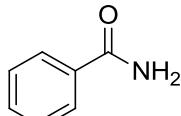


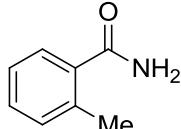
Supporting Information

Benzamide



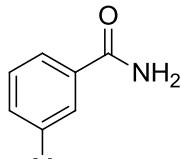
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.40 (s, 1H), 7.47-7.59 (m, 3H), 7.92-7.95 (m, 2H), 8.10 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 127.9 (2CH₂), 128.6 (2CH₂), 131.7 (CH), 134.7 (C), 168.4 (CO).

2-Methylbenzamide



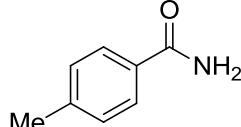
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 2.67 (s, 3H), 7.32-7.52 (m, 2H), 7.64-7.65 (m, 2H), 8.32-8.44 (m, 2H); **¹³CNMR** (DMSO-d₆): δ = 22.2 (CH₃), 126.6, 131.5, 131.7, 132.2 (CH), 133.6, 135.8 (C), 169.6 (CO).

3-Methylbenzamide



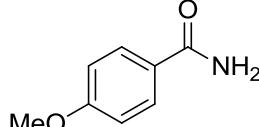
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 2.37 (s, 3H), 7.32-7.39 (m, 2H), 7.67-7.73 (m, 2H), 7.95 (s, 1H), 8.00-8.05 (m, 1H); **¹³CNMR** (DMSO-d₆): δ = 22.2 (CH₃), 126.6, 131.5, 131.7, 132.2 (CH), 133.6, 135.8 (C), 169.6 (CO).

4-Methylbenzamide



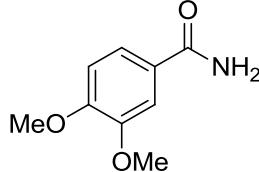
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 2.36 (s, 3H), 7.27 (d, J = 8.23 Hz, 2H), 7.34 (s, 1H, NH₂), 7.83 (d, J = 8.23 Hz, 2H), 7.96 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 21.9 (CH₃), 128.5 (2CH), 129.7 (2CH), 132.2 (C), 142.0 (C), 168.9 (CO).

4-Methoxybenzamide



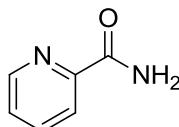
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 3.38 (s, 1H), 3.87 (s, 3H), 7.04 (d, J = 9.18 Hz, 2H), 7.92 (d, J = 9.18 Hz, 2H), 12.6 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 56.3 (OCH₃), 114.6 (2CH), 123.8 (C), 132.2 (2CH), 163.7 (C), 167.8 (CO).

3,4-Dimethoxybenzamide



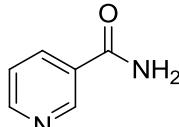
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 3.82 (s, 3H), 3.84 (s, 3H), 6.99-7.05 (m, 2H), 7.23 (s, 1H), 7.47-7.56 (m, 2H), 7.90 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 56.4, 56.5 (2OCH₃), 111.6, 121.6, 127.5 (CH), 149.1, 152.2 (C), 168.4 (CO).

Picolinamide



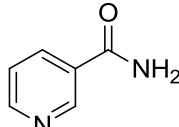
ethyl acetate/hexane (2:1); yield: (%); R_f = ; **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.58-7.62 (m, 1H), 7.71 (s, 1H, NH₂), 7.98-8.03 (m, 1H), 8.07 (s, 1H, NH₂), 8.07-8.09 (m, 1H), 8.17 (s, 1H, NH₂), 8.64-8.66 (m, 1H); **¹³CNMR** (DMSO-d₆): δ = 122.7 (CH), 127.3 (CH), 138.5 (CH), 149.3 (CH), 151.1 (C), 166.9 (CO).

Nicotinamide



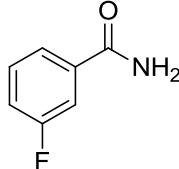
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.48-7.53 (m, 1H), 7.69 (s, 1H, NH₂), 8.24-8.30 (m, 2H), 8.71-8.76 (m, 2H); **¹³CNMR** (DMSO-d₆): δ = 124.4, 130.6, 136.2, 149.7 (CH), 152.9 (C), 167.6 (CO).

Isonicotinamide



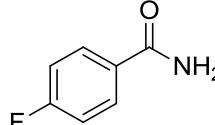
¹H NMR (300 MHz, DMSO-d₆): δ = 7.79-7.84 (m, 3H), 8.72-8.75 (s, 3H); **¹³CNMR** (DMSO-d₆): δ = 122.5 (2CH), 123.9 (C), 142.3 (C), 151.2 (2CH), 167.5 (CO).

3-Fluorobenzamide



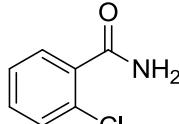
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.38-7.42 (m, 1H), 7.51-7.57 (m, 2H), 7.67-7.71 (m, 1H), 7.74-7.77 (m, 1H), 8.09 (s, 1H, NH₂); **¹³CNMR** (DMSO-d₆): δ = 114.0 (d, J_{3CF} = 24.4 Hz, CH), 118.1 (d, J_{3CF} = 22.2 Hz, CH), 123.6 (d, J_{4CF} = 3.01 Hz, CH), 130.3 (d, J_{3CF} = 8.15 Hz, CH), 136.7 (d, J_{3CF} = 6.91 Hz, C), 161.9 (d, J_{CF} = 244.1 Hz, CF), 166.4 (CO).

4-Fluorobenzamide



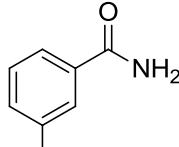
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.29-7.35 (m, 2H), 7.43 (s, 1H), 7.96-8.00 (m, 2H), 8.03 (s, 1H, NH₂); **¹³CNMR** (DMSO-d₆): δ = 115.8 (d, J = 23.8 Hz, 2CH), 130.9 (d, J = 9.40 Hz, 2H), 131.6 (C), 164.7 (d, J = 248.8 Hz, CF), 167.6 (CO).

2-Chlorobenzamide



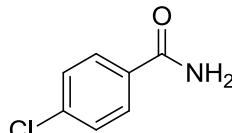
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 6.83-8.13 (m, 6H); **¹³CNMR** (DMSO-d₆): δ = 127.9 (d, J = 13.5 Hz), 129.7, 130.6, 130.9 (CH), 138.0 (C), 169.3 (CO).

3-Chlorobenzamide



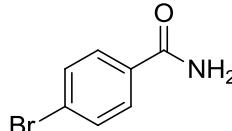
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.50-7.55 (m, 1H), 7.57 (s, 1H, NH₂), 7.60-7.64 (m, 1H), 7.85-7.89 (m, 1H), 7.94-7.96 (m, 1H), 8.12 (s, 1H, NH₂); **¹³CNMR** (DMSO-d₆): δ = 127.0 (CH), 128.1 (CH), 131.1 (CH), 131.9 (CH), 134.0 (C), 137.1 (C), 167.3 (CO)

4-Chlorobenzamide



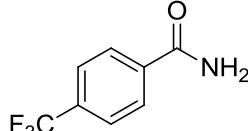
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.49 (s, 1H, NH₂), 7.54-7.58 (m, 2H), 7.91-7.95 (m, 2H), 8.08 (s, 1H, NH₂); **¹³CNMR** (DMSO-d₆): δ = 129.2 (2CH), 130.3 (2CH), 133.9 (C), 137.0 (C), 167.7 (CO).

4-Bromobenzamide



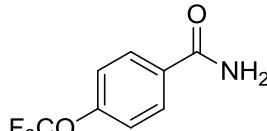
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.53 (s, 1H, NH₂), 7.66-7.71 (m, 2H), 7.83-7.89 (m, 2H), 8.10 (s, 1H, NH₂); **¹³CNMR** (DMSO-d₆): δ = 126.0 (C), 130.5 (2CH), 132.2 (2CH), 134.3 (C), 167.9 (CO).

4-(Trifluoromethyl)benzamide



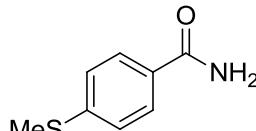
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.69 (s, 1H, NH₂), 7.85-7.89 (m, 2H), 8.10-8.13 (m, 2H), 8.26 (s, 1H, NH₂); **¹³CNMR** (DMSO-d₆): δ = 124.7 (d, *J*_{CF3} = 273.1 Hz, CF₃), 126.2 (2CH), 129.3 (2CH), 132.1 (d, *J* = 31.5 Hz, C), 139.1 (C), 167.7 (CO).

4-(Trifluoromethoxy)benzamide



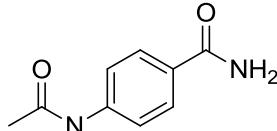
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.43-7.49 (m, 2H), 7.58 (s, 1H), 8.03-8.08 (m, 2H), 8.16 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 121.1 (d, *J*_{CF3} = 253.9 Hz, OCF₃), 121.3 (2CH), 130.6 (2CH), 132.4, 134.2, 151.2 (C), 167.5 (CO).

4-(Thiomethyl)benzamide



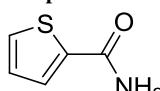
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 3.31 (s, 3H), 7.70 (s, 1H), 8.01-8.09 (m, 2H), 8.10-8.16 (m, 2H), 8.25 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 44.2 (CH₃), 127.9 (2CH), 129.3 (2CH), 139.7, 143.8 (C), 167.5 (CO).

4-Acetamidobenzamide



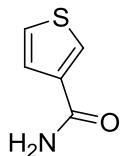
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 2.11 (s, 3H), 7.26 (s, 1H), 7.65-7.68 (m, 2H), 7.83-7.89 (m, 3H), 10.2 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 124.2 (2CH), 129.6 (2CH), 140.7, 149.8 (C), 166.9 (CO).

Thiophene-2-carboxamide



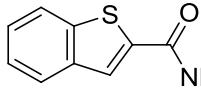
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.15-7.17 (m, 3H), 7.40 (s, 1H), 7.76-7.79 (m, 2H), 7.99 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 128.8, 129.6, 131.9 (CH), 141.3 (C), 163.8 (CO).

Thiophene-3-carboxamide



ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.27 (s, 1H), 7.51-7.53 (m, 1H), 7.58-7.59 (m, 1H), 7.82 (s, 1H), 8.18-8.17 (m, 1H); **¹³CNMR** (DMSO-d₆): δ = 127.4, 128.0, 129.9 (CH), 138.9 (C), 164.6 (CO).

Benzo[b]thiophene-2-carboxamide



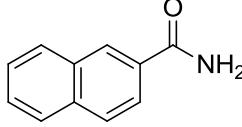
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.43-7.54 (m, 2H), 7.67 (s, 1H), 7.93-7.98 (m, 1H), 8.02-8.07 (m, 1H), 8.09-8.12 (m, 1H), 8.28 (m, 1H); **¹³CNMR** (DMSO-d₆): δ = 123.7, 125.7, 125.9, 126.1, 127.0 (CH), 140.1, 141.2, 141.3 (C), 164.2 (CO).

1-Naphthamide



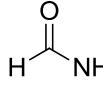
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.54-7.70 (m, 5H), 7.98-8.07 (m, 3H), 8.33-8.38 (m, 1H); **¹³CNMR** (DMSO-d₆): δ = 125.7, 125.9, 126.4, 126.9, 127.3, 128.9, 130.4 (CH), 130.5, 133.7, 133.9, 135.4(C), 171.2 (CO).

2-Naphthamide



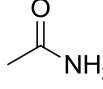
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 7.60-7.77 (m, 3H), 7.79-8.04 (m, 4H), 8.13-8.17 (m, 1H), 8.64 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 126.2, 127.8, 128.7, 129.1, 129.3, 129.4, 130.5 (CH), 131.4, 133.2, 135.9 (C), 168.6 (CO).

Formamide

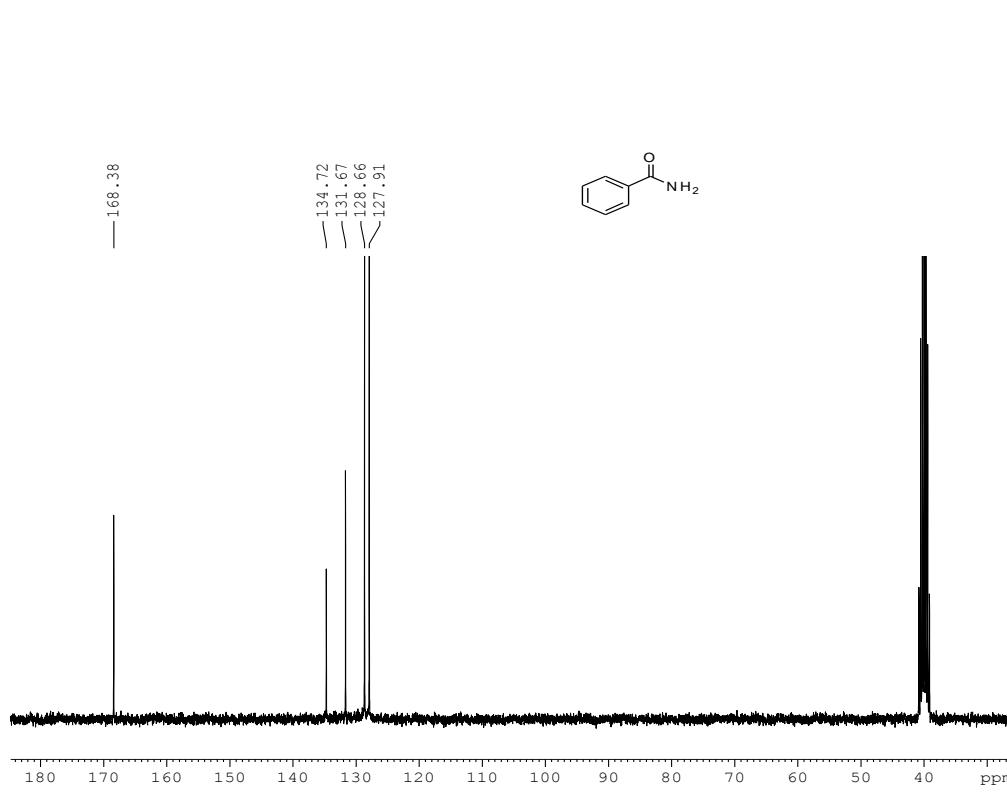
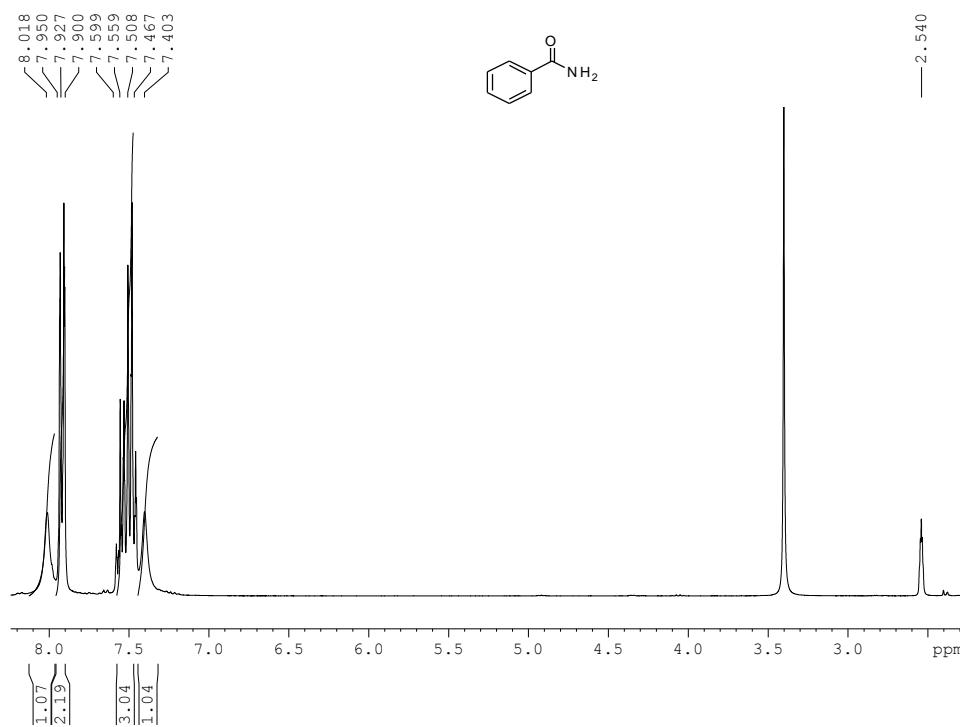


