

Supporting Information for:

Palladium-Catalyzed Direct C–H Silylation and Germanylation of Benzamides and Carboxamides

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General Information. All reactions using oxygen- and moisture-sensitive materials were conducted by Schlenk techniques under an argon atmosphere or in a dry box filled with argon. Flash column chromatography was performed using Kanto Chemical silica gel (spherical, 40–63 µm). Thin layer chromatography (TLC) was performed on Merck Kieselgel 60 F₂₅₄ (0.25 mm) plates. Visualization was accomplished with UV light (254 nm) and/or an alkaline KMnO₄ solution followed by heating.

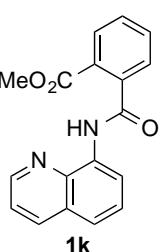
Apparatus. NMR spectra were recorded on JEOL ECX500 (500 MHz for ¹H NMR and 126 MHZ for ¹³C NMR) and JEOL ECS400 (400 MHZ for ¹H NMR and 100 MHz for ¹³C NMR) spectrometers. Proton chemical shifts are reported relative for residual solvent peaks (CDCl₃ at δ 7.26 ppm or C₆D₆ at δ 7.15 ppm). Carbon chemical shifts are reported relative to CDCl₃ at δ 77.0 ppm. ¹H NMR data are reported as follows: chemical shift, multiplicity (s = singlet, d = doublet, t = triplet, q = quartet, br = broad, m = multiplet), coupling constants (Hz), and integration. Infrared (IR) spectra were recorded on a JASCO FT/IR 410 Fourier transform infrared spectrophotometer. High resolution mass spectra (HRMS) were obtained using JEOL JMS-T100LC AccuTOF spectrometer (for ESI+) or JEOL JMS-700 MStation (for FAB+). Microwave experiments were done using InitiatorTM Exp Microwave Synthesis System (Biotage). Preparative recycling gel permeation chromatography (GPC) and preparative recycling silica gel chromatography were performed with a *JAI LC-908 chromatograph* equipped with *JAIGEL-H-P* and *2H* (chloroform as an eluent). Gas chromatography (GC) analysis was performed on a Shimadzu GCMS-QP2010 Ultra equipped with an ENV-1 column

(Kanto Chemical, 30 m x 0.25 mm, pressure = 31.7 kPa, detector = FID, 290 °C) with helium gas as a carrier.

Chemicals. Unless otherwise noted, commercially available chemicals were used without further purification or degassing. Anhydrous 1,4-dioxane was purchased from Kanto Chemical. Pd(OAc)₂ was purchased from Wako Pure Chemical Industries. Ag₂CO₃ was purchased from Sigma-Aldrich Corporation. Disilanes and hexamethyldigermane were purchased either from Sigma-Aldrich Corporation or Tokyo Chemical Industry (TCI). *N*-(8-Quinolinyl)benzamide (**1a**)¹, 2-methyl-*N*-(8-quinolinyl)benzamide (**1b**)¹, 3-methyl-*N*-(8-quinolinyl)benzamide (**1c**)¹, 4-methyl-*N*-(8-quinolinyl)benzamide (**1d**)¹, 3-(benzyloxy)-*N*-(8-quinolinyl)benzamide (**1e**)², 4-methoxy-*N*-(8-quinolinyl)benzamide (**1f**)², 3-((*tert*-butyldimethylsilyl)oxy)-*N*-(8-quinolinyl)benzamide (**1g**)², 4-fluoro-2-methyl-*N*-(8-quinolinyl)benzamide (**1h**)², 2-methyl-3-(8-quinolinylcarbamoyl)phenyl acetate (**1i**)², *N*-(8-quinolinyl)-4-(trifluoromethyl)benzamide (**1j**)³, 3-acetyl-*N*-(8-quinolinyl)benzamide (**1l**)², *N*-(8-quinolinyl)-1-naphthamide (**1n**)¹, *N*-(8-quinolinyl)thiophene-2-carboxamide (**1o**)⁴, *N*-(8-quinolinyl)cyclohexanecarboxamide (**1q**)⁵, *N*-(8-quinolinyl)cyclopentanecarboxamide (**1r**)⁵, 3-phenyl-*N*-(8-quinolinyl)propanamide (**1s**)⁵ were prepared according to the respective literature procedures.

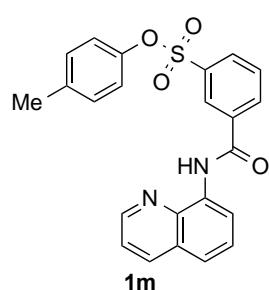
Synthesis of methyl 2-(8-quinolinylcarbamoyl)benzoate (**1k**).

8-Aminoquinoline (0.640 g, 4.40 mmol), 2-(methoxycarbonyl)benzoic acid (1.00 g, 5.60 mmol, 1.3 equiv), and 1-ethyl-3-(3-dimethylaminopropyl)carbodiimide hydrochloride (2.10 g, 11.1 mmol, 2.5 equiv) were measured into an oven-dried 100 mL round-bottomed flask. The flask was evacuated and filled with argon three times. The mixture was dissolved into CH₂Cl₂ (50 mL) and cooled to 0 °C.

 Pyridine (10 mL) was added to the solution, then the solution was warmed gradually to ambient temperature. After stirring overnight (ca. 15 h), the reaction system was quenched with 1.0 M aq. HCl, and extracted with CH₂Cl₂. The combined organic layer was washed with sat. aq. NaHCO₃ and brine, then dried over Na₂SO₄. The clear solution was filtered and evaporated in vacuo. The obtained crude amide was purified using flash column chromatography on silica gel (eluent: hexane/ethyl acetate = 4/1) to afford the desired amide (1.30 g, 4.20 mmol, 95%) as a yellow solid, R_f = 0.38 (hexane/ethyl acetate = 4/1). ¹H NMR (500 MHz, CDCl₃) δ 10.10 (s, 1H), 8.93 (d, J = 7.5 Hz, 1H), 8.76 (dd, J = 4.0, 1.7 Hz, 1H), 8.18

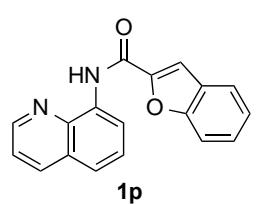
(dd, $J = 8.0, 1.7$ Hz, 1H), 7.98 (dd, $J = 8.0, 1.2$ Hz, 1H), 7.71 (dd, $J = 7.5, 1.2$ Hz, 1H), 7.68–7.54 (m, 4H), 7.45 (dd, $J = 8.6, 4.0$ Hz, 1H), 3.81 (s, 3H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 167.3, 167.0, 148.1, 138.38, 138.37, 136.2, 134.5, 132.0, 130.1, 129.9, 129.3, 127.7, 127.4, 127.3, 121.8, 121.6, 116.6, 52.5; IR (neat, ν / cm^{-1}) 3819, 3018, 1725, 1674, 1526, 1484, 1425, 1388, 1328, 1279, 1217, 771, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{18}\text{H}_{14}\text{N}_2\text{NaO}_3$ [M+Na] $^+$ 329.0902, Found 329.0906.

p-Tolyl 3-(8-quinolinylcarbamoyl)benzenesulfonate (1m). Synthesized following the



synthetic procedure of **1l**. Colorless solid, $R_f = 0.30$ (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.76 (s, 1H), 8.90 (dd, $J = 6.9, 2.3$ Hz, 1H), 8.87 (dd, $J = 4.3, 1.4$ Hz, 1H), 8.55 (s, 1H), 8.37 (d, $J = 8.0$ Hz, 1H), 8.22 (dd, $J = 8.3, 1.4$ Hz, 1H), 8.02 (d, $J = 8.0$ Hz, 1H), 7.72 (dd, $J = 8.0, 8.0$ Hz, 1H), 7.66–7.57 (m, 2H), 7.51 (dd, $J = 8.3, 4.3$ Hz, 1H), 7.09 (d, $J = 8.6$ Hz, 2H), 6.92 (d, $J = 8.6$ Hz, 2H), 2.29 (s, 3H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 163.1, 148.5, 147.3, 138.7, 137.3, 136.48, 136.45, 134.0, 132.8, 131.4, 130.3, 129.7, 128.0, 127.4, 127.2, 122.4, 121.91, 121.89, 116.9, 20.8 (one pair of the aromatic carbons is overlapping); IR (neat, ν / cm^{-1}) 3338, 3019, 1675, 1533, 1503, 1425, 1380, 1176, 752, 669; HRMS (ESI $^+$) Calcd for $\text{C}_{23}\text{H}_{18}\text{N}_2\text{NaO}_4\text{S}$ [M+Na] $^+$ 441.0885, Found 441.0884.

N-(8-Quinolinyl)benzofuran-2-carboxamide (1p). Yellow solid, $R_f = 0.46$ (hexane/ethyl



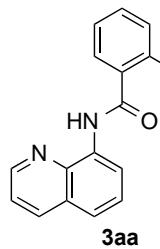
acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 11.03 (s, 1H), 8.96 (dd, $J = 4.1, 1.8$ Hz, 1H), 8.94 (dd, $J = 6.9, 2.3$ Hz, 1H), 8.21 (dd, $J = 8.1, 1.8$ Hz, 1H), 7.73 (d, $J = 7.9$ Hz, 1H), 7.71 (d, $J = 8.1$ Hz, 1H), 7.66 (s, 1H), 7.65–7.56 (m, 2H), 7.52 (dd, $J = 8.1, 4.1$ Hz, 1H), 7.48 (ddd, $J = 7.9, 7.9, 1.3$ Hz, 1H), 7.34 (dd, $J = 7.5, 7.5$ Hz, 1H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 156.9, 155.1, 149.1, 148.5, 138.7, 136.3, 134.0, 128.0, 127.7, 127.3, 127.1, 123.8, 122.7, 122.2, 121.8, 117.0, 112.2, 111.2; IR (neat, ν / cm^{-1}) 3336, 3019, 1673, 1575, 1530, 1487, 1425, 1330, 1279, 1216, 771, 669; HRMS (ESI $^+$) Calcd for $\text{C}_{18}\text{H}_{12}\text{N}_2\text{NaO}_2$ [M+Na] $^+$ 311.0796, Found 311.0809.

A General Procedure for Palladium-Catalyzed C(sp²)-H Silylation and Germanylation of Benzamides. Benzamide (**1**, 0.200 mmol), $\text{Pd}(\text{OAc})_2$ (4.5 mg, 20 μmol), Ag_2CO_3 (110 mg,

0.400 mmol), and CaSO₄ (54.4 mg, 0.400 mmol) were measured into a 15 mL oven-dried vial. The vial was evacuated and refilled with argon three times, then taken into a dry box. 1,4-Dioxane (1.0 mL) and disilane or hexamethyldigermane (1.00 mmol) were added into the vial successively. The vial was then capped, taken out from the dry box, and heated. The temperature and reaction time are specified in Tables 1, Scheme 1, and Scheme 5. The resulting mixture was filtered through a silica gel pad, concentrated in vacuo, and purified by flash column chromatography on silica gel to give the corresponding products. The yields are listed in Table 1, Scheme 1, and Scheme 5.

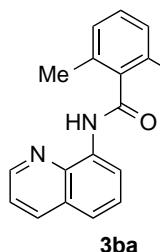
Spectroscopic Data for Silylated Benzamides

N-(8-Quinolinyl)-2-(trimethylsilyl)benzamide (3aa). Colorless solid, R_f = 0.60 (hexane/ethyl



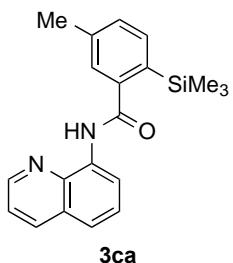
acetate = 4/1). ¹H NMR (500 MHz, CDCl₃) δ 10.36 (s, 1H), 8.98 (d, J = 8.0 Hz, 1H), 8.79 (dd, J = 4.1, 1.7 Hz, 1H), 8.17 (dd, J = 8.6, 1.7 Hz, 1H), 7.82–7.75 (m, 2H), 7.62 (dd, J = 7.7, 7.7 Hz, 1H), 7.55 (dd, J = 8.3, 1.4 Hz, 1H), 7.54–7.51 (m, 2H), 7.45 (dd, J = 8.0, 4.1 Hz, 1H), 0.41 (s, 9H); ¹³C NMR (CDCl₃, 126 MHz) δ 169.2, 148.2, 142.7, 140.0, 138.5, 136.3, 135.5, 134.7, 129.7, 128.9, 127.9, 127.4, 126.3, 121.7, 121.6, 116.4, 0.2; IR (neat, ν / cm⁻¹) 3346, 3018, 2975, 2400, 1672, 1524, 1487, 1424, 1386, 1327, 1216, 843, 769, 669; HRMS (ESI⁺) Calcd for C₁₉H₂₀N₂NaOSi [M+Na]⁺ 343.1243, Found 343.1242.

2-Methyl-N-(8-quinolinyl)-6-(trimethylsilyl)benzamide (3ba). Colorless solid, R_f = 0.52



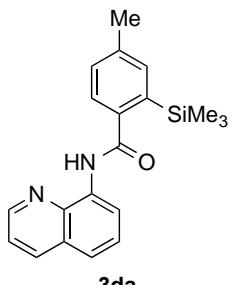
(hexane/ethyl acetate = 4/1). ¹H NMR (500 MHz, CDCl₃) δ 9.96 (s, 1H), 8.99 (d, J = 7.5 Hz, 1H), 8.73 (d, J = 2.9 Hz, 1H), 8.18 (d, J = 6.9 Hz, 1H), 7.62 (dd, J = 7.7, 7.7 Hz, 1H), 7.57 (d, J = 7.5 Hz, 1H), 7.49 (d, J = 8.4, 4.3 Hz, 1H), 7.44 (dd, J = 8.4, 4.3 Hz, 1H), 7.35 (dd, J = 7.5, 7.5 Hz, 1H), 7.29 (d, J = 7.5 Hz, 1H), 2.36 (s, 3H), 0.28 (s, 9H); ¹³C NMR (CDCl₃, 126 MHz) δ 169.9, 148.2, 143.3, 138.5, 137.3, 136.3, 134.5, 134.0, 132.2, 131.1, 128.5, 128.0, 127.5, 121.9, 121.7, 116.8, 19.6, -0.1; IR (neat, ν / cm⁻¹) 3343, 3019, 2975, 2898, 1672, 1523, 1484, 1425, 1386, 1326, 1216, 767, 669; HRMS (ESI⁺) Calcd for C₂₀H₂₂N₂NaOSi [M+Na]⁺ 357.1399, Found 357.1395.

5-Methyl-N-(8-quinolinyl)-2-(trimethylsilyl)benzamide (3ca). Colorless solid, $R_f = 0.52$



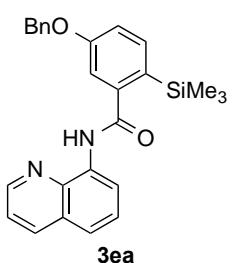
(hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.27 (s, 1H), 8.95 (d, $J = 7.2$ Hz, 1H), 8.80 (d, $J = 4.4$ Hz, 1H), 8.18 (d, $J = 8.1$ Hz, 1H), 7.66–7.51 (m, 4H), 7.46 (dd, $J = 8.1, 4.4$ Hz, 1H), 7.33 (d, $J = 7.2$ Hz, 1H), 2.45 (s, 3H), 0.35 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 169.5, 148.2, 143.0, 139.0, 138.6, 136.3, 136.2, 135.5, 134.8, 130.4, 128.0, 127.5, 127.1, 121.6, 116.5, 21.3, 0.17 (one pair of the aromatic carbons is overlapping); IR (neat, ν / cm^{-1}) 3349, 3014, 2954, 1670, 1523, 1487, 1385, 1327, 1250, 1216, 755, 667; HRMS (ESI $^+$) Calcd for $\text{C}_{20}\text{H}_{22}\text{N}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 357.1399, Found 357.1391.

4-Methyl-N-(8-quinolinyl)-2-(trimethylsilyl)benzamide (3da). Colorless solid, $R_f = 0.50$



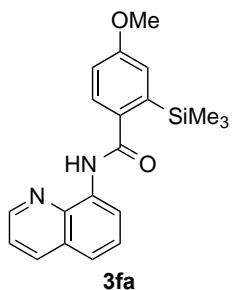
(hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.32 (s, 1H), 8.93 (dd, $J = 7.7, 1.5$ Hz, 1H), 8.79 (dd, $J = 4.2, 1.5$ Hz, 1H), 8.20 (dd, $J = 8.6, 1.5$ Hz, 1H), 7.70 (d, $J = 7.7$ Hz, 1H), 7.60 (dd, $J = 7.7, 7.7$ Hz, 1H), 7.55 (dd, $J = 8.2, 1.5$ Hz, 1H), 7.52 (s, 1H), 7.47 (dd, $J = 8.2, 4.2$ Hz, 1H), 7.31 (dd, $J = 6.9, 1.5$ Hz, 1H), 2.44 (s, 3H), 0.35 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 169.2, 148.2, 140.2, 139.9, 139.7, 138.6, 136.3, 134.9, 129.5, 128.0, 127.5, 126.5, 121.6, 121.5, 116.4, 21.6, 0.2 (one pair of the aromatic carbons is overlapping); IR (neat, ν / cm^{-1}) 3347, 3019, 2976, 1669, 1524, 1487, 1424, 1387, 1327, 1216, 1045, 771, 669; HRMS (ESI $^+$) Calcd for $\text{C}_{20}\text{H}_{22}\text{N}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 357.1399, Found 357.1394.

5-(Benzylxy)-N-(8-quinolinyl)-2-(trimethylsilyl)benzamide (3ea). Colorless oil, $R_f = 0.48$



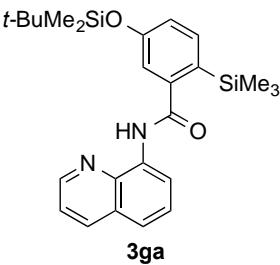
(hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.26 (s, 1H), 8.94 (d, $J = 3.7$ Hz, 1H), 8.79 (d, $J = 3.7$ Hz, 1H), 8.19 (d, $J = 9.2$ Hz, 1H), 7.63 (d, $J = 8.0$ Hz, 1H), 7.60 (d, $J = 7.2$ Hz, 1H), 7.56 (d, $J = 8.0$ Hz, 1H), 7.51–7.44 (m, 3H), 7.43–7.37 (m, 3H), 7.35 (dd, $J = 7.2, 7.2$ Hz, 1H), 7.11 (dd, $J = 8.0, 2.3$ Hz, 1H), 5.16 (s, 2H), 0.33 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 169.0, 159.3, 148.3, 144.4, 138.6, 137.0, 136.5, 136.3, 134.7, 130.9, 128.7, 128.1, 128.0, 127.6, 127.4, 121.7, 121.6, 116.6, 115.7, 113.6, 70.1, 0.3; IR (neat, ν / cm^{-1}) 3344, 3017, 2954, 1733, 1671, 1593, 1523, 1488, 1424, 1385, 1327, 1216, 1085, 755, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{26}\text{H}_{26}\text{N}_2\text{NaO}_2\text{Si} [\text{M}+\text{Na}]^+$ 449.1661, Found 449.1640.

4-Methoxy-N-(8-quinolinyl)-2-(trimethylsilyl)benzamide (3fa). Colorless solid, $R_f = 0.61$



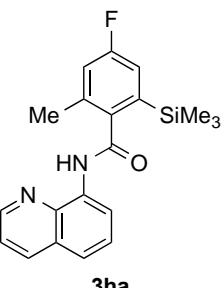
(hexane/ethyl acetate = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 10.32 (s, 1H), 8.91 (dd, $J = 7.8, 1.6$ Hz, 1H), 8.80 (dd, $J = 4.1, 1.6$ Hz, 1H), 8.18 (dd, $J = 8.5, 1.6$ Hz, 1H), 7.80 (d, $J = 8.7$ Hz, 1H), 7.60 (dd, $J = 8.0, 8.0$ Hz, 1H), 7.53 (dd, $J = 8.5, 1.6$ Hz, 1H), 7.45 (dd, $J = 8.5, 4.1$ Hz, 1H), 7.26 (s, 1H), 6.98 (dd, $J = 8.7, 2.7$ Hz, 1H), 3.89 (s, 3H), 0.36 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 168.5, 160.6, 148.2, 143.1, 138.6, 136.3, 135.0, 134.8, 128.2, 128.0, 127.5, 122.0, 121.6, 121.4, 116.3, 112.7, 55.3, 0.3; IR (neat, ν / cm^{-1}) 3353, 3019, 1669, 1525, 1479, 1424, 1386, 1328, 1215, 770, 669; HRMS (ESI $^+$) Calcd for $\text{C}_{20}\text{H}_{22}\text{N}_2\text{NaO}_2\text{Si} [\text{M}+\text{Na}]^+$ 373.1348, Found 373.1344.

5-(tert-Butyldimethylsilyloxy)-N-(8-quinolinyl)-2-(trimethylsilyl)benzamide (3ga).



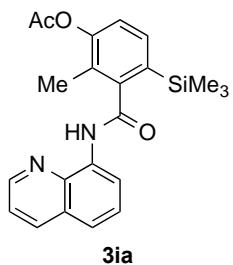
Colorless solid, $R_f = 0.60$ (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.30 (s, 1H), 8.93 (d, $J = 8.0$ Hz, 1H), 8.79 (dd, $J = 4.2, 1.6$ Hz, 1H), 8.20 (dd, $J = 8.5, 1.6$ Hz, 1H), 7.53–7.51 (m, 3H), 7.47 (dd, $J = 8.5, 4.2$ Hz, 1H), 7.26 (d, $J = 9.2$ Hz, 1H), 6.99 (dd, $J = 8.5, 2.3$ Hz, 1H), 1.03 (s, 9H), 0.34 (s, 9H), 0.29 (s, 6H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 168.7, 156.1, 147.9, 144.0, 138.4, 136.8, 136.1, 134.5, 131.2, 127.8, 127.2, 121.40, 121.38, 121.1, 118.1, 116.2, 25.4, 18.0, 0.0, -4.6; IR (neat, ν / cm^{-1}) 3346, 3015, 2955, 2931, 2859, 1672, 1590, 1524, 1473, 1298, 1252, 1216, 758, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{25}\text{H}_{34}\text{N}_2\text{NaO}_2\text{Si}_2 [\text{M}+\text{Na}]^+$ 473.2057; Found 473.2033.

4-Fluoro-2-methyl-N-(8-quinolinyl)-6-(trimethylsilyl)benzamide (3ha). Colorless solid, $R_f =$



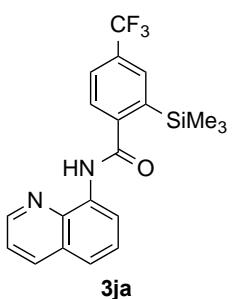
0.68 (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 9.94 (s, 1H), 8.98 (d, $J = 7.5$ Hz, 1H), 8.75 (d, $J = 4.5$ Hz, 1H), 8.19 (d, $J = 8.5$ Hz, 1H), 7.70–7.55 (m, 2H), 7.45 (dd, $J = 8.5, 4.5$ Hz, 1H), 7.16 (d, $J = 9.2$ Hz, 1H), 6.97 (d, $J = 9.2$ Hz, 1H), 2.36 (s, 3H), 0.27 (s, 9H); ^{13}C NMR (CDCl_3 , 100 MHz) δ 169.2, 162.4 (d, $J_{\text{F}} = 255$ Hz), 148.3, 141.0 (d, $J_{\text{F}} = 3.8$ Hz), 139.3 (d, $J_{\text{F}} = 2.9$ Hz), 138.4, 137.3 (d, $J_{\text{F}} = 7.7$ Hz), 136.4, 134.3, 128.0, 127.5, 122.0, 121.7, 118.5 (d, $J_{\text{F}} = 20.0$ Hz), 117.6 (d, $J_{\text{F}} = 20.9$ Hz), 116.8, 19.7, -0.3; IR (neat, ν / cm^{-1}) 3336, 3018, 1697, 1684, 1653, 1558, 1541, 1507, 1215, 757, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{20}\text{H}_{21}\text{FN}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 375.1305, Found 375.1298.

2-Methyl-3-(8-quinolinylcarbamoyl)-4-(trimethylsilyl)phenyl ethanoate (3ia). Colorless



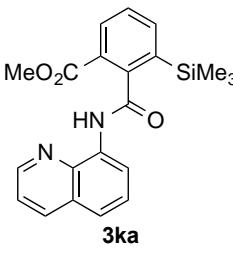
solid, $R_f = 0.34$ (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 9.99 (s, 1H), 8.96 (dd, $J = 5.7, 1.5$ Hz, 1H), 8.74 (dd, $J = 4.0, 1.5$ Hz, 1H), 8.17 (d, $J = 8.6$ Hz, 1H), 7.64–7.54 (m, 2H), 7.51 (d, $J = 8.0$ Hz, 1H), 7.45 (dd, $J = 8.0, 4.0$ Hz, 1H), 7.12 (d, $J = 8.0$ Hz, 1H), 2.35 (s, 3H), 2.27 (s, 3H), 0.26 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 169.2, 168.7, 150.1, 148.3, 145.1, 138.4, 136.2, 135.2, 134.3, 133.5, 130.0, 127.4, 126.7, 122.1, 122.0, 121.7, 116.8, 20.8, 13.0, -0.2; IR (neat, ν / cm^{-1}) 3339, 3019, 2957, 2898, 1755, 1671, 1521, 1486, 1425, 1371, 1327, 1252, 1214, 1152, 1130, 1084, 889, 841, 764, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{22}\text{H}_{24}\text{N}_2\text{NaO}_3\text{Si} [\text{M}+\text{H}]^+$ 415.1454, Found 415.1442.

N-(8-Quinolinyl)-4-(trifluoromethyl)-2-(trimethylsilyl)benzamide (3ja). Colorless solid, $R_f =$



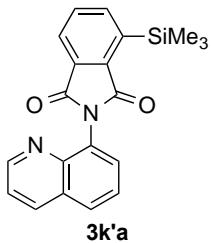
0.65 (hexane/ethyl acetate = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 10.31 (s, 1H), 8.91 (dd, $J = 7.0, 2.1$ Hz, 1H), 8.79 (dd, $J = 4.4, 1.6$ Hz, 1H), 8.21 (dd, $J = 8.5, 1.6$ Hz, 1H), 7.94 (s, 1H), 7.85 (d, $J = 8.5$ Hz, 1H), 7.75 (dd, $J = 8.5, 1.6$ Hz, 1H), 7.66–7.56 (m, 2H), 7.48 (dd, $J = 8.5, 4.4$ Hz, 1H), 0.38 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 168.1, 148.4, 145.9, 141.6, 138.5, 136.4, 134.4, 132.5 (q, $J_F = 3.6$ Hz), 131.4 (q, $J_F = 32.8$ Hz), 128.0, 127.4, 126.5, 125.9 (q, $J_F = 3.6$ Hz), 124.0 (q, $J_F = 272$ Hz), 122.2, 121.8, 116.7, -0.1; IR (neat, ν / cm^{-1}) 3340, 3019, 2958, 2899, 1676, 1524, 1489, 1425, 1387, 1325, 1216, 1131, 1079, 756, 669; HRMS (ESI $^+$) Calcd for $\text{C}_{20}\text{H}_{19}\text{F}_3\text{N}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 411.1116, Found 411.1108.

Methyl 2-(8-quinolinylcarbamoyl)-3-(trimethylsilyl)benzoate (3ka). Colorless solid, $R_f =$



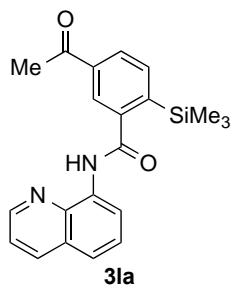
0.33 (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 9.60 (s, 1H), 8.66 (d, $J = 7.5$ Hz, 1H), 8.42 (d, $J = 4.0$ Hz, 1H), 7.87 (d, $J = 8.0$ Hz, 1H), 7.73 (d, $J = 8.1$ Hz, 1H), 7.54 (d, $J = 7.5$ Hz, 1H), 7.32 (d, $J = 8.1$ Hz, 1H), 7.27–7.19 (m, 2H), 7.13 (dd, $J = 8.0, 4.0$ Hz, 1H), 3.73 (s, 3H), 0.29 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 168.8, 166.9, 148.2, 144.1, 139.1, 138.8, 138.4, 136.3, 134.7, 130.9, 128.5, 128.1, 127.9, 127.6, 121.8, 121.6, 116.9, 52.4, -0.1; IR (neat, ν / cm^{-1}) 3343, 3018, 2954, 1719, 1677, 1524, 1486, 1386, 1327, 1253, 1216, 1113, 755, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{21}\text{H}_{22}\text{N}_2\text{NaO}_3\text{Si} [\text{M}+\text{Na}]^+$ 401.1297, Found 401.1288.

2-(8-Quinolinyl)-4-(trimethylsilyl)isoindoline-1,3-dione (3k'a). Colorless solid, $R_f = 0.37$



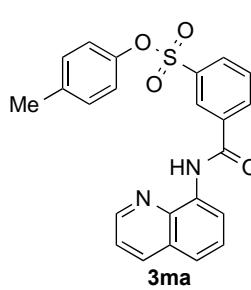
(hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 8.86 (dd, $J = 4.2, 1.8$ Hz, 1H), 8.22 (dd, $J = 8.5, 1.8$ Hz, 1H), 7.99 (d, $J = 7.6$ Hz, 1H), 7.97–7.91 (m, 2H), 7.79–7.71 (m, 2H), 7.68 (d, $J = 7.6$ Hz, 1H), 7.44 (dd, $J = 8.5, 4.2$ Hz, 1H), 0.42 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 169.1, 168.3, 150.8, 144.3, 140.4, 140.0, 136.8, 136.1, 132.8, 132.6, 130.2, 129.9, 129.4, 129.3, 126.1, 124.2, 121.9, -1.2; IR (neat, ν / cm^{-1}) 3019, 2352, 2321, 1772, 1717, 1506, 1396, 1215, 1113; HRMS (ESI $^+$) Calcd for $\text{C}_{20}\text{H}_{18}\text{N}_2\text{NaO}_2\text{Si}$ [M+Na] $^+$ 369.1035, Found 369.1032.

5-Ethanoyl-N-(8-quinolinyl)-2-(trimethylsilyl)benzamide (3la). Colorless solid, $R_f = 0.35$



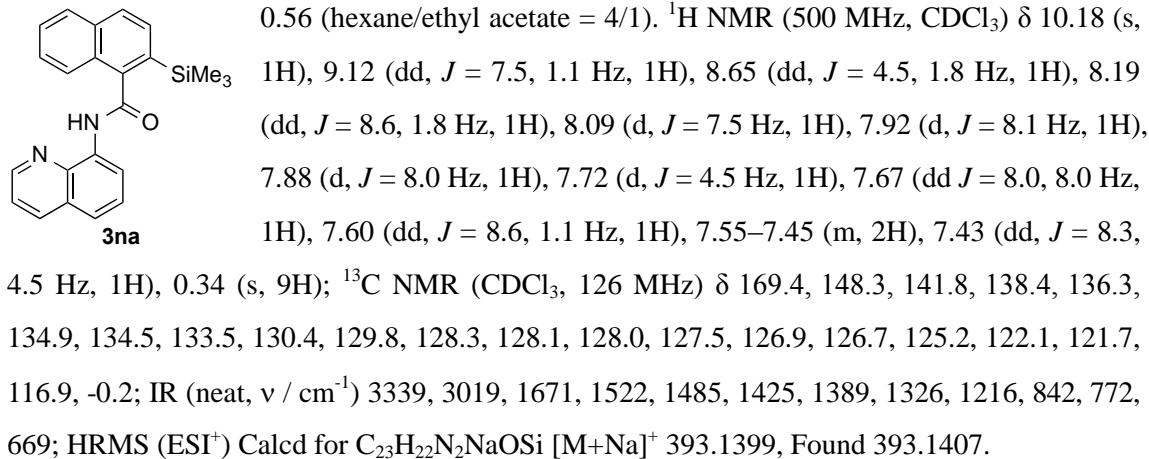
(hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.33 (s, 1H), 8.92 (d, $J = 7.5$ Hz, 1H), 8.79 (dd, $J = 4.3, 1.7$ Hz, 1H), 8.29 (d, $J = 1.7$ Hz, 1H), 8.19 (dd, $J = 8.2, 1.7$ Hz, 1H), 8.04 (dd, $J = 8.2, 1.7$ Hz, 1H), 7.83 (d, $J = 8.2$ Hz, 1H), 7.65–7.55 (m, 2H), 7.47 (dd, $J = 8.2, 4.3$ Hz, 1H), 2.66 (s, 3H), 0.37 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 197.5, 168.6, 148.4, 146.4, 143.4, 138.5, 137.2, 136.4, 135.9, 134.5, 128.9, 128.0, 127.4, 125.5, 122.0, 121.8, 116.6, 26.7, -0.1; IR (neat, ν / cm^{-1}) 3344, 3013, 2957, 2360, 1671, 1523, 1485, 1425, 1385, 1326, 1270, 1253, 1216, 1140, 757, 667; HRMS (ESI $^+$) Calcd for $\text{C}_{21}\text{H}_{22}\text{N}_2\text{NaO}_2\text{Si}$ [M+Na] $^+$ 385.1348, Found 385.1348.

p-Tolyl 3-(8-Quinolinylcarbamoyl)-4-(trimethylsilyl)benzenesulfonate (3ma). Colorless

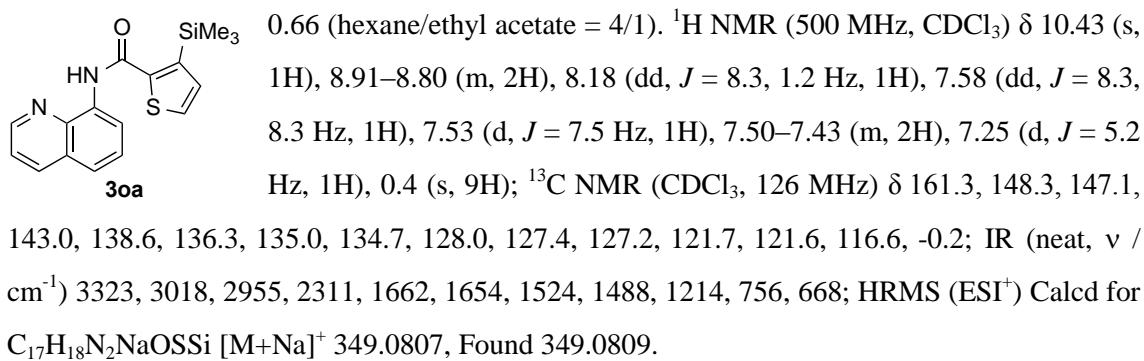


solid, $R_f = 0.48$ (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 10.30 (s, 1H), 8.87 (dd, $J = 6.3, 2.9$ Hz, 1H), 8.81 (dd, $J = 4.6, 1.7$ Hz, 1H), 8.21 (dd, $J = 8.0, 1.7$ Hz, 1H), 8.18 (d, $J = 1.7$ Hz, 1H), 7.90 (dd, $J = 8.0, 1.7$ Hz, 1H), 7.87 (d, $J = 8.1$ Hz, 1H), 7.62–7.59 (m, 2H), 7.49 (dd, $J = 8.1, 8.1$ Hz, 1H), 7.13 (d, $J = 8.1$ Hz, 2H), 7.00–6.94 (m, 2H), 2.30 (s, 3H), 0.38 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 167.2, 148.5, 147.3, 143.7, 138.5, 137.2, 136.4, 136.3, 136.2, 134.2, 130.3, 128.7, 128.0, 127.3, 125.3, 122.4, 121.9, 121.8, 116.7, 20.8, -0.1 (two pairs of the aromatic carbons are overlapping); IR (neat, ν / cm^{-1}) 3336, 3018, 2956, 1675, 1524, 1487, 1379, 1216, 1197, 756, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{26}\text{H}_{26}\text{N}_2\text{NaO}_4\text{SSi}$ [M+Na] $^+$ 513.1280, Found 513.1261.

