>13</sup>C NMR



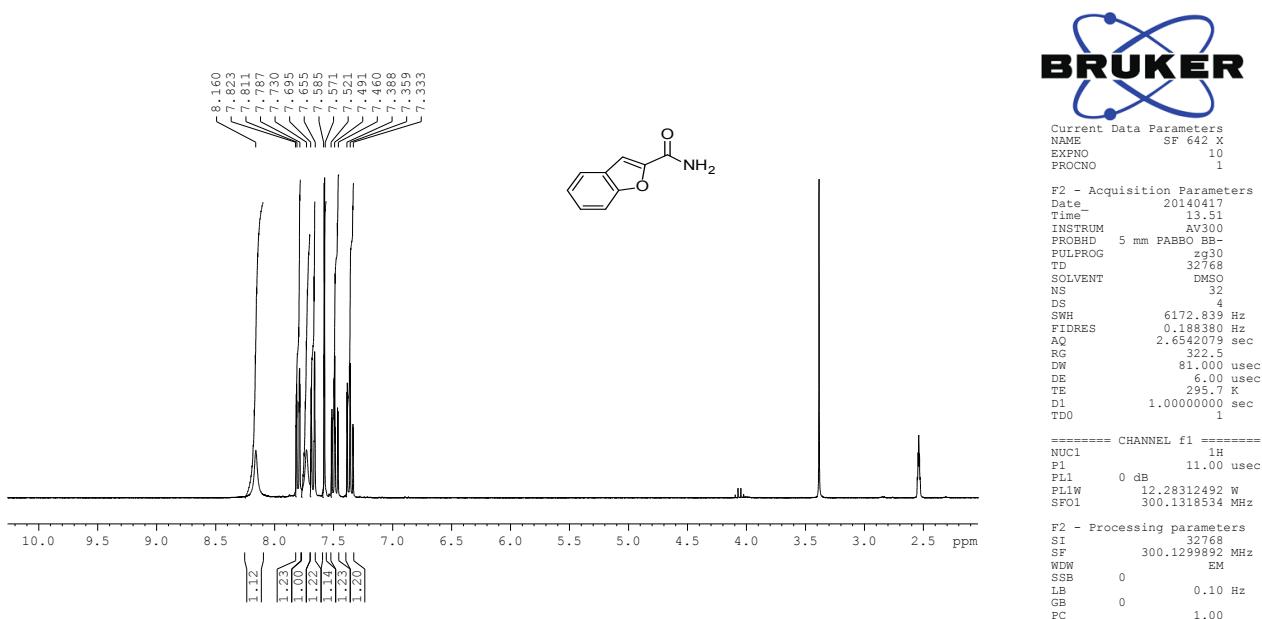
**<sup>1</sup>H NMR**



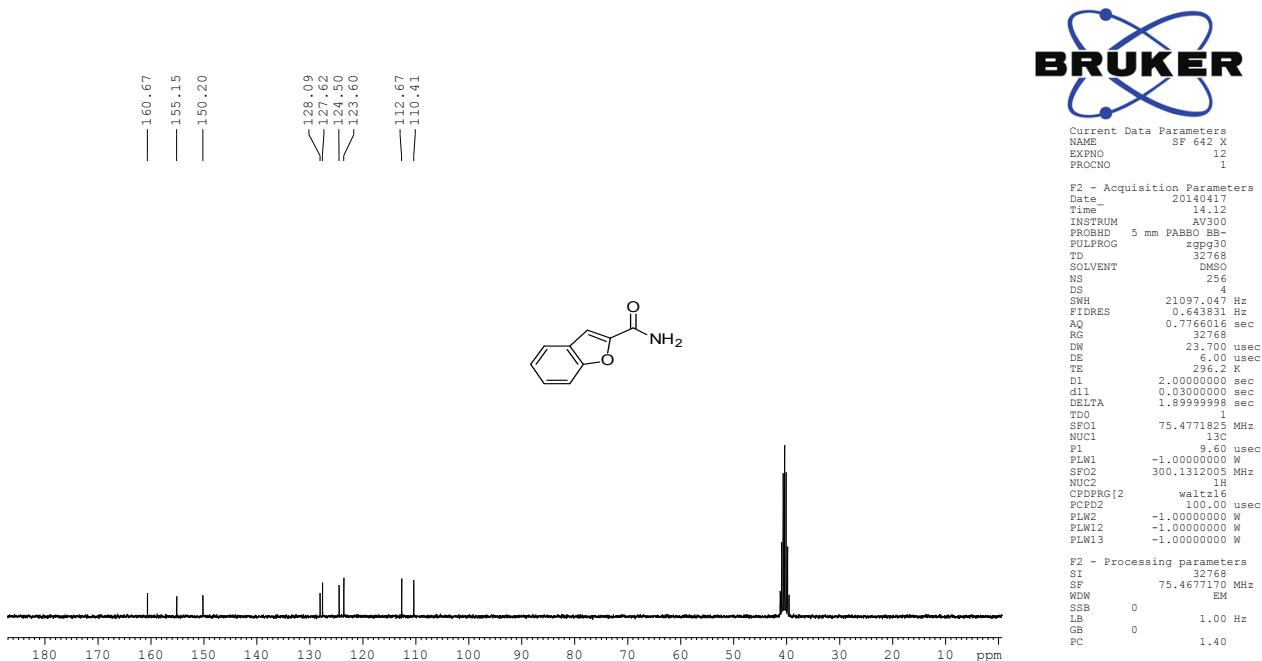
**<sup>13</sup>C NMR**



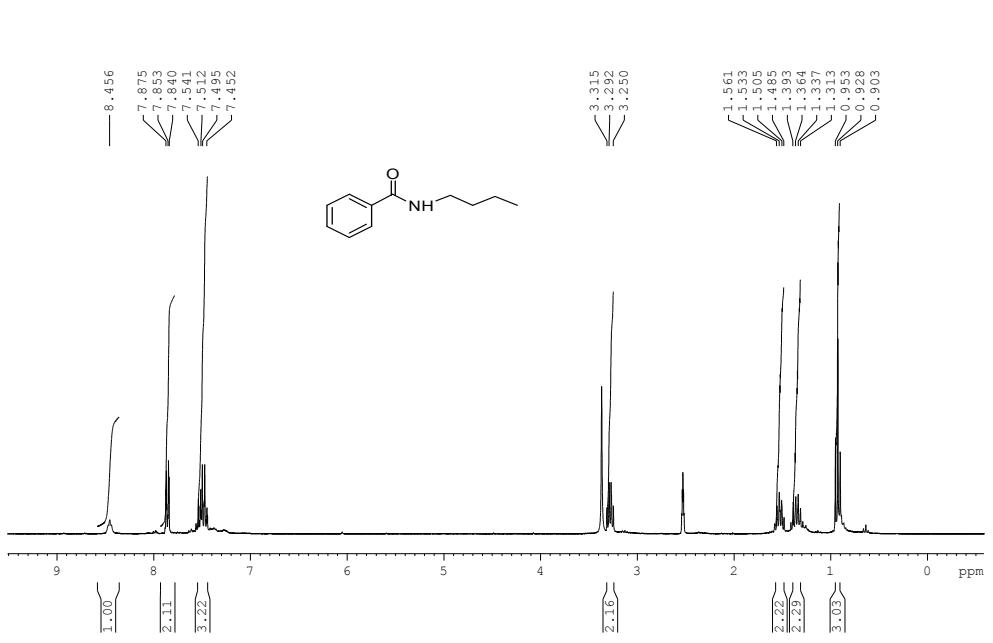
## **<sup>1</sup>H NMR**



## <sup>13</sup>CNMR



### <sup>1</sup>H NMR



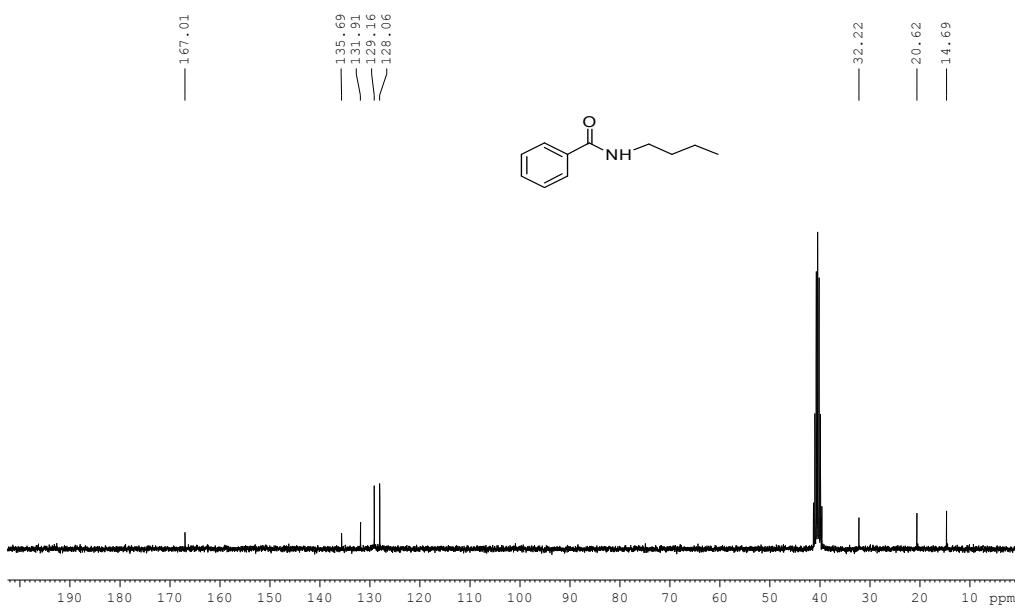
Current Data Parameters  
NAME SF 642 N 2  
EXPNO 10  
PROCNO 1

F2 - Acquisition Parameters  
Date 20140411  
Time 0.48  
INSTRUM AV300  
PROBHD 5 mm PABBO BB-  
PULPROG zg30  
TD 32768  
SOLVENT DMSO  
NS 32  
DS 4  
SWH 6172.82 Hz  
FIDRES 0.188380 Hz  
AQ 2.6542079 sec  
RG 322.5  
DW 81.000 usec  
DE 6.00 usec  
TE 295.6 K  
D1 1.00000000 sec  
TDO 1

===== CHANNEL f1 =====  
NUC1 1H  
P1 11.00 usec  
PL1 0 dB  
PL1W 12.28312492 W  
SF01 300.1318534 MHz

F2 - Processing parameters  
SI 32768  
SF 300.1299922 MHz  
WDW EM  
SSB 0  
LB 0.10 Hz  
GB 0  
PC 1.00