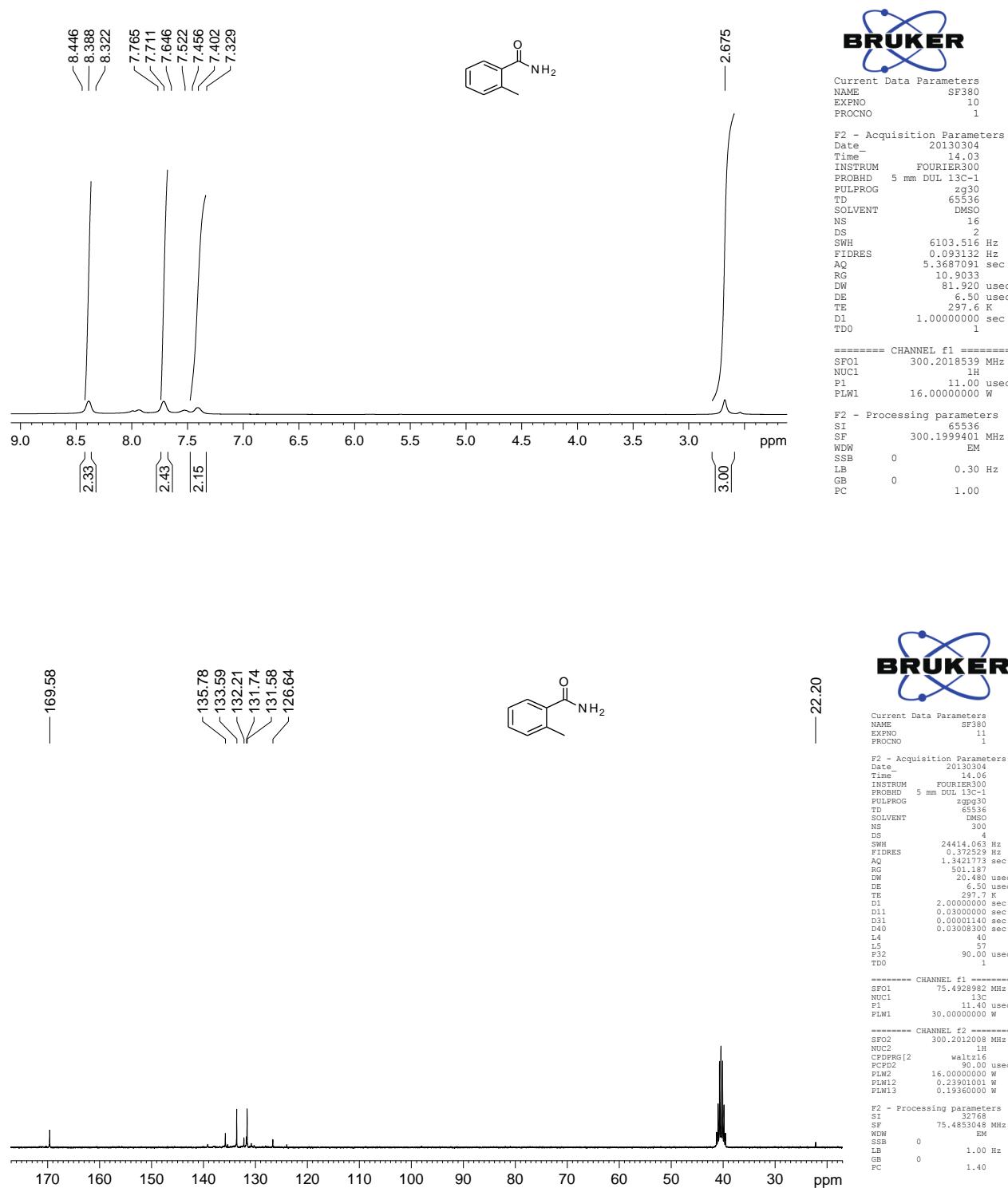
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 5.52 (s, 2H), 7.37 (m, 1H); **¹³CNMR** (DMSO-d₆): δ = 160.6 (CO).

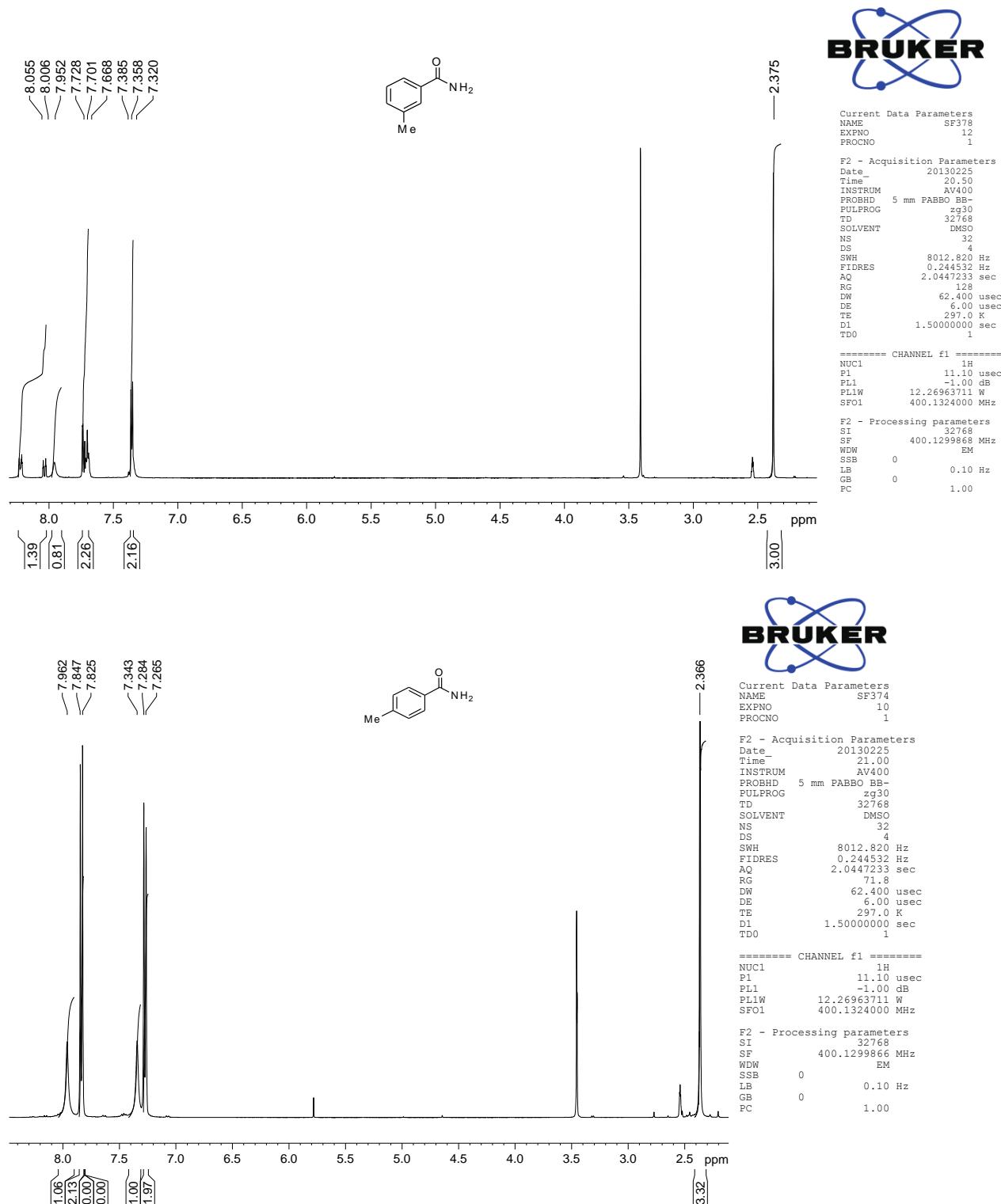
acetamide



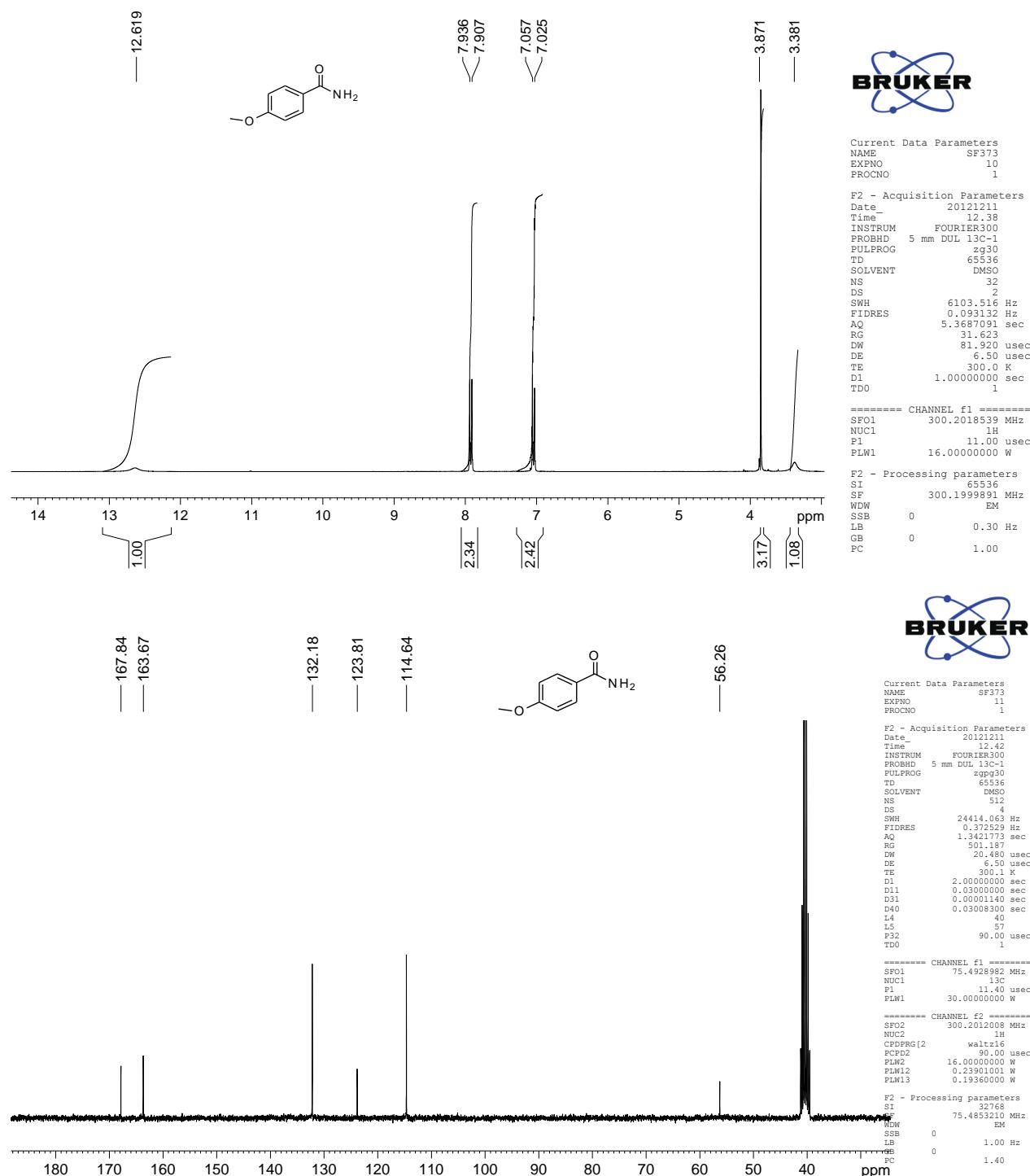
ethyl acetate/hexane (2:1); **¹H NMR** (300 MHz, DMSO-d₆): δ = 1.79 (s, 3H), 6.71 (s, 1H), 7.31 (s, 1H); **¹³CNMR** (DMSO-d₆): δ = 23.4 (CH₃), 160.6 (CO).



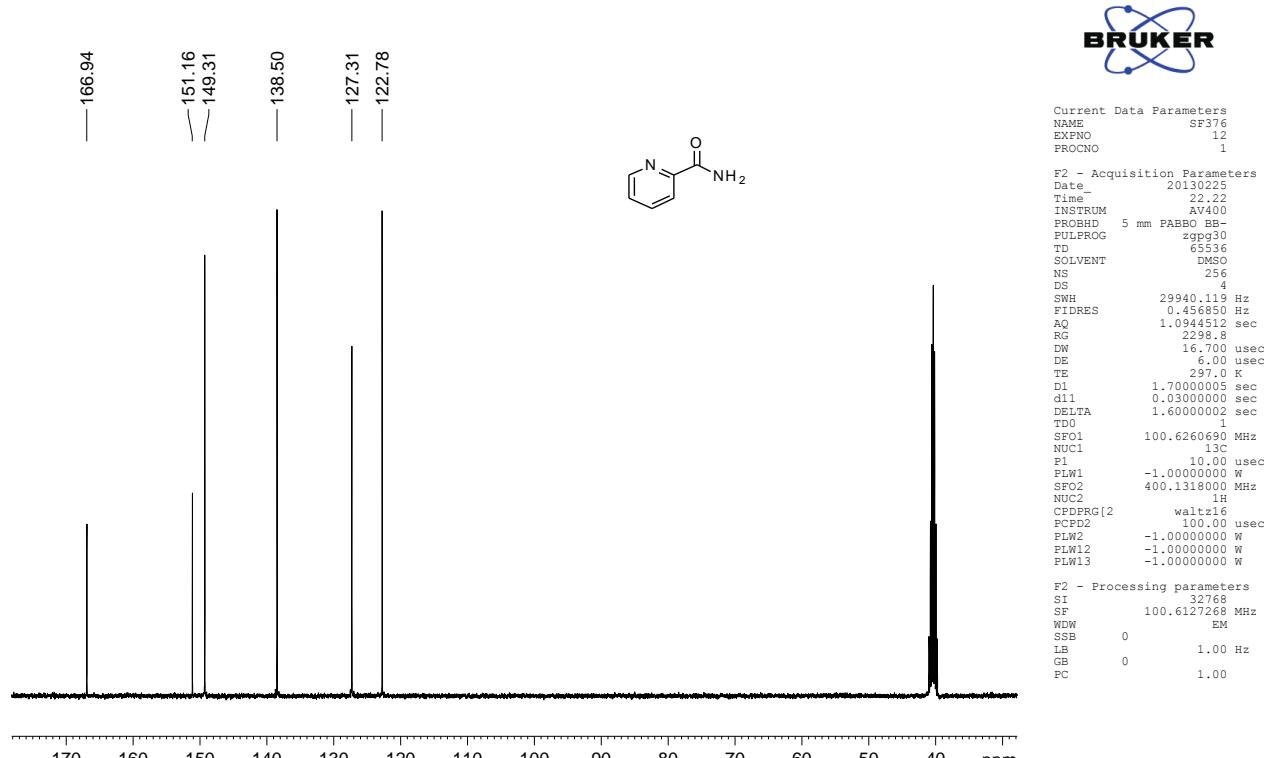
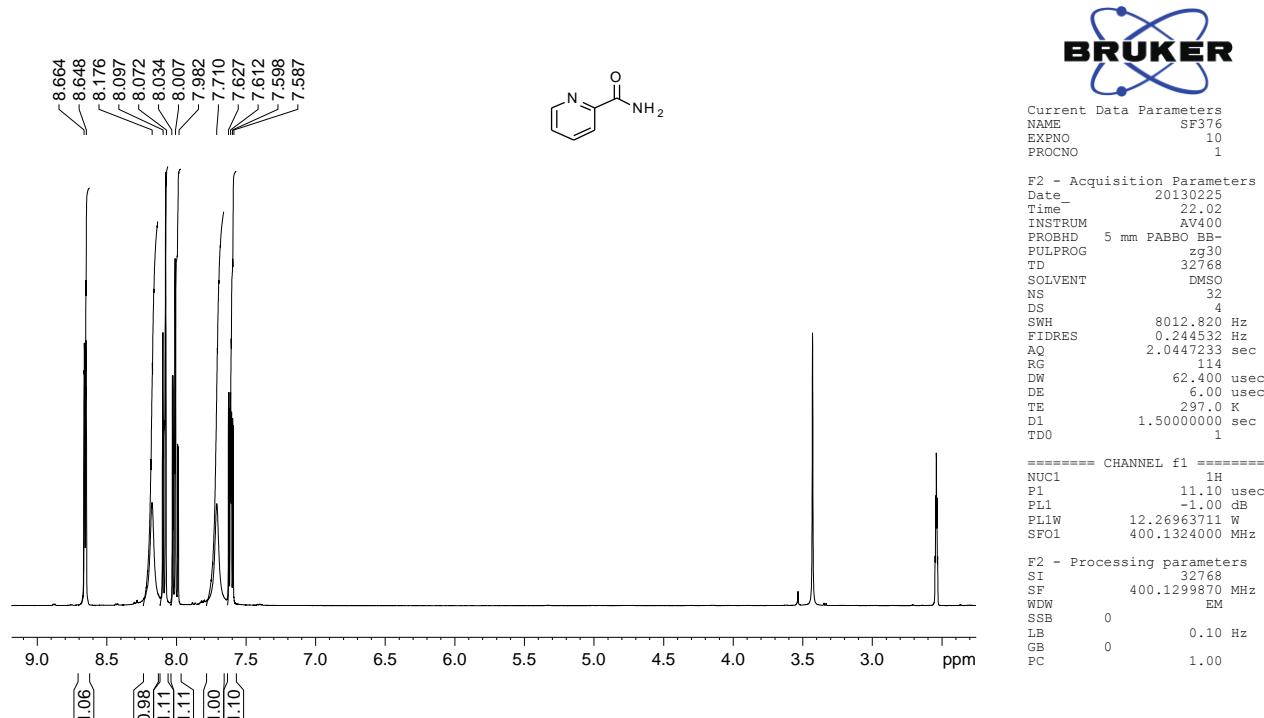