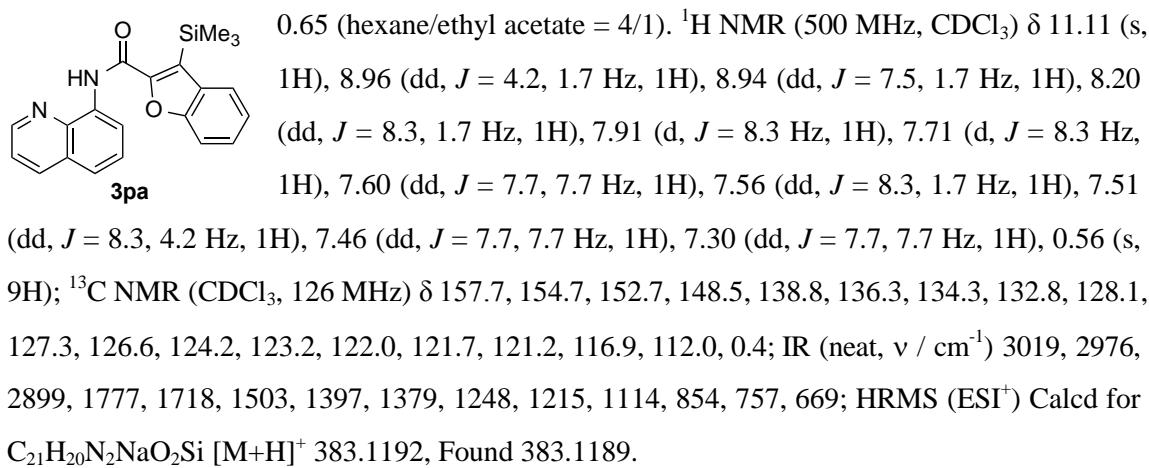
N-(8-Quinolinyl)-2-(trimethylsilyl)naphthalene-1-carboxamide (3na). Colorless solid, $R_f =$



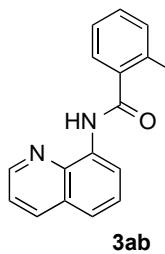
N-(8-Quinolinyl)-3-(trimethylsilyl)thiophene-2-carboxamide (3oa). Colorless solid, $R_f =$



N-(8-Quinolinyl)-3-(trimethylsilyl)benzofuran-2-carboxamide (3pa). Colorless solid, $R_f =$

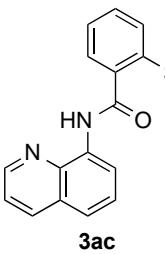


2-(Benzylidemethylsilyl)-N-(8-quinolinyl)benzamide (3ab). Colorless solid, $R_f = 0.53$



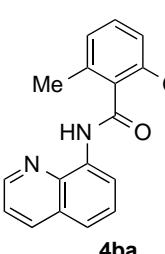
(hexane/ethyl acetate = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 10.39 (s, 1H), 8.99 (d, $J = 8.2$ Hz, 1H), 8.80 (dd, $J = 4.0, 1.6$ Hz, 1H), 8.20 (dd, $J = 8.2, 1.6$ Hz, 1H), 7.82 (d, $J = 7.5$ Hz, 1H), 7.65–7.32 (m, 6H), 7.13 (dd, $J = 7.5, 7.5$ Hz, 2H), 7.04–6.95 (m, 3H), 2.57 (s, 2H), 0.28 (s, 6H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 169.1, 148.3, 142.7, 140.4, 138.63, 138.56, 136.4, 136.1, 134.8, 129.7, 129.1, 128.5, 128.0, 127.9, 127.5, 126.3, 123.8, 121.8, 121.7, 116.5, 26.3, -2.0; IR (neat, ν / cm^{-1}) 3346, 3018, 1670, 1524, 1488, 1215, 756, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{25}\text{H}_{24}\text{N}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 419.1556, Found 419.1540.

2-(Dimethyl(phenyl)silyl)-N-(8-quinolinyl)benzamide (3ac). Colorless solid, $R_f = 0.55$



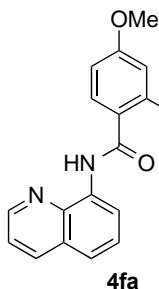
(hexane/ethyl acetate = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 10.12 (s, 1H), 8.73 (dd, $J = 6.9, 1.8$ Hz, 1H), 8.71 (dd, $J = 4.6, 1.8$ Hz, 1H), 8.15 (dd, $J = 8.6, 1.8$ Hz, 1H), 7.74 (d, $J = 7.5$ Hz, 1H), 7.61–7.39 (m, 8H), 7.20–7.10 (m, 3H), 0.64 (s, 6H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 167.0, 148.1, 143.3, 139.1, 138.6, 138.2, 136.8, 136.3, 134.6, 134.2, 129.6, 129.2, 128.5, 127.9, 127.45, 127.36, 126.3, 121.6, 121.5, 116.6, -1.2; IR (neat, ν / cm^{-1}) 3019, 2977, 1524, 1425, 1216, 1046, 929, 768, 669; HRMS (ESI $^+$) Calcd for $\text{C}_{24}\text{H}_{22}\text{N}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 405.1399, Found 405.1392.

2-Methyl-N-(8-quinolinyl)-6-(trimethylgermyl)benzamide (4ba). Colorless solid, $R_f = 0.48$



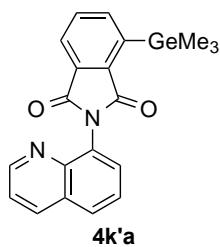
(Hexane/EtOAc = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 9.97 (s, 1H), 9.01 (d, $J = 7.4$ Hz, 1H), 8.74 (d, $J = 2.3$ Hz, 1H), 8.19 (d, $J = 6.9$ Hz, 1H), 7.63 (dd, $J = 8.0, 8.0$ Hz, 1H), 7.58 (d, $J = 7.5$ Hz, 1H), 7.47–7.40 (m, 2H), 7.35 (dd, $J = 7.5, 7.5$ Hz, 1H), 7.27 (d, $J = 4.0$ Hz, 1H), 2.48 (s, 3H), 0.39 (s, 9H); ^{13}C NMR (CDCl_3 , 125 MHz) δ 169.5, 148.2, 143.0, 139.5, 138.5, 136.3, 134.5, 134.3, 131.5, 130.6, 128.7, 128.0, 127.5, 121.9, 121.7, 116.8, 19.7, -0.5; IR (neat, ν / cm^{-1}) 3343, 3018, 2975, 2910, 1671, 1521, 1484, 1425, 1386, 1326, 1215, 1123, 1042, 827, 783, 668; HRMS (FAB $^+$) Calcd for $\text{C}_{20}\text{H}_{22}\text{GeN}_2\text{O} [\text{M}]^+$ 380.0944, Found 380.0942.

4-Methoxy-N-(8-quinoliny)-2-(trimethylgermyl)benzamide (4fa). Colorless solid, $R_f = 0.56$



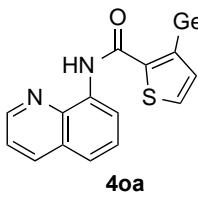
(Hexane/EtOAc = 4/1). Colorless solid. ^1H NMR (400 MHz, CDCl_3) δ 10.4 (s, 1H), 8.92 (d, $J = 7.5$ Hz, 1H), 8.81 (d, $J = 4.4$ Hz, 1H), 8.18 (d, $J = 8.6$ Hz, 1H), 7.84 (d, $J = 8.0$ Hz, 1H), 7.59 (dd, $J = 8.0, 7.5$ Hz, 1H), 7.53 (d, $J = 8.0$ Hz, 1H), 7.46 (dd, $J = 8.6, 4.4$ Hz, 1H), 7.22 (d, $J = 2.9$ Hz, 1H), 6.97 (dd, $J = 8.6, 2.9$ Hz, 1H), 3.90 (s, 3H), 0.48 (s, 9H); ^{13}C NMR (CDCl_3 , 125 MHz) δ 168.1, 160.7, 148.2, 145.6, 138.7, 136.3, 135.0, 134.0, 128.2, 128.0, 127.5, 121.6, 121.3, 121.0, 116.3, 112.6, 55.3, 0.1; IR (neat, ν / cm^{-1}) 3354, 3019, 2975, 1668, 1590, 1525, 1480, 1424, 1386, 1328, 1216, 757, 669; HRMS (FAB+) Calcd for $\text{C}_{20}\text{H}_{23}\text{GeN}_2\text{O}_2$ $[\text{M}+\text{H}]^+$ 393.0971, Found 393.0970.

2-(8-Quinolinyl)-4-(trimethylgermyl)isoindoline-1,3-dione (4k'a). Colorless solid, $R_f = 0.25$



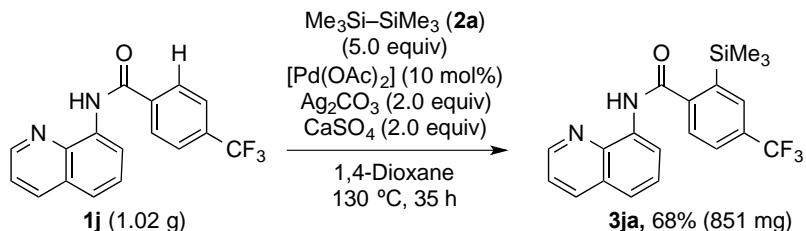
(Hexane/EtOAc = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 8.85 (d, $J = 2.9$ Hz, 1H), 8.21 (d, $J = 6.7$ Hz, 1H), 8.00–7.92 (m, 2H), 7.90 (d, $J = 7.5$ Hz, 1H), 7.77 (d, $J = 6.9$ Hz, 1H), 7.72 (dd, $J = 7.5, 7.5$ Hz, 1H), 7.67 (dd, $J = 7.7, 7.7$ Hz, 1H), 7.43 (dd, $J = 8.6, 4.1$ Hz, 1H), 0.54 (s, 9H); ^{13}C NMR (CDCl_3 , 125 MHz) δ 169.1, 168.3, 150.8, 144.3, 142.4, 139.3, 136.4, 136.1, 132.8, 132.6, 130.1, 130.0, 129.4, 129.3, 126.1, 123.8, 121.9, -1.4; IR (neat, ν / cm^{-1}) 3019, 2976, 1718, 1523, 1475, 1426, 1397, 1216, 1046, 929, 770, 669; HRMS (FAB+) Calcd for $\text{C}_{20}\text{H}_{19}\text{GeN}_2\text{O}_2$ $[\text{M}+\text{H}]^+$ 393.0658, Found 393.0661.

N-(8-Quinolyl)-3-(trimethylgermyl)thiophene-2-carboxamide (4oa). Colorless solid, $R_f =$



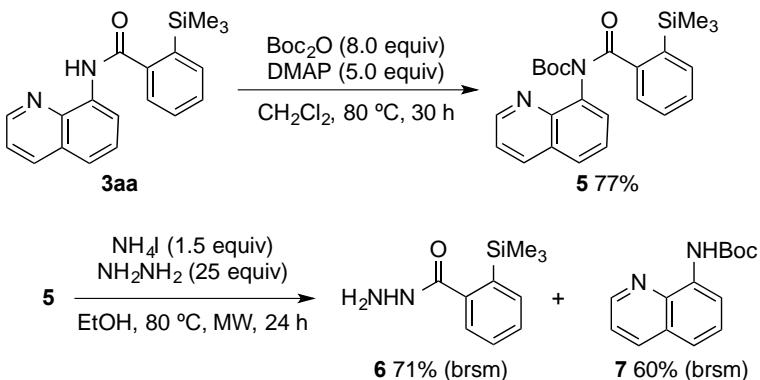
0.68 (Hexane/EtOAc = 4/1). ^1H NMR (400 MHz, CDCl_3) δ 10.42 (s, 1H), 8.86–8.83 (m, 2H), 8.17 (dd, $J = 8.6, 1.5$ Hz, 1H), 7.58 (dd, $J = 7.7, 7.7$ Hz, 1H), 7.53 (dd, $J = 8.0, 1.5$ Hz, 1H), 7.51 (d, $J = 4.9$ Hz, 1H), 7.47 (dd, $J = 8.3, 4.3$ Hz, 1H), 7.23 (d, $J = 4.9$ Hz, 1H), 0.53 (s, 9H); ^{13}C NMR (CDCl_3 , 125 MHz) δ 161.1, 148.6, 148.3, 141.2, 138.6, 136.3, 134.7, 134.0, 128.0, 127.43, 127.37, 121.7, 121.5, 116.6, -0.3; IR (neat, ν / cm^{-1}) 3322, 3019, 1661, 1524, 1484, 1424, 1329, 1215, 758, 669; HRMS (FAB+) Calcd for $\text{C}_{17}\text{H}_{19}\text{GeN}_2\text{OS}$ $[\text{M}+\text{H}]^+$ 373.0430, Found 373.0420.

Large-Scale Preparation of **3ja**.



N-(8-Quinolinyl)-4-(trifluoromethyl)benzamide (**1j**, 1.02 g, 3.23 mmol), $\text{Pd}(\text{OAc})_2$ (72.5 mg, 0.330 mmol), Ag_2CO_3 (1.78 g, 6.45 mmol), and CaSO_4 (0.878 g, 6.45 mmol) were measured into a 100 mL oven-dried vial. The vial was evacuated and refilled with argon three times, then taken into a dry box. 1,4-Dioxane (15 mL) and hexamethyldisilane (2.36 g, 16.1 mmol) were added into the vial. The vial was then capped, taken out from the dry box, and heated at 130°C for 35 h. The resulting mixture was filtered through a silica gel pad, concentrated in vacuo, and purified by flash column chromatography on silica gel (hexane/ethyl acetate = 4/1) to give the corresponding silytated product **3ja** in 68% yield (851 mg).

Removal and Recovery of Directing Group



tert-Butyl (8-quinolinyl)(2-(trimethylsilyl)benzoyl)carbamate (**5**) was prepared following a reported procedure:⁶ *N*-(8-quinolinyl)-2-(trimethylsilyl)benzamide (**3aa**, 150 mg, 0.470 mmol) was placed in a 50 mL round-bottomed flask, and the flask was flushed with nitrogen. Dichloromethane (5.0 mL) and Boc_2O (0.825 g, 3.78 mmol) were added. *N,N*-Dimethyl-4-aminopyridine (0.287 g, 2.35 mmol) was added in one portion and the resulting solution was refluxed for 30 h. The reaction mixture was concentrated in vacuo, followed by column chromatography on silica gel (hexane/ethyl acetate = 4/1) gave *tert*-butyl (8-quinolinyl)(2-(trimethylsilyl)benzoyl)carbamate (**5**, 152 mg, 0.361 mmol, 77%) as a white

solid.

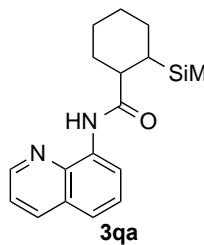
Next, following a modified procedure,⁷ compound **5** (24.1 mg, 57.0 μ mol) and NH₄I (24.1 mg, 85.5 μ mol) were placed in a 5 mL microwave-vial. Ethanol (1.0 mL) and hydrazine monohydrate (71.6 mg, 1.43 mmol) were added into the vial, then the reaction was carried out under microwave irradiation at 80 °C for 24 h. The resulting solution was then concentrated in vacuo, followed by column chromatography on silica gel (hexane/ethyl acetate = 4/1) to give 2-(trimethylsilyl)benzohydrazide (**6**, 5.1 mg, 25 μ mol) and *tert*-butyl 8-quinolinylcarbamate (**7**, 5.0 mg, 21 μ mol) in 71% and 60% yields (based on recovered starting material), respectively.

Spectral data of **6.** Colorless liquid, R_f = 0.28 (hexane/ethyl acetate = 1/1). ¹H NMR (500 MHz, CDCl₃) δ 7.65 (d, J = 7.4 Hz, 1H), 7.44 (ddd, J = 7.5, 7.5, 1.7 Hz, 1H), 7.40 (dd, J = 7.5, 1.7 Hz, 1H), 7.37 (ddd, J = 8.0, 8.0, 1.4 Hz, 1H), 7.12 (s, 1H), 4.06 (br, 2H), 0.32 (s, 9H); ¹³C NMR (CDCl₃, 126 MHz) δ 171.9, 140.0, 139.8, 135.5, 129.9, 128.8, 126.2, -0.1; IR (neat, ν / cm⁻¹) 3287, 3053, 2952, 1655, 1584, 1521, 1459, 1324, 1255, 1245, 1129, 1067, 945, 841, 742; LRMS (ESI) m/z 231.02, [M+Na]⁺; HRMS (ESI⁺) Calcd for C₁₀H₁₆N₂NaOSi [M+Na]⁺ 231.0930; Found 231.0930.

The spectral data of compound **7** matched that of the reported compound.⁸

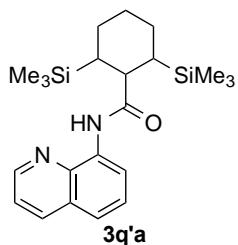
A General Procedure of Palladium-Catalyzed C(sp³)-H Silylation of Carboxyamides. Carboxyamide (**1**, 0.200 mmol), Pd(OAc)₂ (4.5 mg, 20 μ mol), and Ag₂CO₃ (110 mg, 0.400 mmol) were measured into a 15 mL oven-dried vial. The vial was evacuated and refilled with argon three times, then taken into a dry box. 1,4-Dioxane (1.0 mL), dimethylformamide (50.0 μ L, 0.640 μ mol), and hexamethyl disilane (87.8 mg, 0.600 mmol) were successively added into the vial. The vial was then capped, taken out from the dry box, and heated at 130 °C for 35 h (Scheme 4). The resulting mixture was filtered through a silica gel pad, concentrated in vacuo, and purified by flash column chromatography on silica gel to give the corresponding products. The yields are listed in Scheme 6. A mixture of **3qa** and **3q'a** (or **3ra** and **3r'a**) was further separated by preparative recycling GPC.

N-(8-Quinolinyl)-2-(trimethylsilyl)cyclohexane-1-carboxamide (3qa). Colorless liquid, $R_f =$



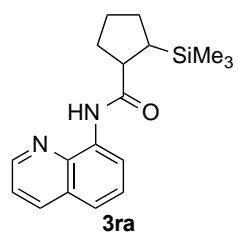
0.70 (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 9.88 (br, 1H), 8.80 (dd, $J = 4.2, 1.5$ Hz, 1H), 8.79 (dd, $J = 7.6, 1.5$ Hz, 1H), 8.15 (dd, $J = 7.6, 1.5$ Hz, 1H), 7.53 (dd, $J = 8.6, 8.6$ Hz, 1H), 7.48 (dd, $J = 8.3, 1.5$ Hz, 1H), 7.44 (dd, $J = 8.3, 4.2$ Hz, 1H), 2.88 (ddd, $J = 4.8, 4.8, 4.8$ Hz, 1H), 2.13–2.02 (m, 2H), 1.95–1.85 (m, 1H); 1.83–1.69 (m, 2H), 1.66–1.60 (m, 1H), 1.58–1.50 (m, 1H), 1.45–1.35 (m, 1H), 1.12 (ddd, $J = 8.6, 4.6, 4.6$ Hz, 1H), 0.03 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 174.5, 148.1, 138.5, 136.3, 134.8, 128.0, 127.5, 121.5, 121.0, 116.3, 44.7, 30.1, 29.0, 26.7, 24.8, 23.3, -1.4; IR (neat, ν / cm^{-1}) 3357, 2925, 2853, 1686, 1523, 1485, 1424, 1381, 1324, 1247, 1159, 842, 826, 756; HRMS (ESI $^+$) Calcd for $\text{C}_{19}\text{H}_{26}\text{N}_2\text{NaOSi} [\text{M}+\text{Na}]^+$ 349.1712; Found 349.1714; (According to NMR and GC-MS, only one product was observed. The stereochemistry was not determined).

N-(8-Quinolinyl)-2,6-bis(trimethylsilyl)cyclohexane-1-carboxamide (3q'a). Colorless liquid,



$R_f = 0.75$ (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 9.93 (s, 1H), 8.81 (dd, $J = 4.5, 1.6$ Hz, 1H), 8.76 (dd, $J = 7.7, 1.6$ Hz, 1H), 8.16 (dd, $J = 7.7, 1.6$ Hz, 1H), 7.52 (dd, $J = 7.7, 7.7$ Hz, 1H), 7.48 (dd, $J = 8.3, 1.6$ Hz, 1H), 7.45 (dd, $J = 8.3, 4.5$ Hz, 1H), 2.82 (t, $J = 4.0$ Hz, 1H), 2.19 (dddd, $J = 13.2, 13.2, 13.2, 4.0$ Hz, 2H), 2.01 (ddd, $J = 13.2, 3.6, 3.6$ Hz, 1H), 1.65–1.56 (m, 2H), 1.32 (dtt, $J = 13.2, 13.2, 3.6$ Hz, 1H), 0.92 (ddd, $J = 13.2, 3.6, 3.6$ Hz, 2H), -0.04 (s, 18H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 174.1, 158.6, 150.0, 136.4, 134.7, 134.0, 128.0, 126.7, 121.5, 121.1, 52.5, 36.2, 31.0, 27.2, -1.8; IR (neat, ν / cm^{-1}) 3358, 3019, 2922, 1681, 1521, 1483, 1424, 1250, 1215, 756, 668; HRMS (ESI $^+$) Calcd for $\text{C}_{22}\text{H}_{34}\text{N}_2\text{NaOSi}_2 [\text{M}+\text{Na}]^+$ 421.2107; Found 421.2094; (According to NMR and GC-MS, only one product was observed. The stereochemistry was not determined).

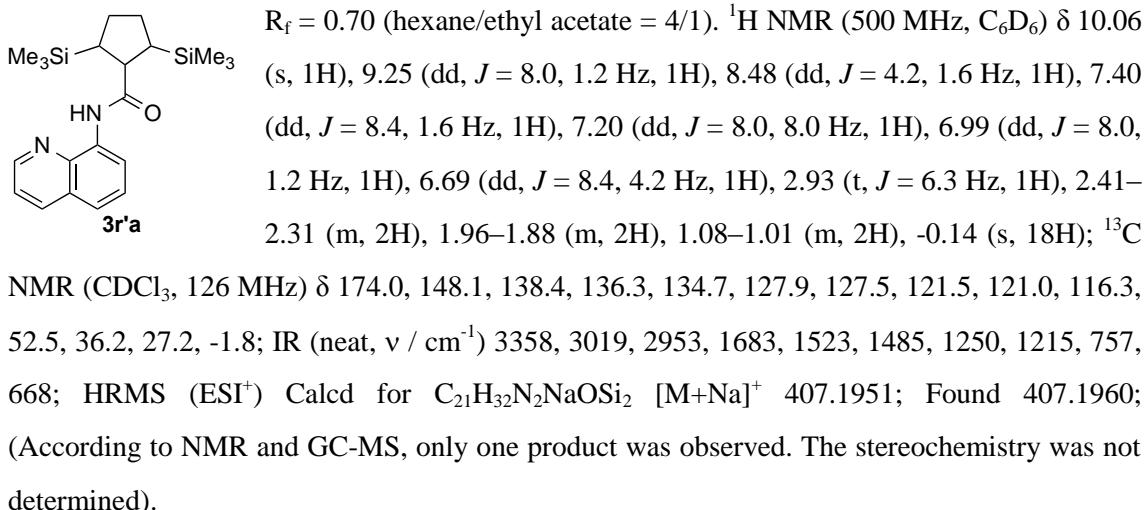
N-(8-Quinolinyl)-2-(trimethylsilyl)cyclopentane-1-carboxamide (3ra). Colorless liquid, $R_f =$



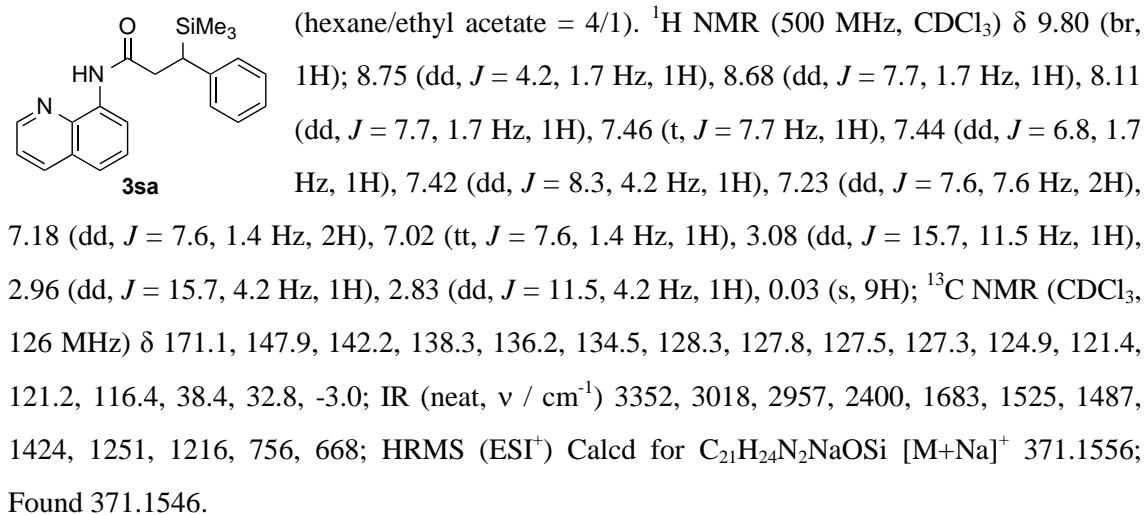
0.65 (hexane/ethyl acetate = 4/1). ^1H NMR (500 MHz, CDCl_3) δ 9.86 (br, 1H), 8.80 (dd, $J = 4.3, 1.5$ Hz, 1H), 8.78 (dd, $J = 7.7, 1.5$ Hz, 1H), 8.16 (dd, $J = 8.3, 1.5$ Hz, 1H), 7.53 (dd, $J = 7.7, 7.7$ Hz, 1H), 7.48 (dd, $J = 8.3, 1.5$ Hz, 1H), 7.45 (dd, $J = 8.3, 4.3$ Hz, 1H), 3.14 (ddd, $J = 7.8, 7.8, 4.3$ Hz, 1H), 2.14–2.02 (m, 2H), 2.00–1.80 (m, 3H), 1.72–1.64 (m, 1H), 1.29 (ddd, $J = 11.5, 11.5, 8.1$ Hz, 1H), 0.01 (s, 9H); ^{13}C NMR (CDCl_3 , 126 MHz) δ 175.2, 148.0, 138.4,

136.3, 134.7, 127.9, 127.5, 121.5, 121.1, 116.3, 49.5, 33.3, 31.6, 28.0, 26.2, -1.7; IR (neat, ν / cm^{-1}) 3357, 3012, 2951, 2864, 1683, 1523, 1485, 1424, 1388, 1324, 1249, 1162, 837, 755, 666; HRMS (ESI $^+$) Calcd for $\text{C}_{18}\text{H}_{24}\text{N}_2\text{NaOSi}$ [M+Na] $^+$ 335.1556; Found 335.1561; (According to NMR and GC-MS, only one product was observed. The stereochemistry was not determined).

N-(8-Quinolinyl)-2,5-bis(trimethylsilyl)cyclopentane-1-carboxamide (3r'a). Colorless liquid,

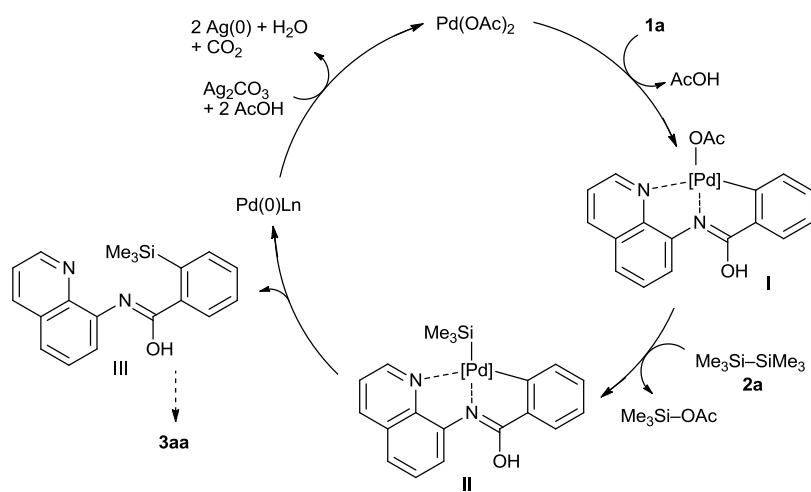


3-Phenyl-N-(8-quinolinyl)-3-(trimethylsilyl)propanamide (3sa). Colorless liquid, $R_f = 0.53$



An alternative mechanism for the silylation reaction is shown below: Activation of *ortho*-C–H bond of a benzamide by Pd(OAc)₂ may form intermediate **I**, followed by transmetallation between intermediate **I** and disilane **2a** to form Ar–Pd–SiMe₃ complex **II** (Scheme S1). Subsequent C–Si reductive elimination would afford Pd(0) species and intermediate **III**, which tautomerizes to the desired product **3aa**. Oxidation of the Pd(0) species by AgCO₃ then regenerates Pd(OAc)₂.

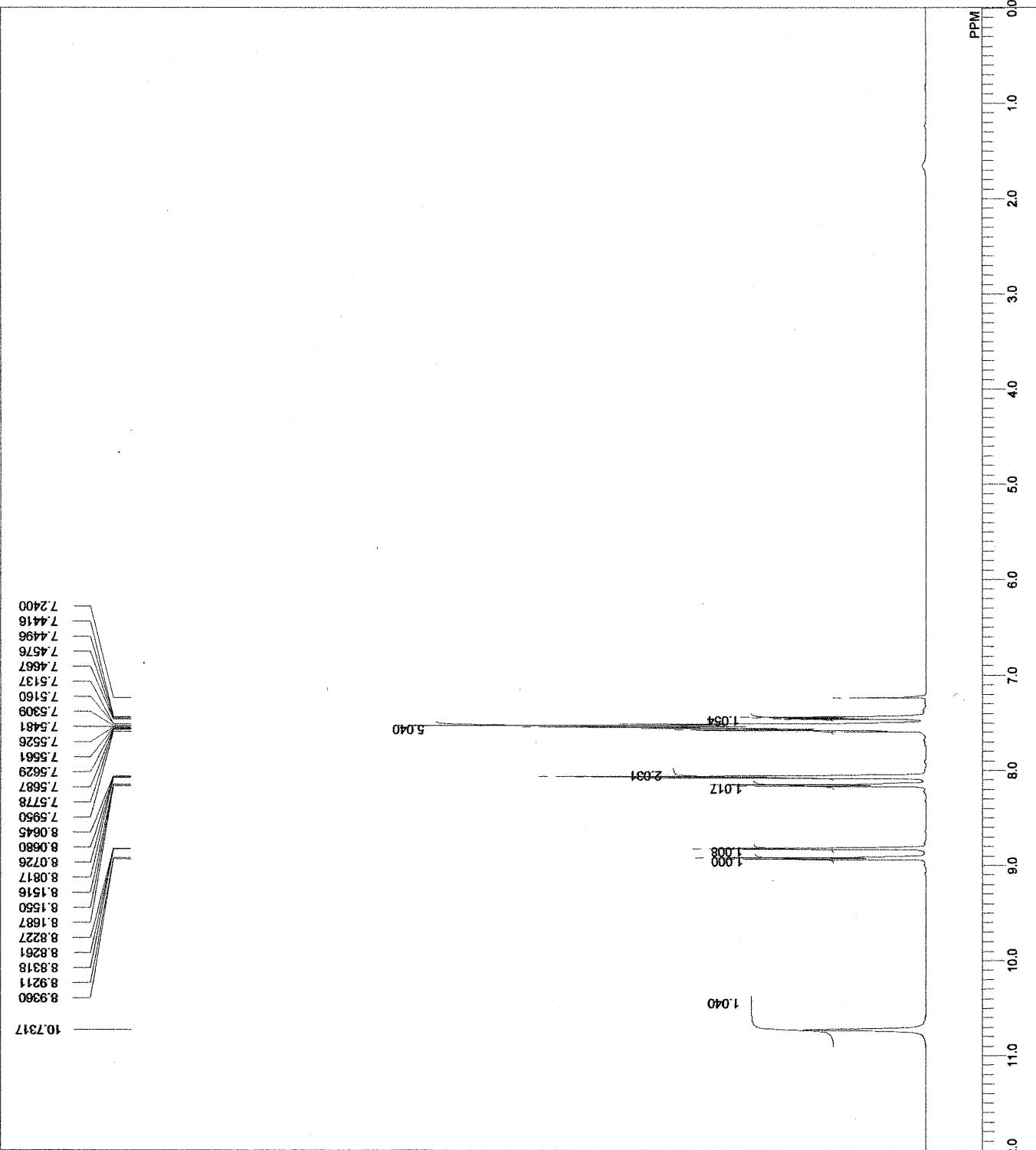
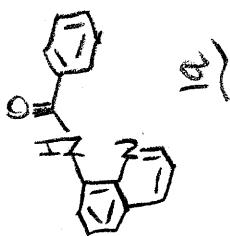
Scheme S1. An Alternative Mechanism for the Silylation Reaction