### <sup>13</sup>C NMR

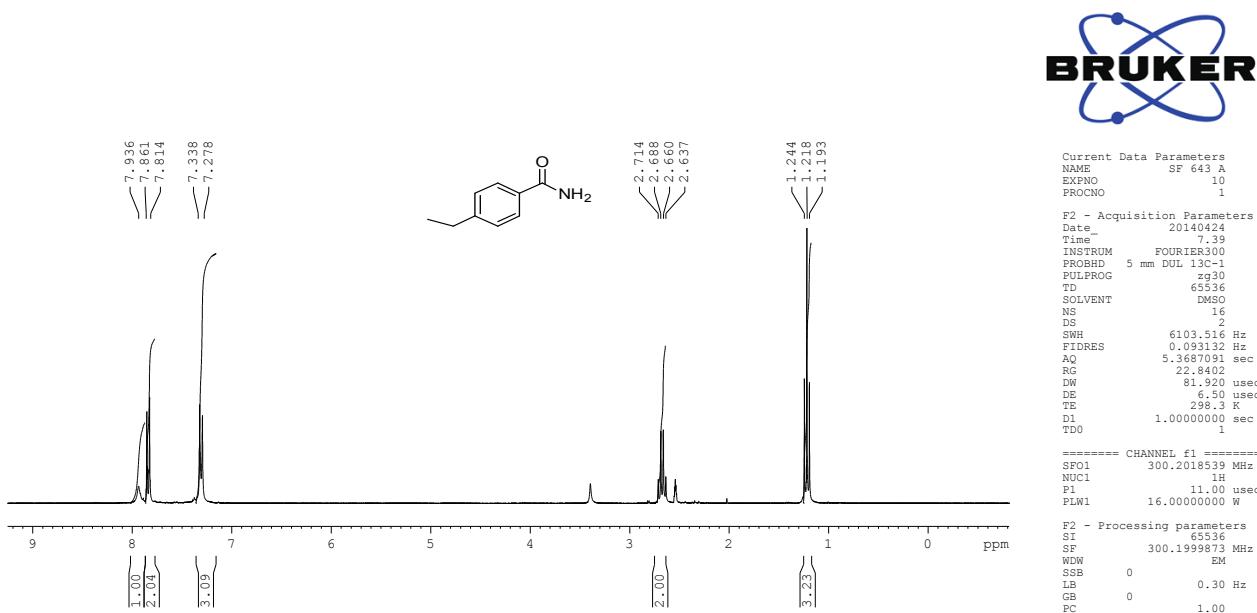


Current Data Parameters  
NAME SF 642 N 2  
EXPNO 11  
PROCNO 1

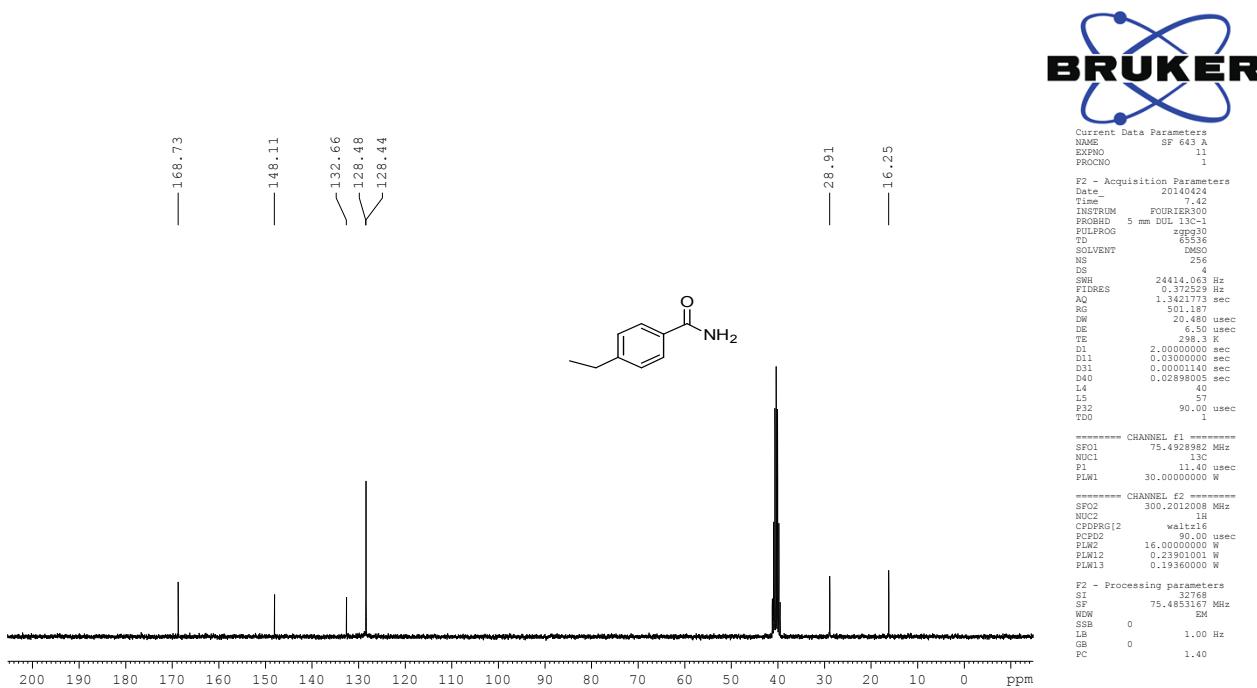
F2 - Acquisition Parameters  
Date 20140411  
Time 1.02  
INSTRUM AV300  
PROBHD 5 mm PABBO BB-  
PULPROG zg30  
TD 32768  
SOLVENT DMSO  
NS 256  
DS 4  
SWH 21097.047 Hz  
FIDRES 0.643831 Hz  
AQ 0.7766016 sec  
RG 32768  
DW 23.700 usec  
DE 6.00 usec  
TE 296.1 K  
D1 2.00000000 sec  
d11 0.03000000 sec  
DELTA 1.89993998 sec  
TDO 1  
SF01 75.4771825 MHz  
NUC1 13C  
P1 9.60 usec  
PLW1 -1.00000000 W  
SF02 300.1312005 MHz  
NUC2 1H  
CPDPBG12 waltz16  
PCPD2 100.00 usec  
PLW2 -1.00000000 W  
PLW12 -1.00000000 W  
PLW13 -1.00000000 W