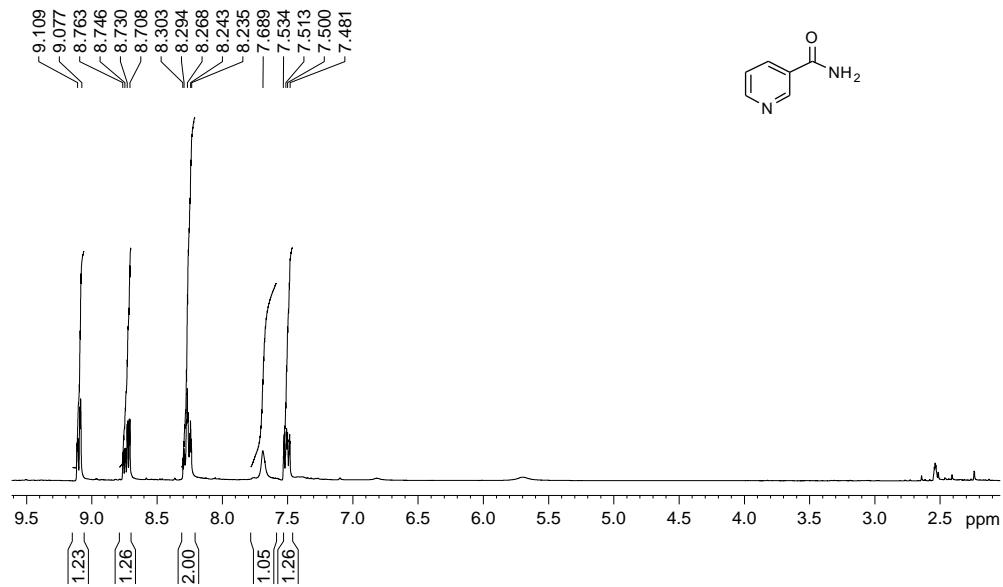
4-methoxybenzamide



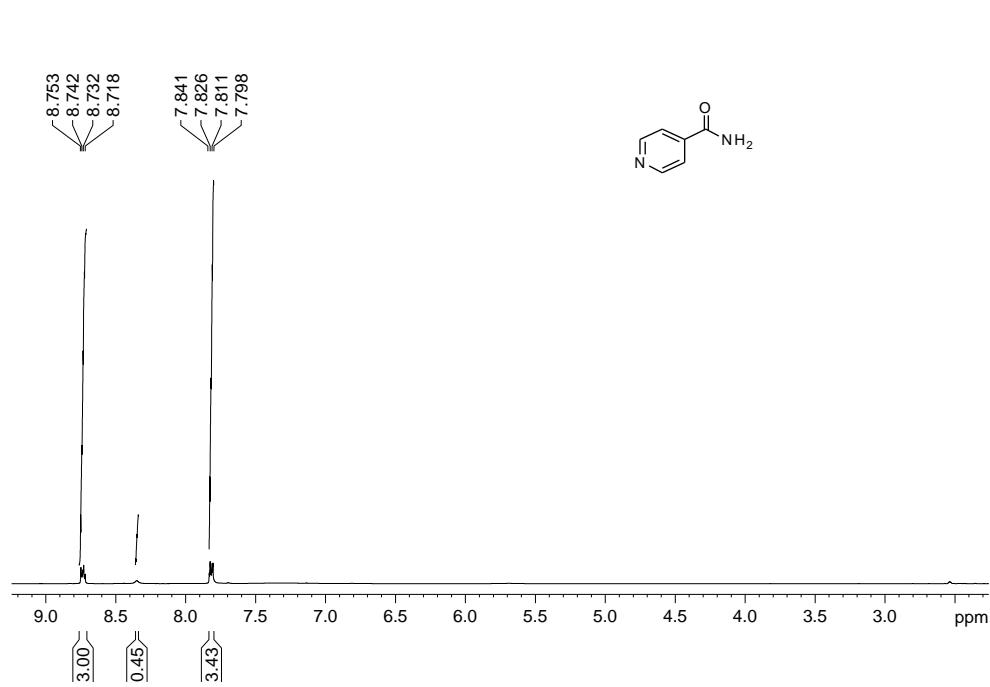
picolinamide

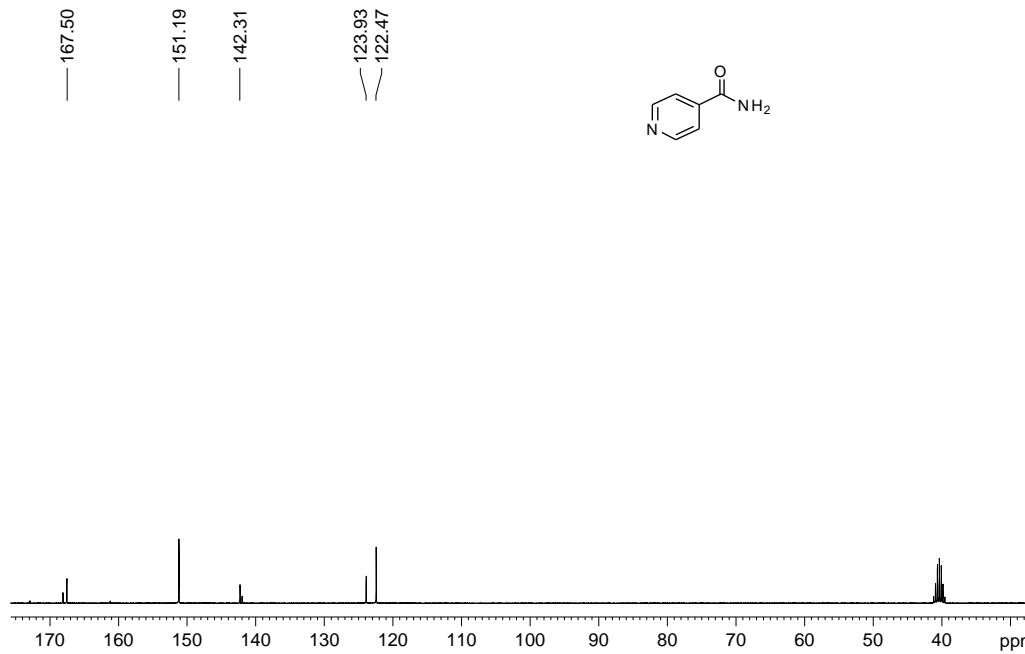


nicotinamide

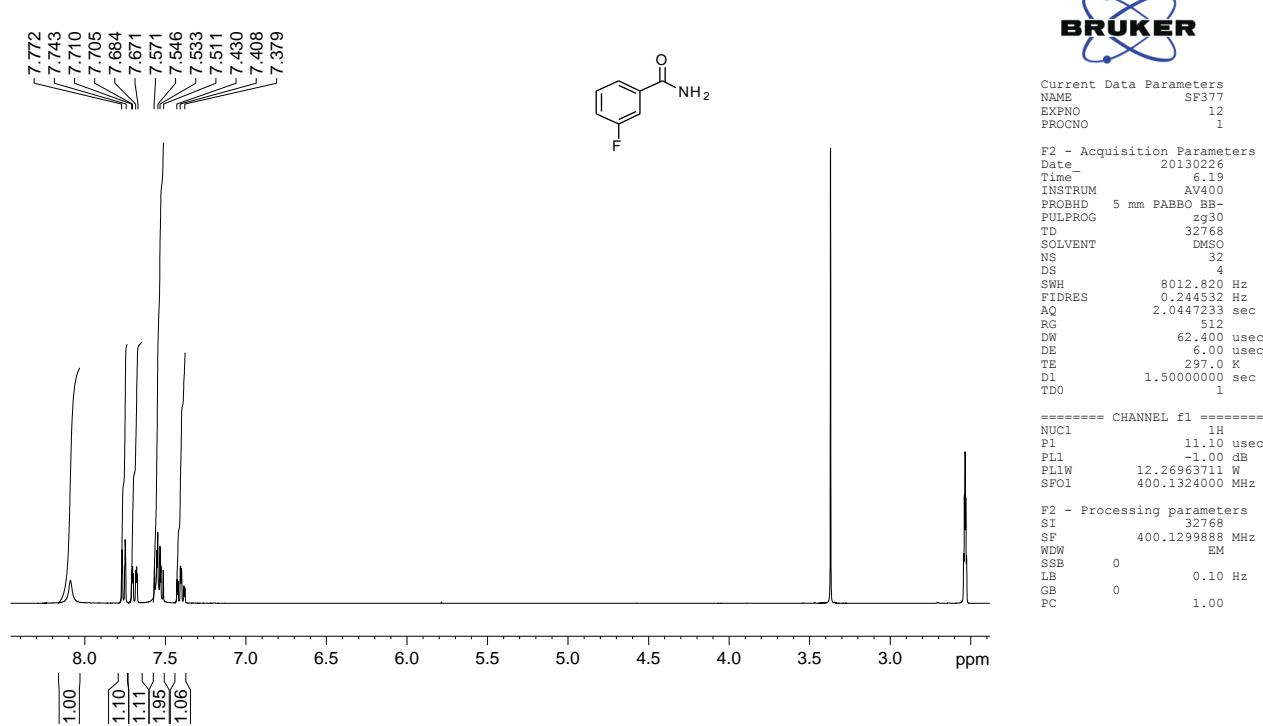


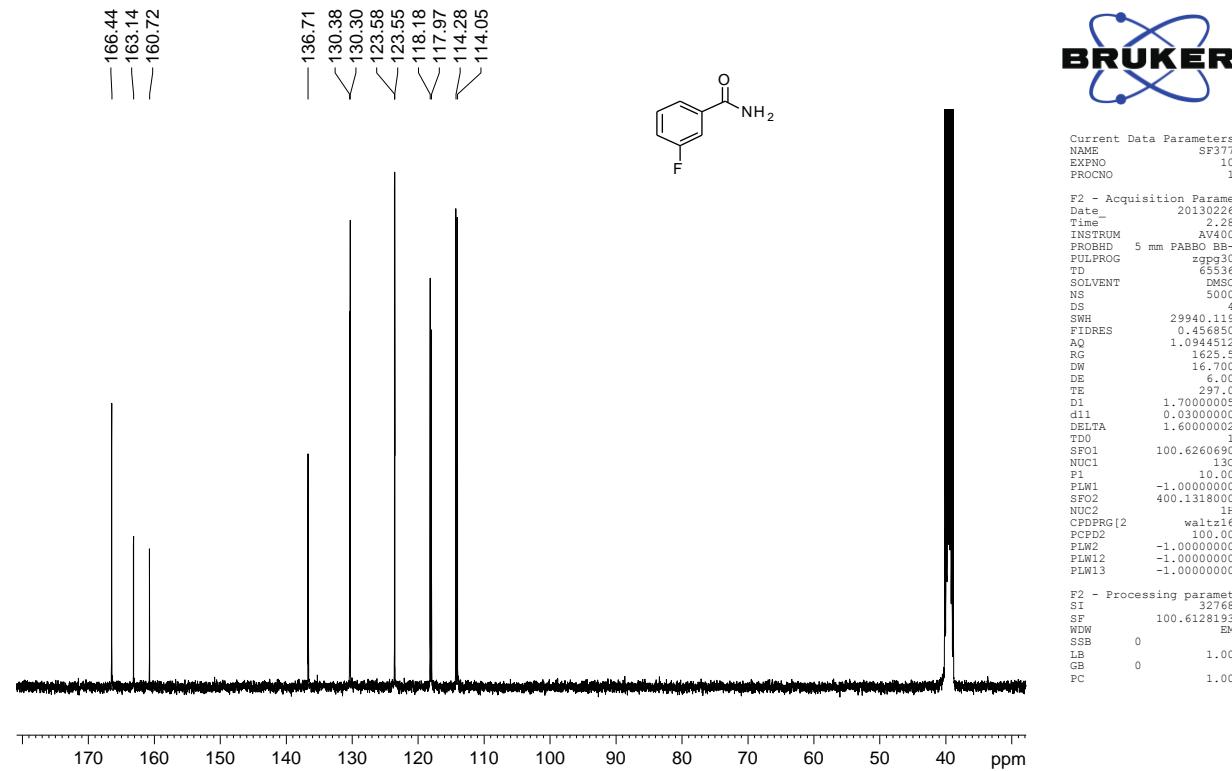
isonicotinamide



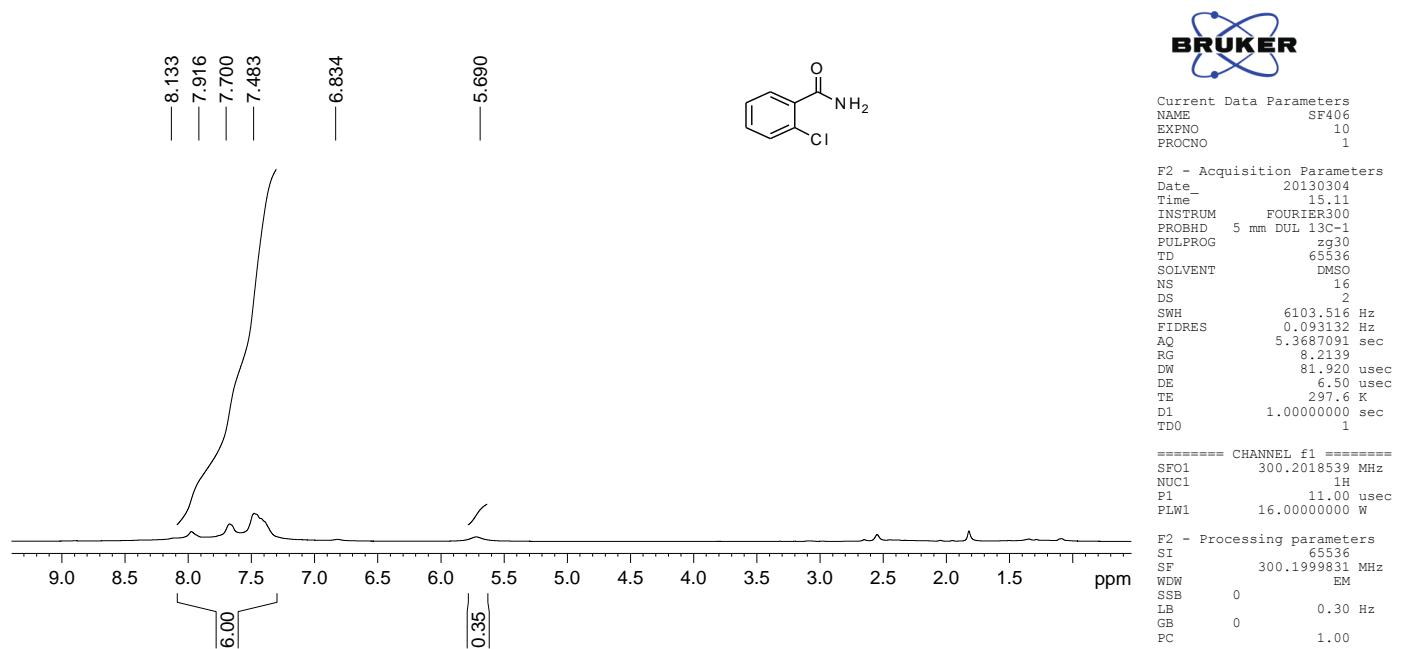


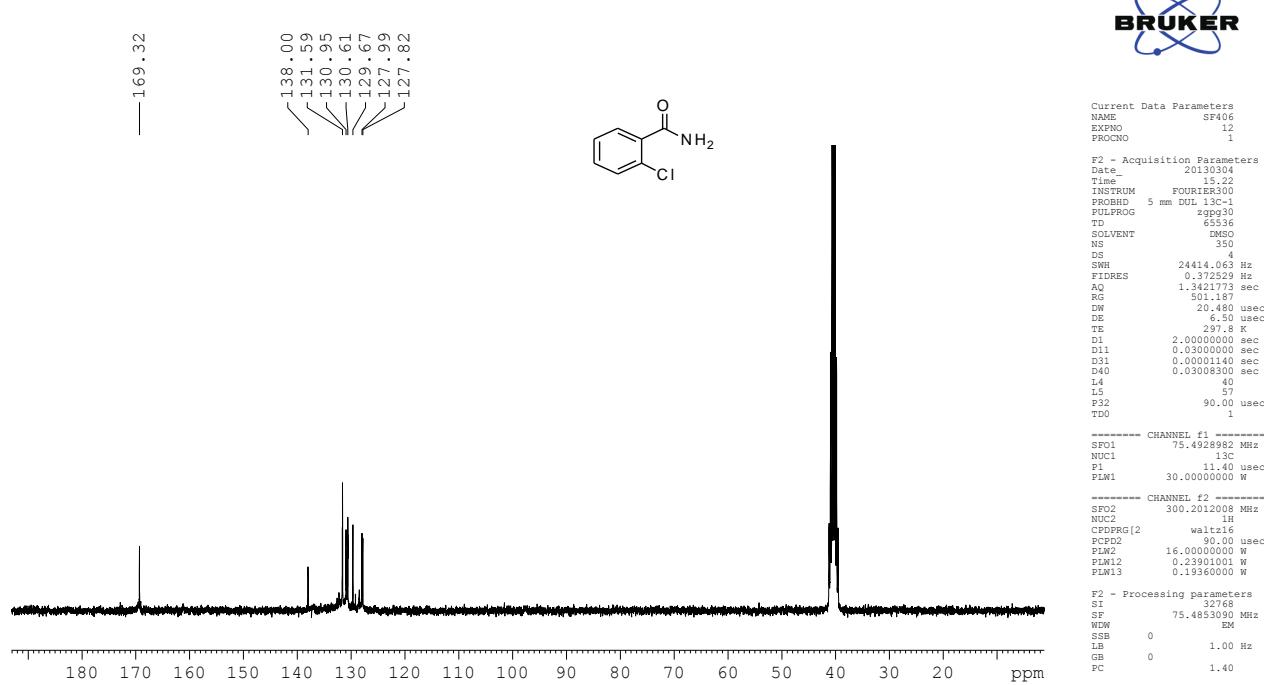