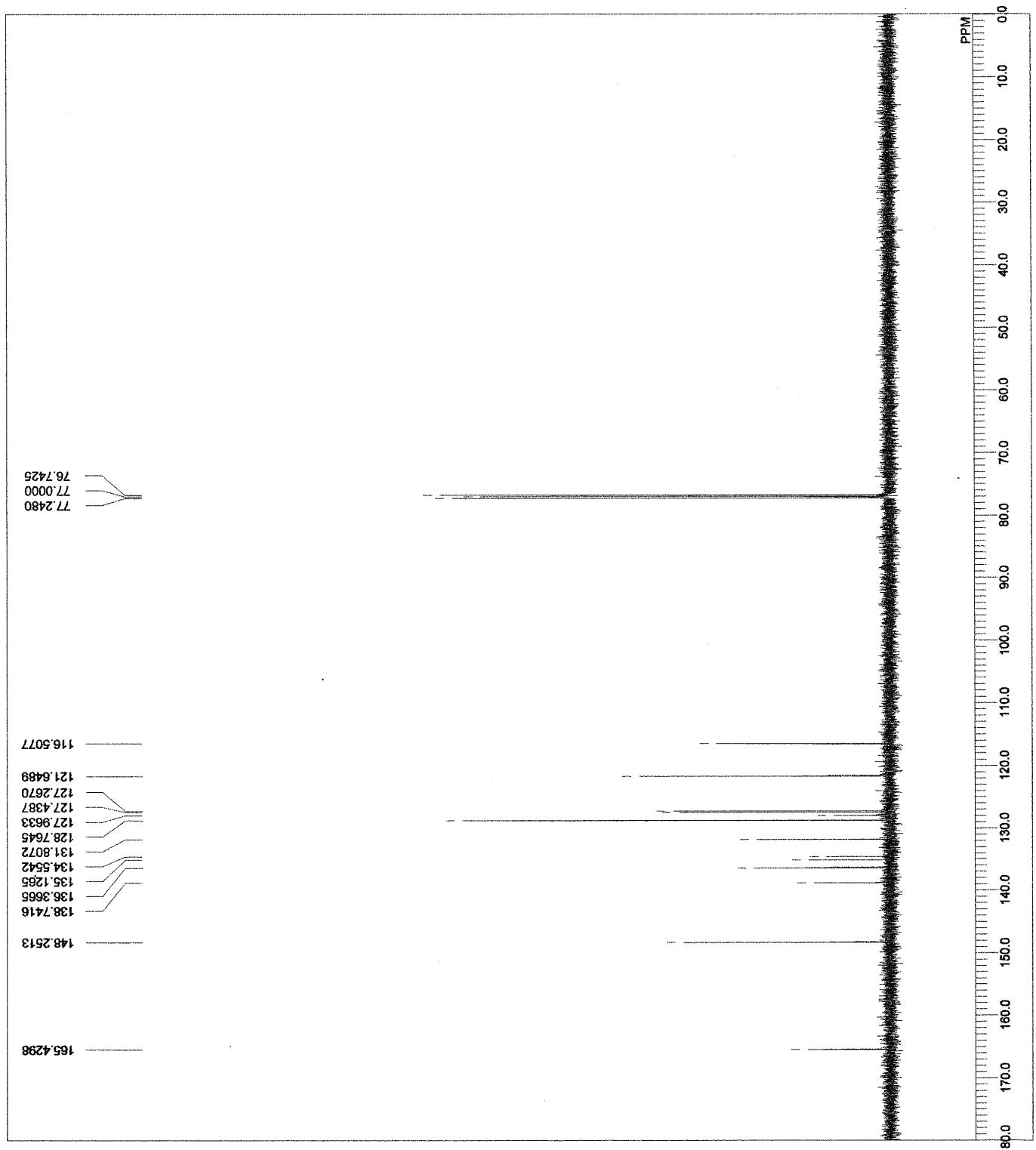


References

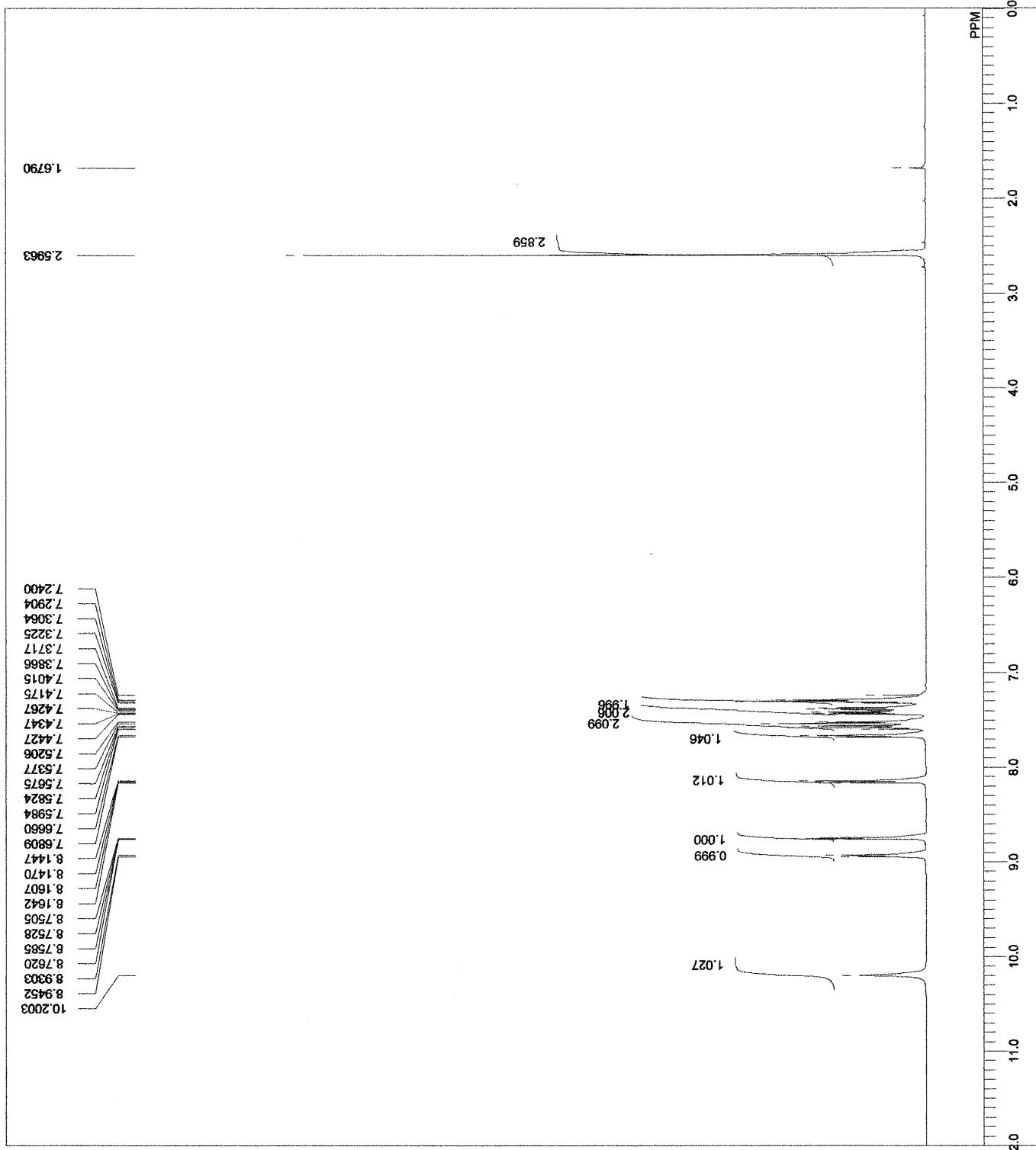
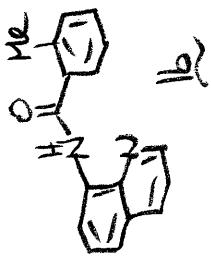
1. Gou, F.-R.; Wang, X.-C.; Huo, P.-F.; Bi, H.-P.; Guan, Z.-H.; Liang, Y.-M. *Org. Lett.* **2009**, *11*, 5726.
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4. Aihara, Y.; Chatani, N. *Chem. Sci.* **2013**, *4*, 664.
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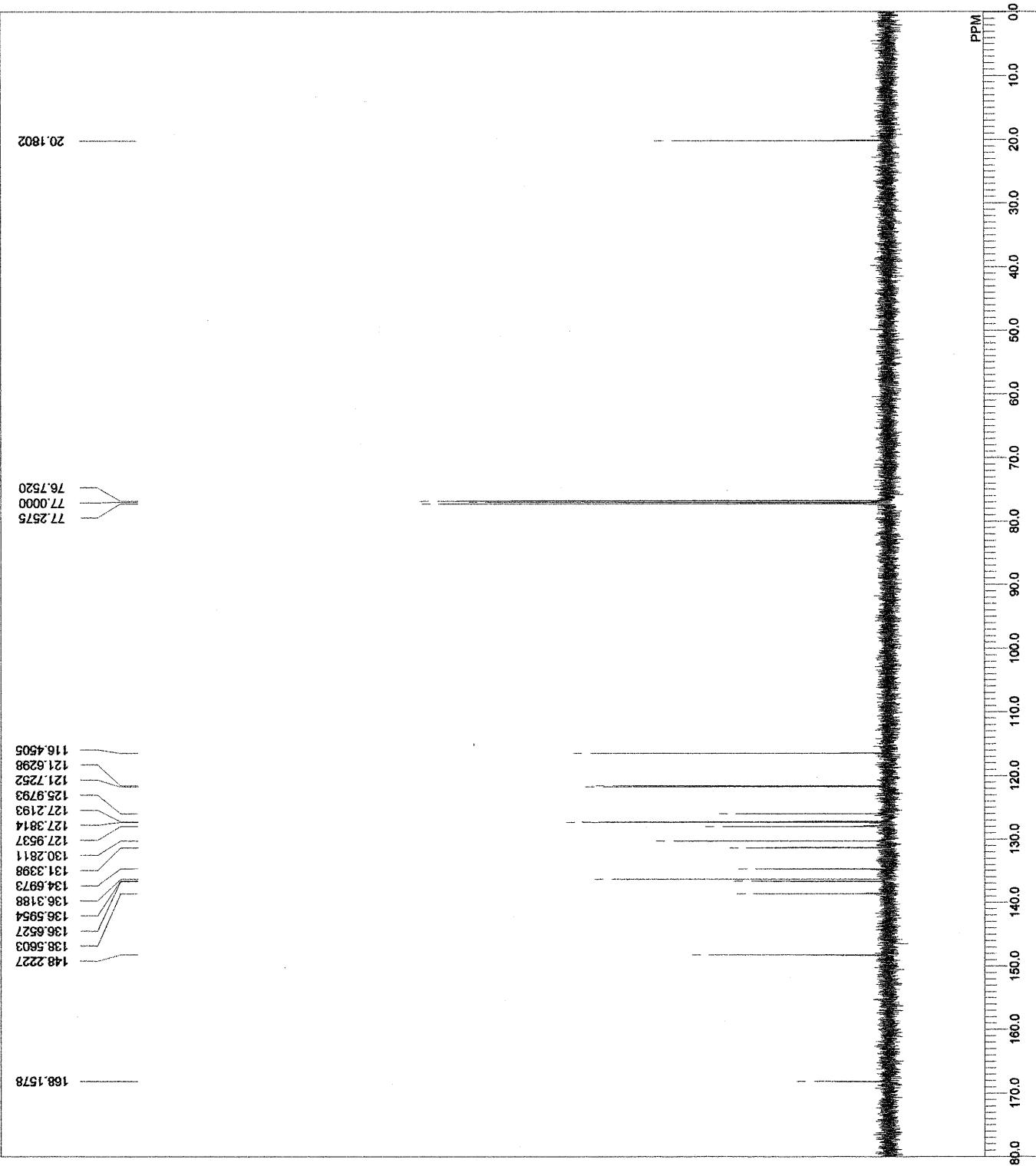
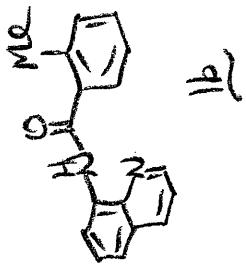




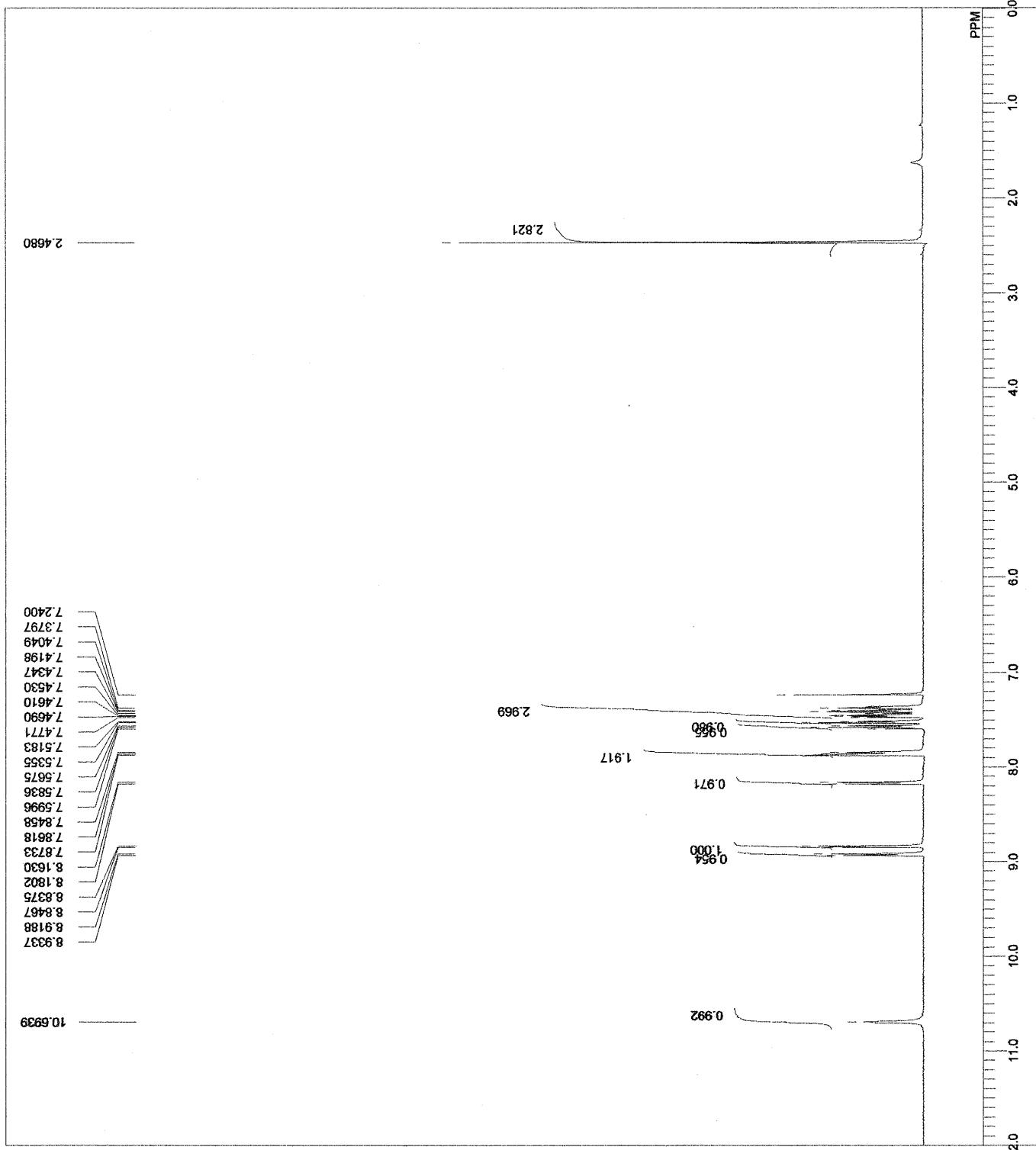
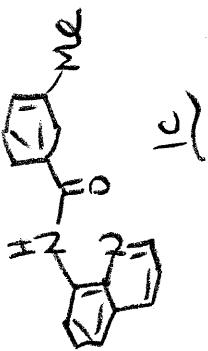
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EXMOD

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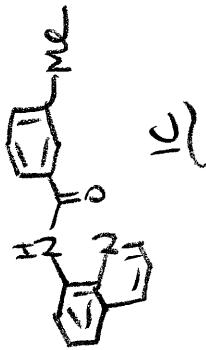
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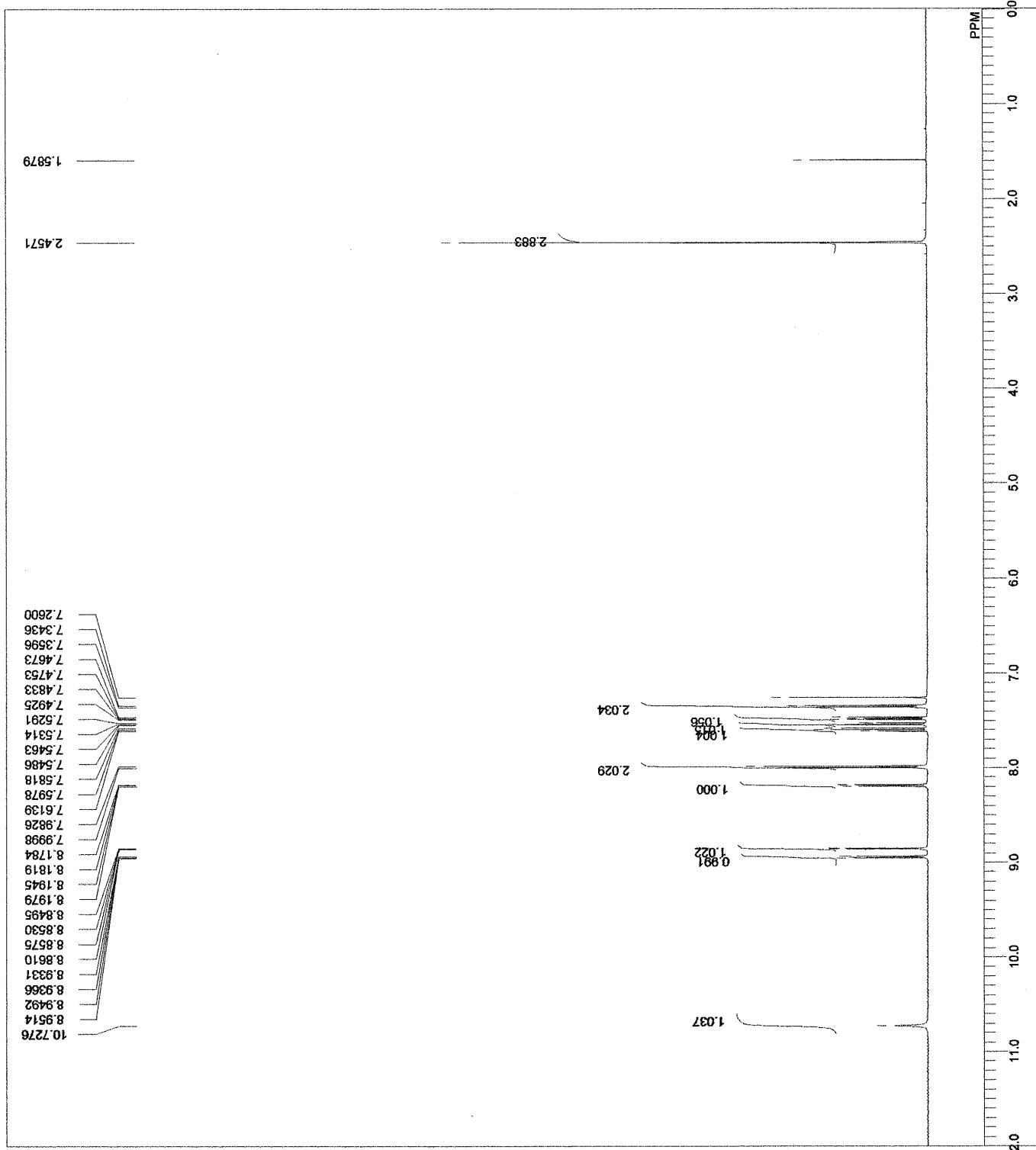
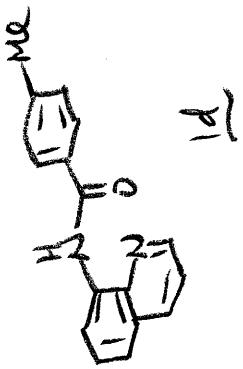
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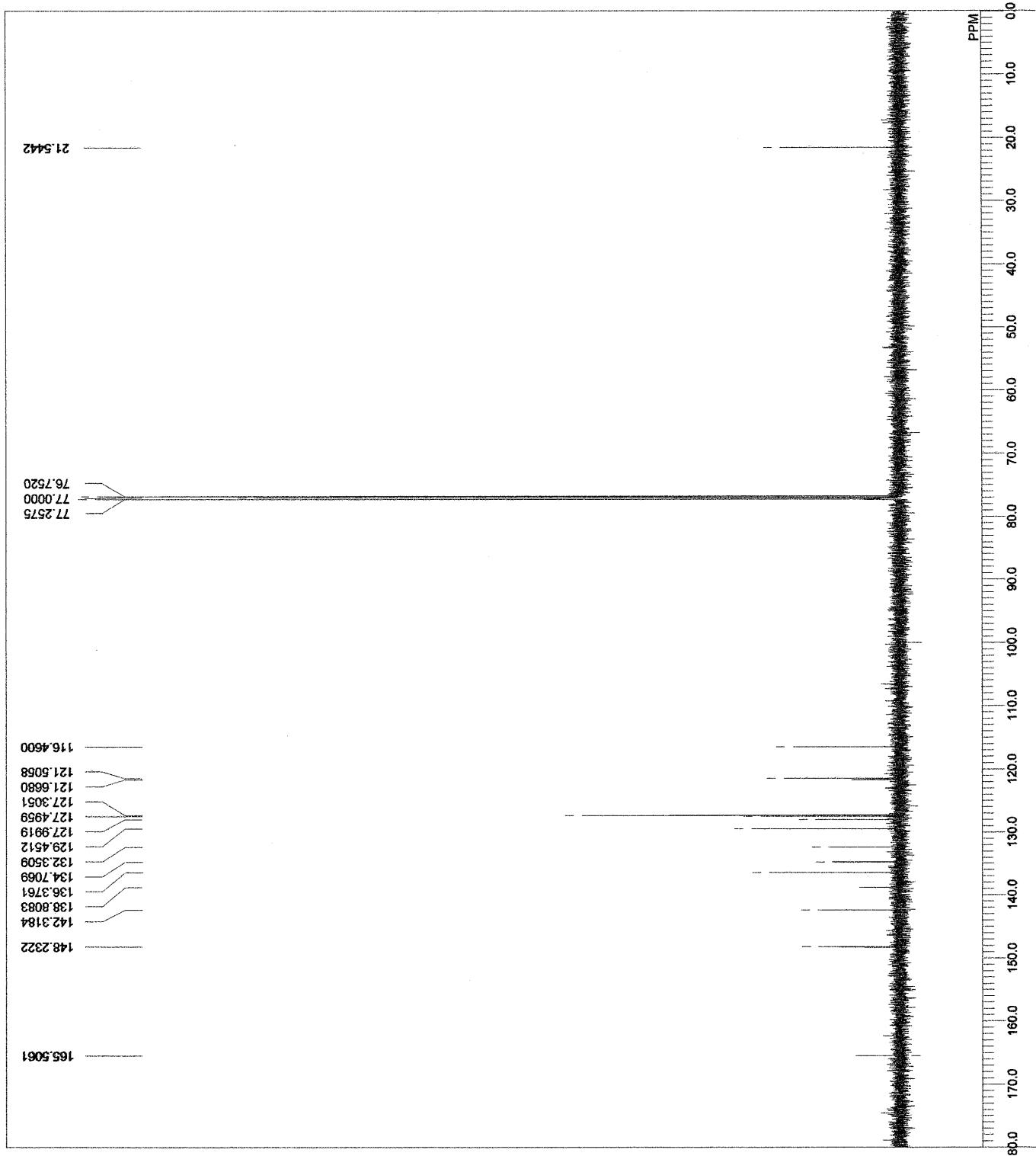
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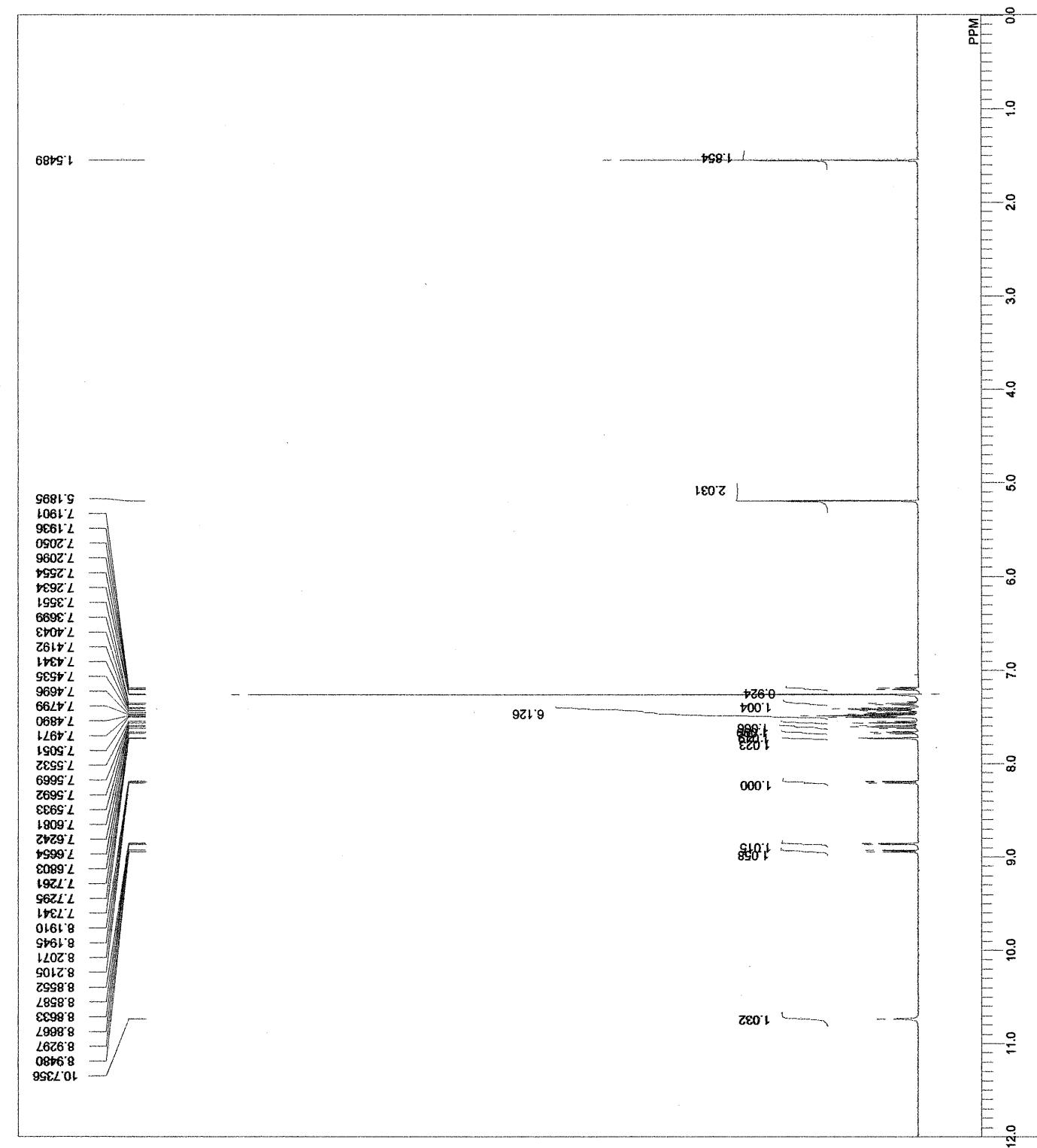


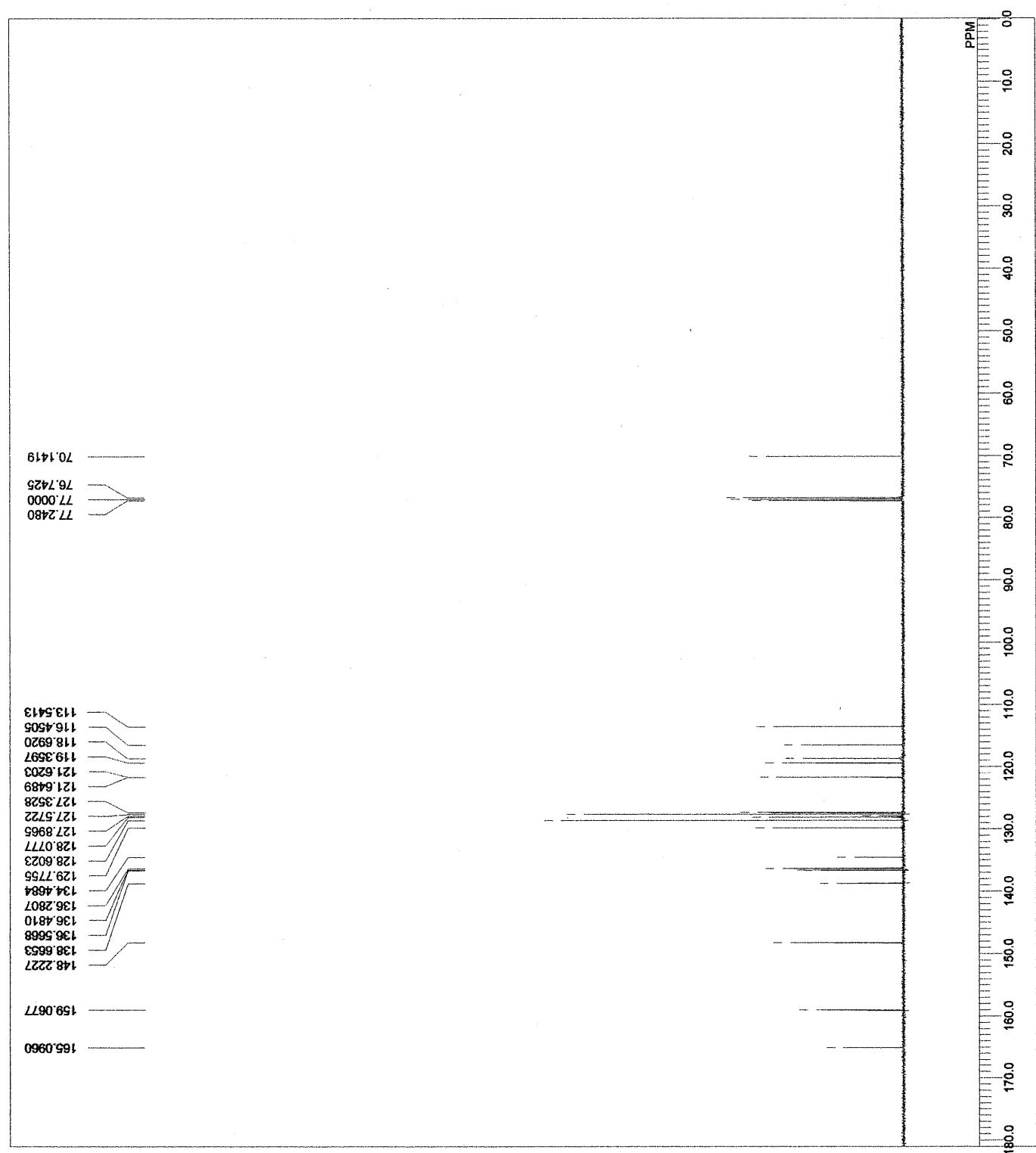
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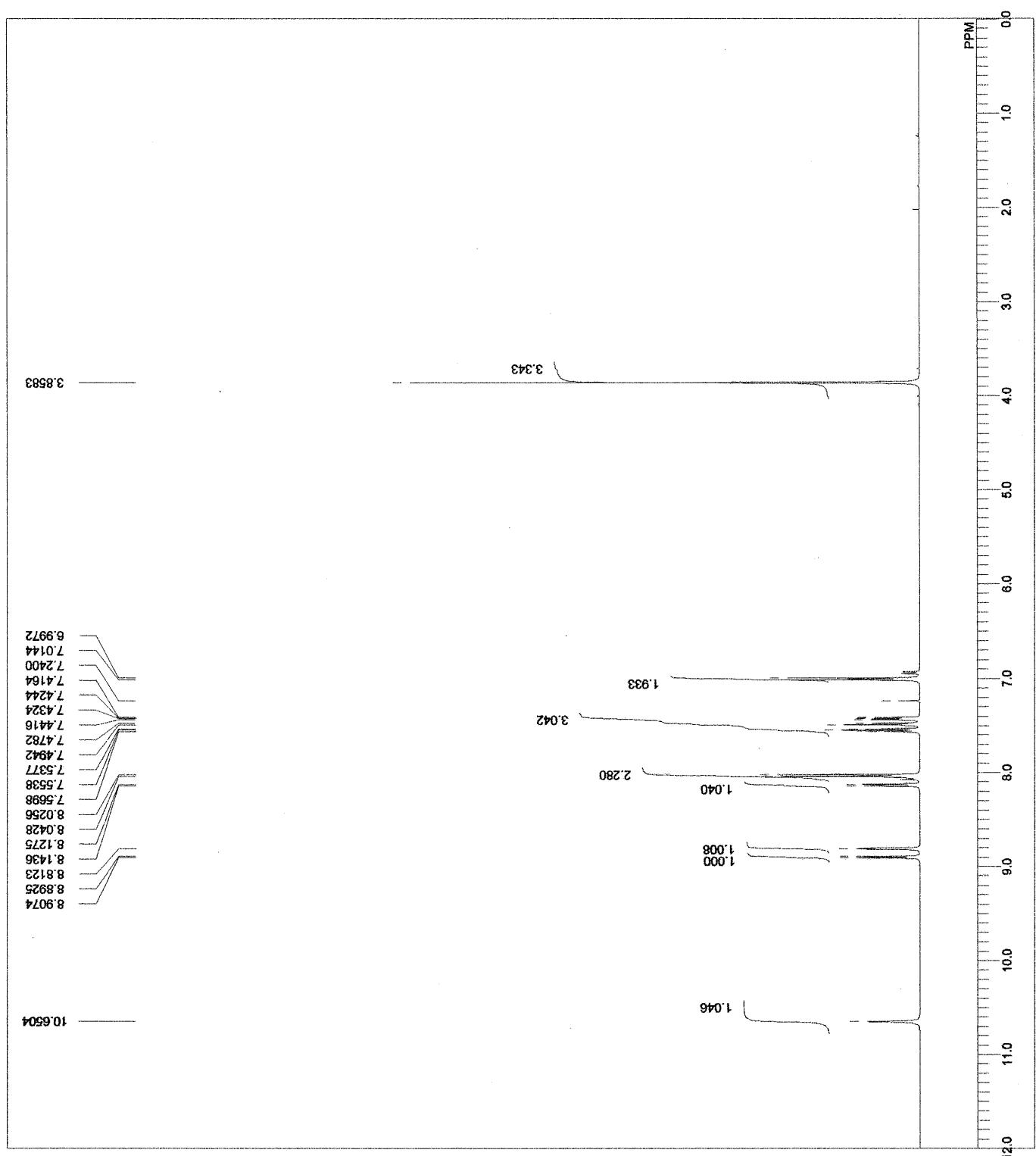
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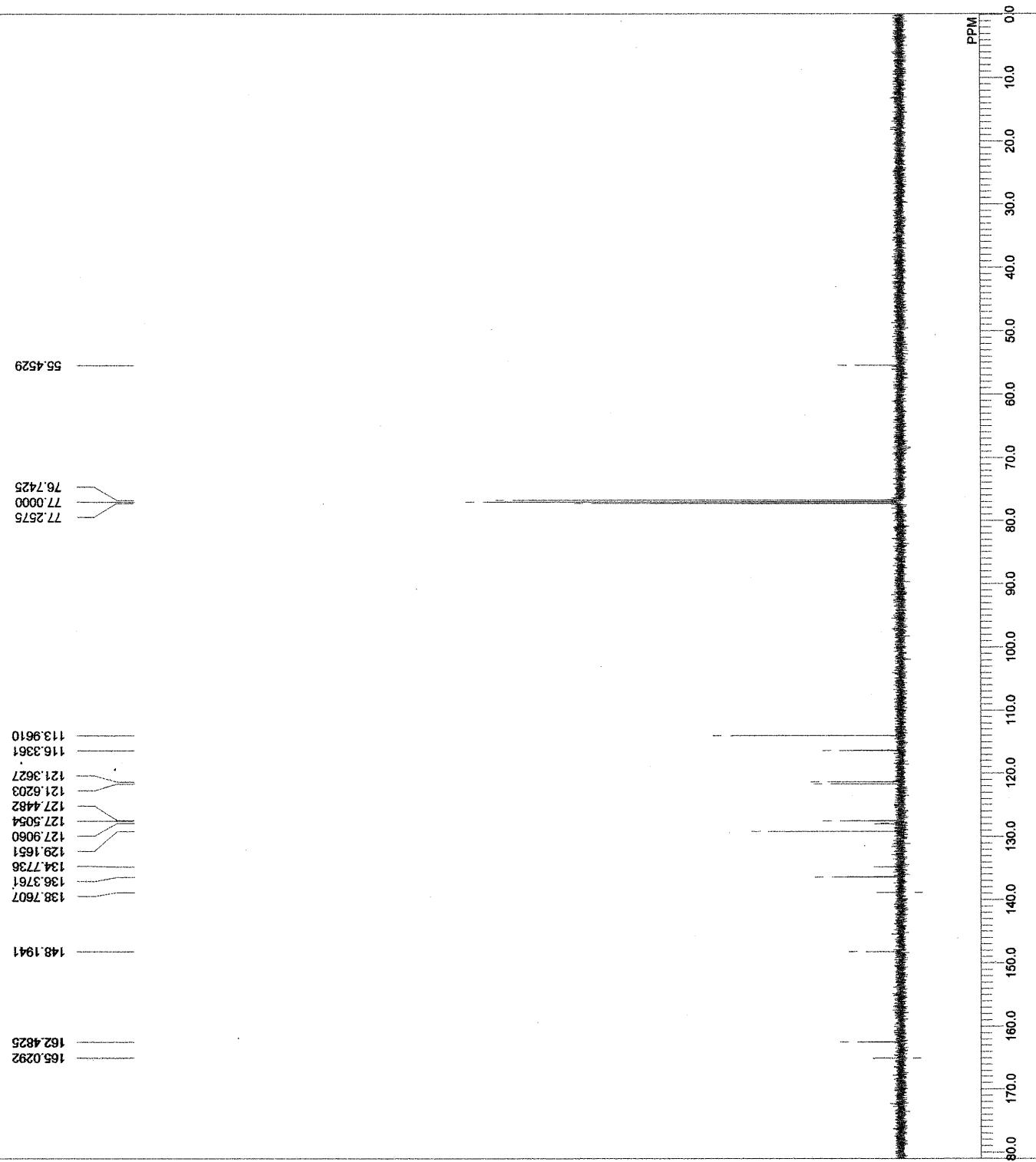
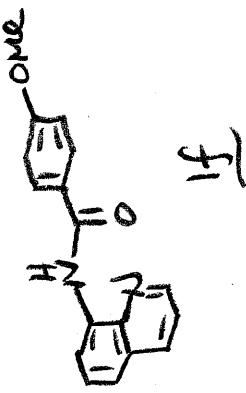