F2 - Processing parameters  
SI 32768  
SF 75.4677113 MHz  
WDW EM  
SSB 0  
LB 1.00 Hz  
GB 0  
PC 1.40

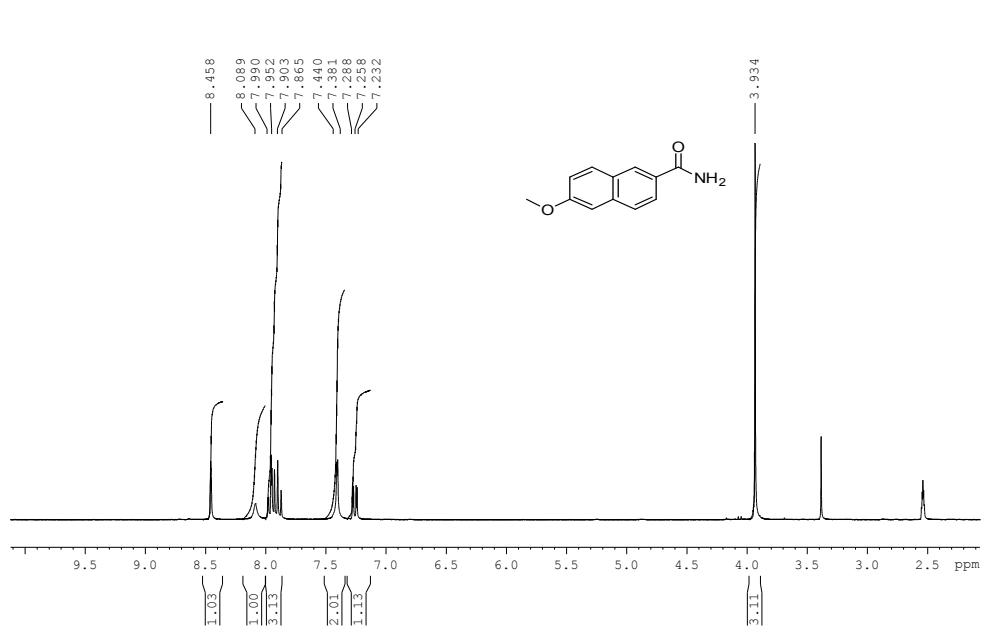
## <sup>1</sup>H NMR



## <sup>13</sup>C