3-fluorobenzamide



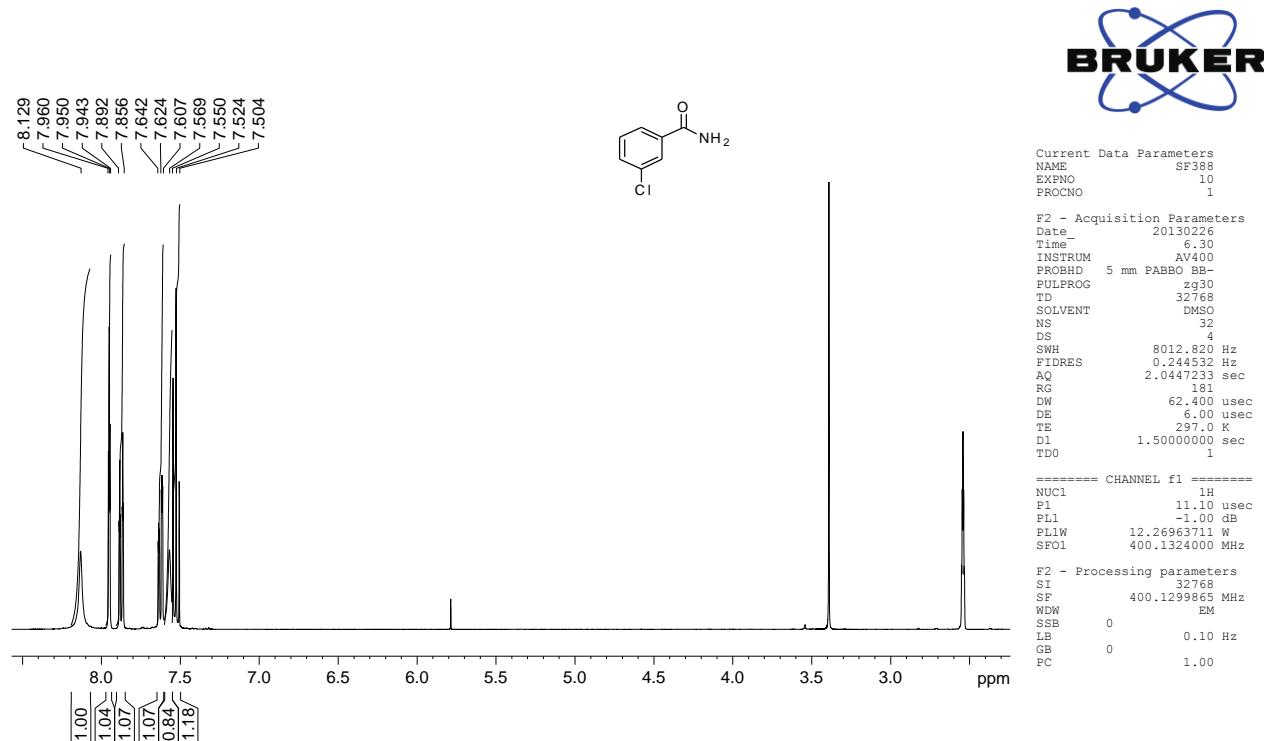


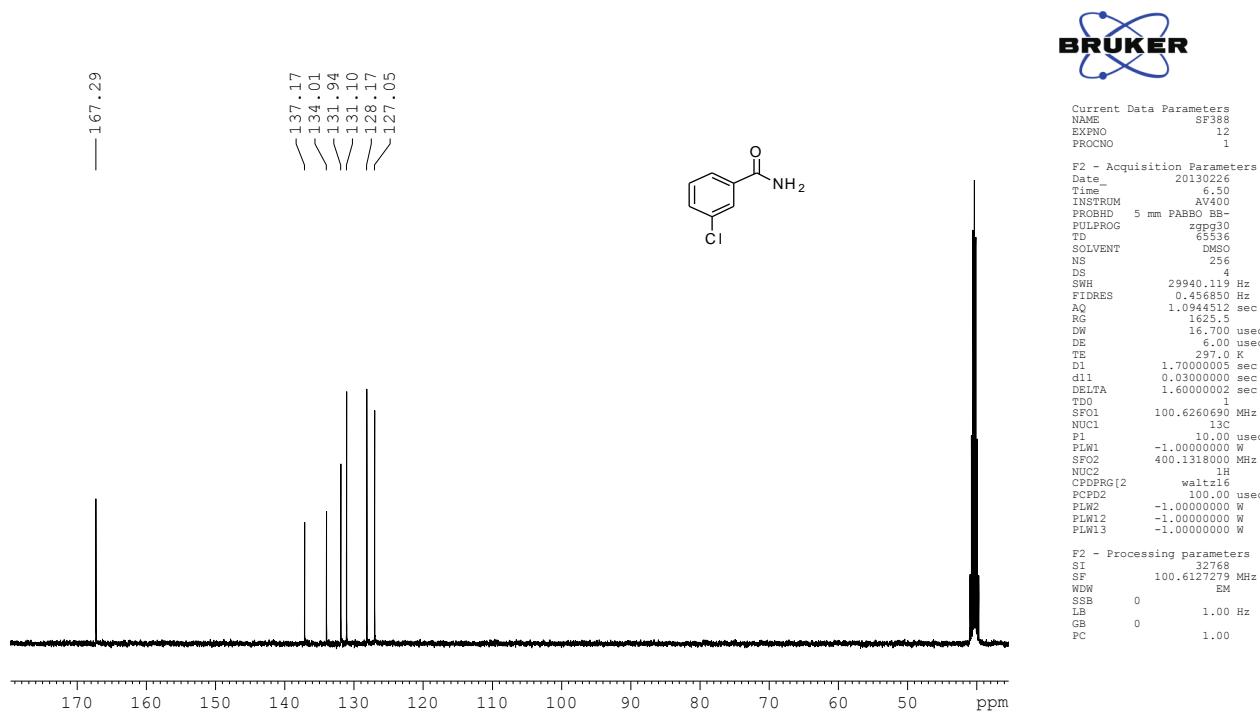
2-chlorobenzamide



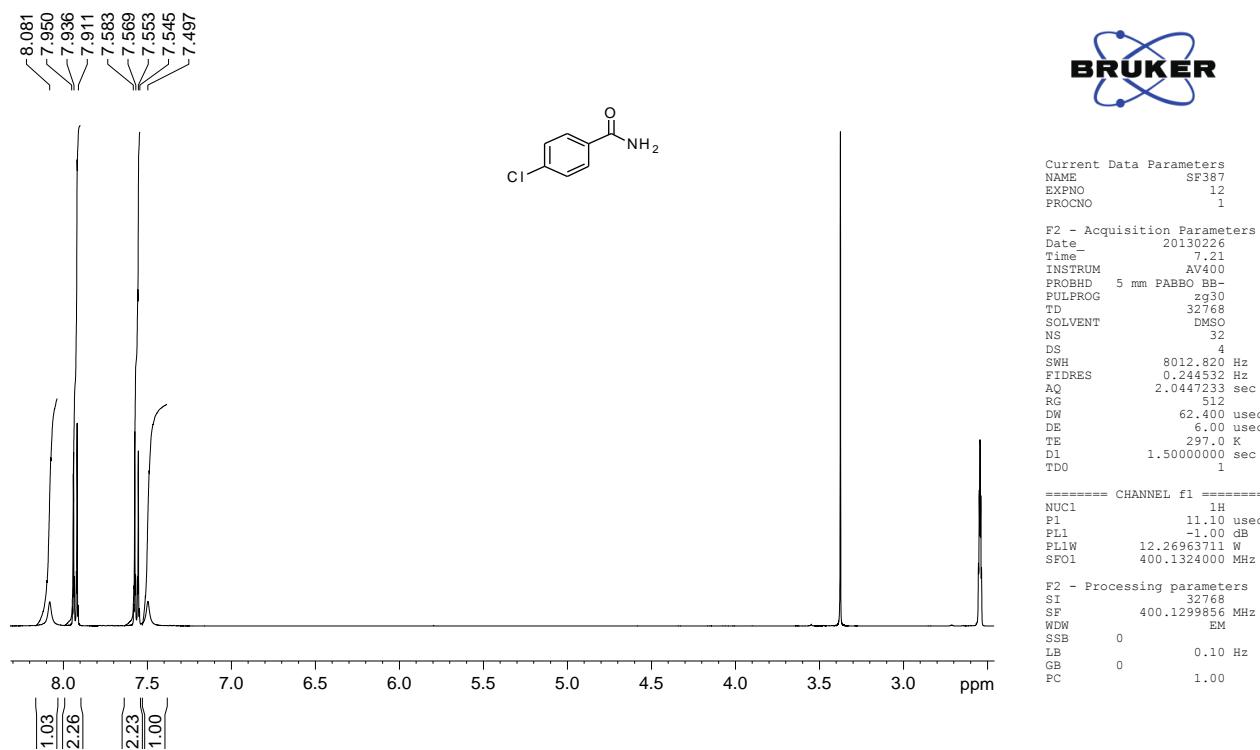


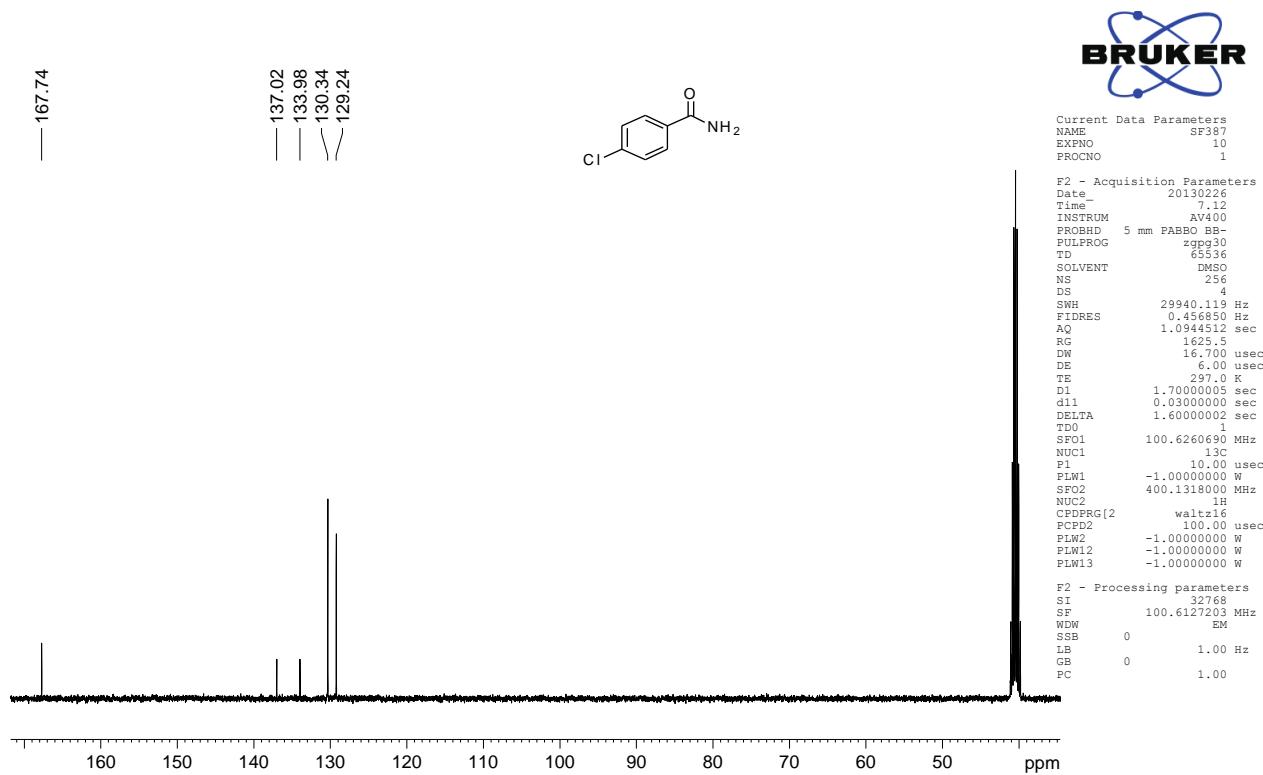
3-chlorobenzamide



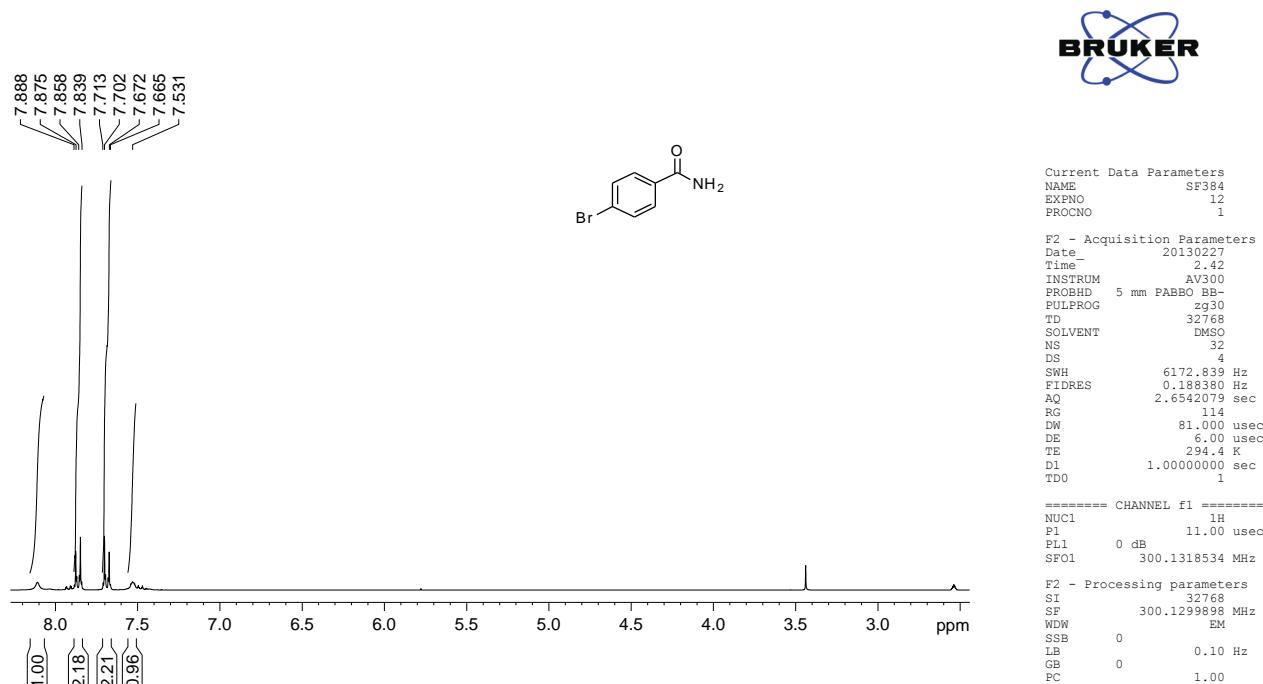


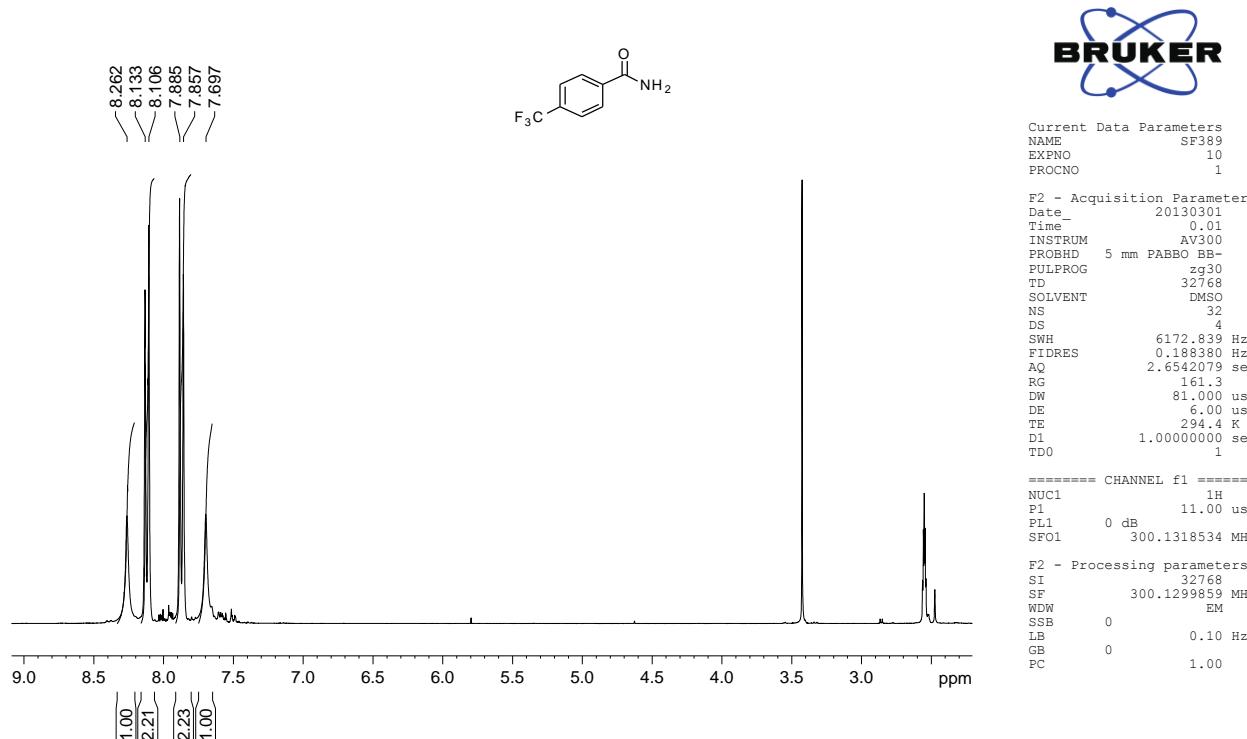
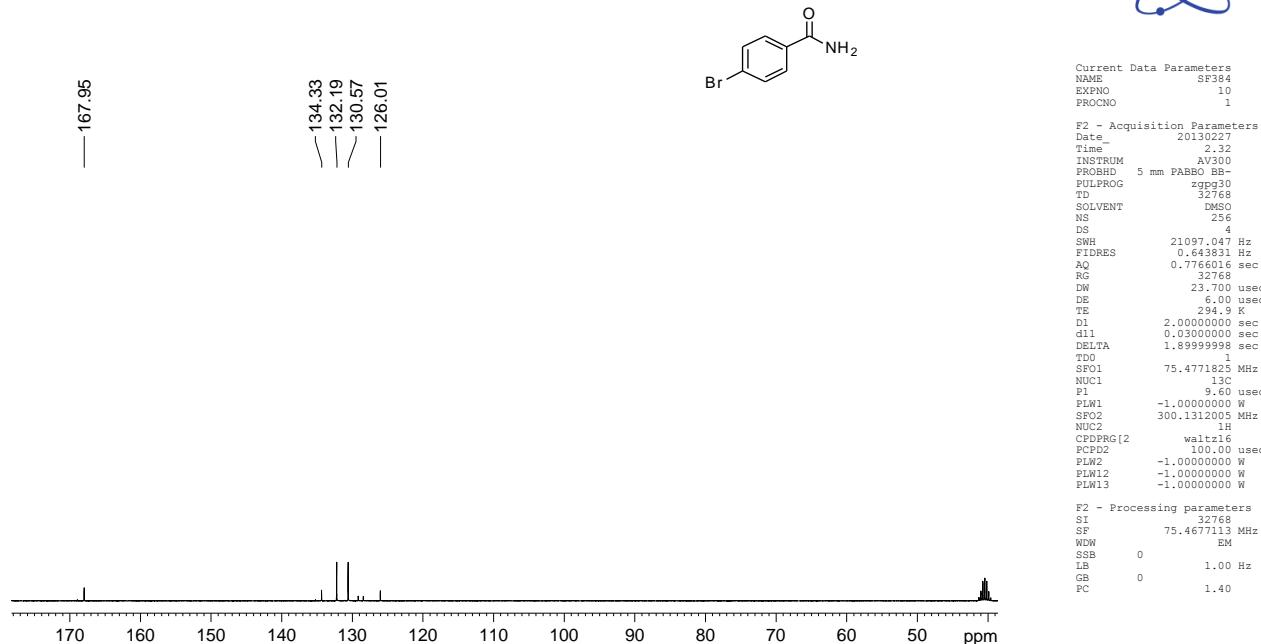
4-chlorobenzamide