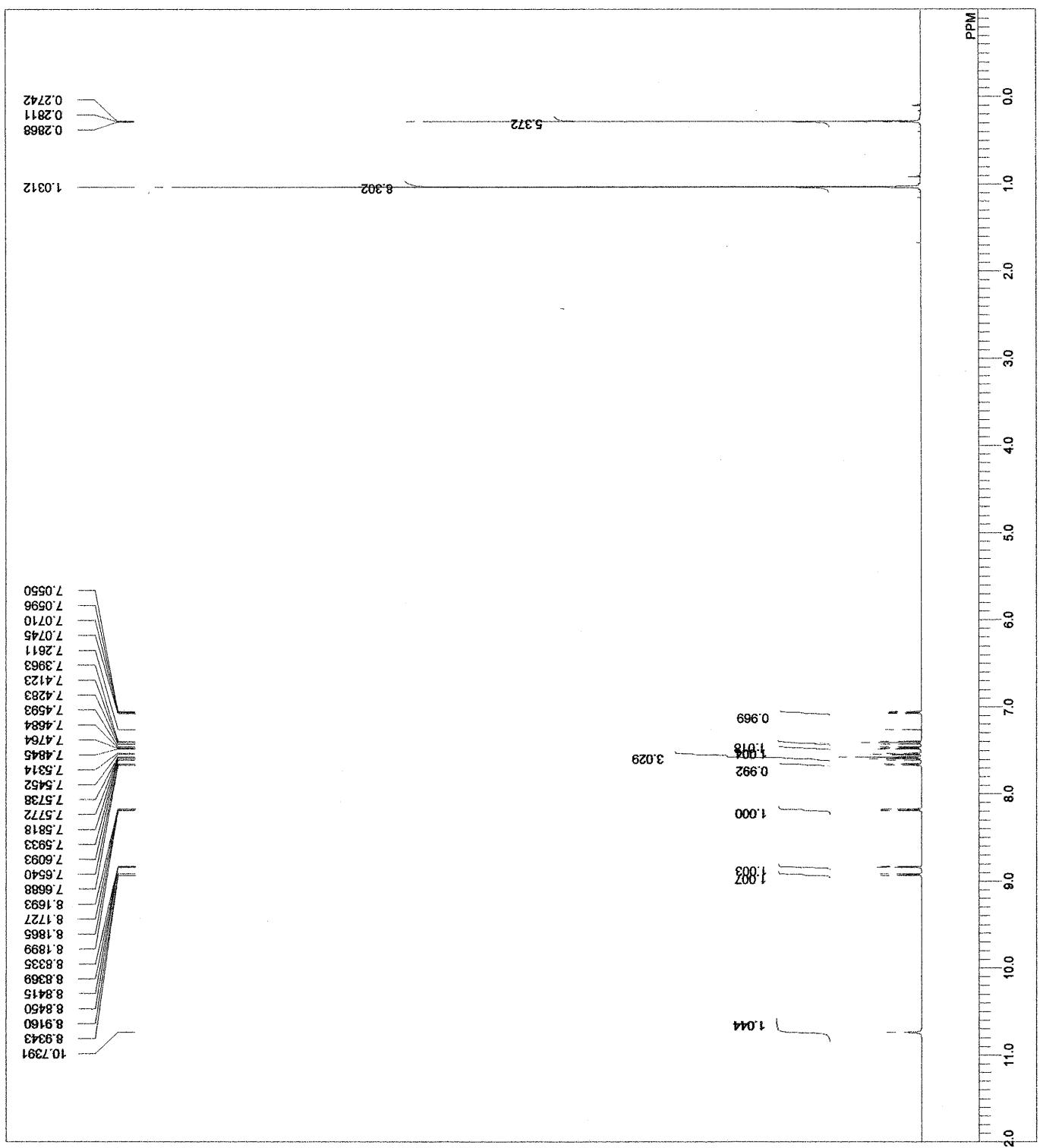


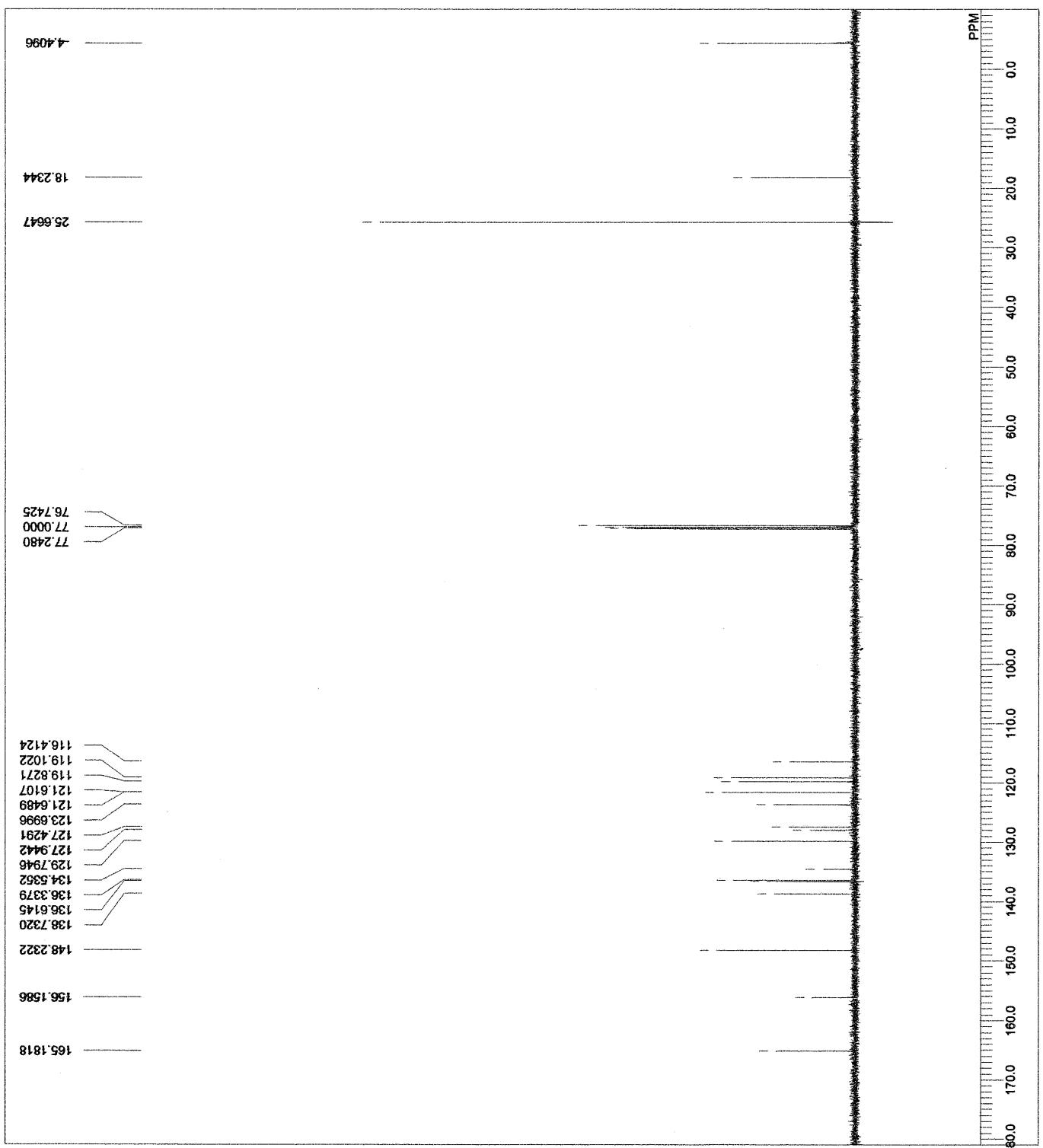
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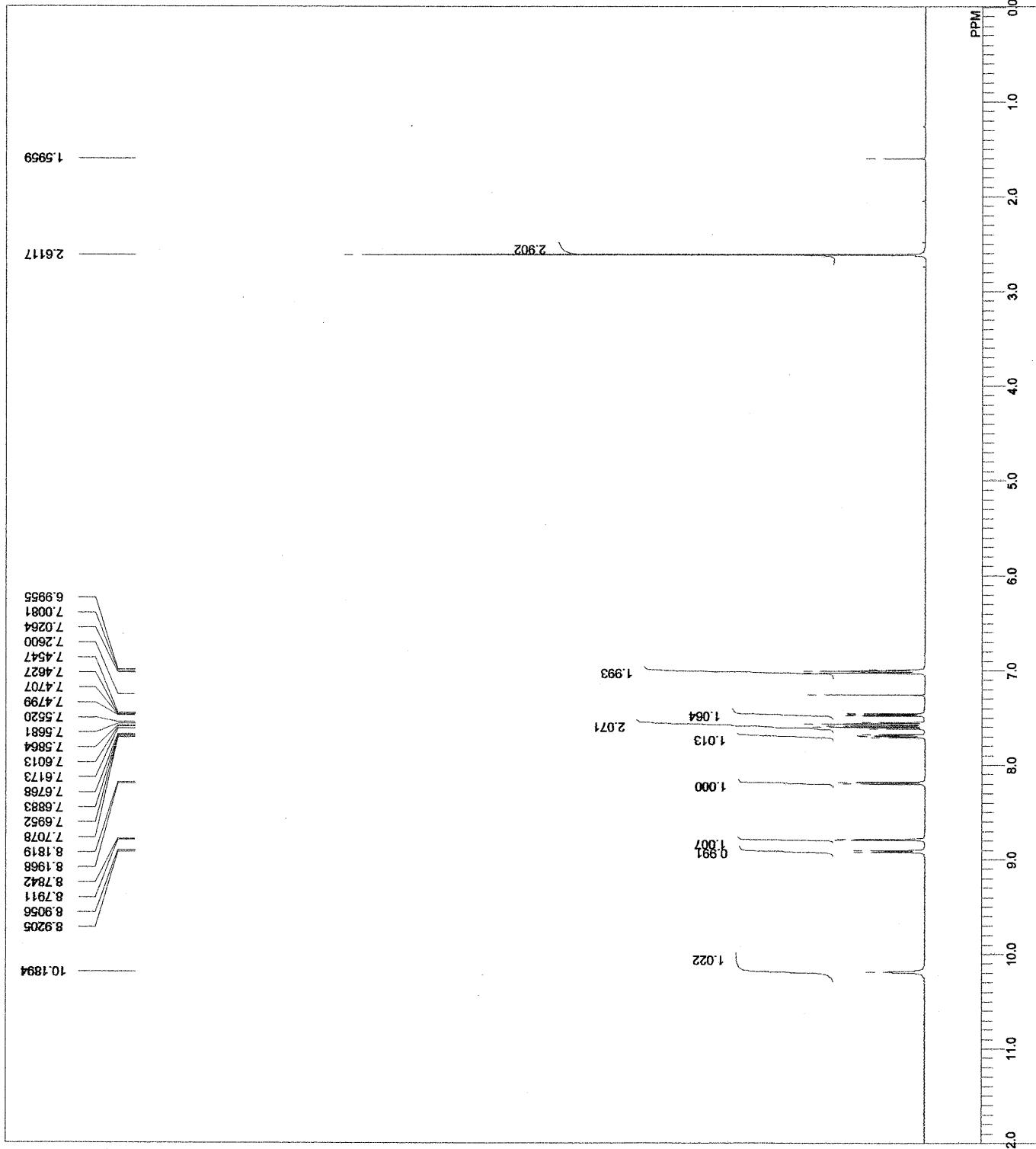
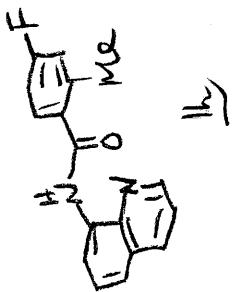
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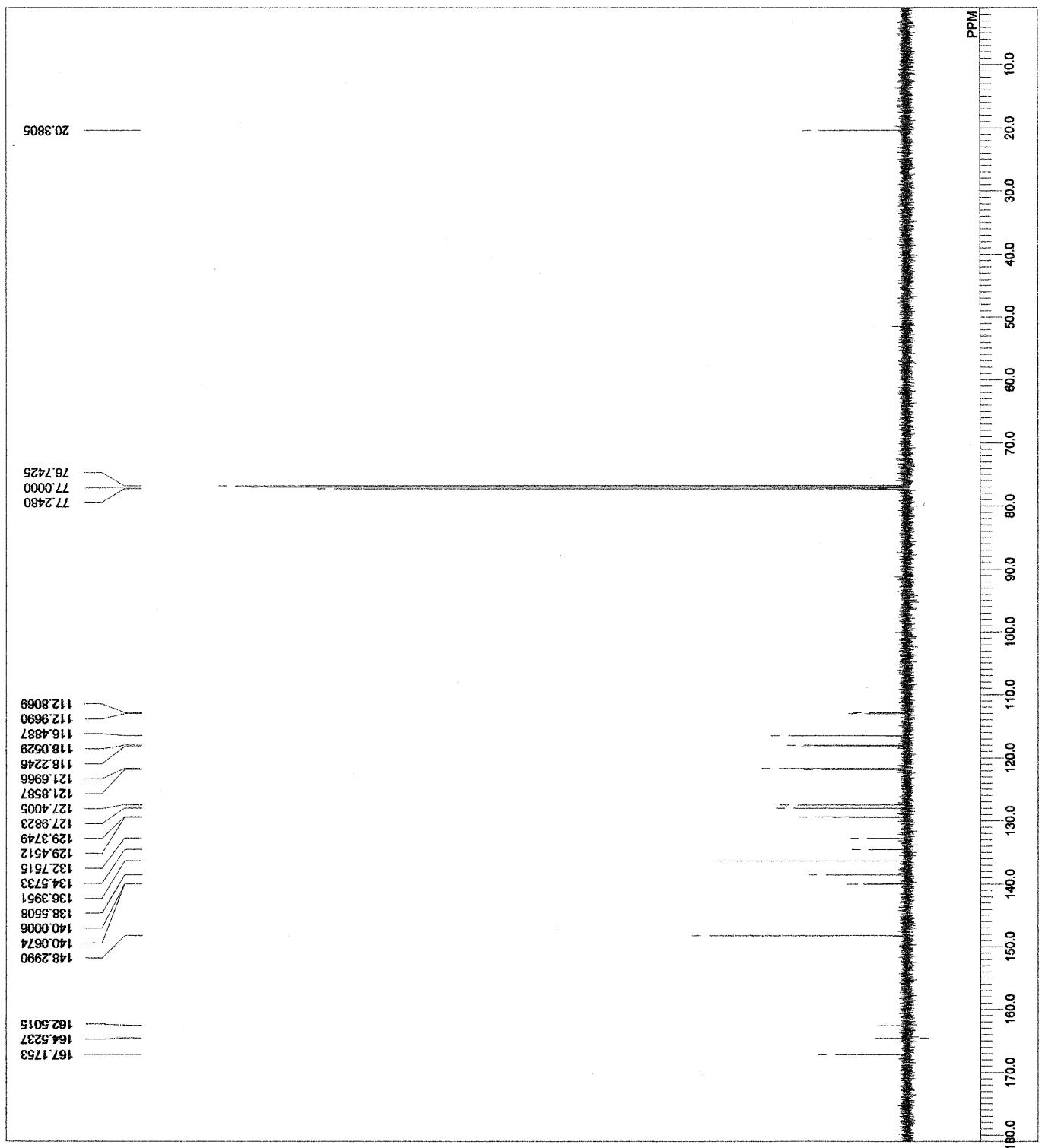






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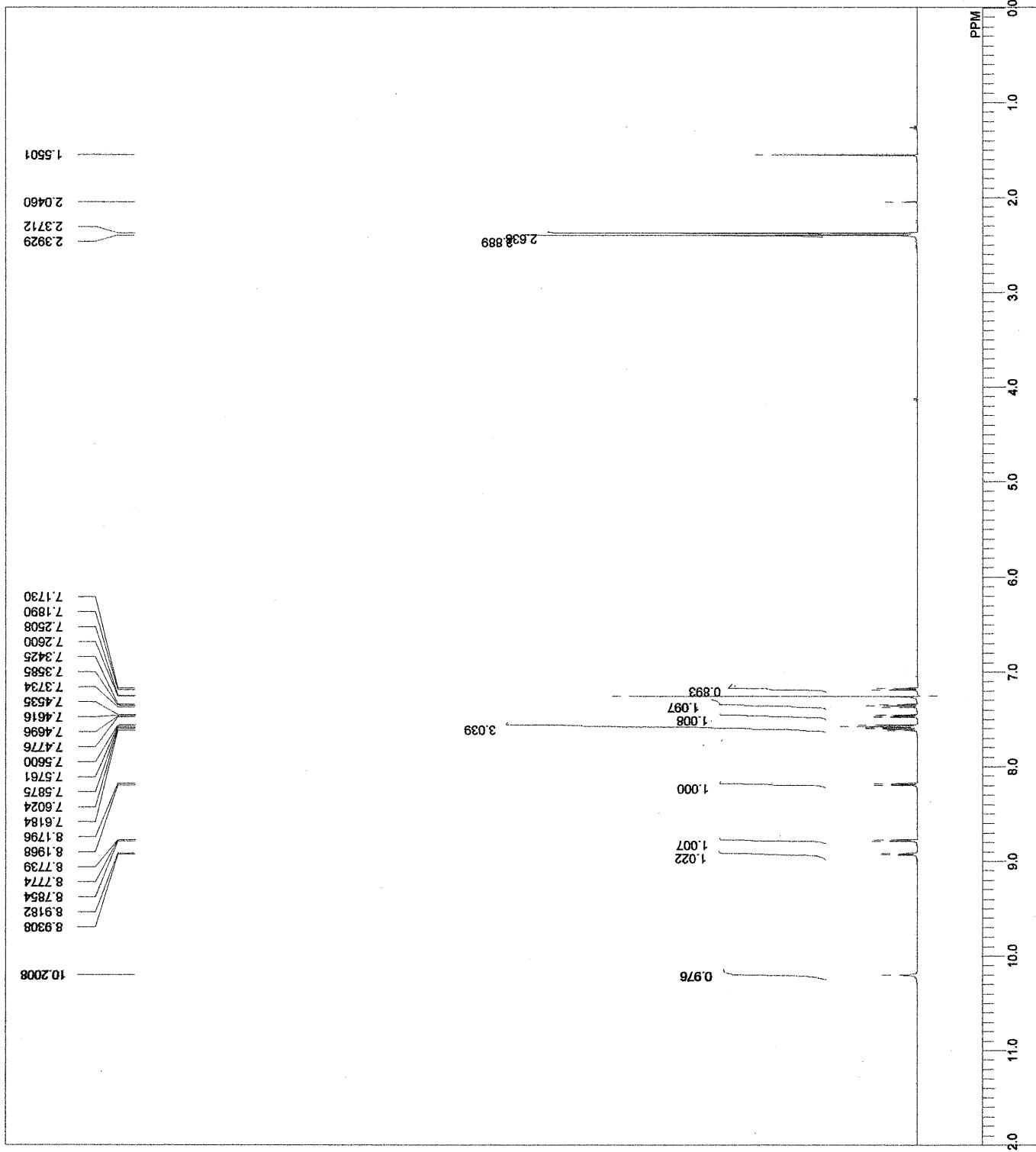
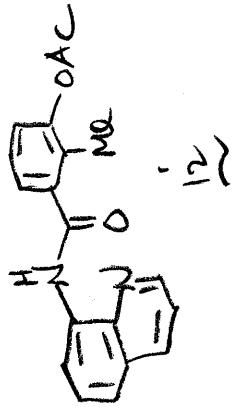


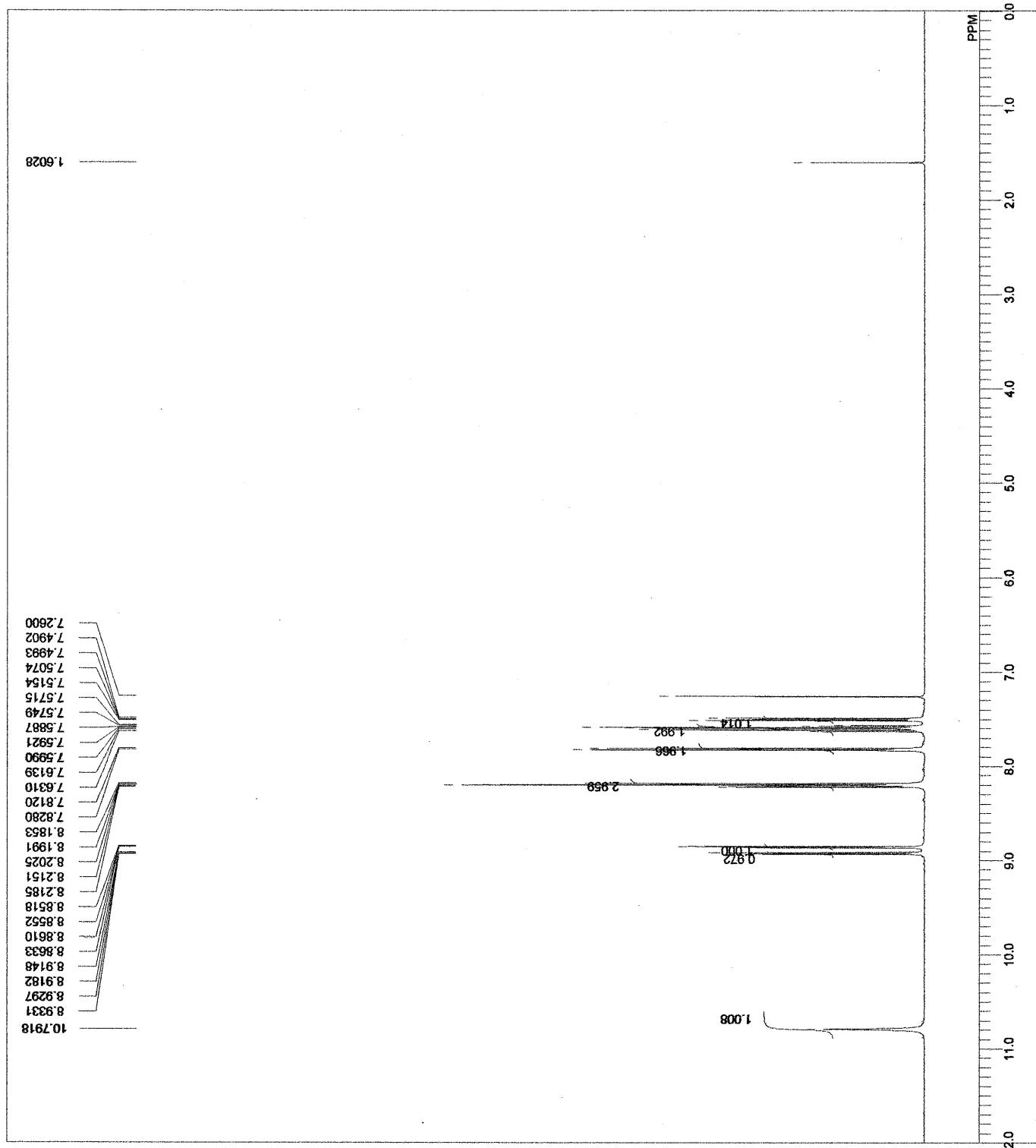


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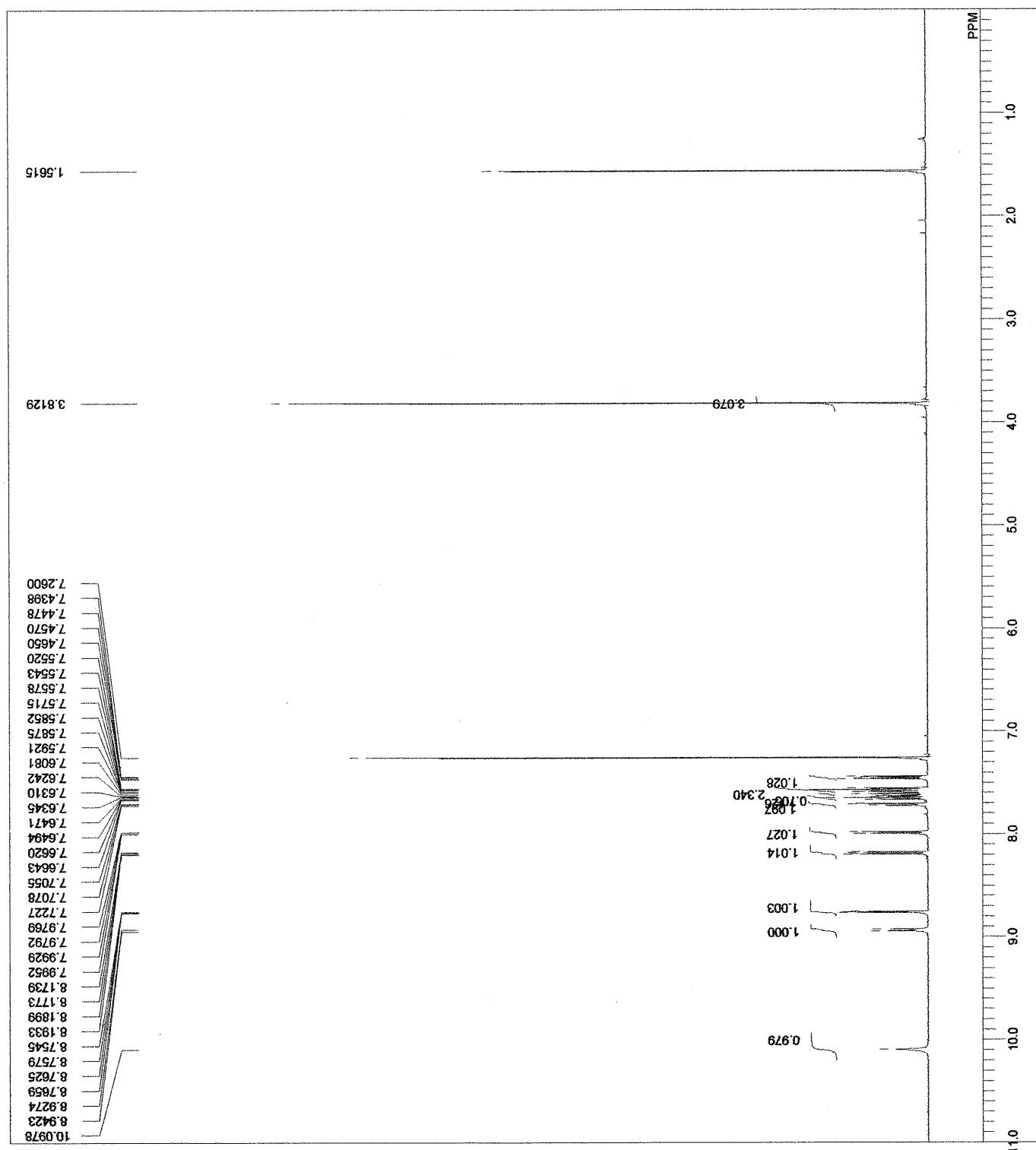
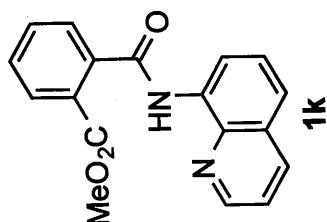
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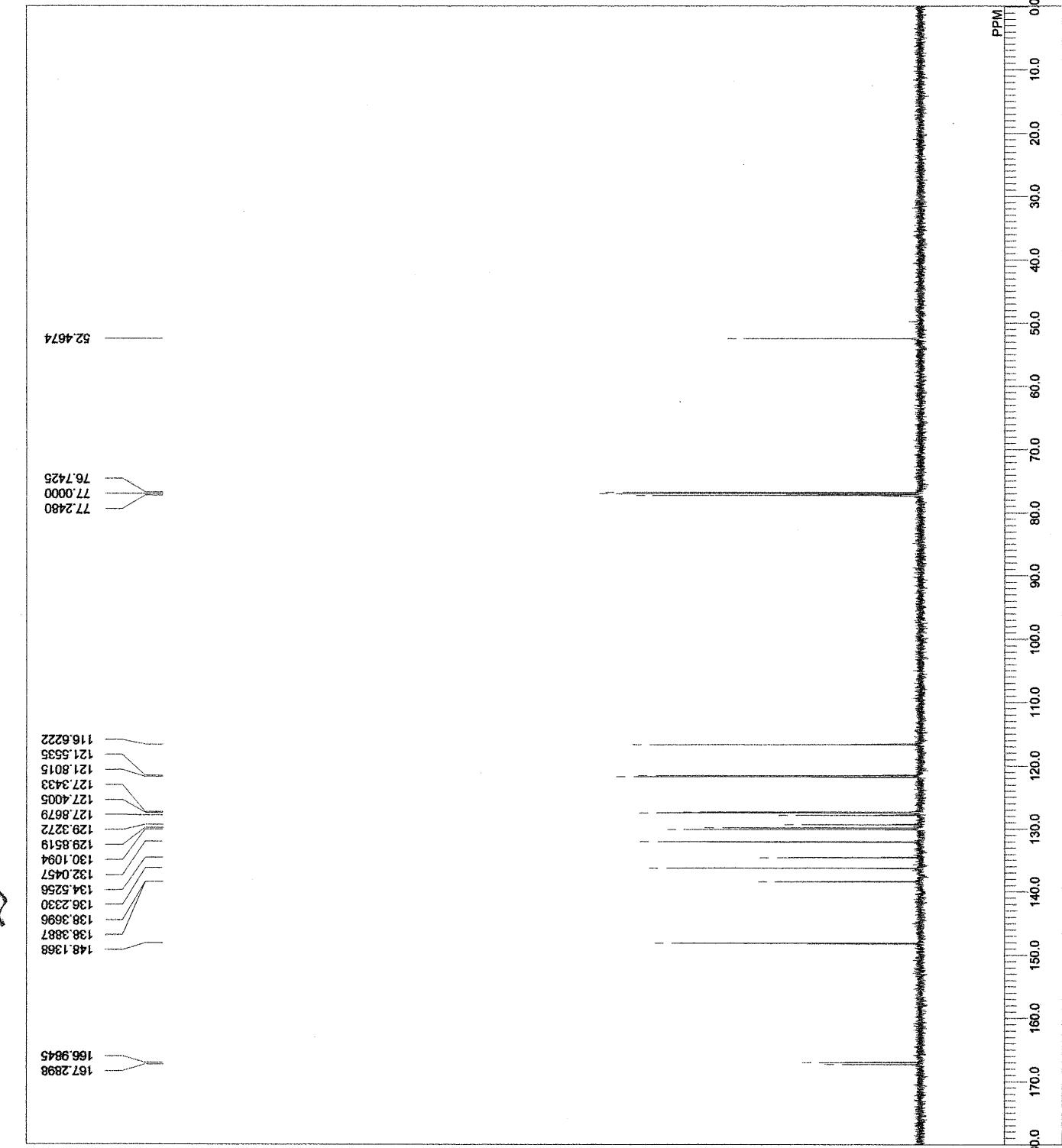
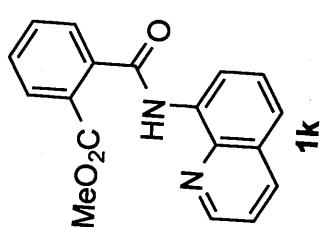
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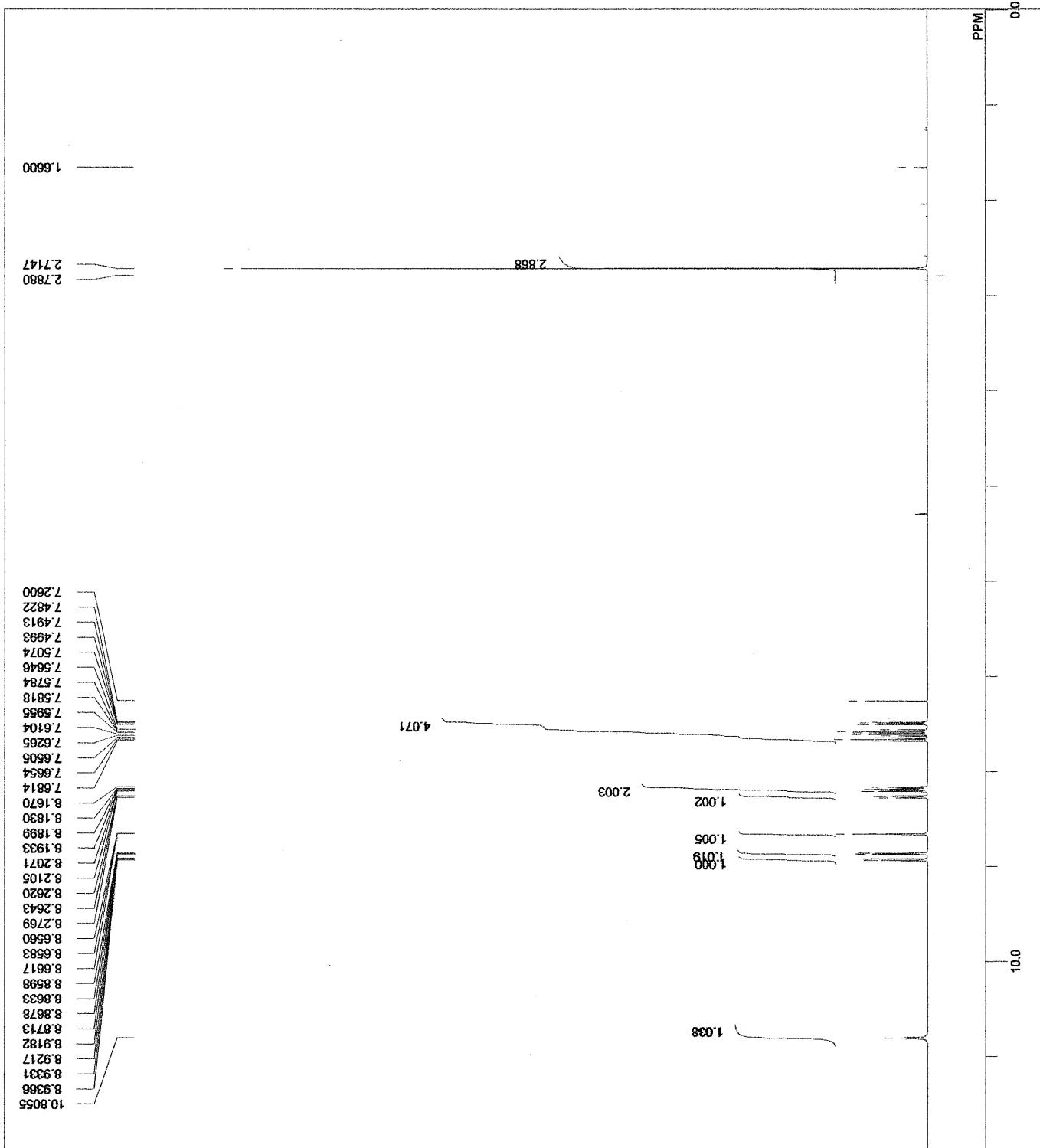
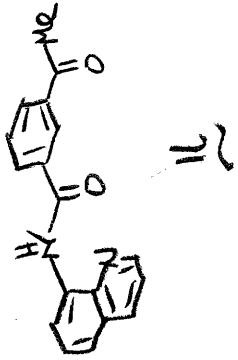
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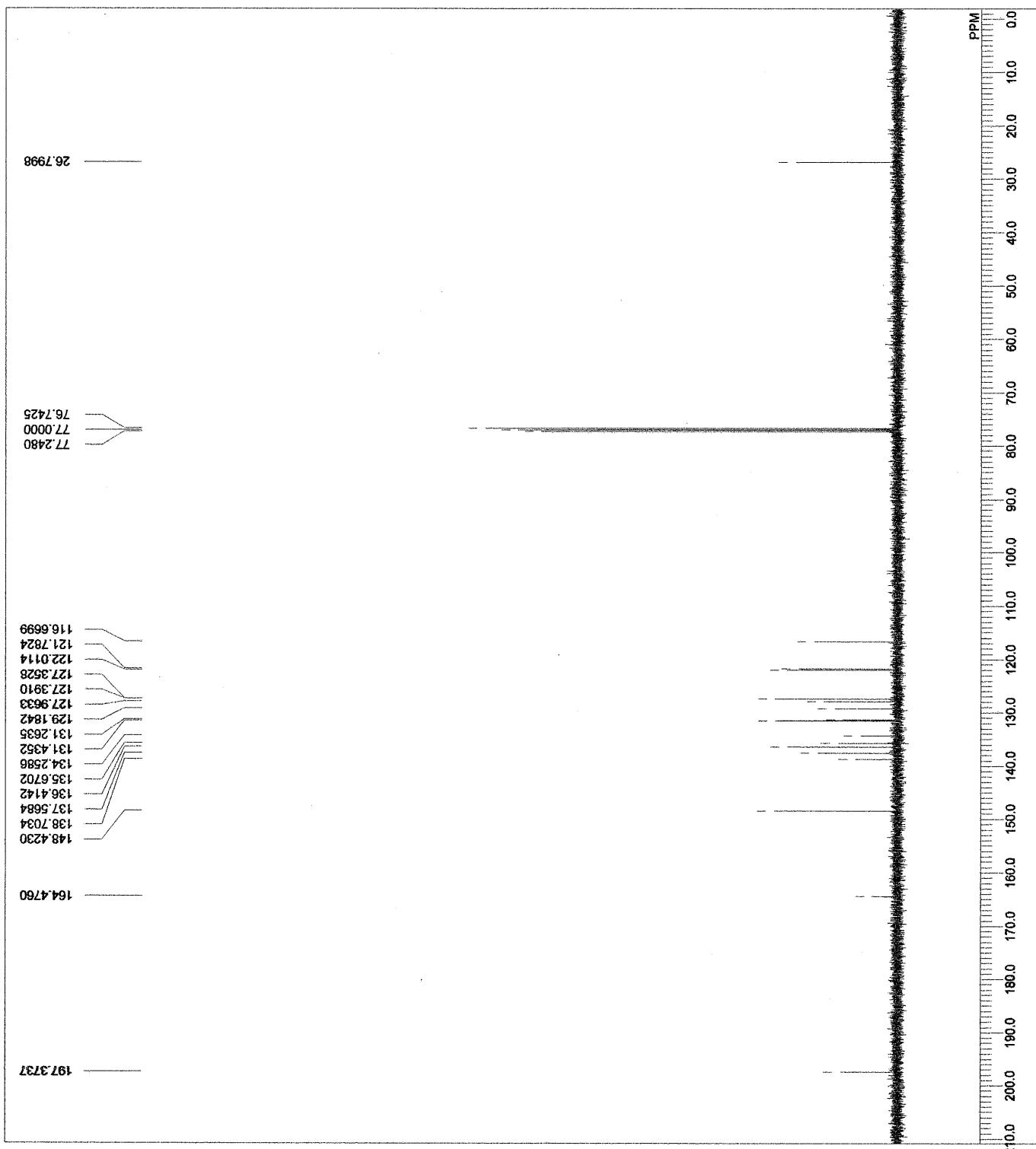
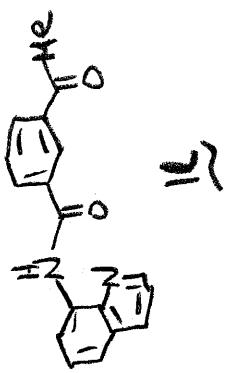
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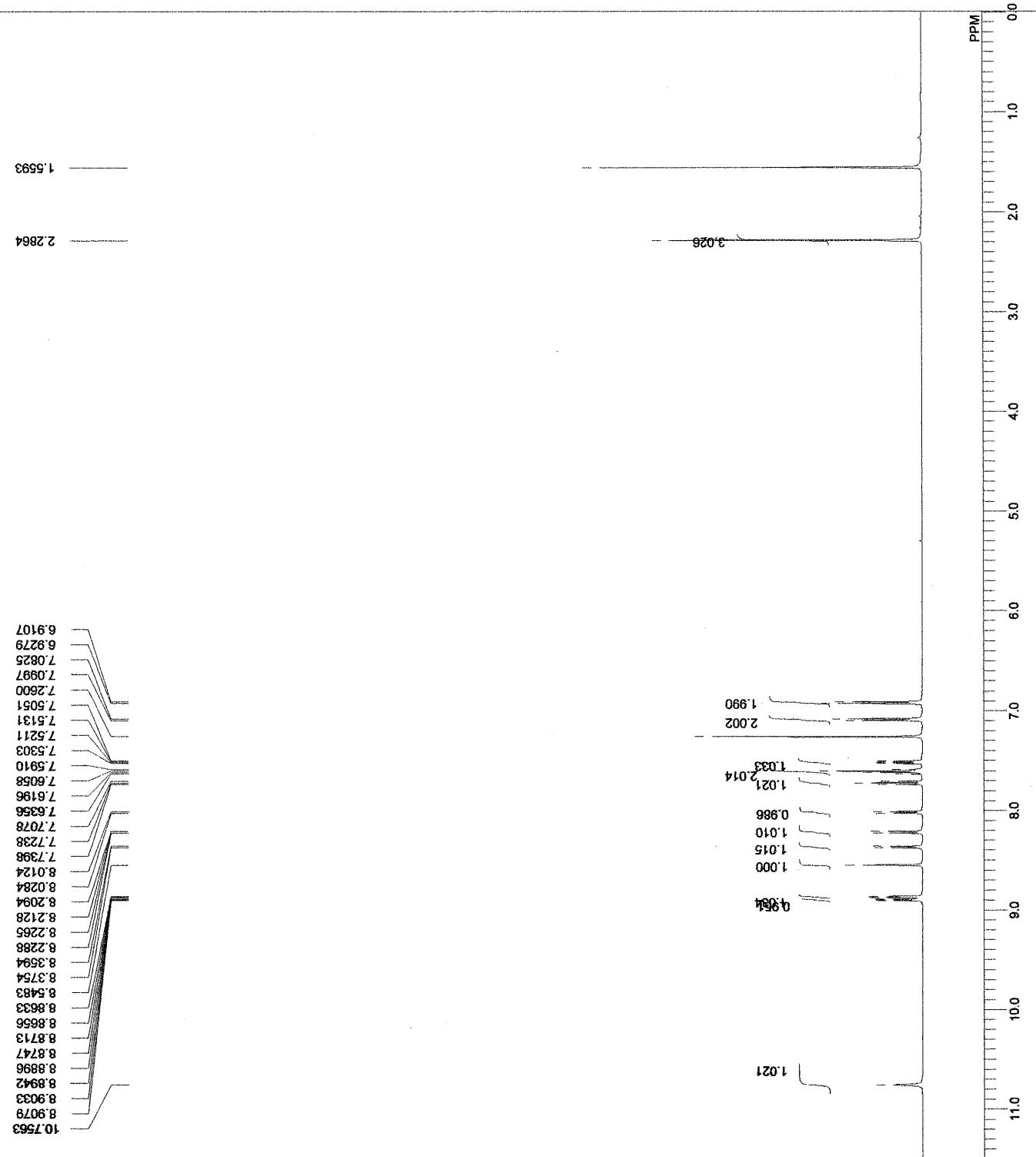
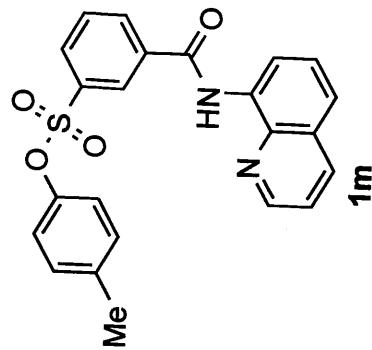
skk-sm-11-C(ArCOMe)-1-1.jdf
skk-sm-11-C(ArCOMe)
2014-02-18 16:43:09