NMR

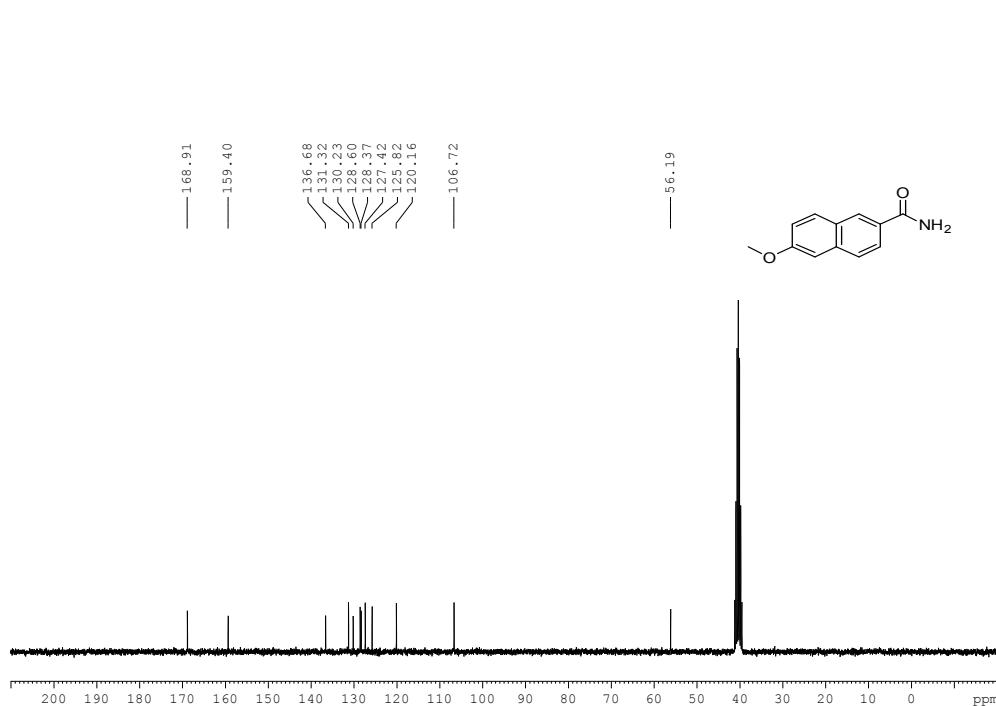


<sup>1</sup>H NMR



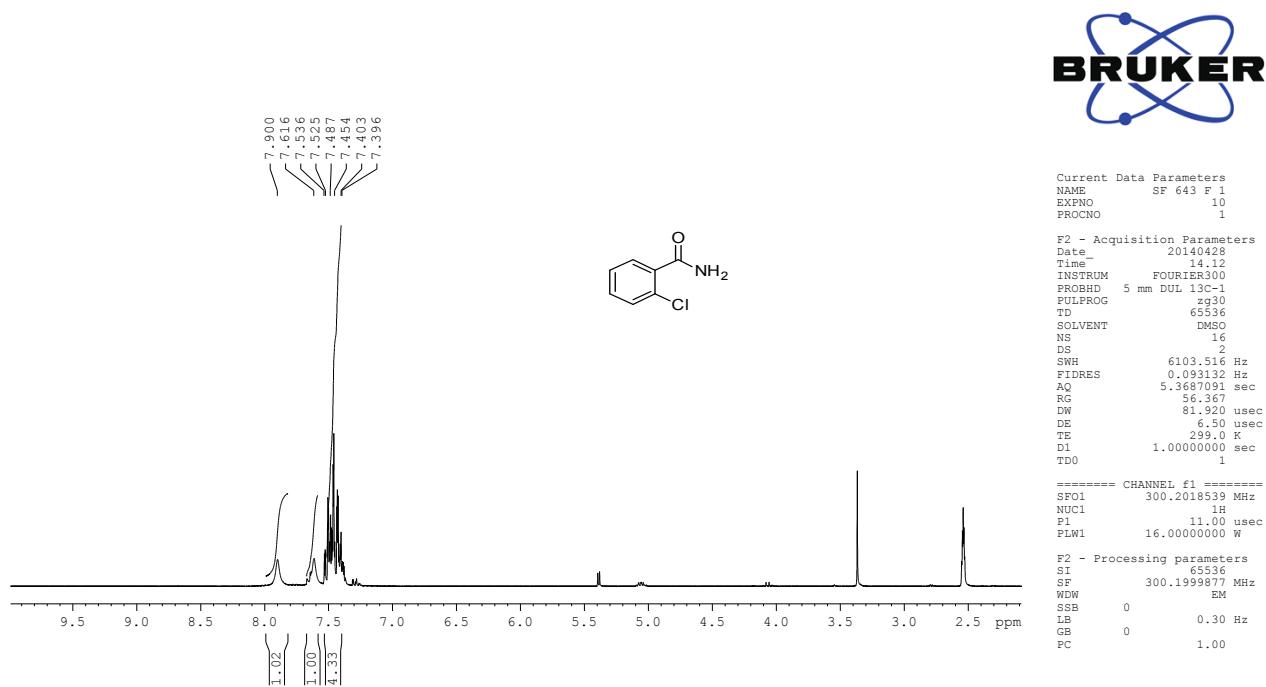
Current Data Parameters  
NAME SF 643 D 2  
EXPNO 10  
PROCNO 1  
  