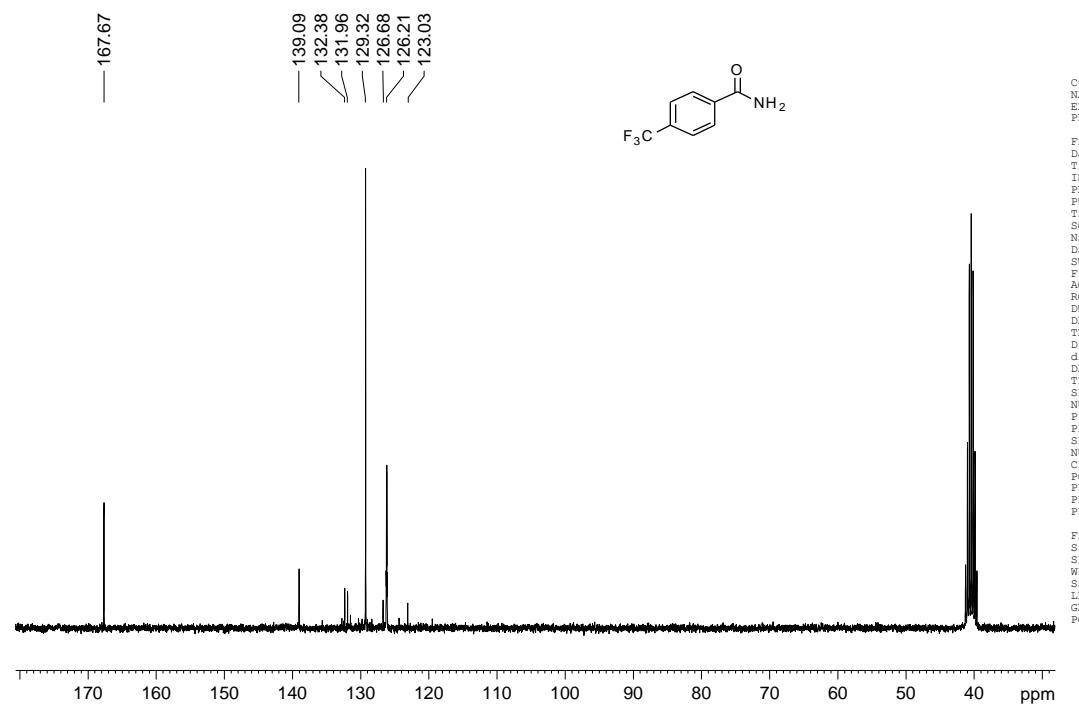




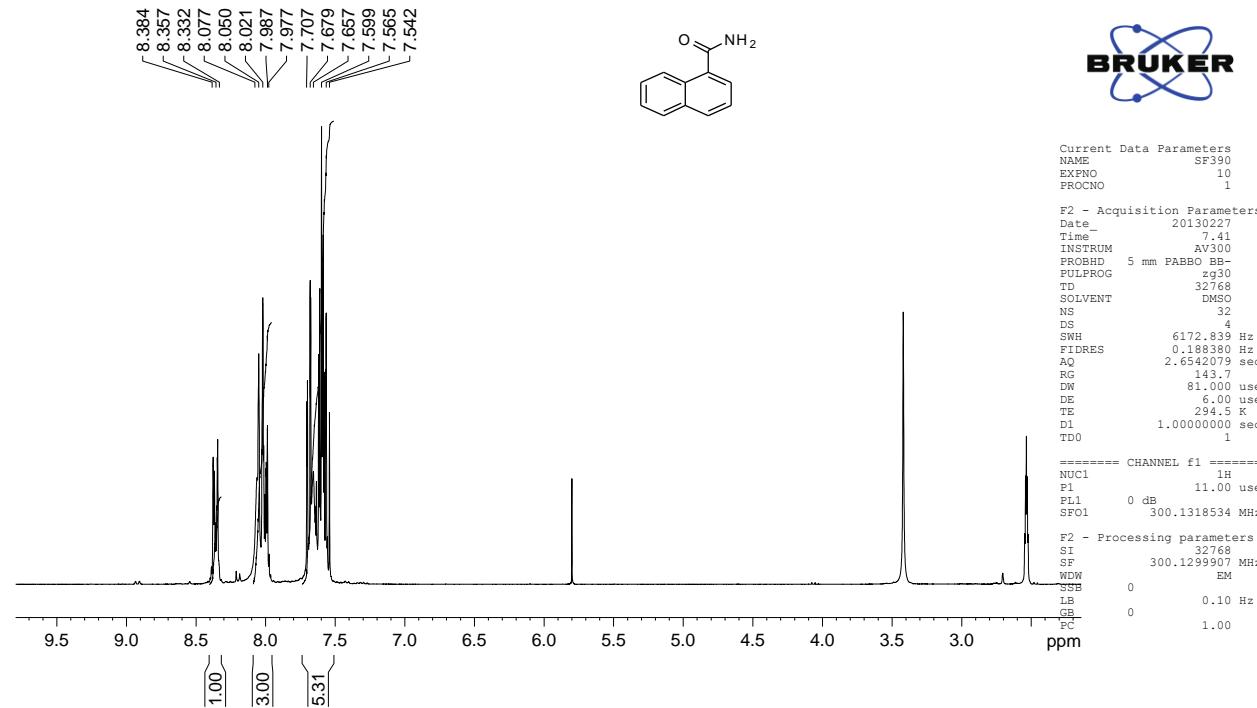
4-bromobenzamide

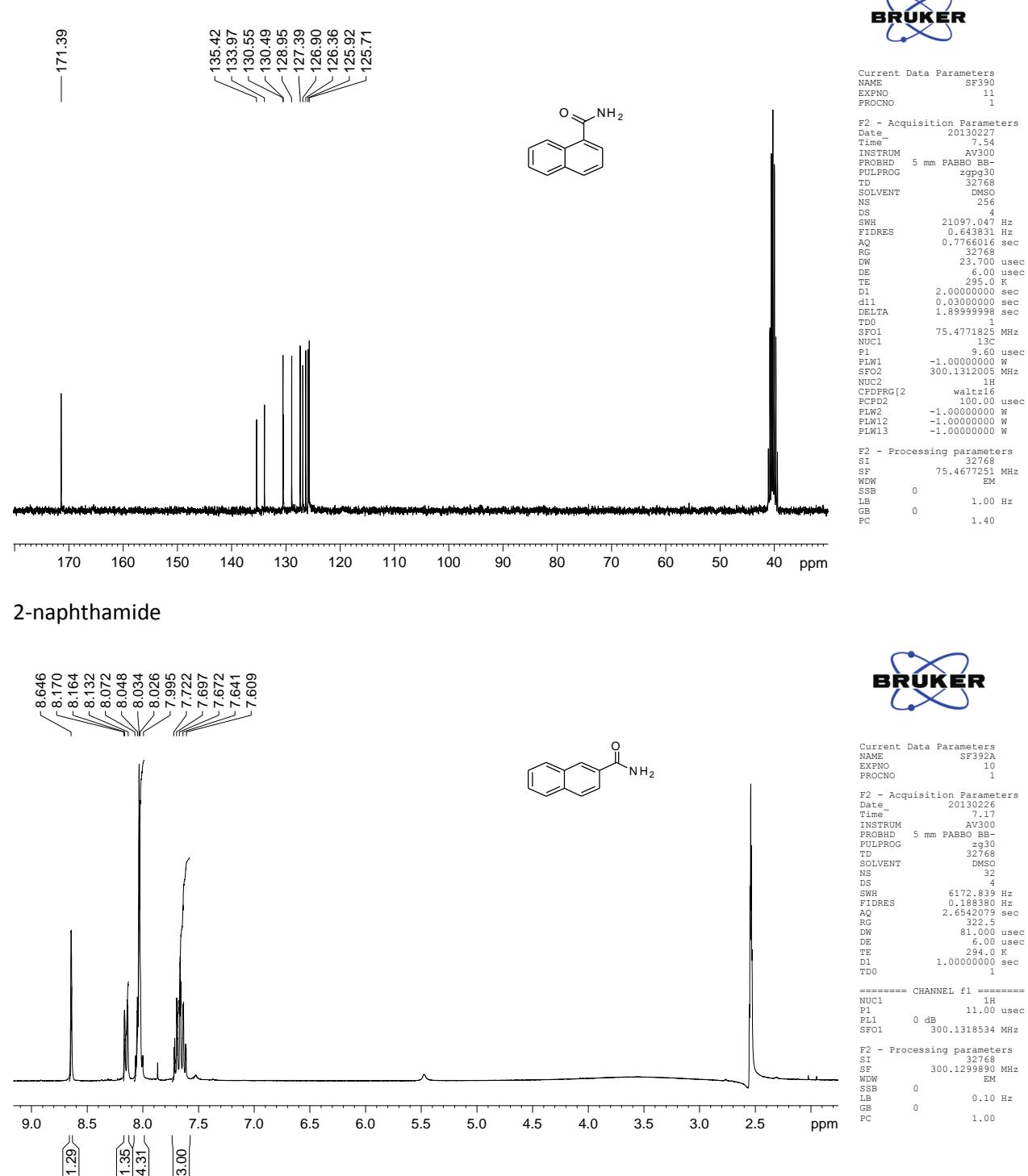


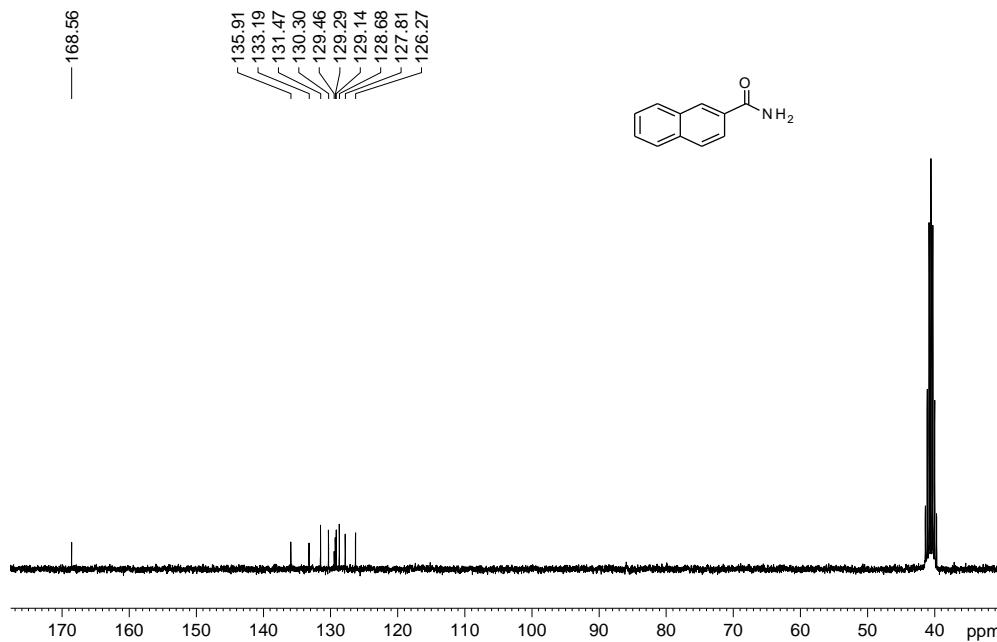




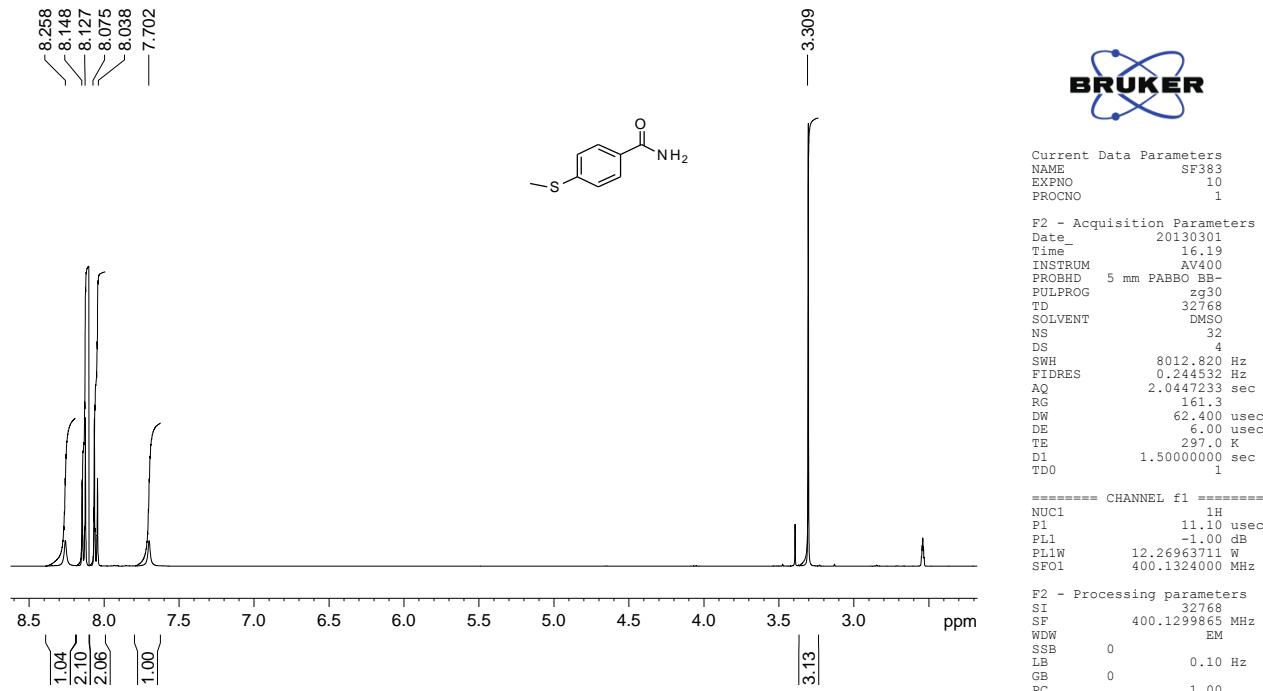
1-naphthamide

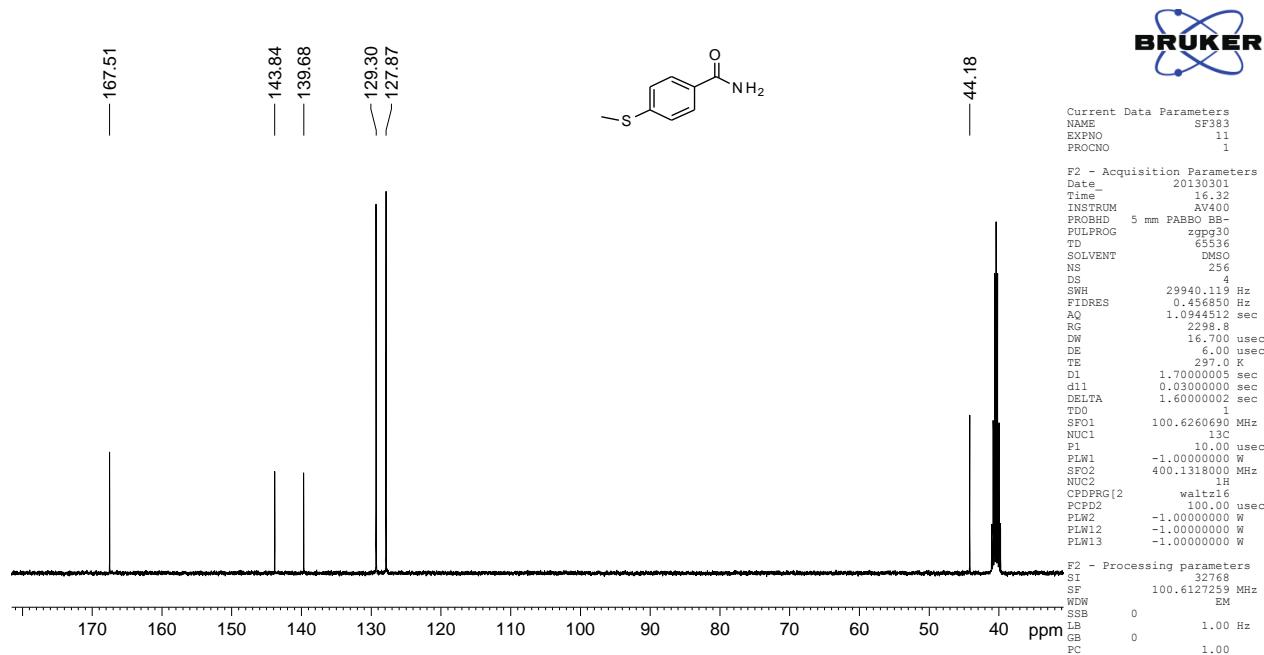