DFILE
COMNT
DATIM
OBNUC
EXMOD
OBFRQ
OBSET
OBFIN
POINT
FREQU
SCANS
ACQTM
PD
PW1
IRNUC
CTEMP
SLVNT
EXREF
BF
RGAIN

13C
carbon-13D
125.77 MHz
7.87 kHz
4.21 Hz
32767
36308.18 Hz
105
0.8336 sec
3.0000 sec
3.40 usec
1H
CDCL3
77.00 ppm
0.12 Hz
60

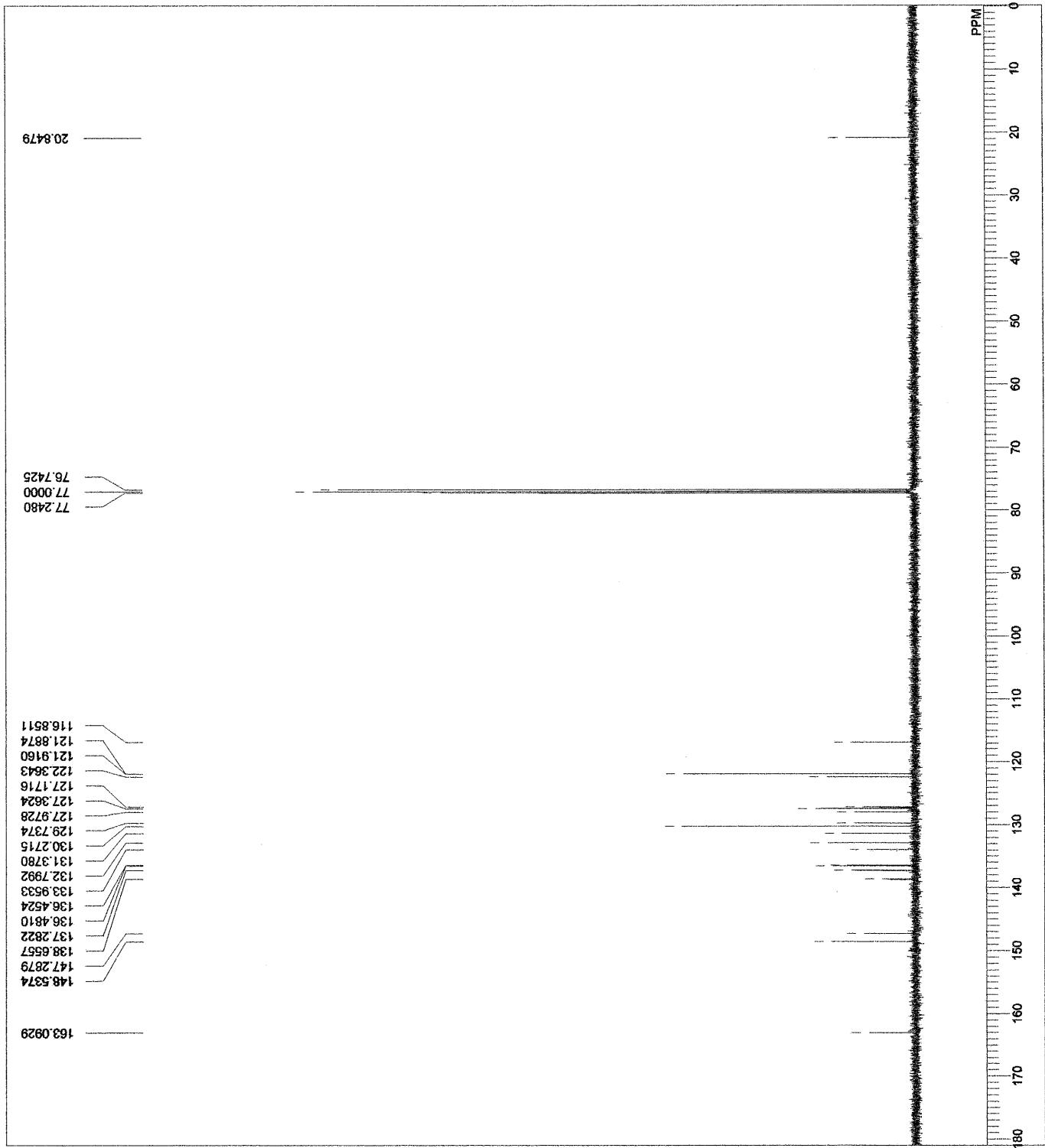


skk-p50-SMsyn-ArSO₃Ar-H

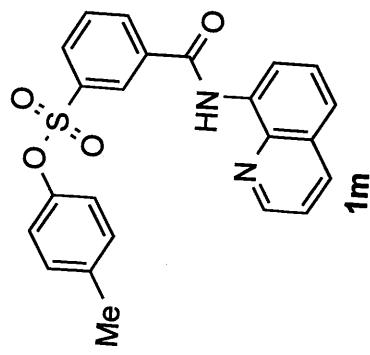




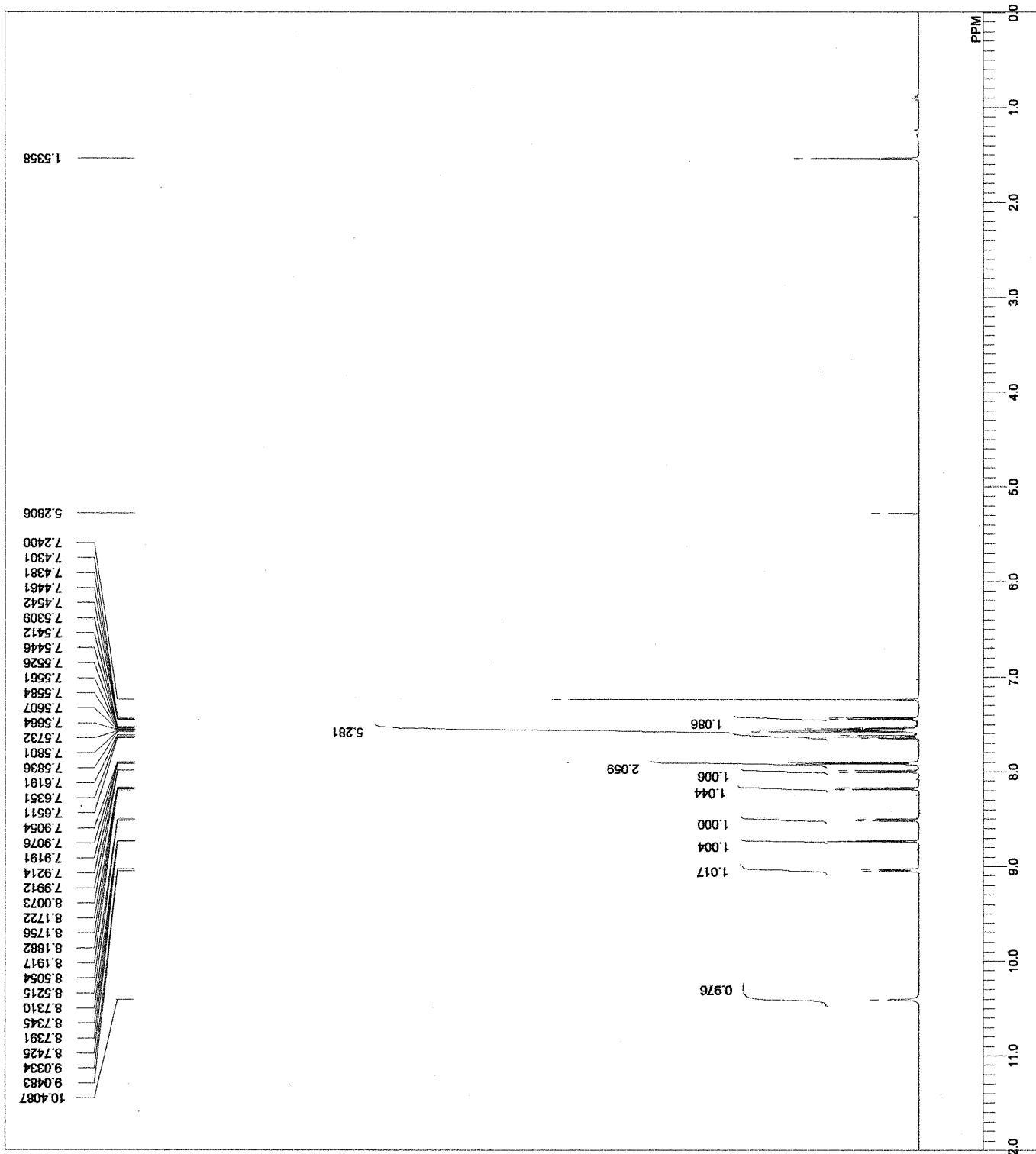
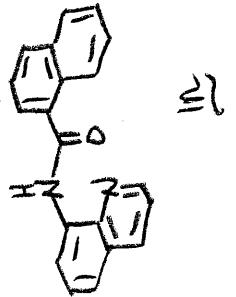
163.0929
148.5374
147.2879
138.6657
137.2822
136.4610
136.4524
133.9533
132.7992
131.3780
130.2715
129.7374
127.9728
127.7324
127.7116
122.7143
121.9160
121.8874
116.8511
77.2480
77.0000
76.7425
20.8479



skk-p55-SMsyn-8AQSO3Ar-C-1-1.d1f
skk-p55-SMsyn-8AQSO3Ar-C
2013-10-28 15:06:11
13C
carbon.jdx
EXMOD
OBFRQ 125.77 MHz
OBSET 7.87 kHz
OBFIN 4.21 Hz
POINT 32767
FREQU 39308.18 Hz
SCANS 357
ACQTM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usec
IRNUC 1H
CTEMP 24.2 c
SLVNT CDCl₃
EXREF 77.00 ppm
BF 0.12 Hz
RGAIN 60



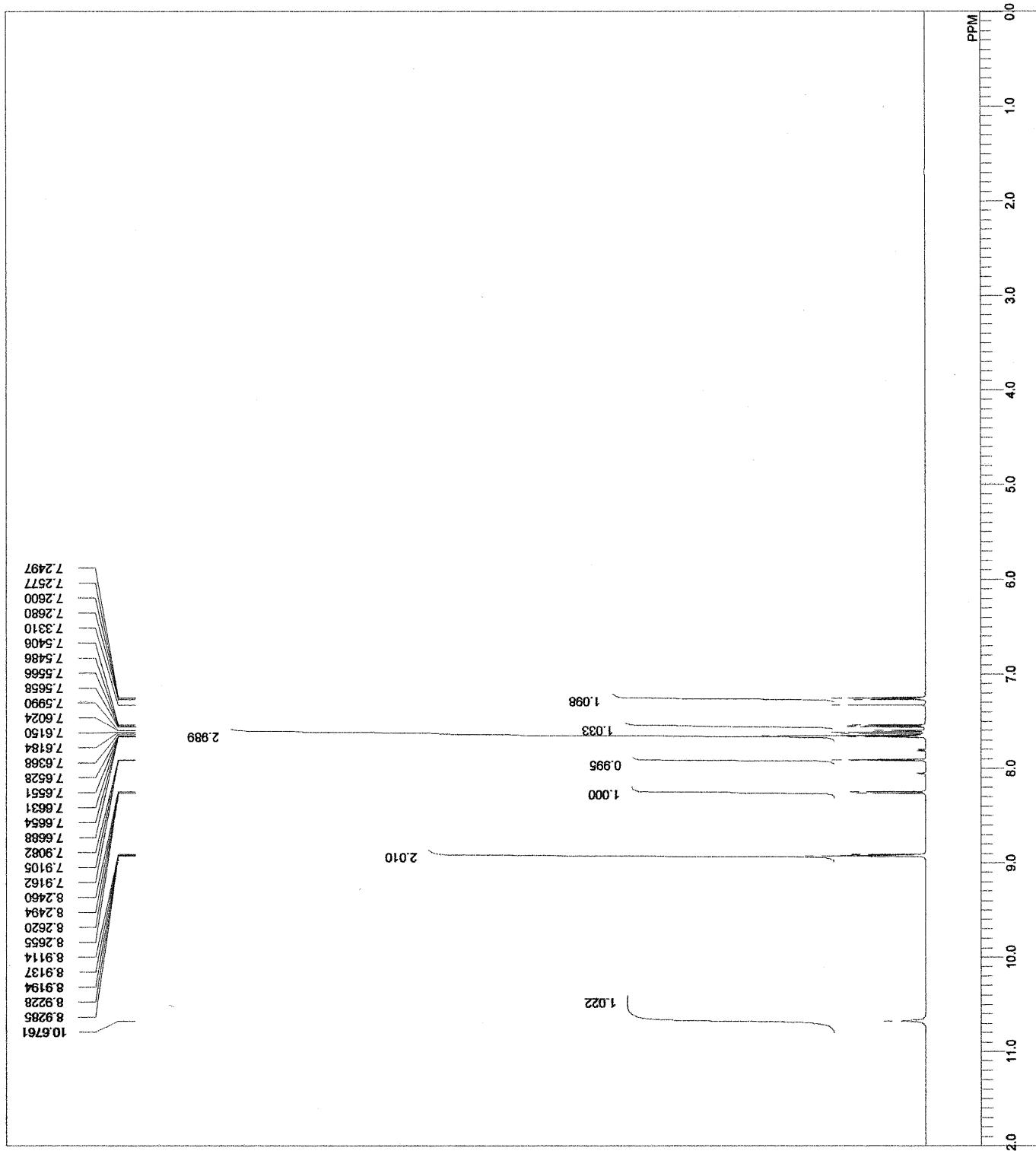
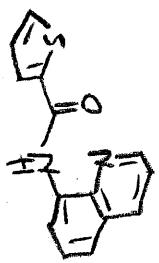
DFILE	skk-smr-1m-H-4.1.jdf
COMMENT	skk-smr-1m-H
DATIM	2014-02-18 18:38:43
OBNUC	1H
EXMOD	proton
BFRQ	500.00 MHz
OBSET	2.41 kHz
OBFIN	6.01 Hz
POINT	16384
FREEF	9384.38 Hz
SCANS	8
ACQTM	1.7459 sec
PD	5,0000 sec
PWM	5.55 usec
IRNUC	1H
CTEMP	21.3 c
SLVNT	CDCL3
EXREF	7.24 ppm
BF	0.12 Hz
RGAIN	40

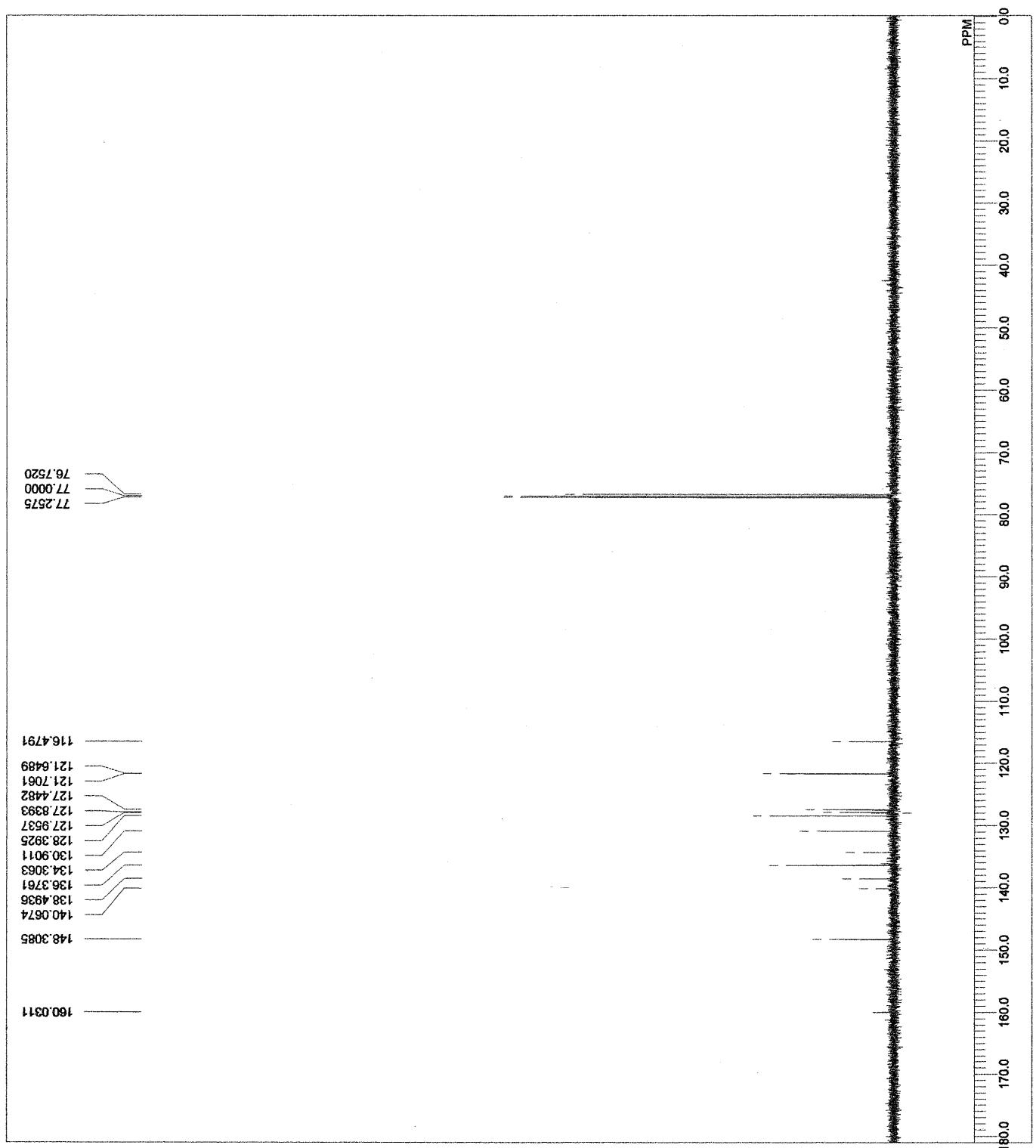


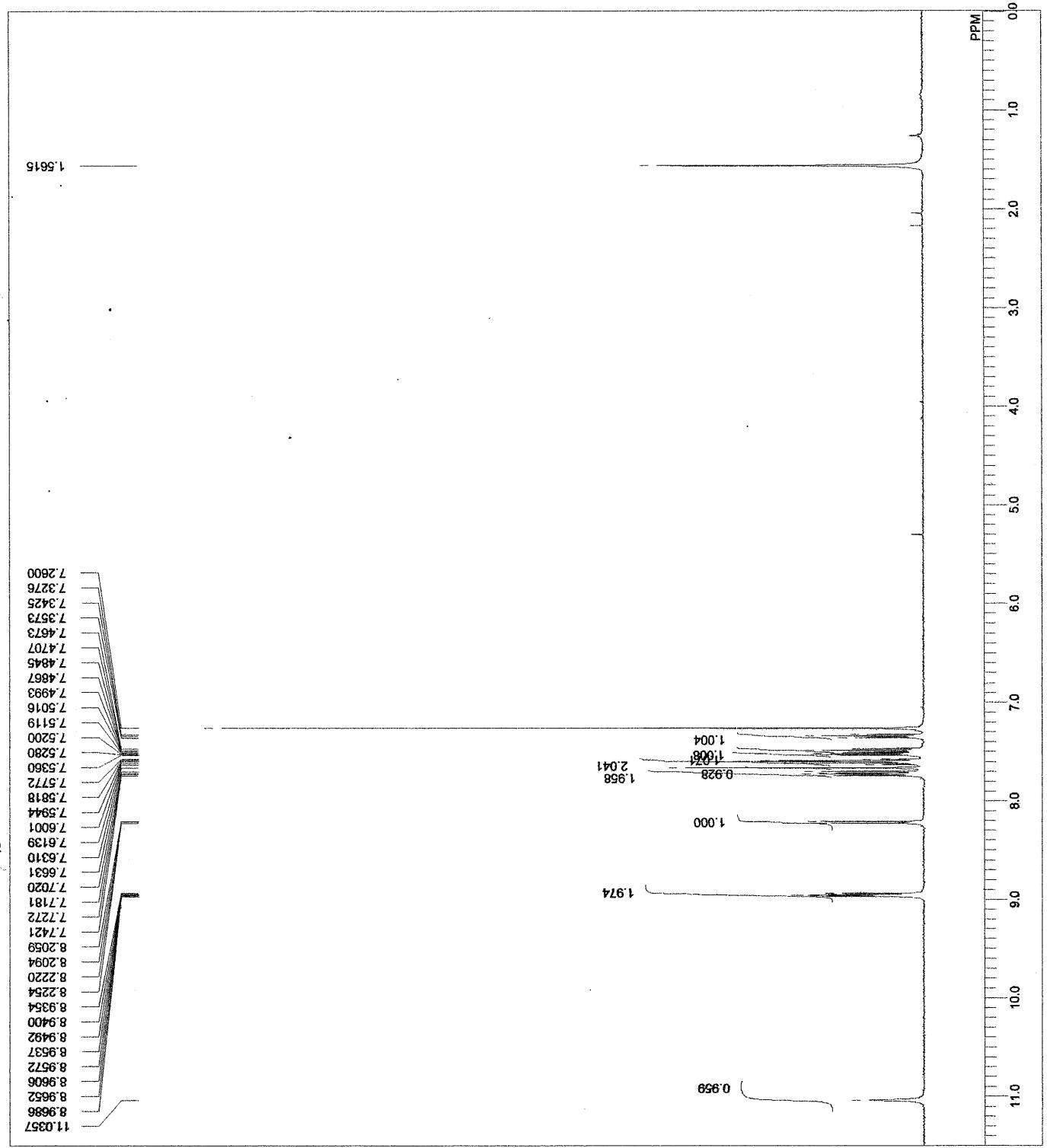
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sik-sm-11-(thioP)-1-1.b3
sik-sm-11-(thioP)
COMMNT
DATIM 2014-02-18 16:52:49
1H
proton1D
500.16 MHz
2.41 kHz
6.01 Hz
16384
9384.38 Hz
8
ACQTIME
5.0000 sec
PD
5.35 ussec
PW1
1H
IRNUC
CTEMP
SLVNT
CDCL3
21.7 c
EXREF
7.26 ppm
BF
0.12 Hz
40

```



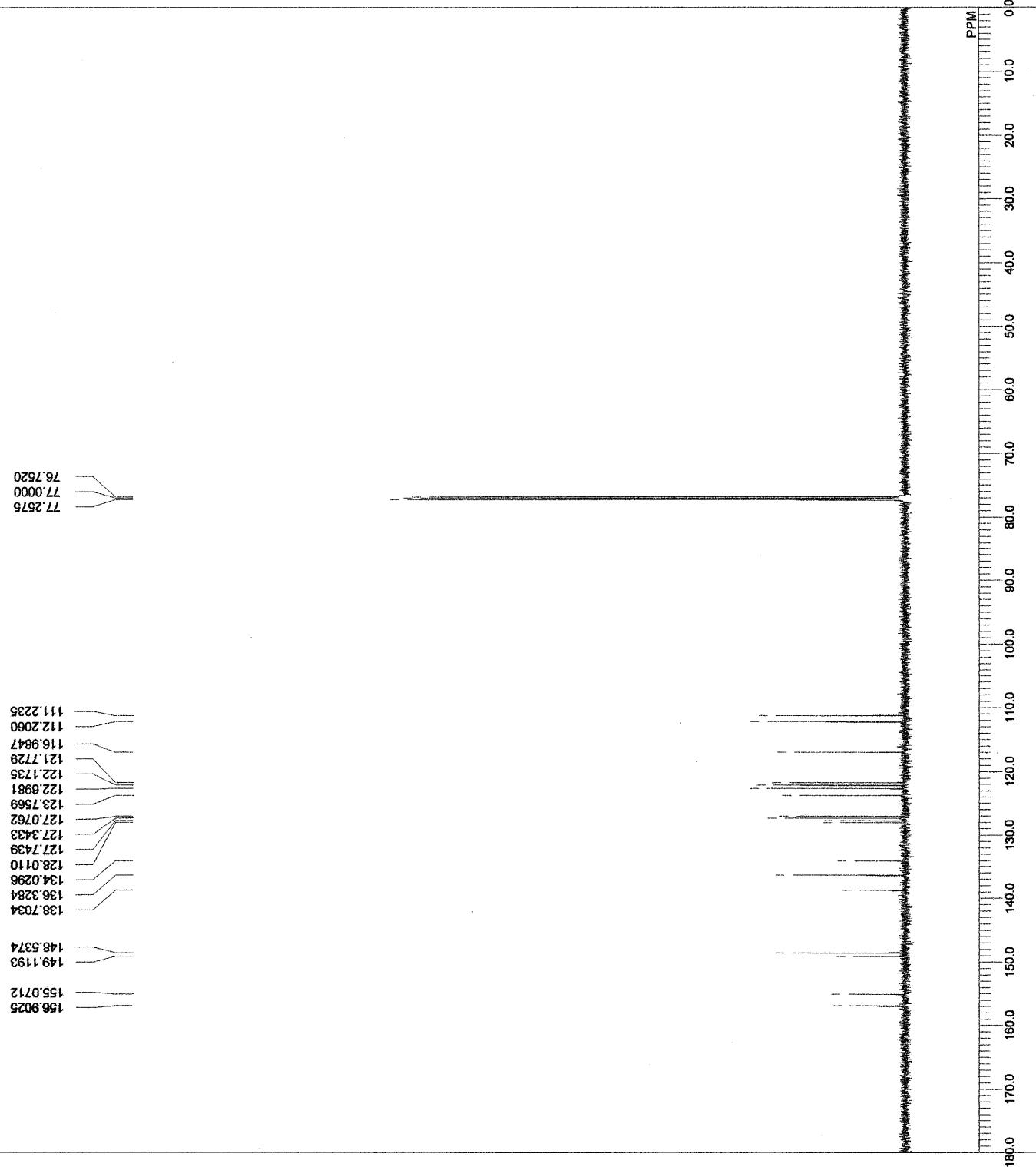
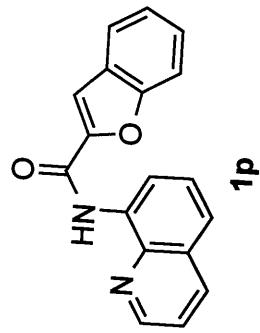




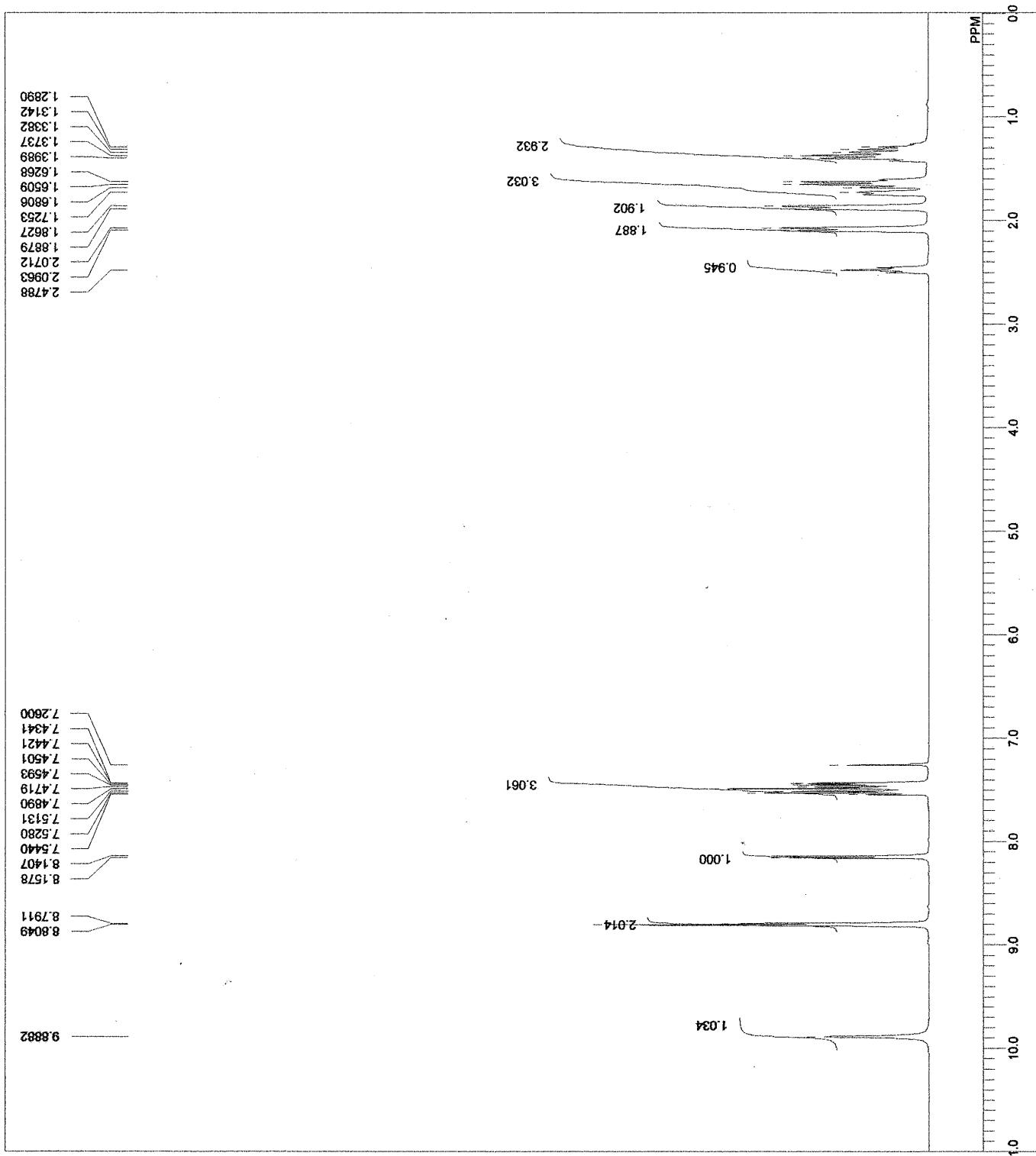
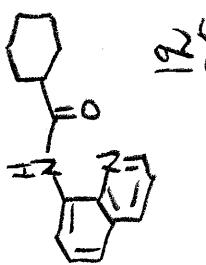
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DFILE    shk-p55-SMsyn-Bfurian-C-1-a1s
COMMENT
DATIM   2013-10-28 16:20:01
OBNUC   13C
EXMOD   carbon13C
OBRQ   125.77 MHz
OBSET   7.87 kHz
OBEIN   4.21 Hz
POINT   32/167
FREQU   3980.18 Hz
SCANS   170
ACQTIM  0.8336 sec
PD      3.0000 sec
PW1     3.40 usec
IRNUC   1H
CTEMP   24.0 c
SLNT    CDCL3
EXREF   77.00 ppm
BF      1.20 Hz
RGAIN   60