F2 - Acquisition Parameters  
Date 20140428  
Time 13.38  
INSTRUM FOURIER300  
PROBHD 5 mm DUL 13C-1  
PULPROG zg30  
TD 65536  
SOLVENT DMSO  
NS 16  
DS 2  
SWH 6103.516 Hz  
FIDRES 0.093132 Hz  
AQ 5.368708 sec  
RG 34.8032  
DW 81.920 usec  
DE 6.50 usec  
TE 299.2 K  
D1 1.0000000 sec  
TDO 1  
  
===== CHANNEL f1 =====  
SF01 300.2018539 MHz  
NUC1 1H  
P1 11.00 usec  
PLW1 16.0000000 W  
  
F2 - Processing parameters  
SI 65536  
SF 300.1999877 MHz  
NDW 0 EM  
SSB 0 0.30 Hz  
LB 0  
GB 0 1.00  
PC

<sup>13</sup>C NMR

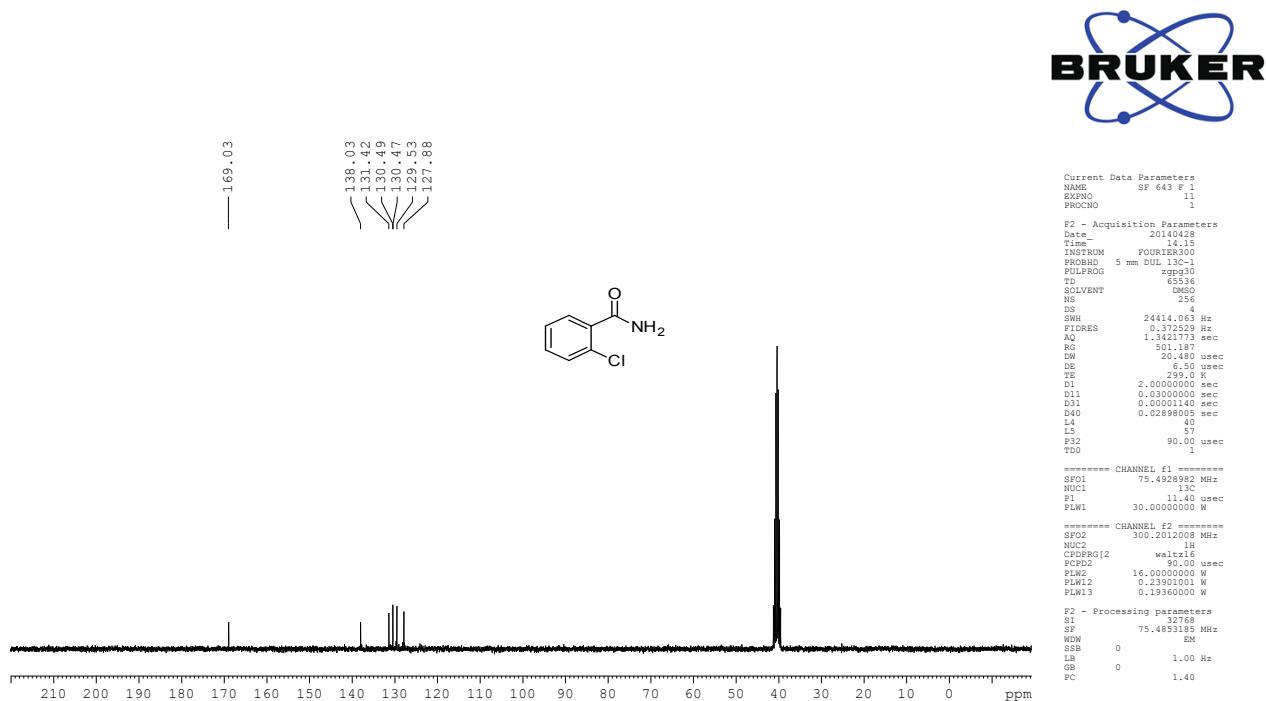


Current Data Parameters  
NAME SF 643 D 2  
EXPNO 11  
PROCNO 1  
  
F2 - Acquisition Parameters  
Date 20140428  
Time 13.41  
INSTRUM FOURIER300  
PROBHD 5 mm DUL 13C-1  
PULPROG zg30  
TD 65536  
SOLVENT DMSO  
NS 256  
DS 4  
SWH 24414.063 Hz  
FIDRES 0.1321773 sec  
AQ 1.3421773 sec  
RG 501.187  
DW 20.480 usec  
DE 5 usec  
TE 299.2 K  
D1 2.0000000 sec  
D11 0.0000000 sec  
D2 0.00001140 sec  
D40 0.02899005 sec  
L4 40  
LS 57  
P22 90.00 usec  
TDO 1  
  