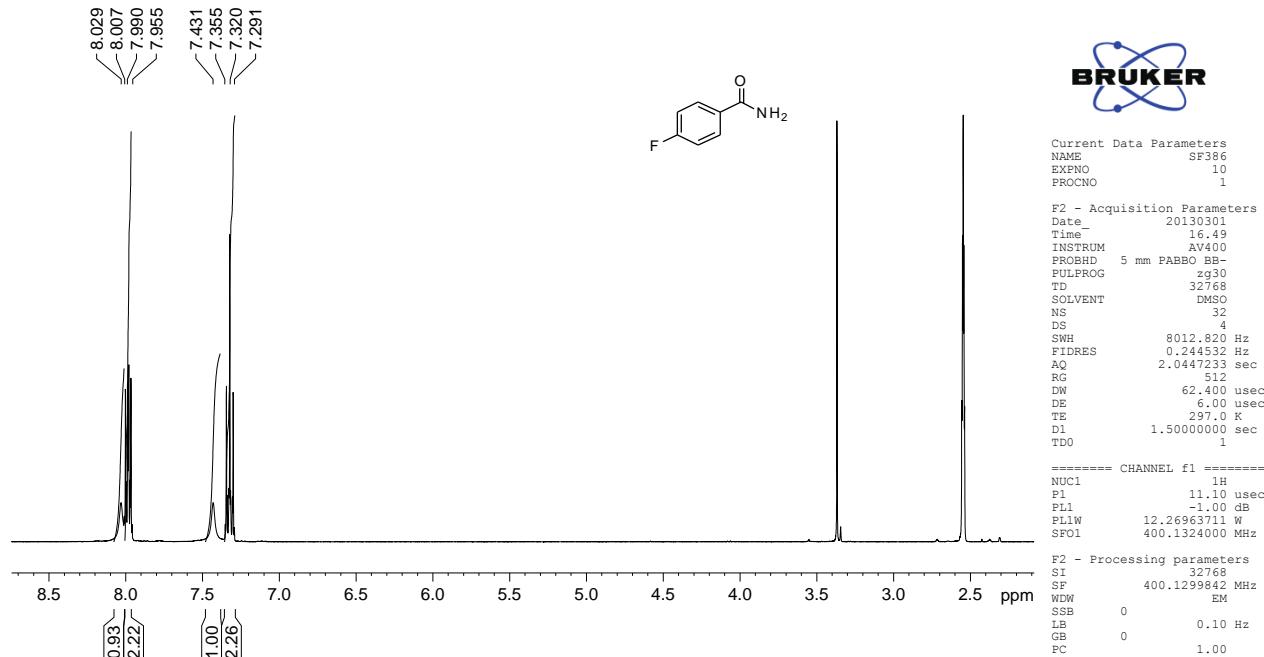


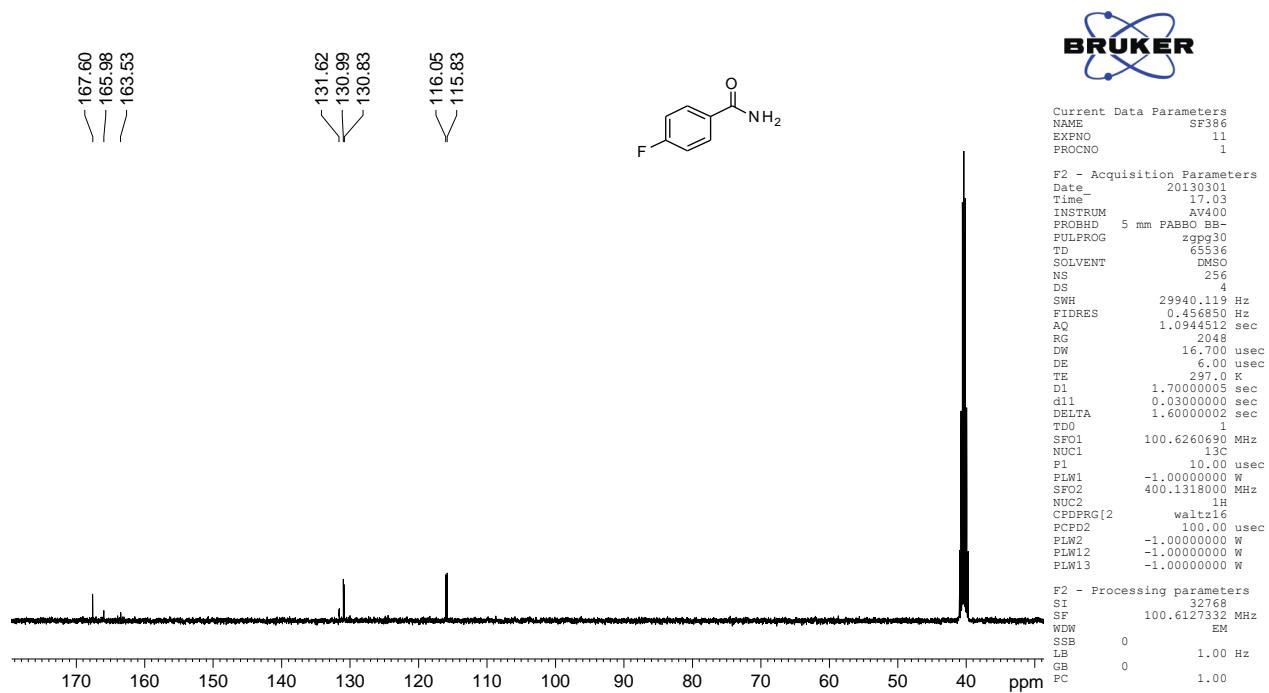
4-(methylthio)benzamide



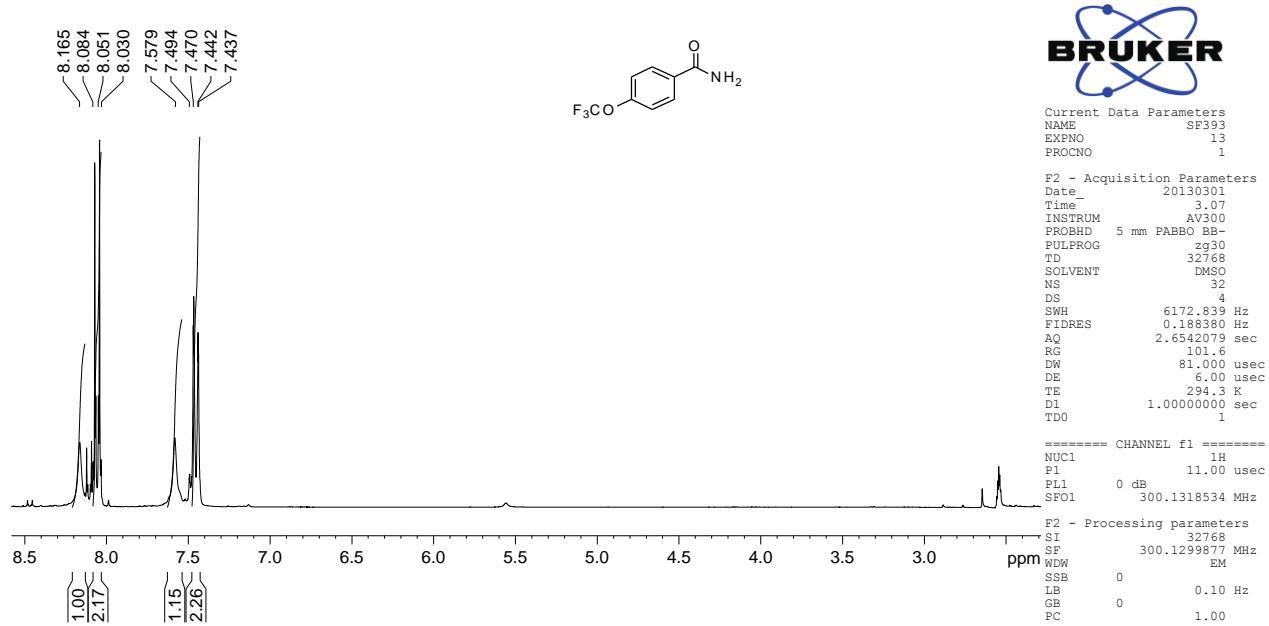


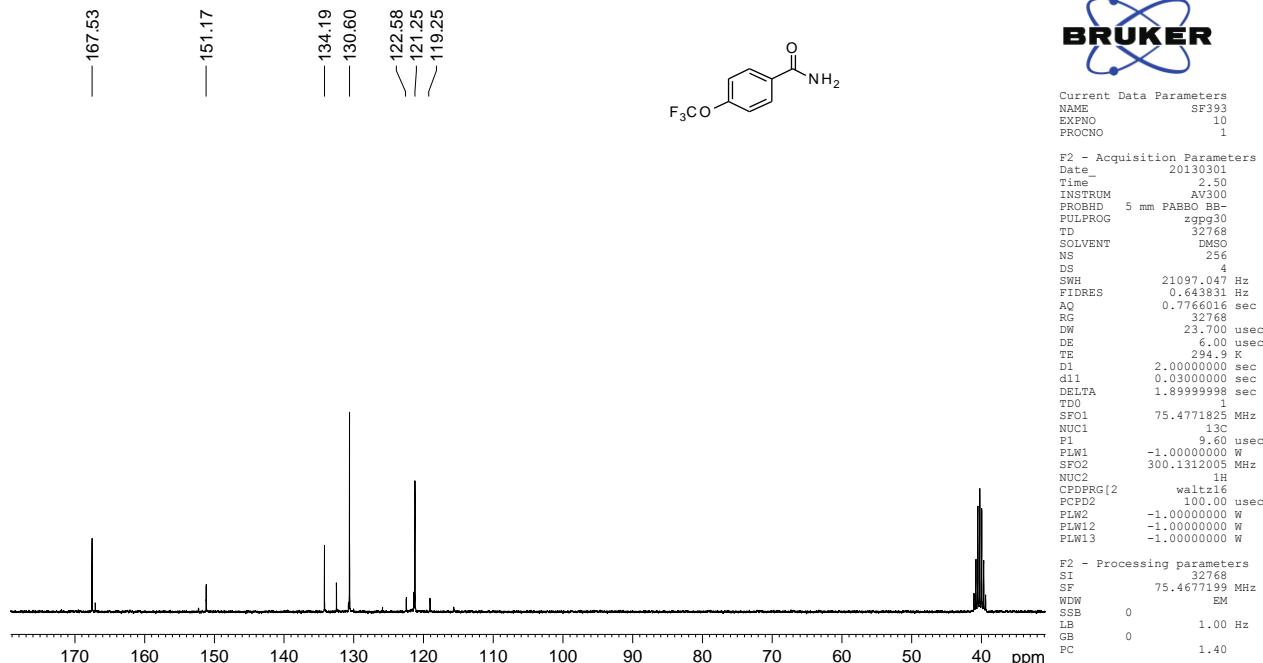
4-fluorobenzamide



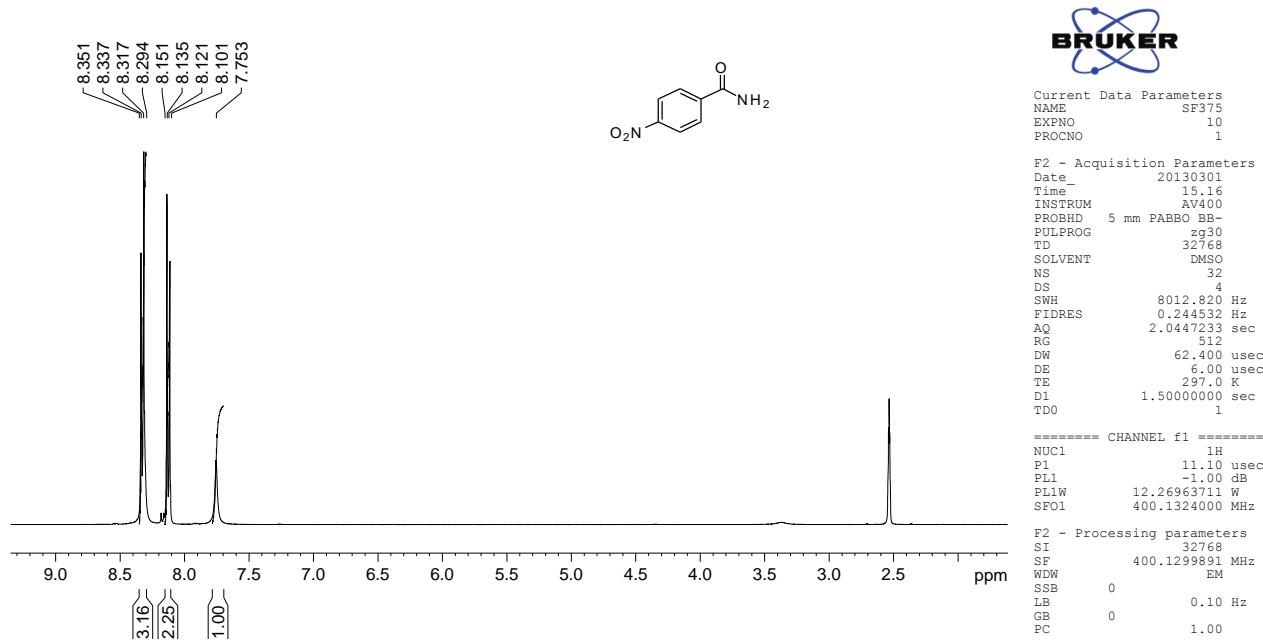


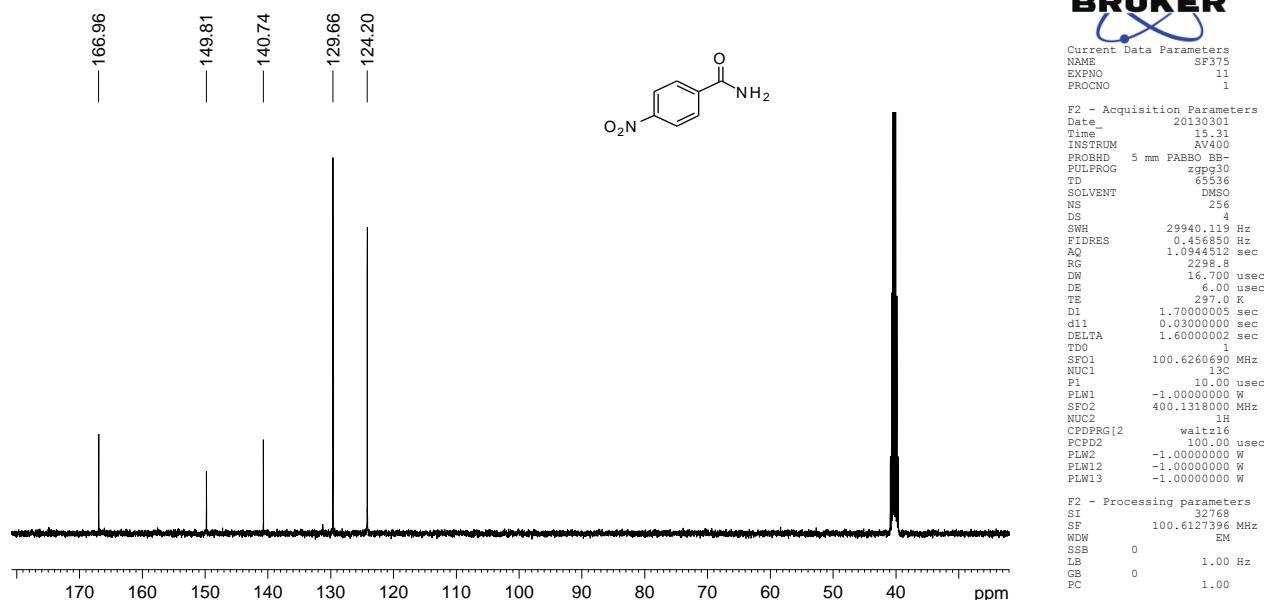