```



DFILE skk-sm-1r-H-1-1.jdf
 COMNT
 DATIM 2014-02-18 18:06:01
 OBNUC H
 EXMOD proton.jdp
 OBFRQ 500.16 MHz
 OBSET 2.41 kHz
 OBFIN 6.01 Hz
 POINT 16384
 FREQU 9384.38 Hz
 SCANS 8
 ACQTM 1.7459 sec
 PD 5.0000 sec
 PW1 5.55 usec
 INUC 1H
 CTEMP 21.1 C
 SLVNT CDCL₃
 EXREF 7.26 ppm
 BF 1.00 Hz
 RGAIN 36



skk-sm-1r-C-1-1.jdf

DFILE

COMNT

DATIM

OBNUC

EXMOD

OBFRQ

OBSET

OBFIN

POINT

FREQU

SCANS

ACQTM

PD

PW1

IRNUC

CTEMP

SLVNT

EXREF

BF

RGAIN

carbon, kp

125.77 MHz

7.87 kHz

4.24 Hz

32.76

39308.18 Hz

0.8336 sec

3.0000 sec

3.40 usec

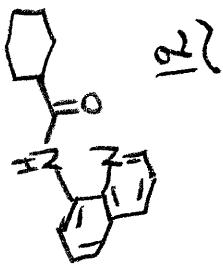
1H

CDCL₃

77.00 ppm

1.00 Hz

60



29.7376



46.8875



77.2575



116.3742



121.2101



127.5154



127.4387



127.9156



134.6401



136.3379



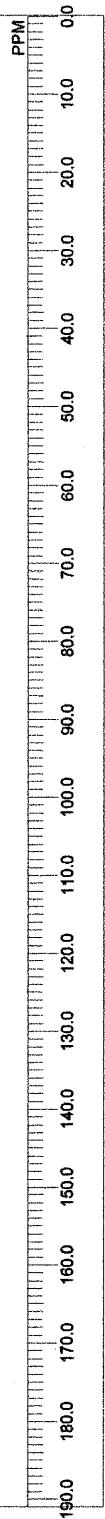
138.4745



148.0701



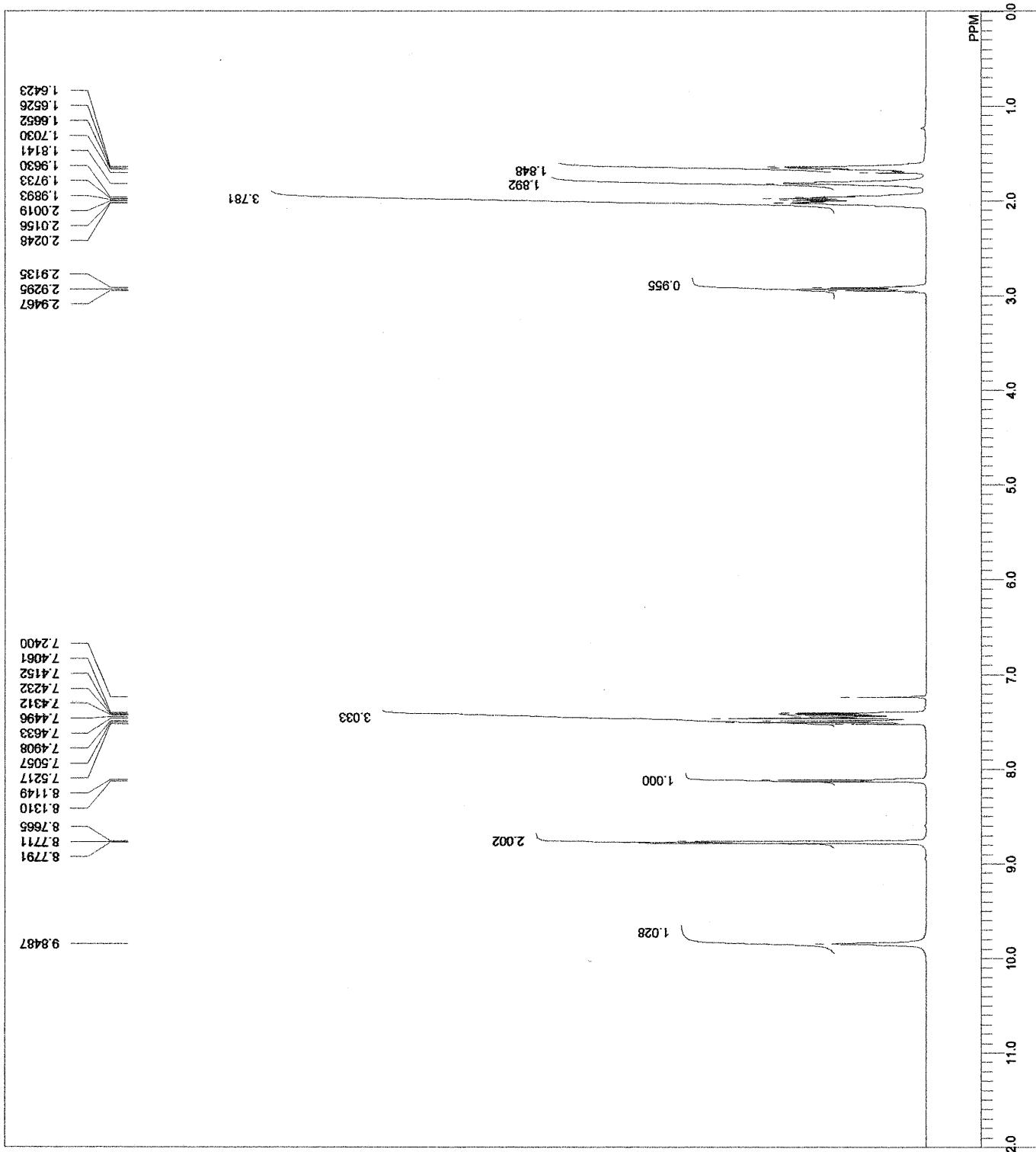
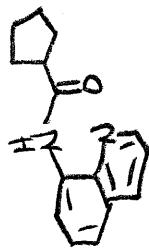
174.8632



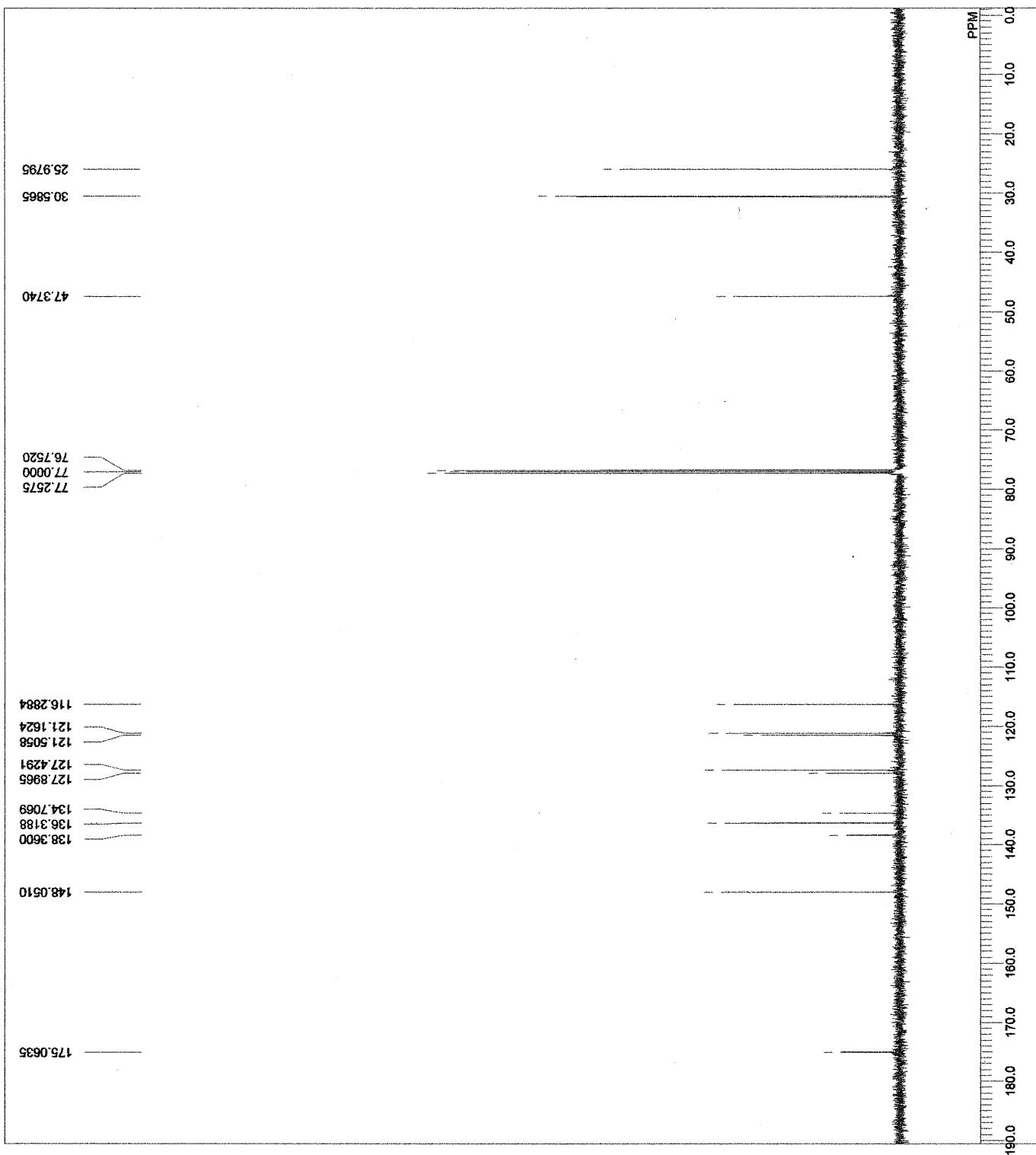
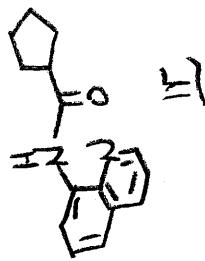
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DFILE      skk-sm-1t-H-1.jdf
COMMENT    skk-sm-1t-H
DATIM     2014-02-18 18:26:30
OBNUC
EXMOD
OBFRQ   500.16 MHz
OBSET   2.41 kHz
OBFIN   6.01 Hz
POINT   16384
FREQU   9344.38 Hz
SCANS   8
ACQTM   1.7455 sec
PD      5.0000 sec
PW1     5.55 usec
IRNUC
CTEMP
SLVNT
EXREF
BF
RGAIN
CDCL3
21.5 c
7.24 ppm
1.00 Hz
36

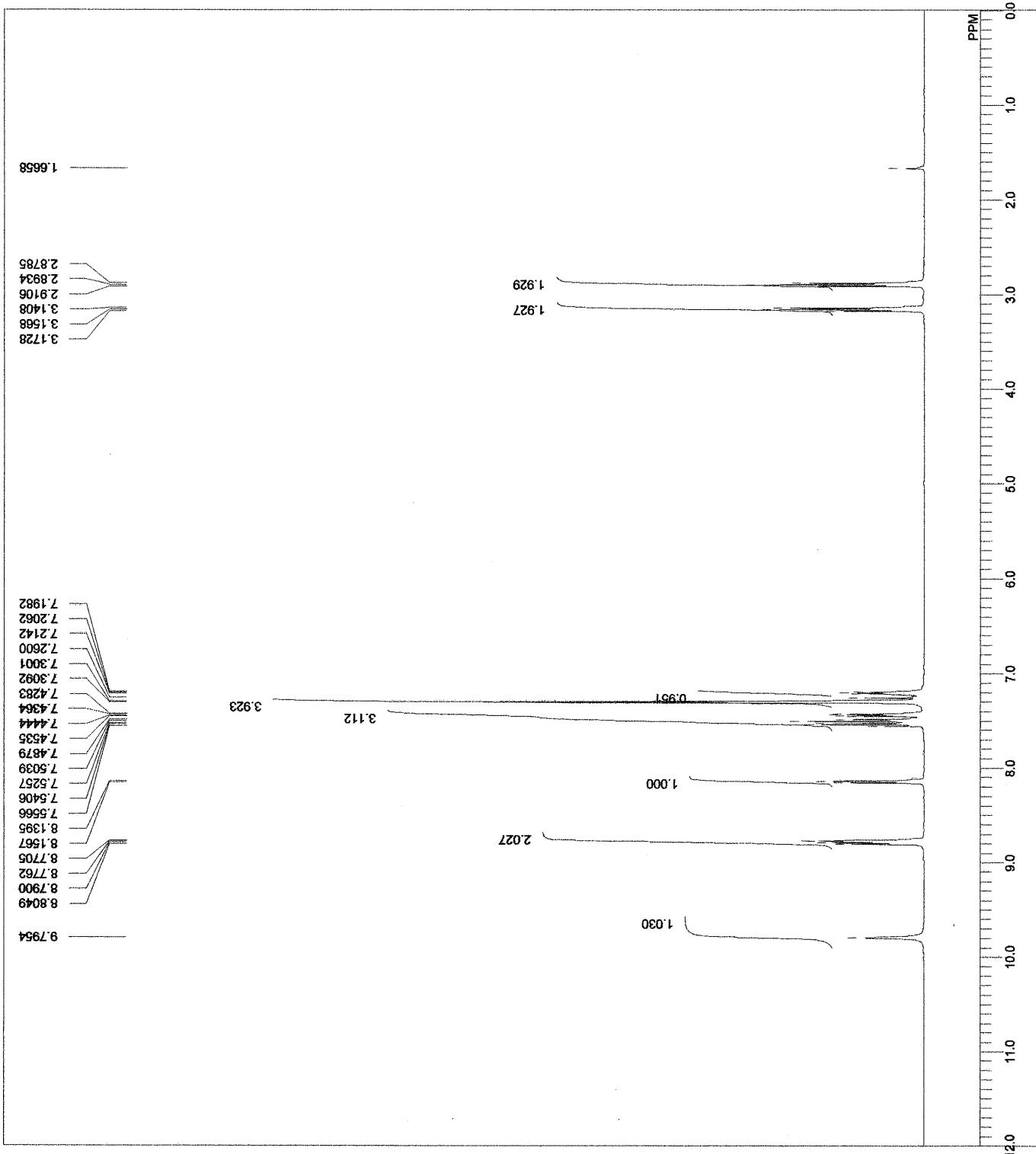
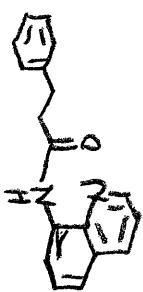
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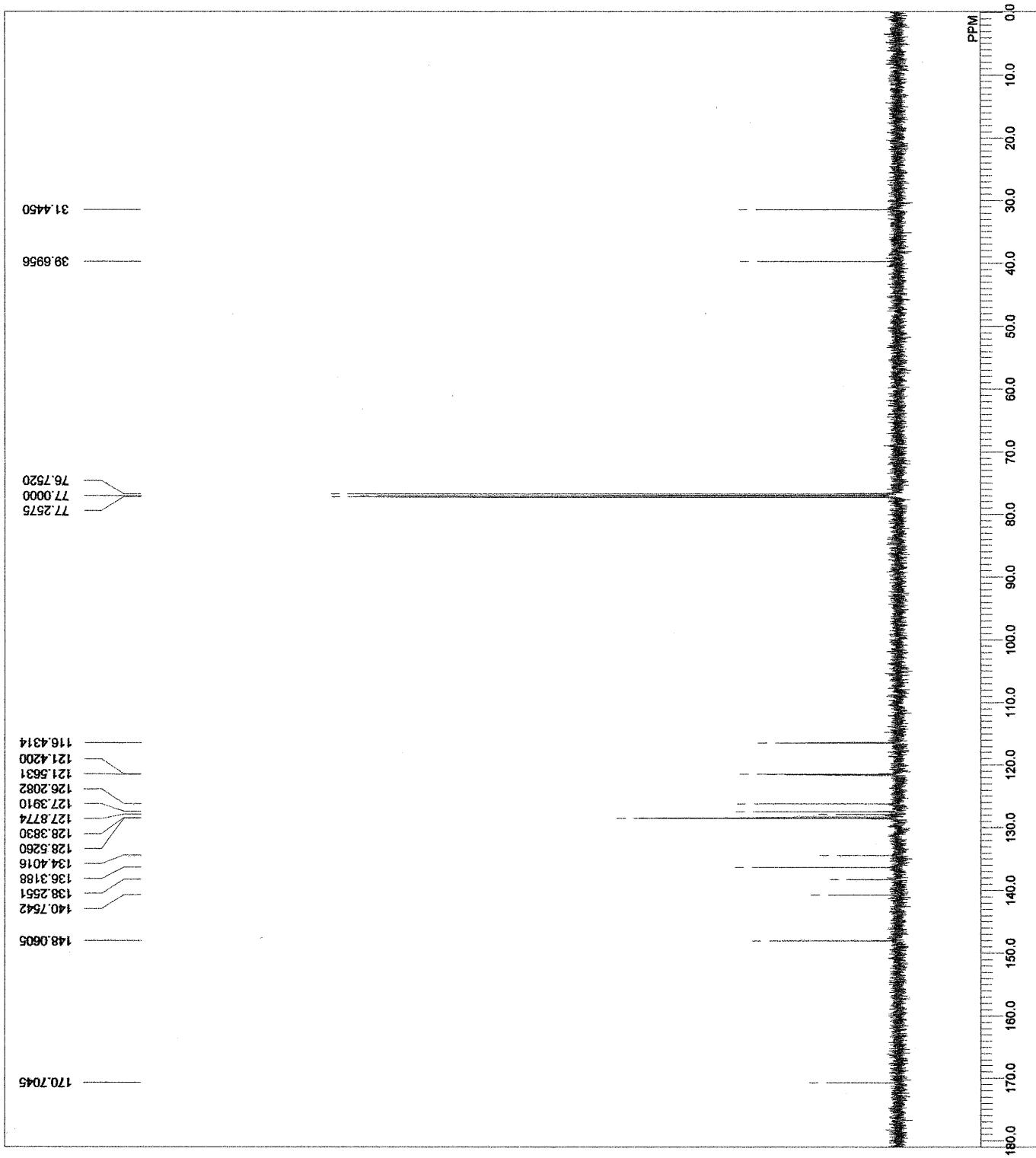
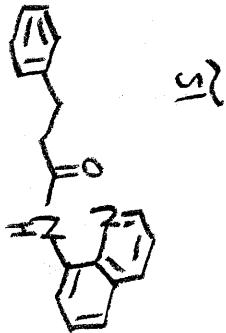
DFILE skk-sm-1f-C-1-1.als
COMNT skk-sm-1f-C
DATM 2014-02-18 18:28:22
OBNUC 13C
EXMOD carbon,kp
OBFRQ 125.77 MHz
OBFIN 7.87 kHz
POINT 4.21 Hz
FREQU 32767
SCANS 91
ACQTM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usec
IRNUC 1H
CTEMP 21.4 C
SLVNT CDCL₃
EXREF 77.00 ppm
BF 1.00 Hz
RGAIN 60



DFILE skk-sm-1s-H-1.als
 COMNT skk-sm-1s-H
 DATIM 2014-02-18 18:17:26
 OBNUC 1H
 EXMOD proton,JP
 OBFRQ 500.16 MHz
 OBSET 2.41 kHz
 OBFIN 6.01 Hz
 POINT 16384
 FREQU 9384.38 Hz
 SCANS 8
 ACQTM 1.7459 sec
 PD 5.0000 sec
 PW1 5.55 usec
 IRNUC 1H
 CTEMP 21.4 c
 SILVNT CDCL₃
 EXREF 7.26 ppm
 BF 1.00 Hz
 RGAIN 36

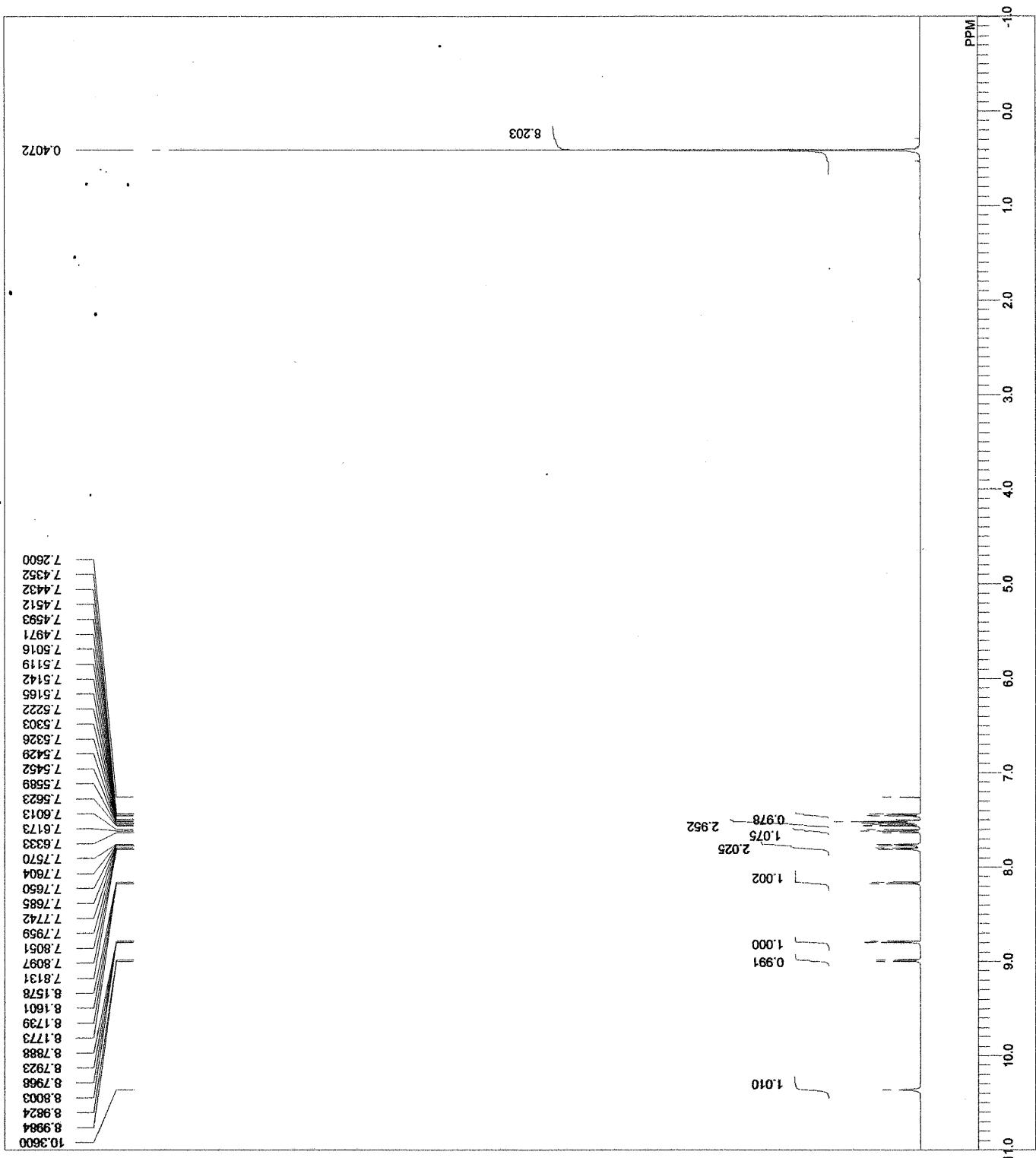


DFILE skk-sm-1s-C-1.als
 COMINT
 DATIM 2014-02-18 18:19:22
 OBNUC 13C
 EXMOD carbon,13C
 OBFRQ 125.77 MHz
 OBSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 32767
 FREQU 39308.18 Hz
 SCANS 70
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IRNUC 1H
 CTEMP 21.6 c
 SLVNT CDCL₃
 EXREF 77.00 ppm
 BF 1.00 Hz
 RGAIN 60



skk-p75-ArSiMe3

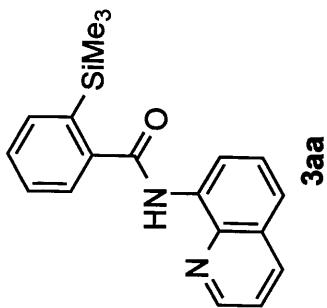
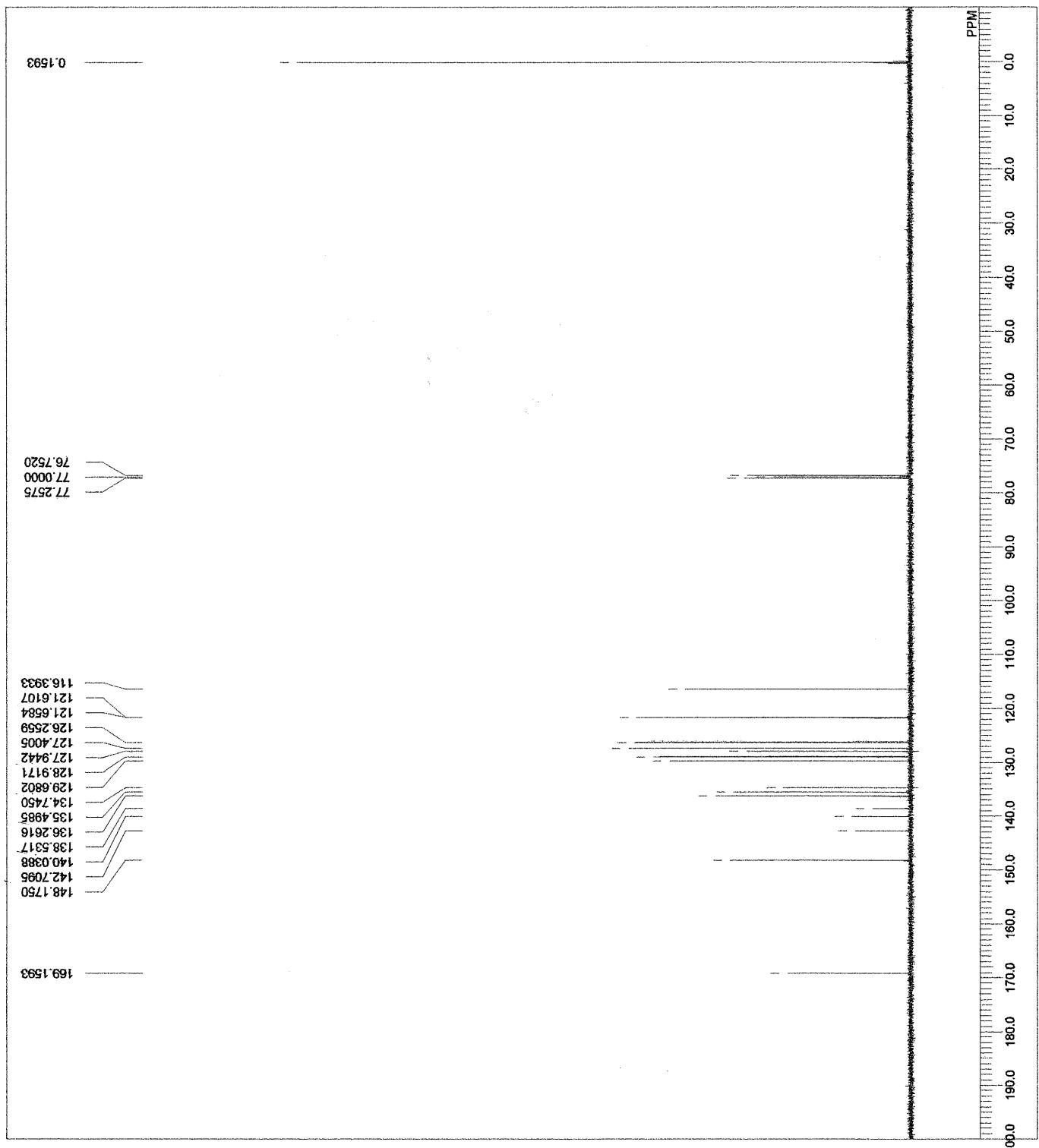
3aa



```

DFILE      skk-p75-ArSiMe3-1-1.jdf
COMMENT    skk-p75-ArSiMe3
DATIM     2013-12-10 17:24:00
OBNUC      1H
EXMOD      proton
QBFRQ     500.16 MHz
OBBET      2.41 kHz
OBIN      6.01 Hz
POINT     16384
FREQU     9384.38 Hz
SCANS      8
ACQTM      1.7459 sec
PD        5.0000 sec
PW1       5.35 usec
IRNUC      1H
CTEMP      21.4 c
SLVNT      CDCl3
EXREF      7.26 ppm
BF        0.12 Hz
RGAIN      28

```

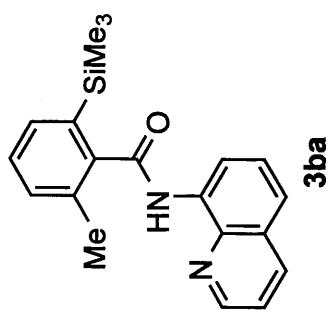


skk-p202-SI-H 3ba

This figure displays a proton NMR spectrum (1H NMR) with the x-axis representing chemical shift in PPM, ranging from 0.0007 to 10.0. The y-axis represents intensity. Key peaks are labeled with their chemical shifts and relative intensities:

- 9.9604 (1.025)
- 9.0007 (1.000)
- 8.7293 (0.996)
- 8.7350 (0.998)
- 8.9858 (0.999)
- 8.1876 (1.000)
- 8.1739 (1.000)
- 7.6379 (1.000)
- 7.6219 (1.000)
- 7.5623 (1.000)
- 7.5772 (1.000)
- 7.6070 (1.000)
- 7.4467 (1.000)
- 7.4558 (1.000)
- 7.4386 (1.000)
- 7.4306 (1.000)
- 7.3688 (1.000)
- 7.3539 (1.000)
- 7.3390 (1.000)
- 7.2955 (1.000)
- 7.2806 (1.000)
- 7.2600 (1.000)
- 2.5567 (2.831)
- 2.4617 (1.6154)
- 0.3728 (0.2778)

The spectrum shows a complex multiplet between 7.0 and 8.5 PPM, characteristic of aromatic protons, and several singlets in the aliphatic region (0.5-2.5 PPM).

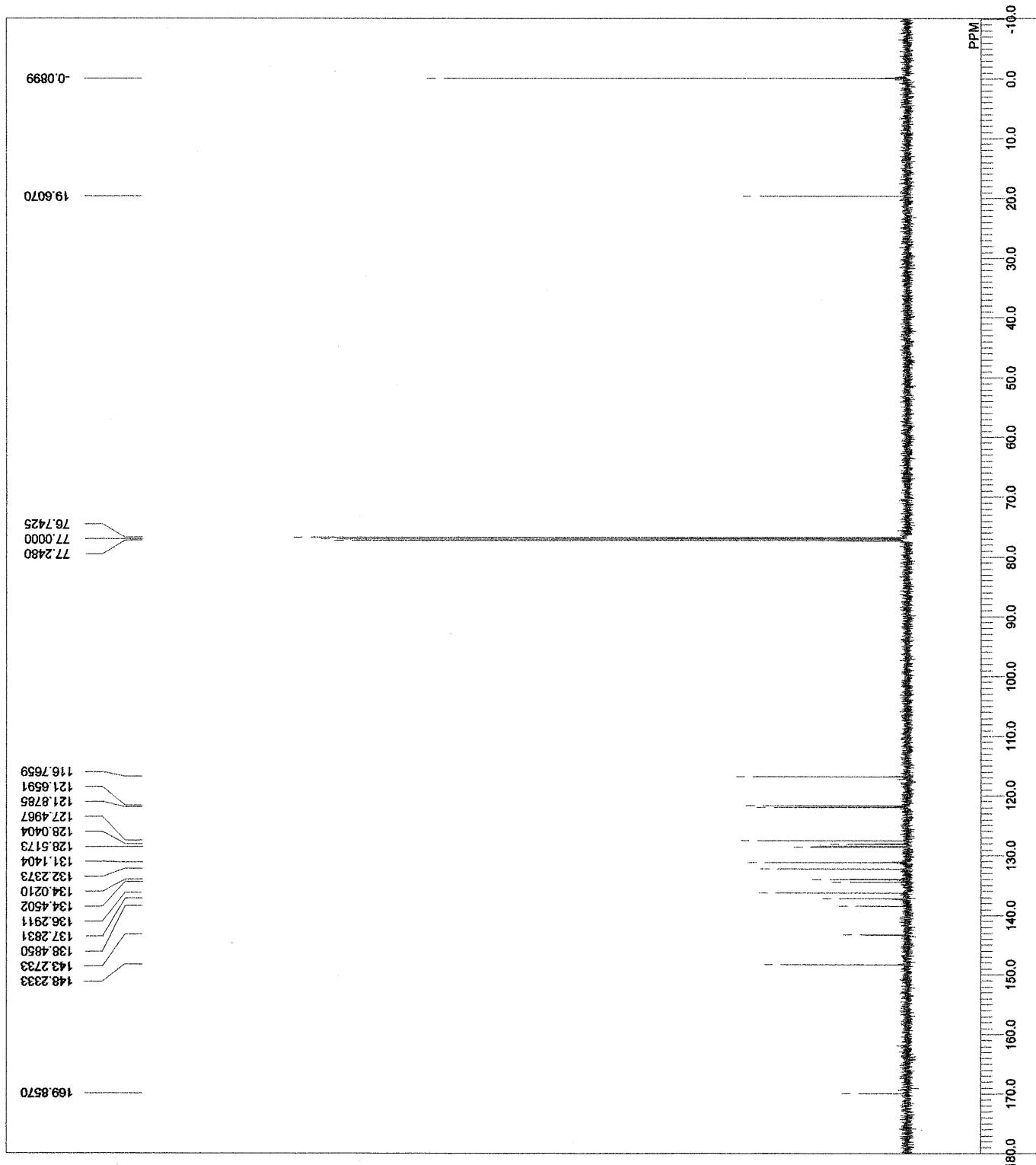
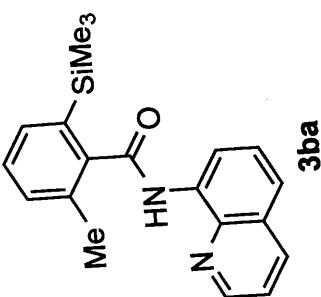