===== CHANNEL f1 =====  
SF01 75.492982 MHz  
NUC1 13C  
P1 11.40 usec  
PLW1 30.0000000 W  
  
===== CHANNEL f2 =====  
SF02 300.2012008 MHz  
NUC2 1H  
CPDPGR2 waltz16  
PCP2 90.00 usec  
PLW2 16.0000000 W  
PLW3 0.2390111 W  
PLW13 0.193860000 W  
  
F2 - Processing parameters  
SI 32768  
SF 75.4853178 MHz  
NDW 0 EM  
SSB 0 1.00 Hz  
LB 0  
GB 0 1.40

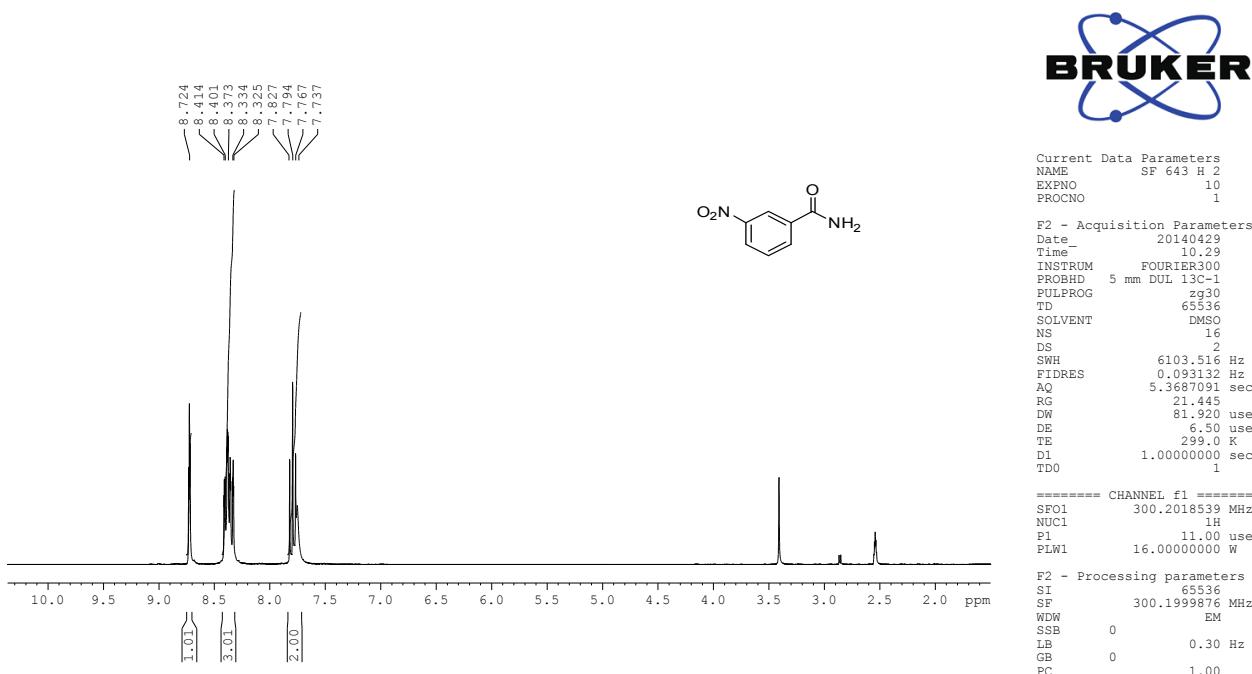
<sup>1</sup>H NMR



<sup>13</sup>C NMR



<sup>1</sup>H NMR



<sup>13</sup>C NMR