4-(trifluoromethoxy)benzamide



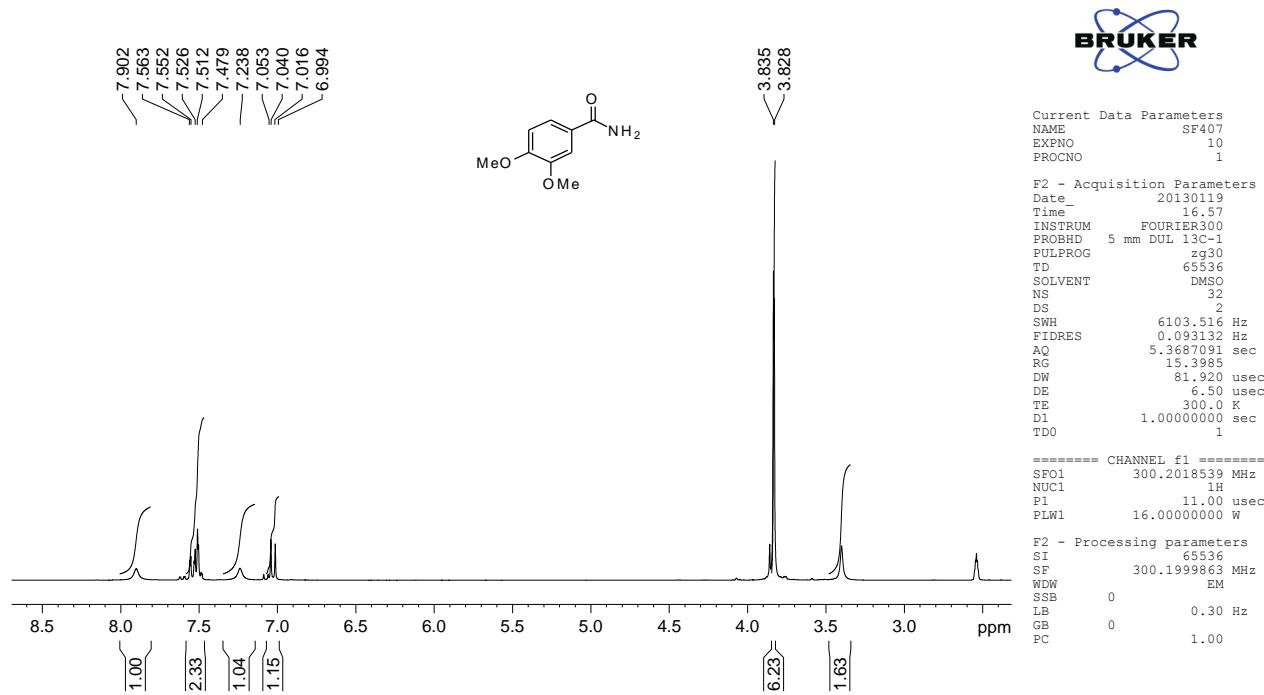


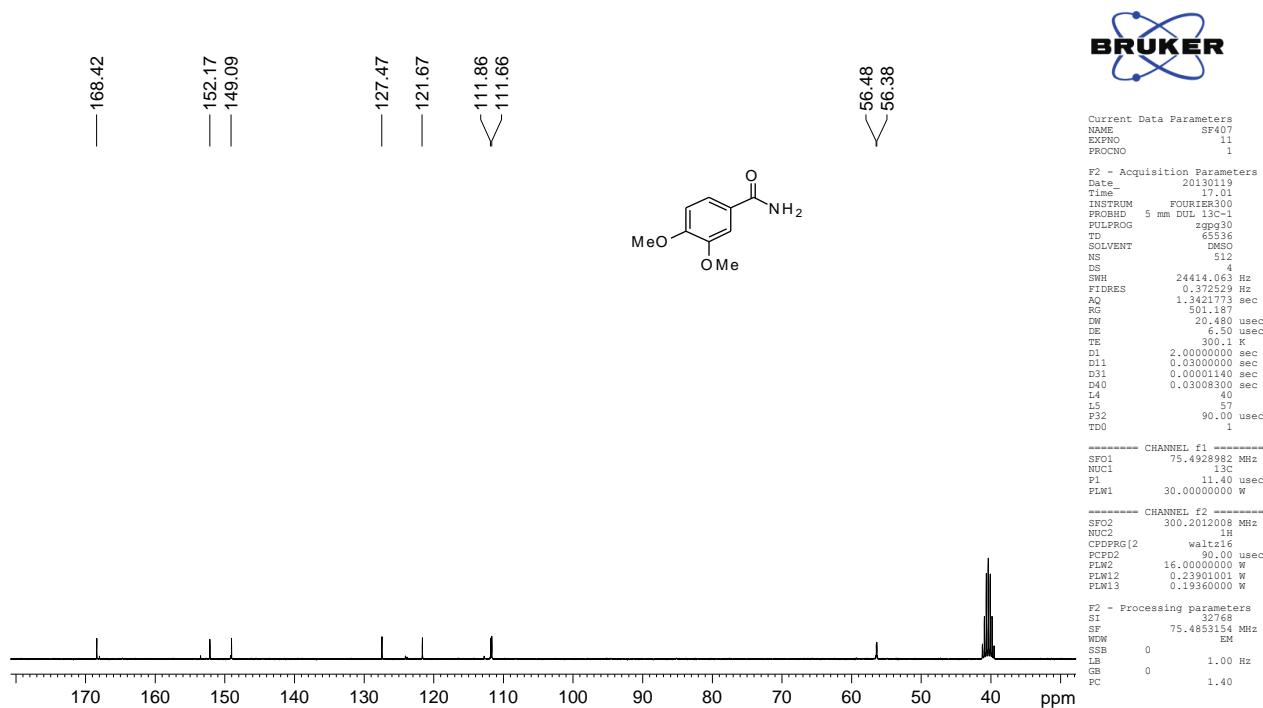
4-nitrobenzamide



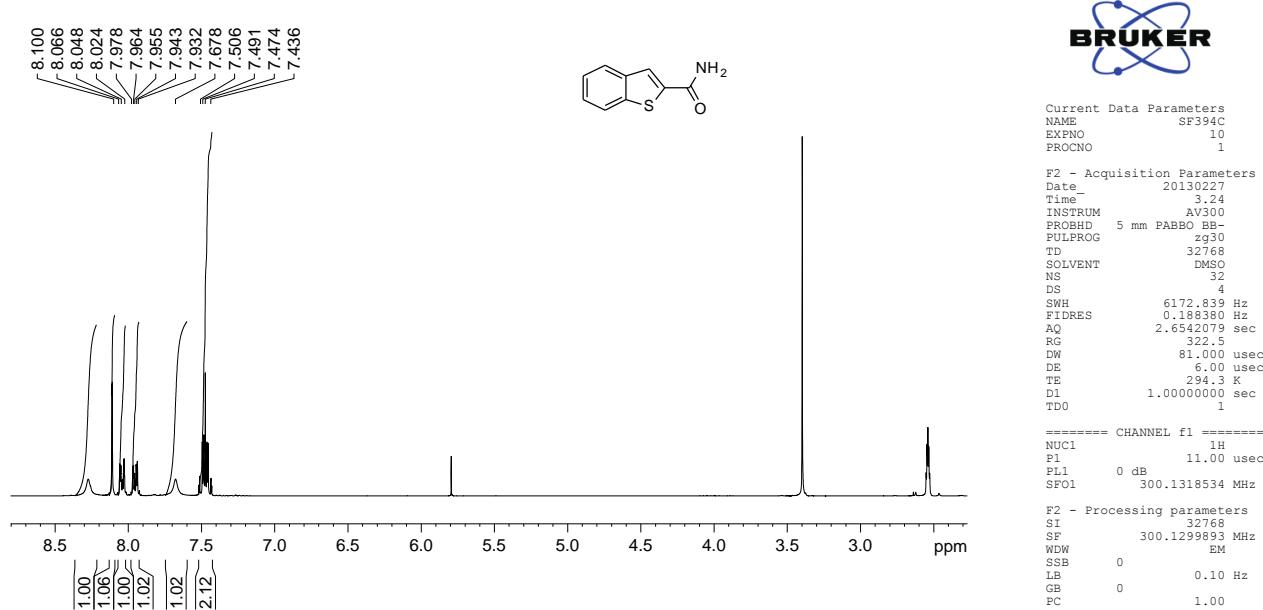


3,4-dimethoxybenzamide

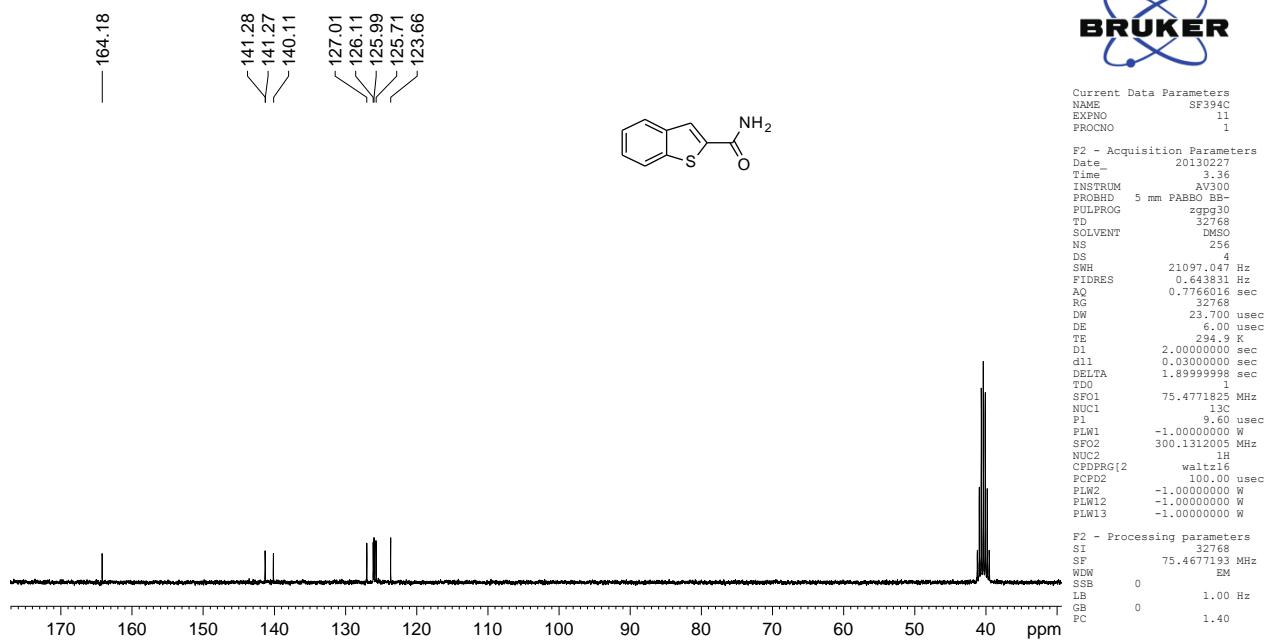




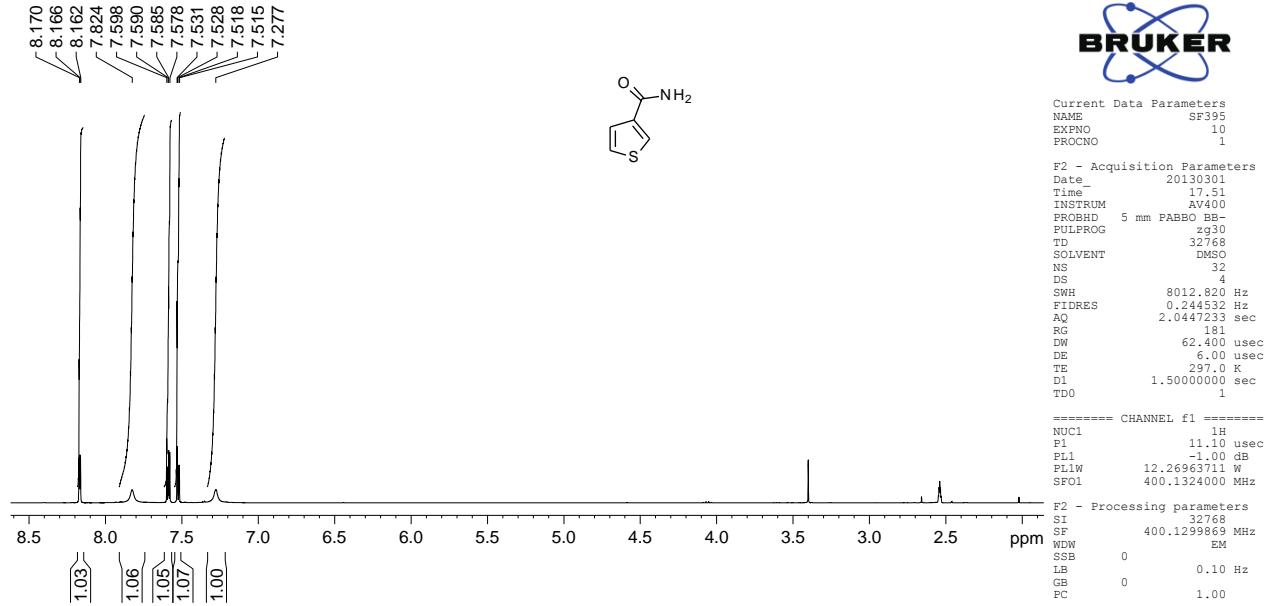