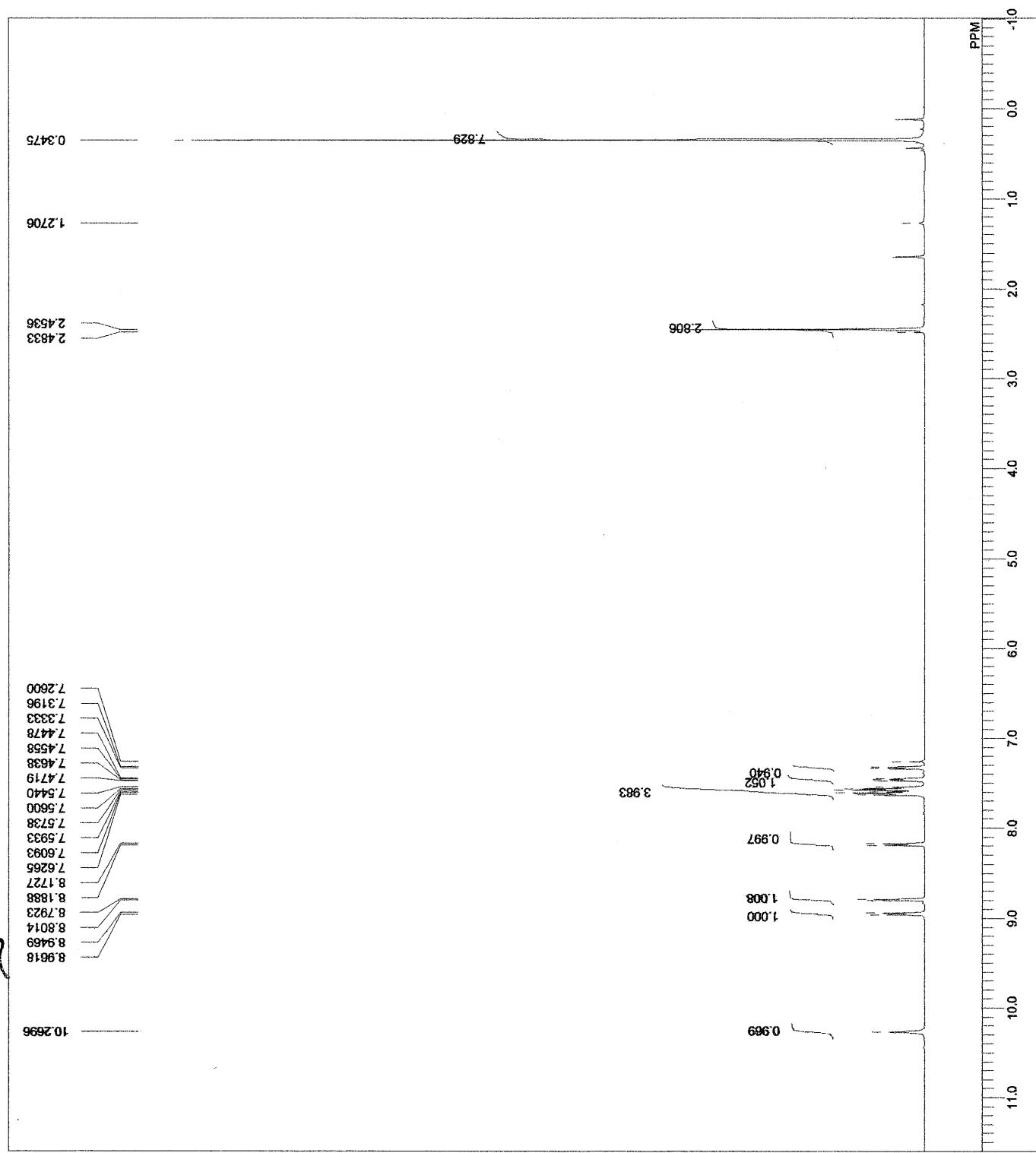
```

DFILE      skk-p202-SH-1-jdf
COMMIT     skk-p202-SH
DATIM     2013-08-12 17:50:53
DBRUC
EXMOD
FREQMOD photon xp
FRQFRQ   500.16 MHz
OFFSET    2.41 kHz
OFIN     6.01 Hz
POINT    16184
FREQU   9384.38 Hz
SCANS
ACQTM
PD      5,000,000 sec
FW1
IRNUC
CTEMP
SLVNT
EXREF
BFREF
RGAIN
CDCL3
7.26 ppm
1.00 Hz
36

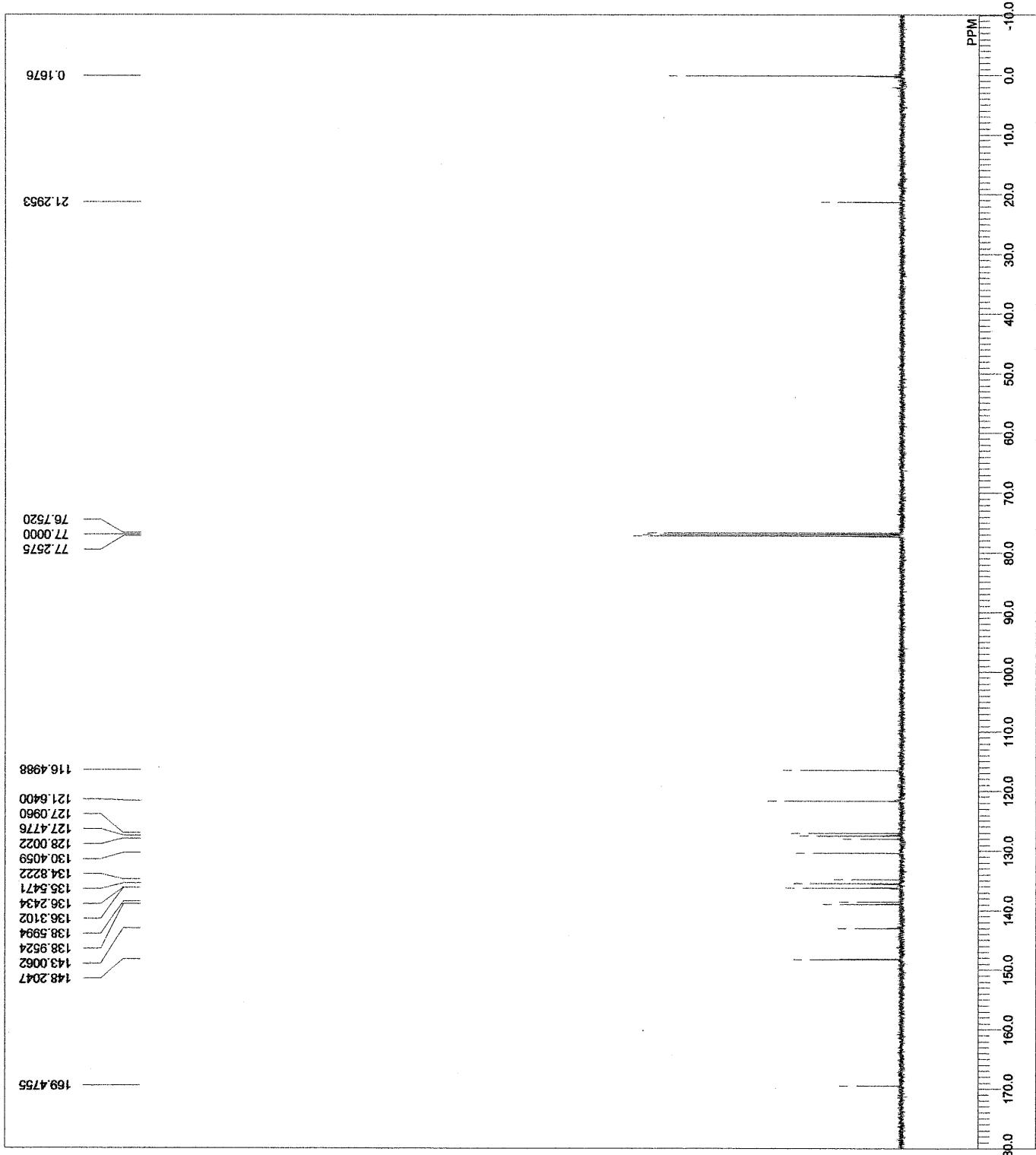
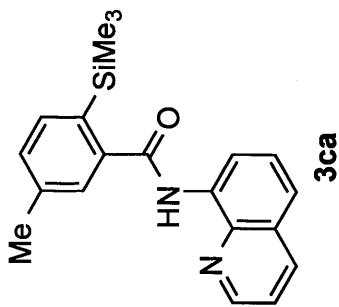
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DFILE skk-p202-SI-C-1.ais
 COMNT skk-p202-SI-C
 DATIM 2013-08-12 17:53:52
 OBNUC 13C
 EXMOD carbon,kp
 OBFRQ 125.77 MHz
 OBSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 26214
 FREQU 31446.54 Hz
 SCANS 169
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IRNUC 1H
 CTEMP 26.5 C
 SLVNT CDCl₃
 EXREF 77.00 ppm
 BF 1.00 Hz
 RGAIN 60

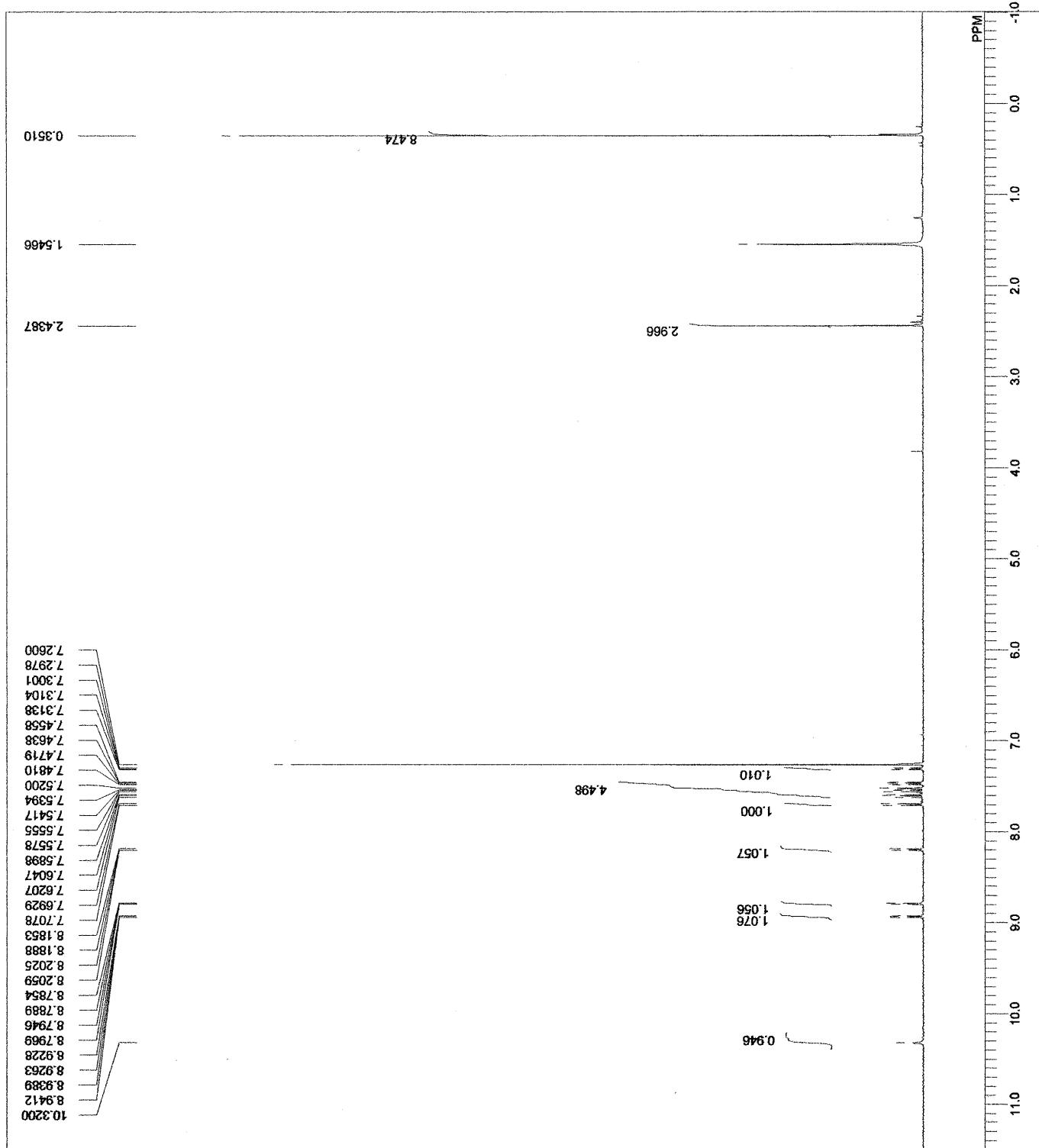
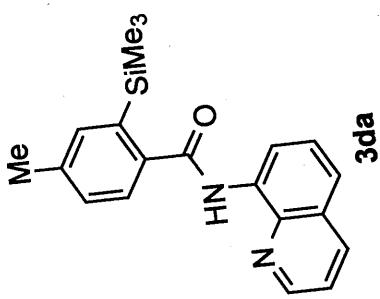




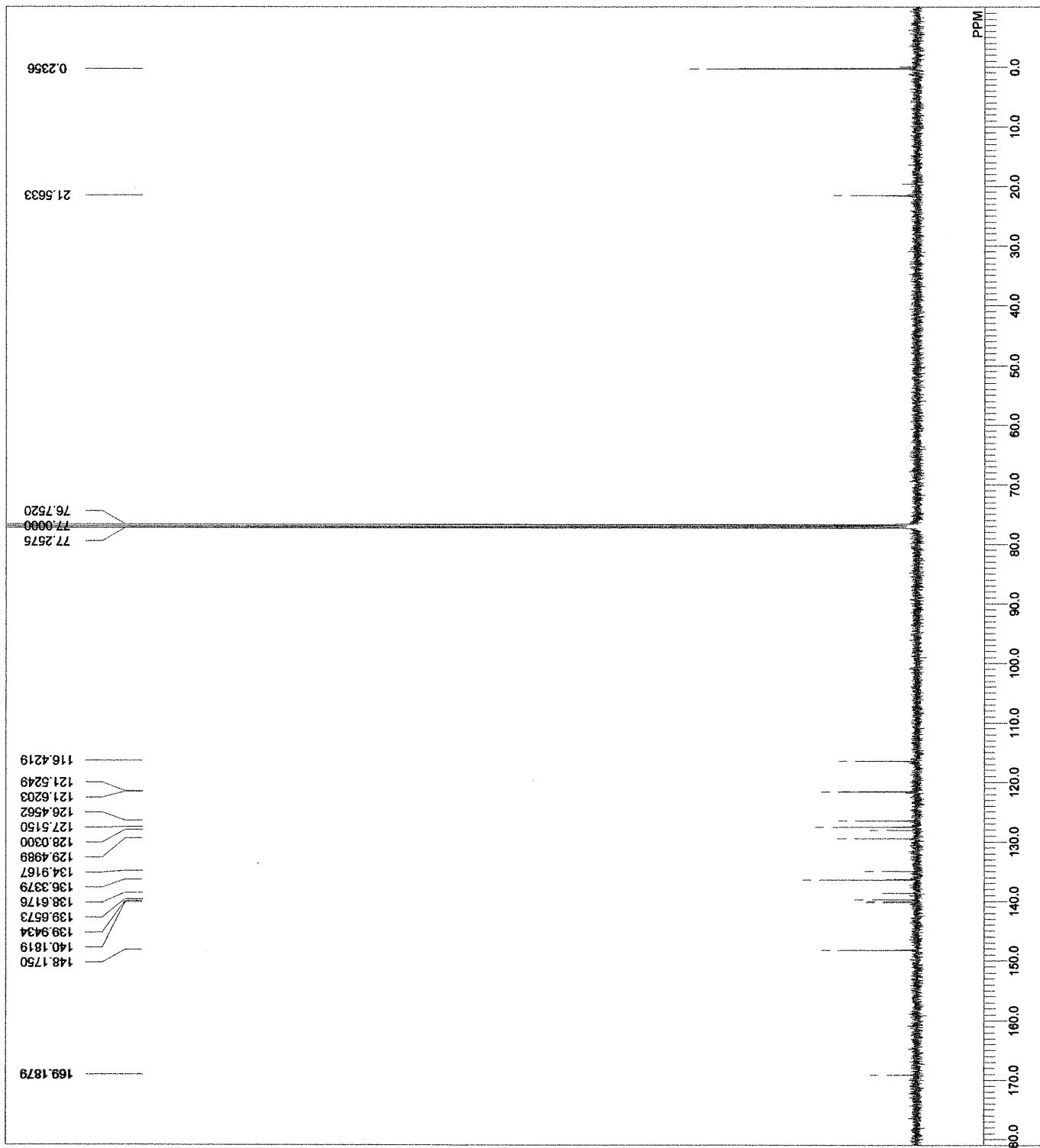
DFILE
 skk-p18-Si-Ar(m)-Me-C-1-1.als
 COMNT
 DATIM 2013-08-16 18:53:47
 OBNUC 13C
 EXMOD carbon,bq
 OBFRQ 125.77 MHz
 OBSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 26214
 FREQU 31446.54 Hz
 SCANS 102
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IRNUC 1H
 CTEMP 23.4 c
 SLVNT CDCl₃
 EXREF 77.00 ppm
 BF 1.00 Hz
 RGAIN 60

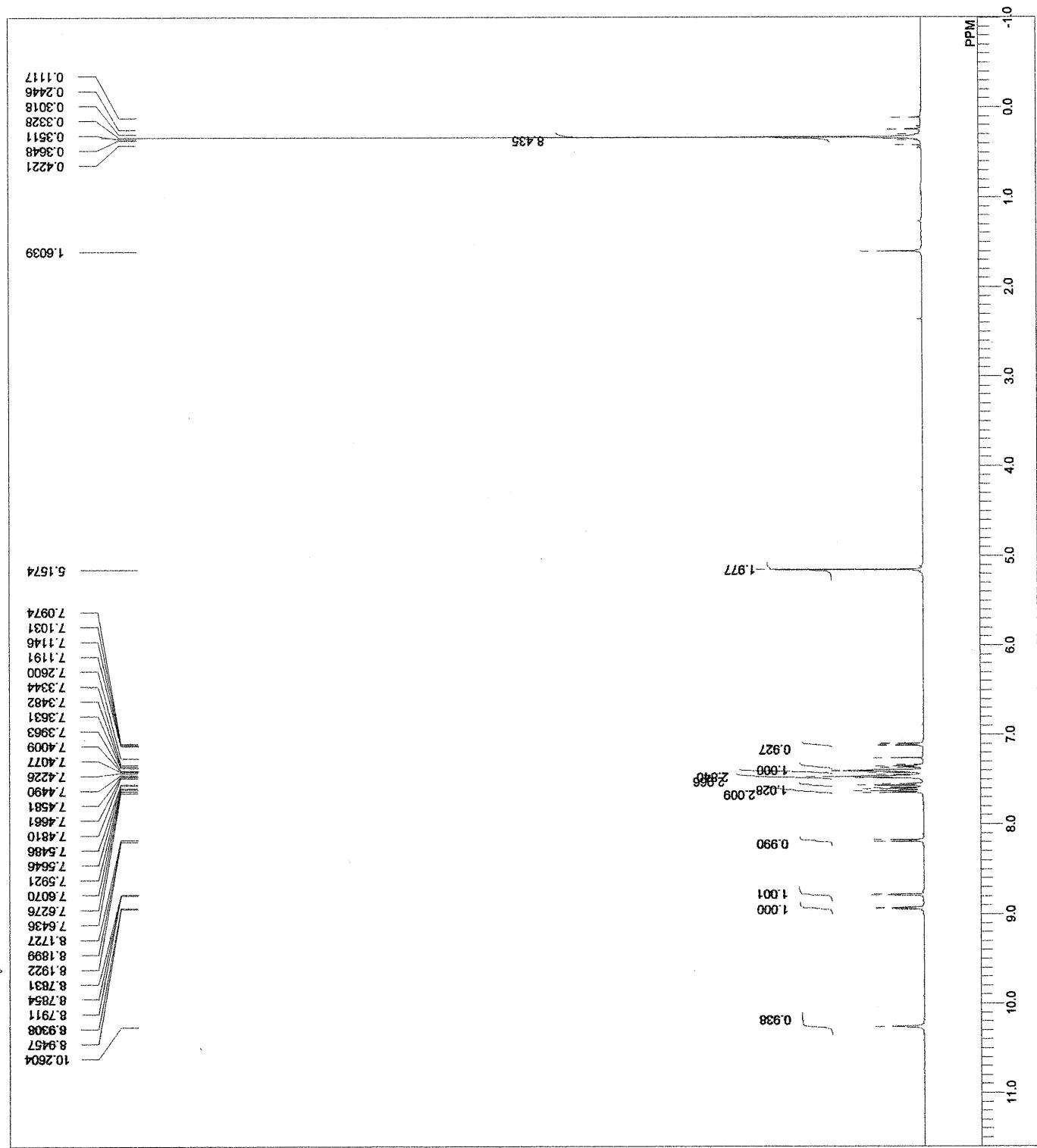


skk-ArpMe-SiMe3-H-1-1.als
 shk-ArpMe-SiMe3-H
 2013-11-18 19:56:52
 1H
 proton NMR
 500.16 MHz
 2.41 kHz
 6.01 Hz
 13107
 7507.51 Hz
 8
 1.7459 sec
 5.0000 sec
 5.55 usec
 1H
 22.4 c
 CDCl₃
 7.26 ppm
 0.12 Hz
 42

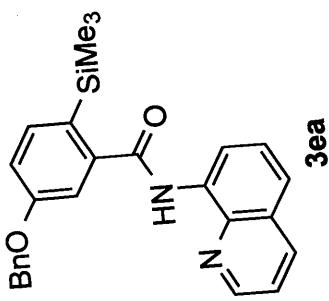


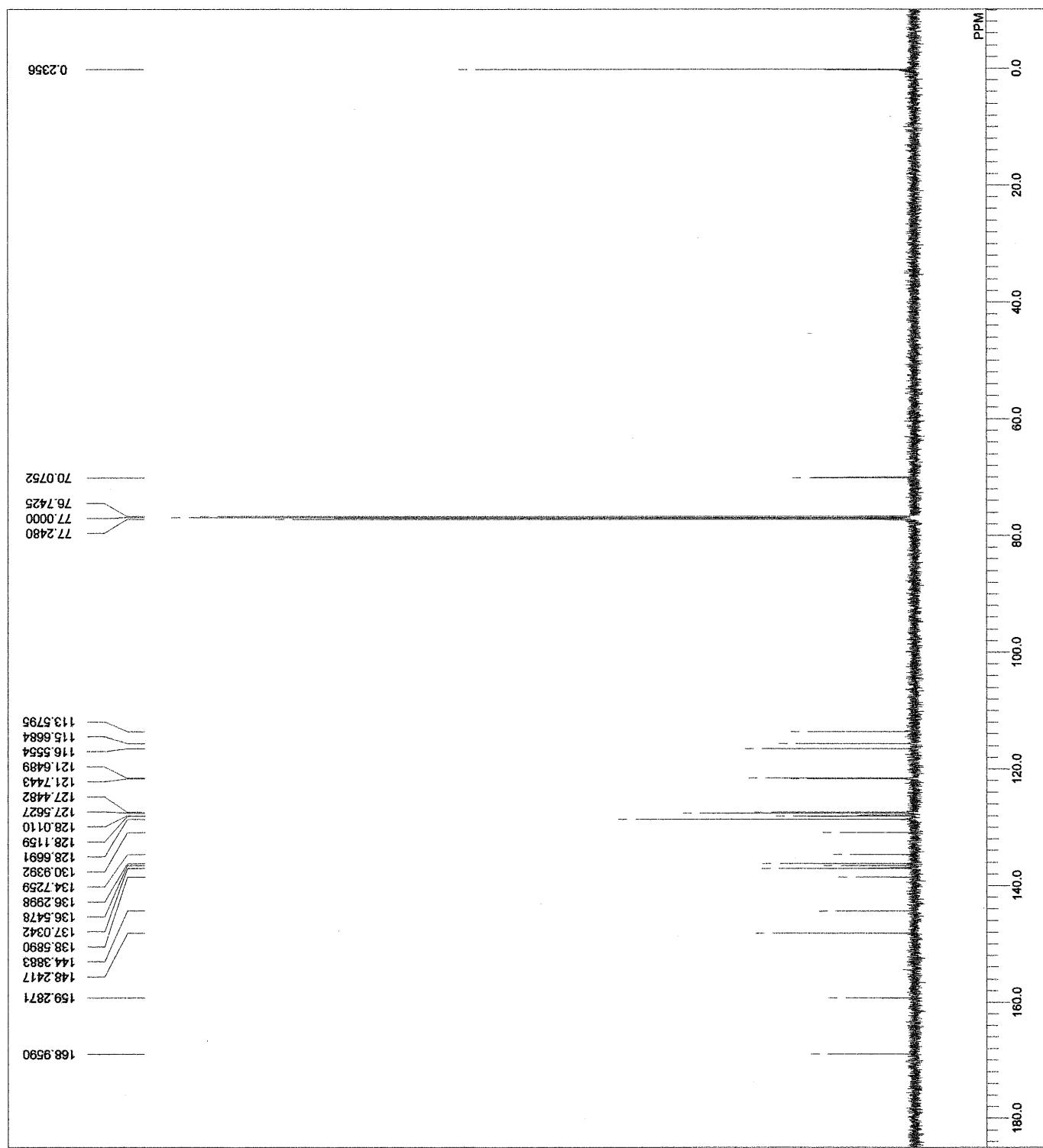
skk-C · *3d^a*



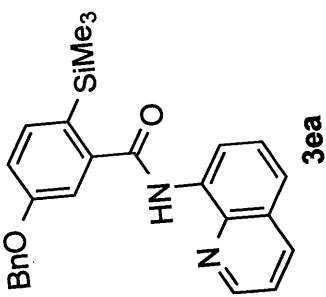


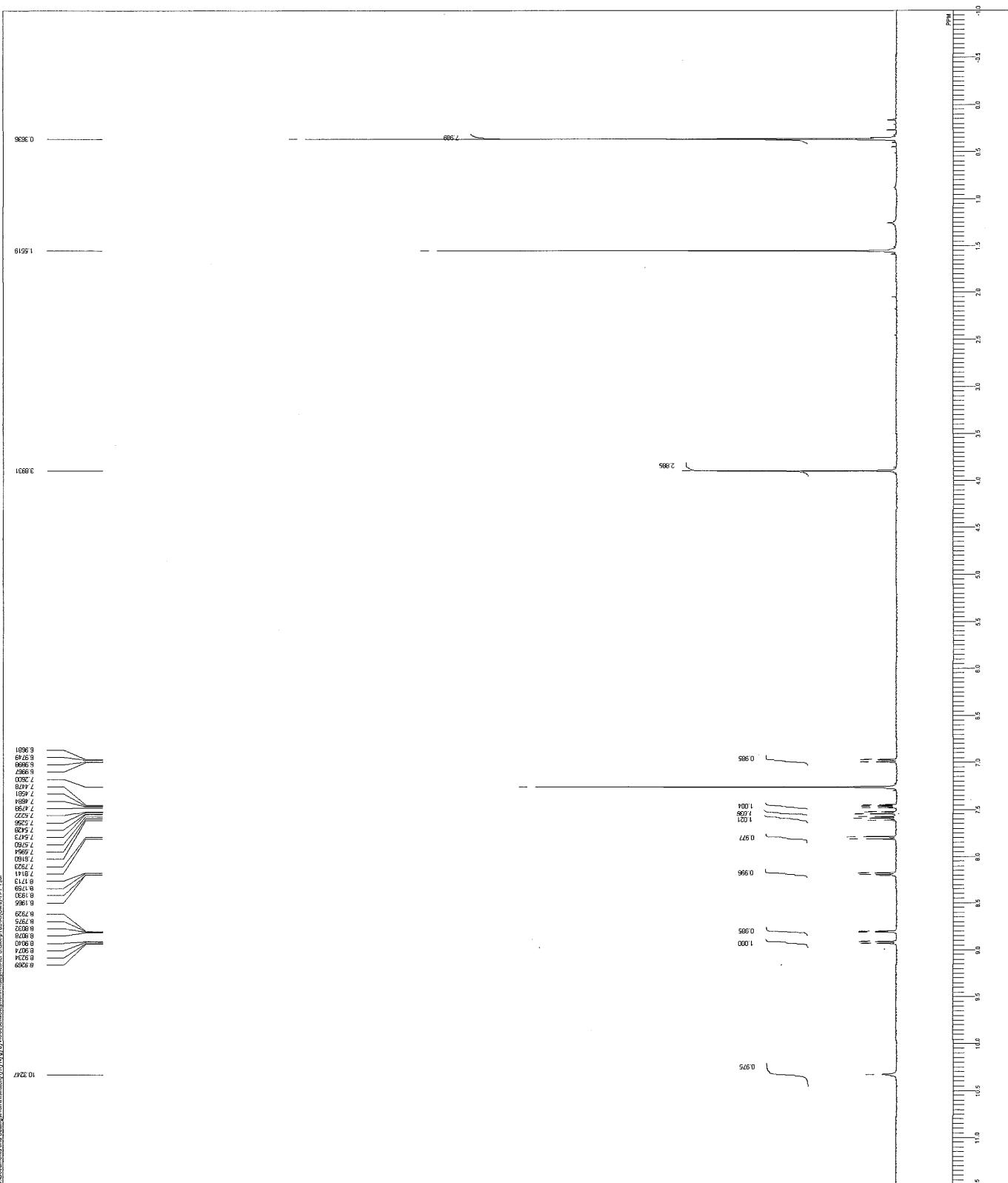
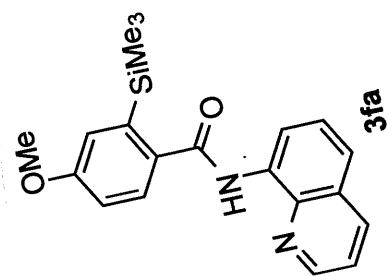
DFILE skk-p52-OBn-H-1-1.dff
COMNT skk-p52-OBn-H
DATIM 2013-10-25 15:32:11
OBNUC 1H
EXMOD proton.kdp
OBFREQ 500.16 MHz
OBSET 2.41 kHz
OBFIN 6.01 Hz
POINT 16384
FREQU 9384.38 Hz
SCANS 8
ACQTM 1.7459 sec
PD 5.0000 sec
PW1 5.55 usec
IRNUC 1H
CTEMP 25.4 c
SLVNT CDCl₃
EXREF 7.26 ppm
BF 1.00 Hz
RGAIN 30





DFILE
skk-p52-OBn-C_copy2-1-1.jdf
COMNT
2013-10-25 15:34:07
DATIM
13C
EXMOD
125.77 MHz
OBFRQ
7.87 kHz
OBSET
4.21 Hz
OBFIN
POINT
32767
FREQU
39308.18 Hz
SCANS
145
ACQTM
0.0000 sec
PD
3.0000 sec
PW1
3.40 usec
IRNUC
1H
CTEMP
25.3 c
CDCL3
77.00 ppm
SLVNT
1.00 Hz
EXREF
BF
RGAIN
60





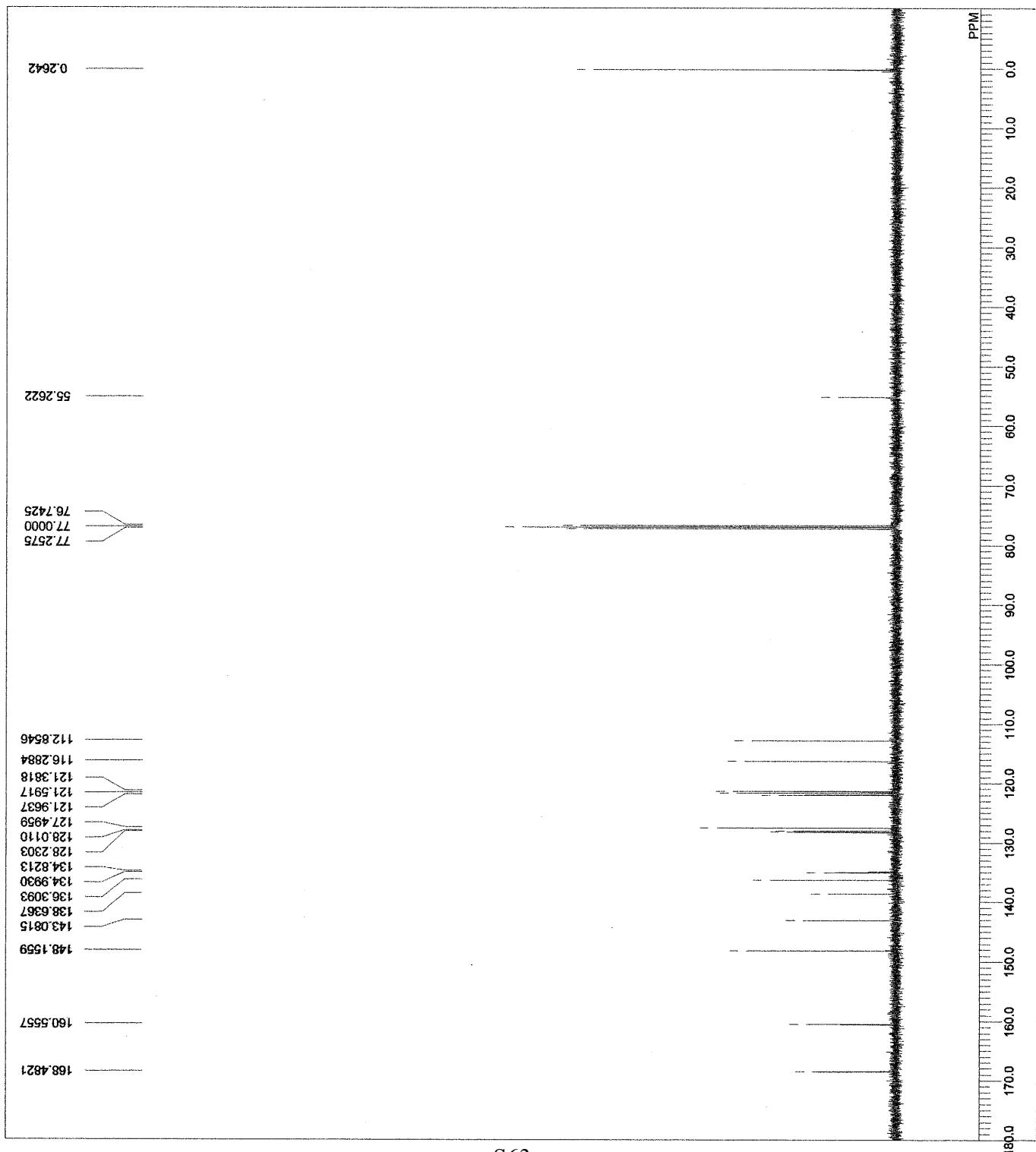
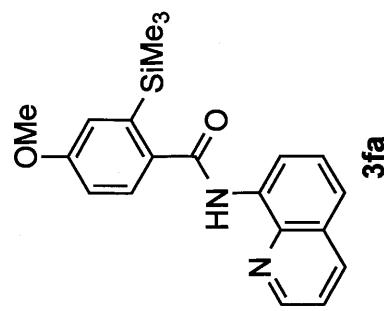
3fa

skk-p169-Ar(OMe)-C-C