benzo[b]thiophene-2-carboxamide

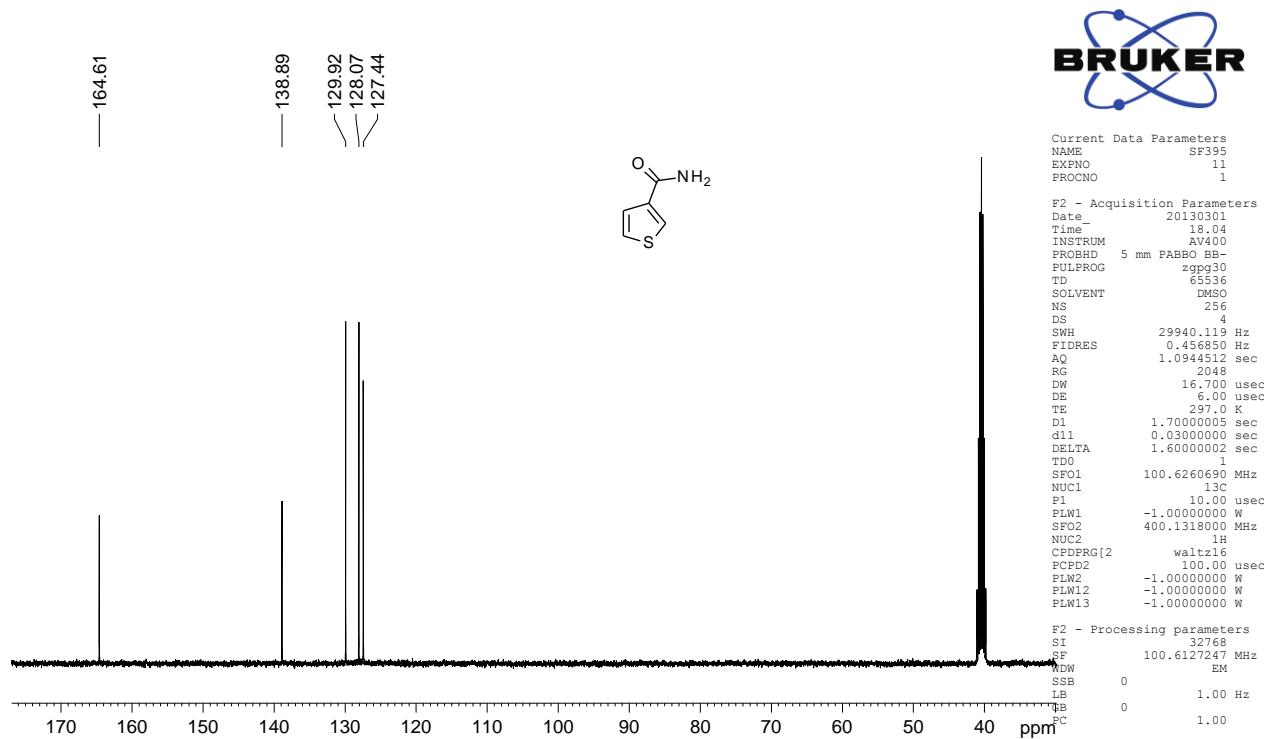


— 164.18 —

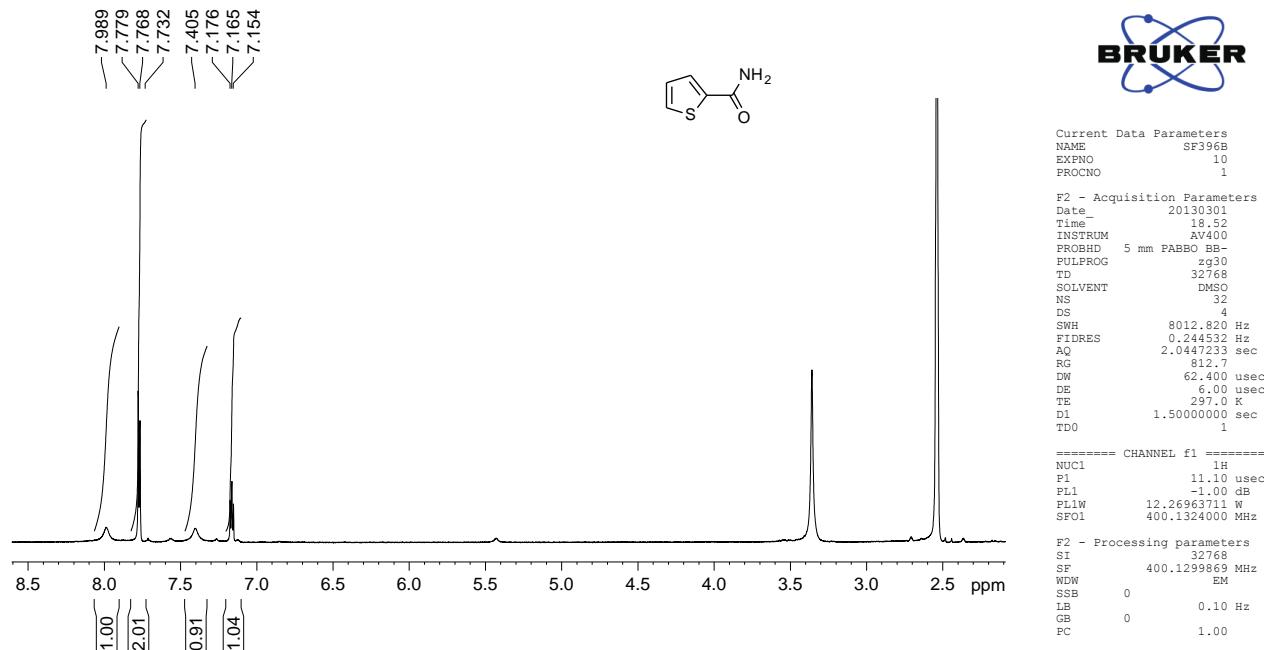


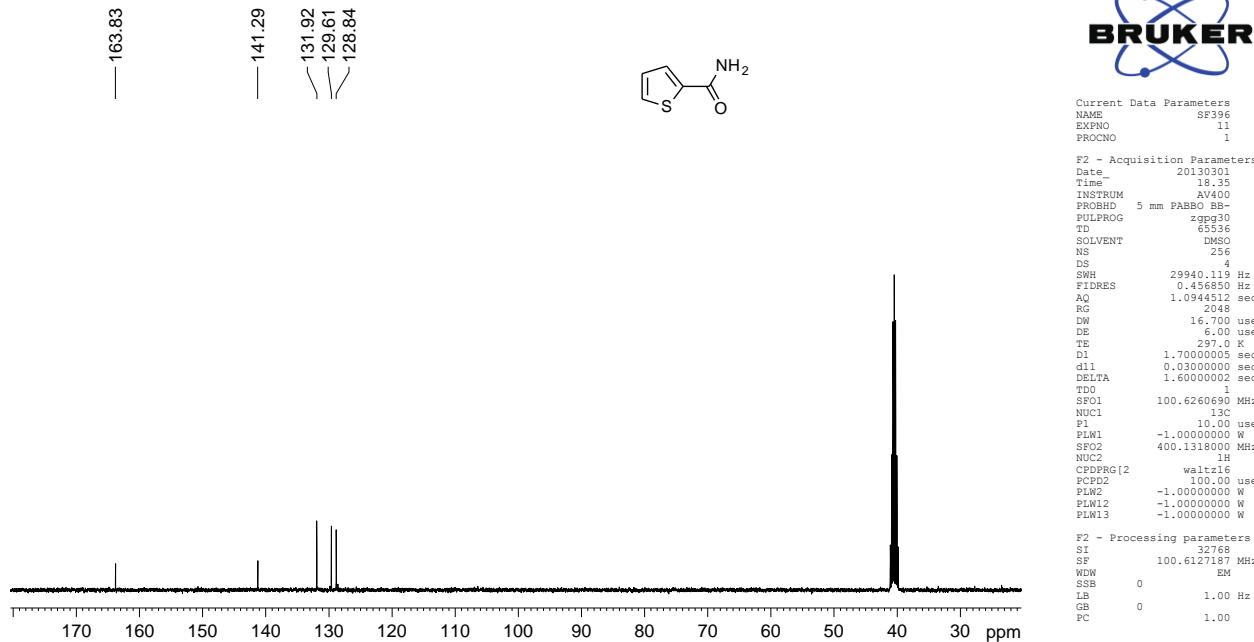
thiophene-3-carboxamide



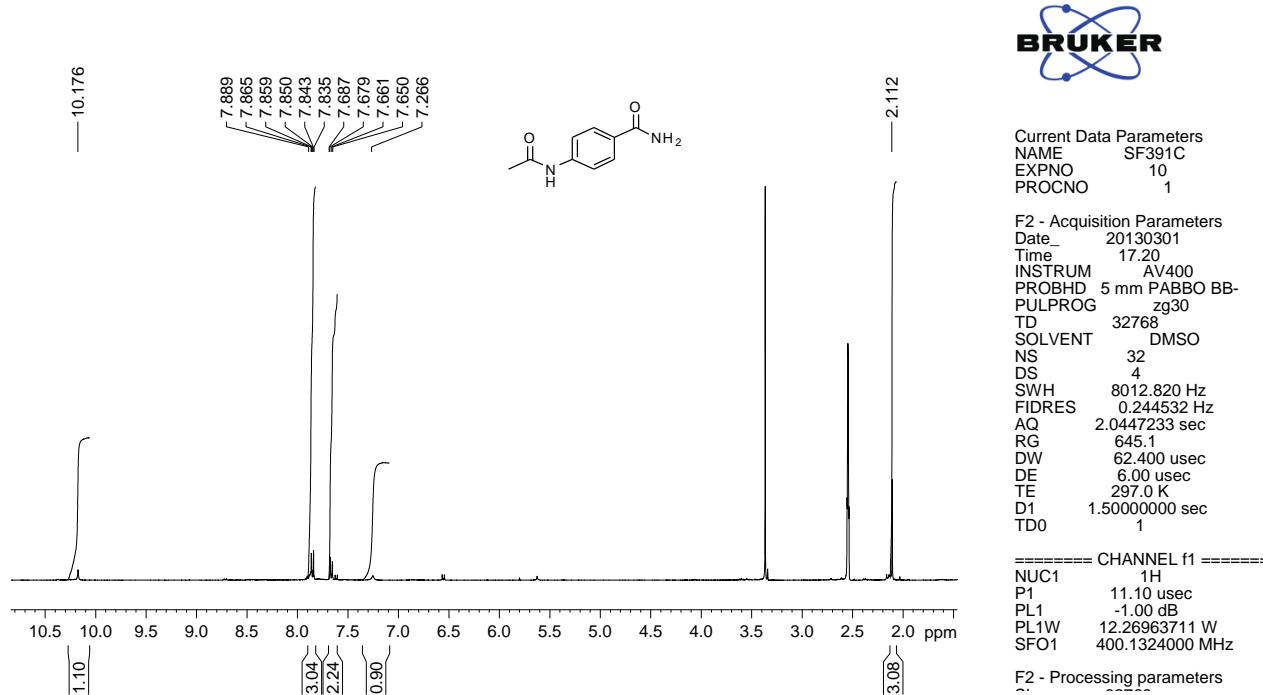


thiophene-2-carboxamide





4-acetamidobenzamide



Acetamide