```

DFILE      skk-p169-Ar(OMe)-C-1-1.jdf
COMNT
DATIM    2013-07-08 18:18:01
13C
EXMOD
OBFRQ   125.77 MHz
OBSET   7.87 kHz
OBFIN   4.21 Hz
POINT   32/167
FREQU   39308.18 Hz
SCANS   103
ACQTM
PD      0.8336 sec
3.0000 sec
PW1    3.40 usec
IRNUC
CTEMP
SLVNT
EXREF
BF      0.12 Hz
RGAIN
CDCl3
77.00 ppm
0.12 Hz
60

```

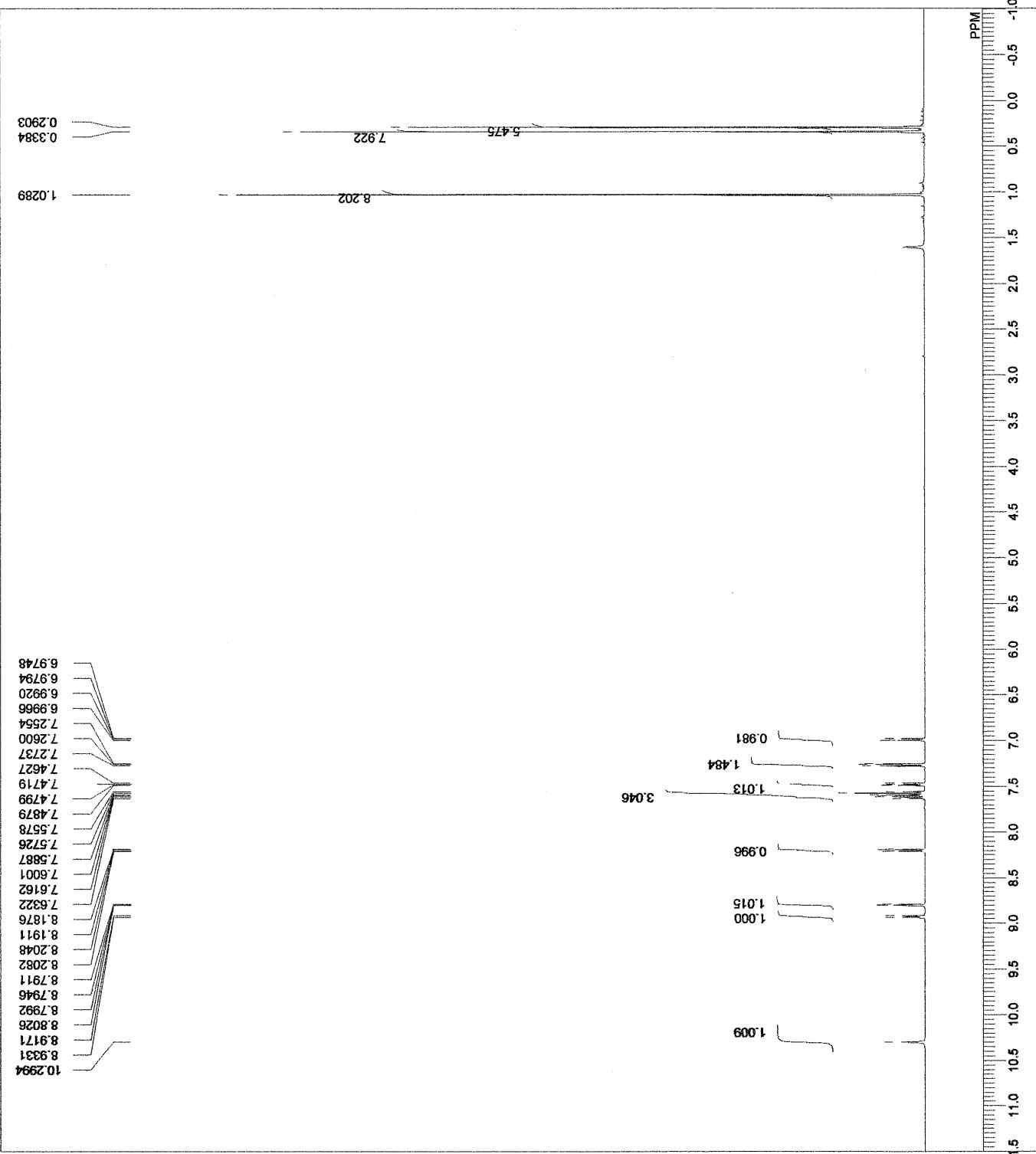
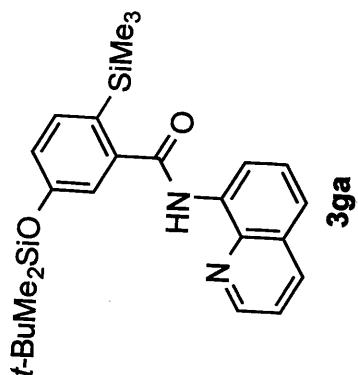


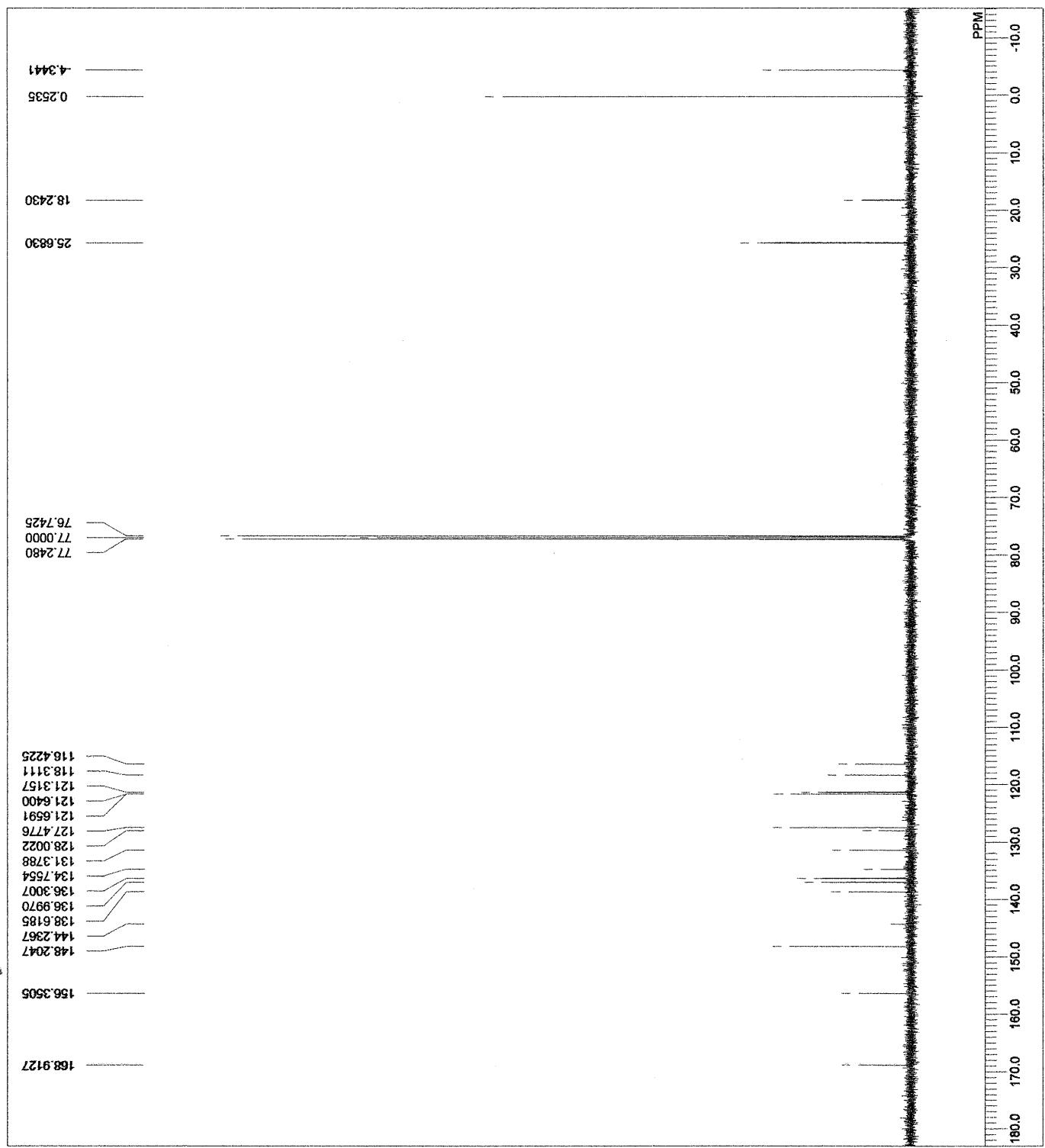
399

skk-173-ArOTBS-H

skk-173-ArOTBS-H-1-1.als

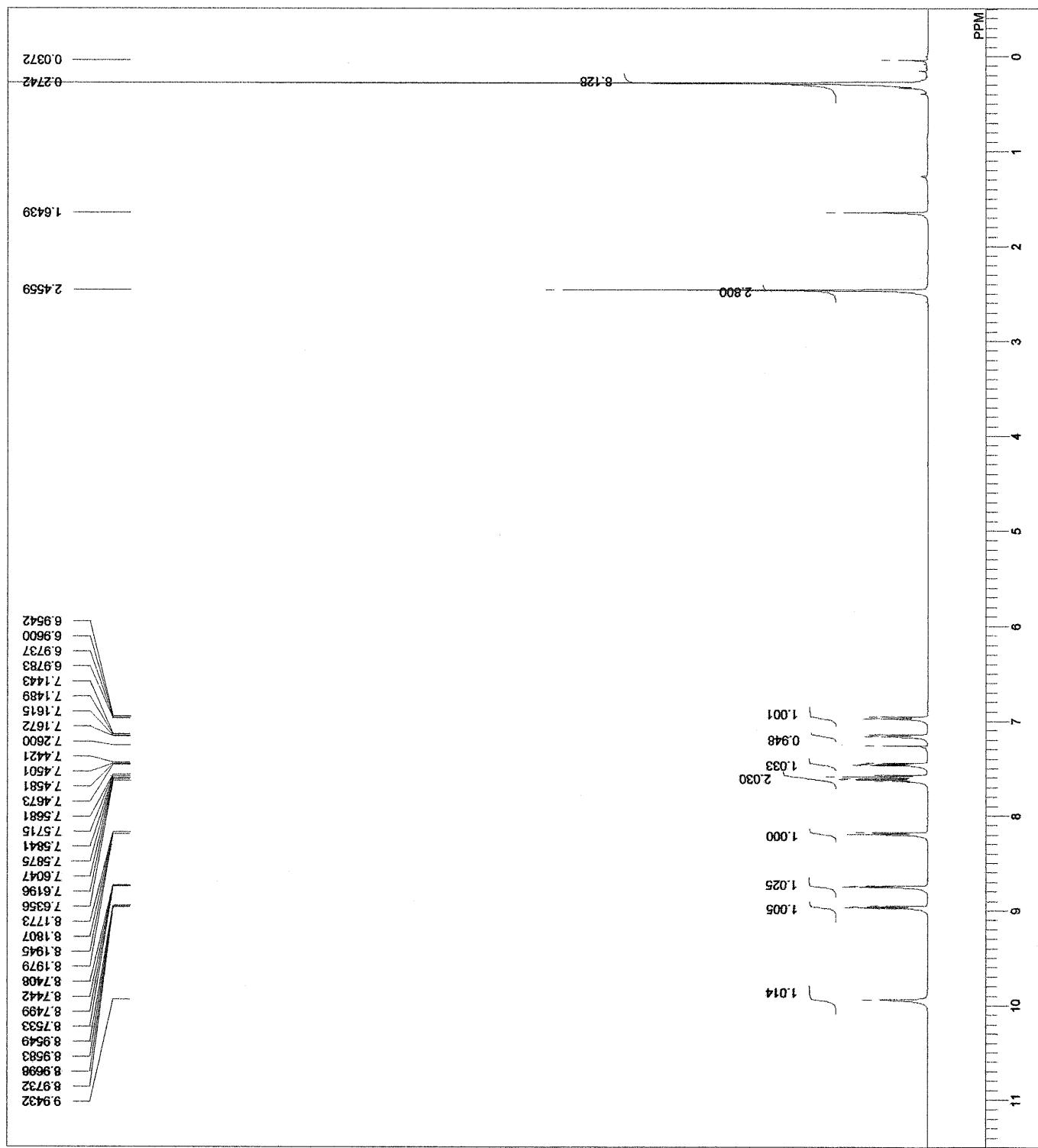
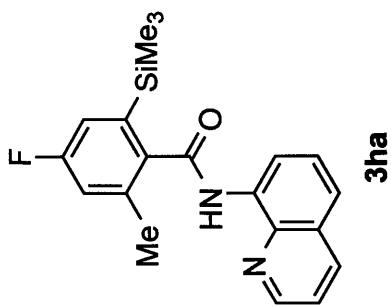
skk-173-ArOTBS-H
1H
OBNUC
EXMOD
proton,1D
500.16 MHz
2.41 kHz
6.01 Hz
13107
7507.51 Hz
16
SCANS
ACQTM
PD
OBSET
OBFIN
POINT
FREQU
SCWV
PW1
IRNUC
CTEMP
SLVNT
CDCL3
1H
1.7459 sec
5.0000 sec
5.55 usec
1H
22.9 c
0.2903
0.384
1.0289
6.9748
6.974
6.9920
6.9966
7.2554
7.2600
7.2737
7.4227
7.4719
7.4999
7.4879
7.5678
7.526
7.5887
7.6001
7.6162
7.6322
8.1876
8.1911
8.2048
8.2082
8.7911
8.7946
8.7992
8.8026
8.9171
8.9331
10.2994



3ga

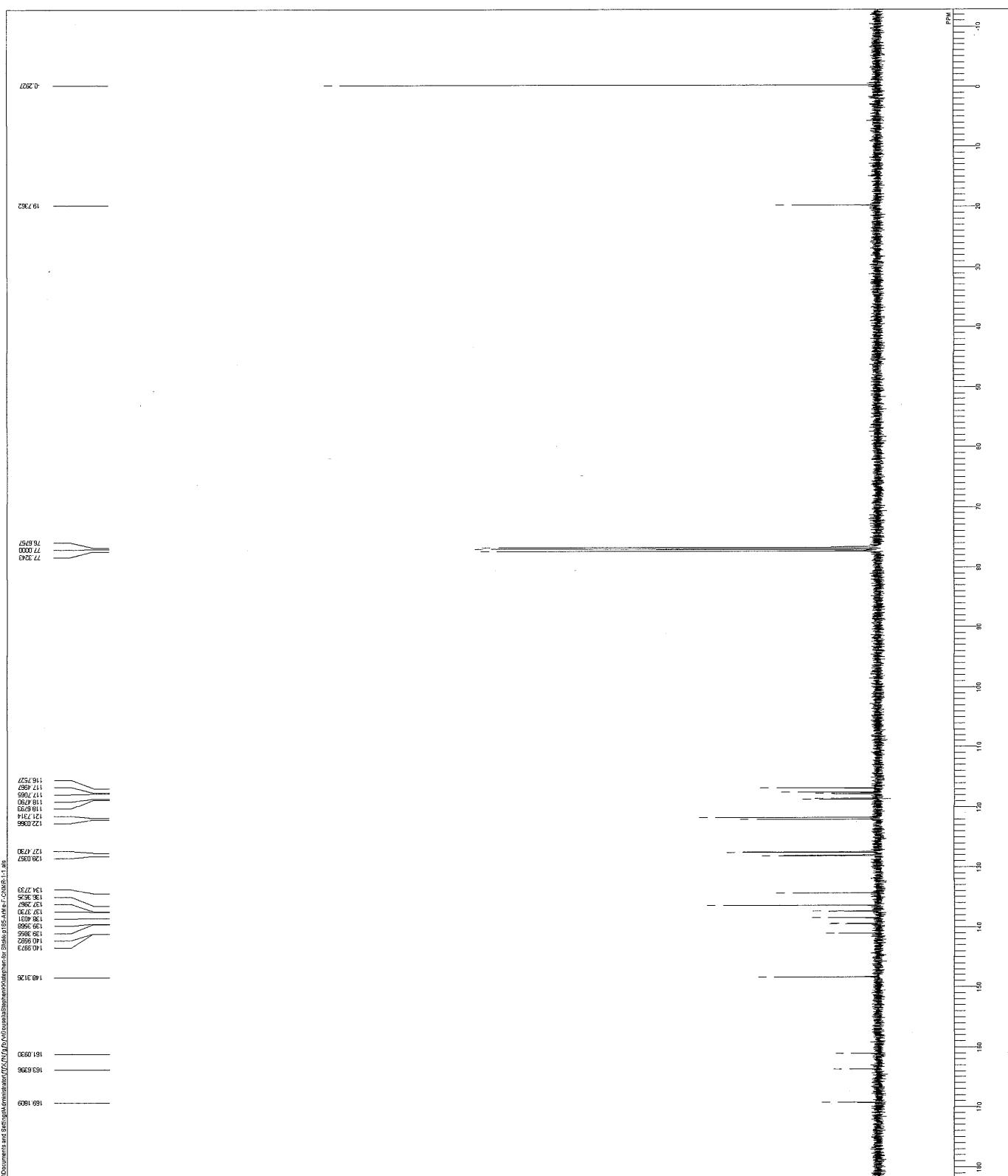
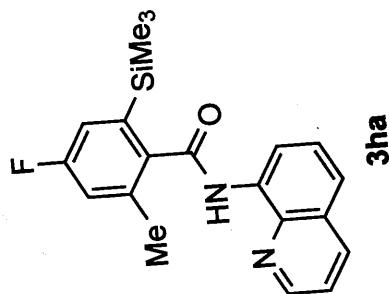
skk-p185-ArCOMe-H-1-1.als
ArneF-SMe3

DFILE
COMNT
DATIM
OBNUC
EXMOD
proton.kp
OBFRQ
2.41 kHz
OBSET
6.01 Hz
POINT
13107
FREQU
SCANS
16
ACQTM
1.7459 sec
PD
5.0000 sec
PW1
5.55 usec
IRNUC
1H
CTEMP
22.9 c
SLVNT
CDCl₃
EXREF
7.26 ppm
BF
0.12 Hz
RGAIN
34



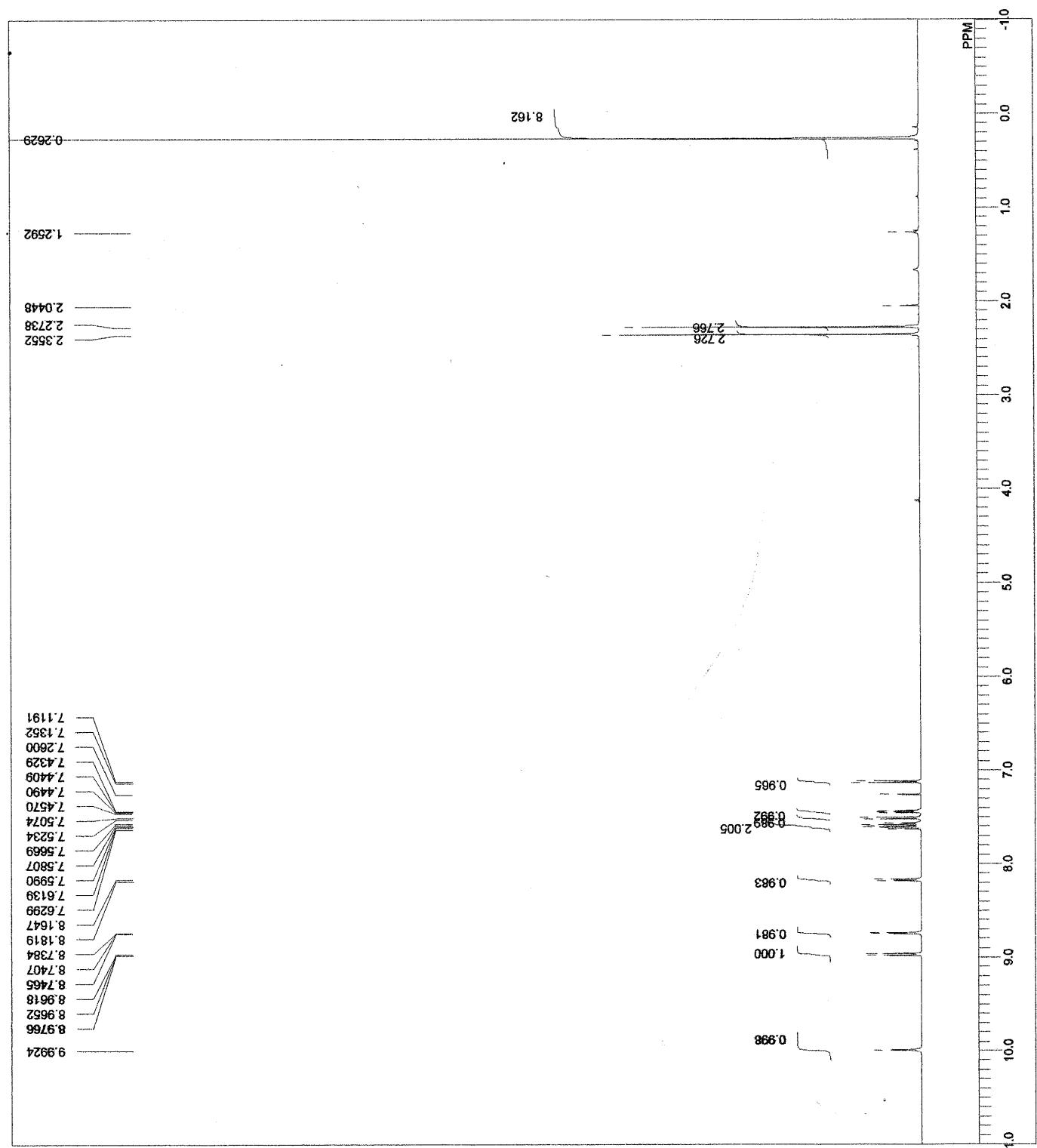
三
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skk-p18b-ArMe-H-CNMR

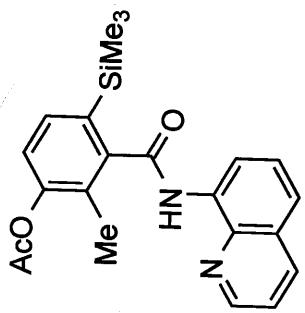


3ia

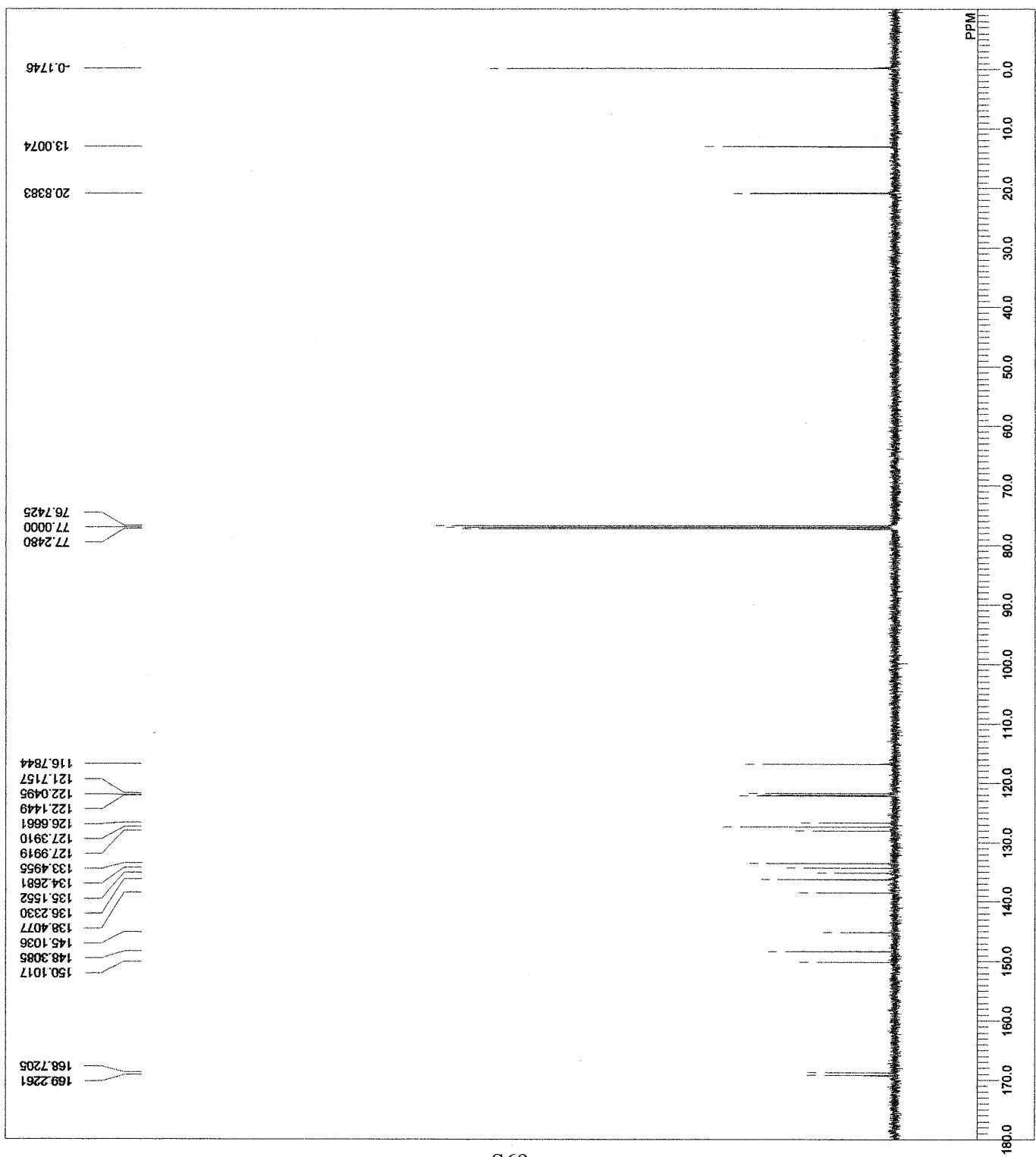
skk-p30-ArMe/OAc-Si-H

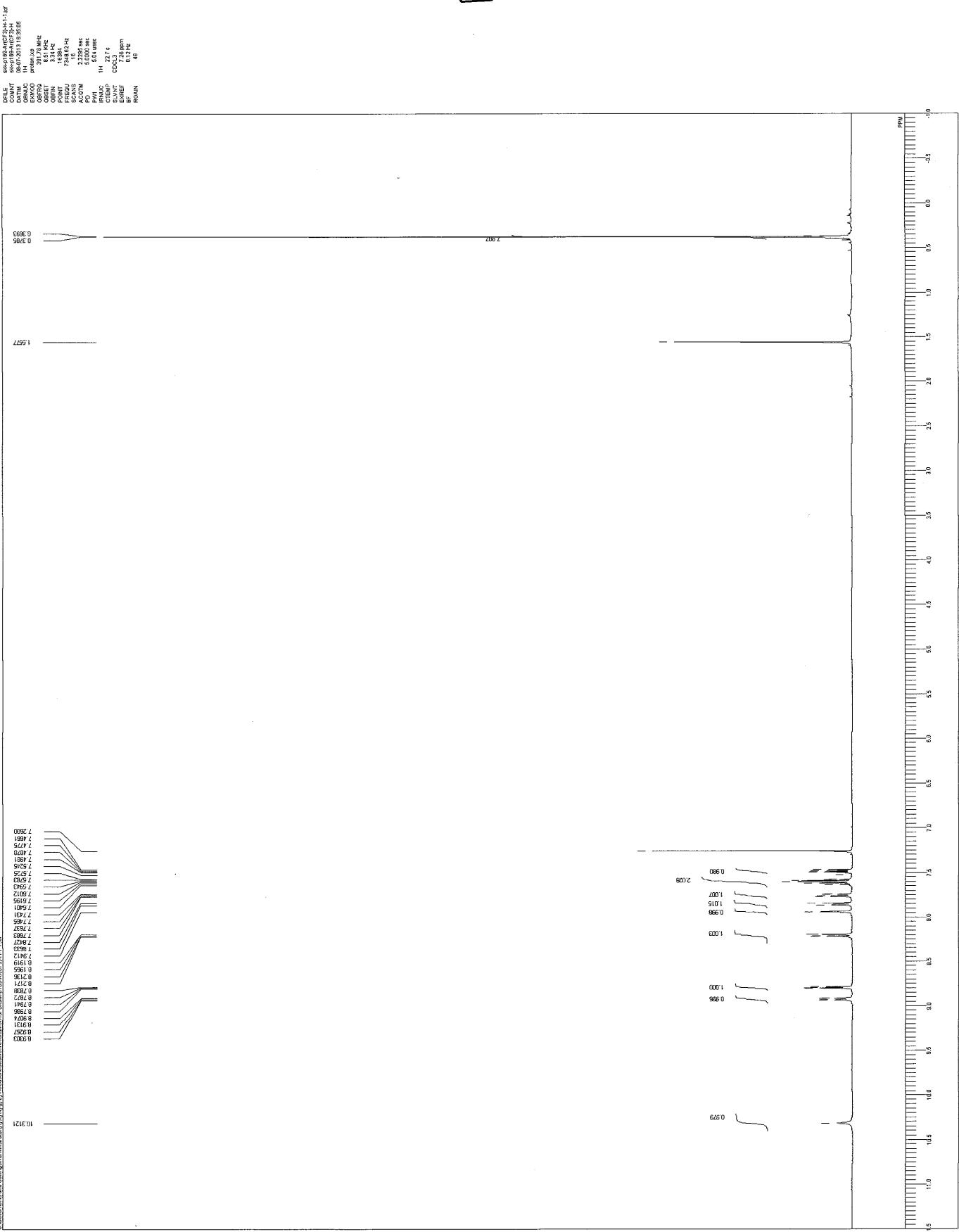
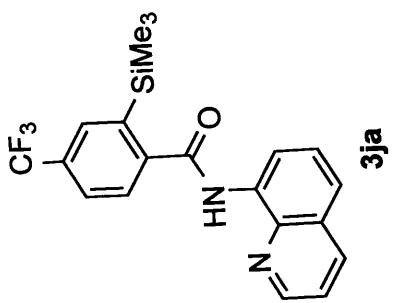


skk-p30-ArMe OAc-SiH-1-1.iqf
skk-p30-ArMe(OAc-Si-H
1H
proton.iqx
500.16 MHz
OBFRQ 2.41 kHz
OBSET 6.01 Hz
POINT 16384
FREQU 9384.38 Hz
SCANS 8
ACQTM 1.7459 sec
PD 5.0000 sec
PW1 5.55 usec
IRNUC 1H
CTEMP 24.7 c
SLVNT CDCl₃
EXREF 7.26 ppm
BF 0.12 Hz
RGAIN 36

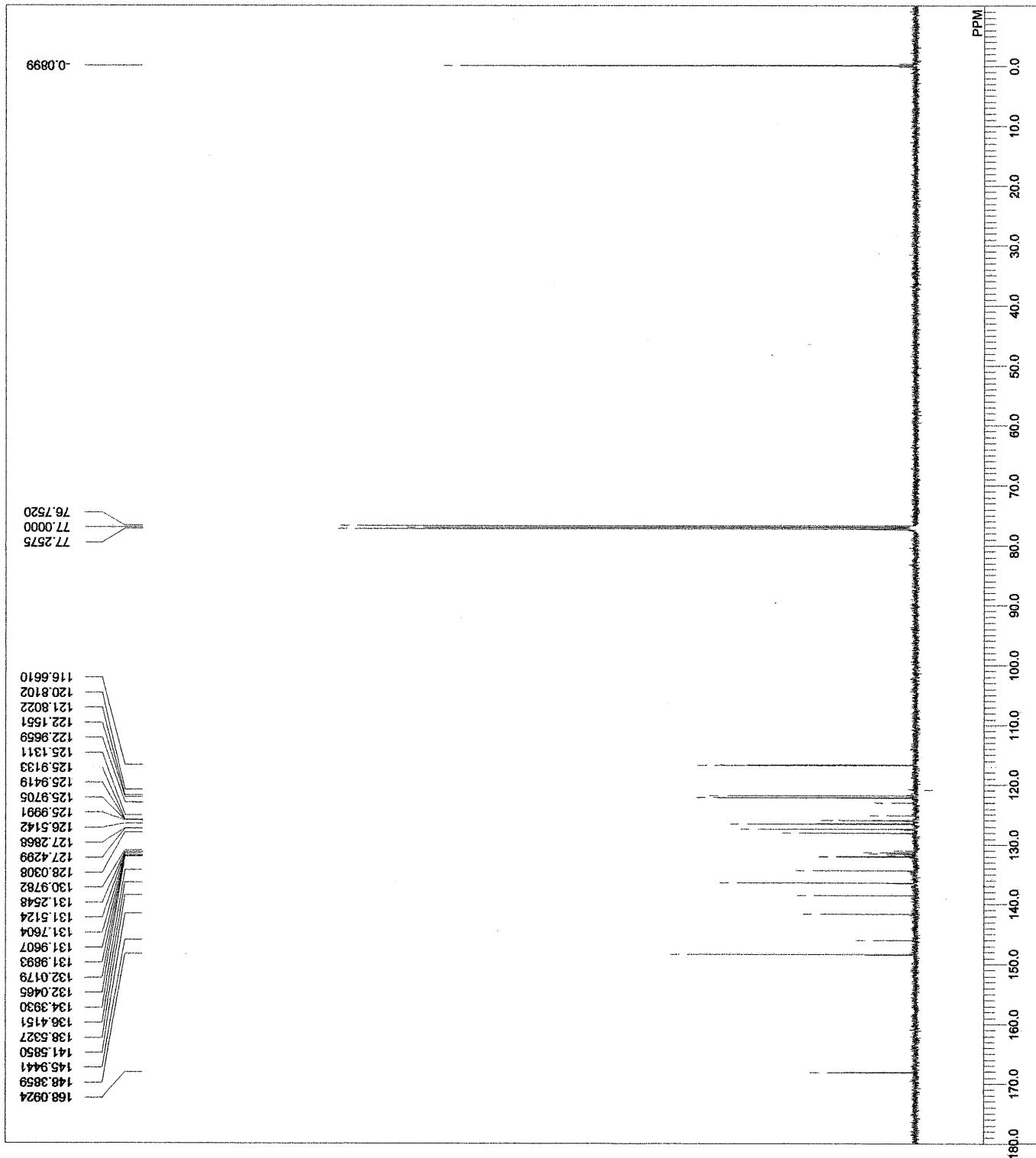
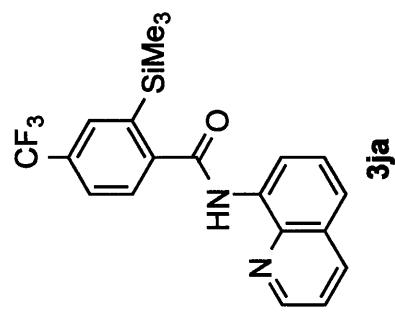


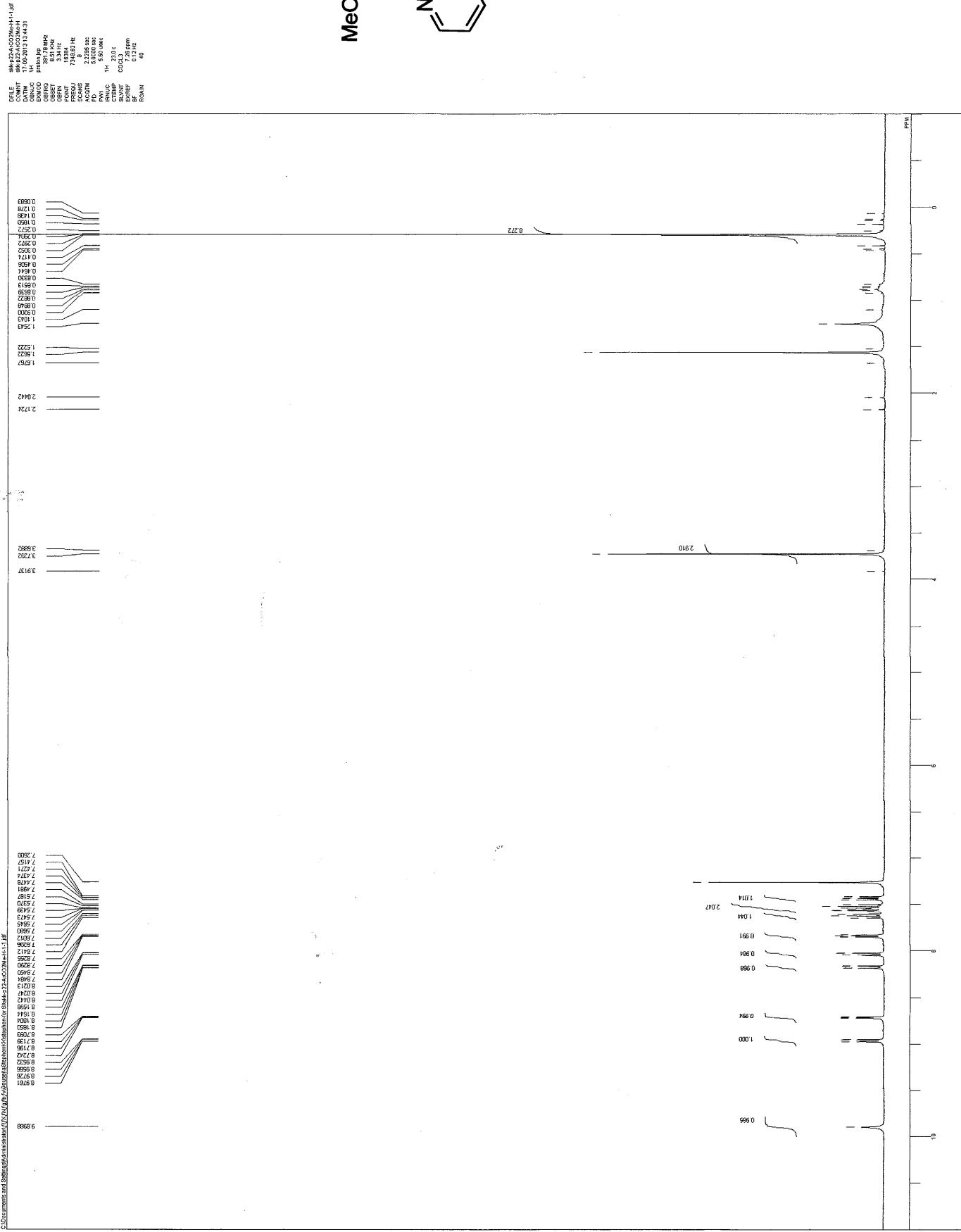
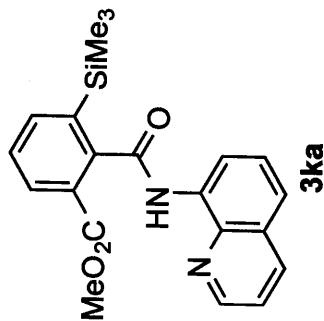
3ia



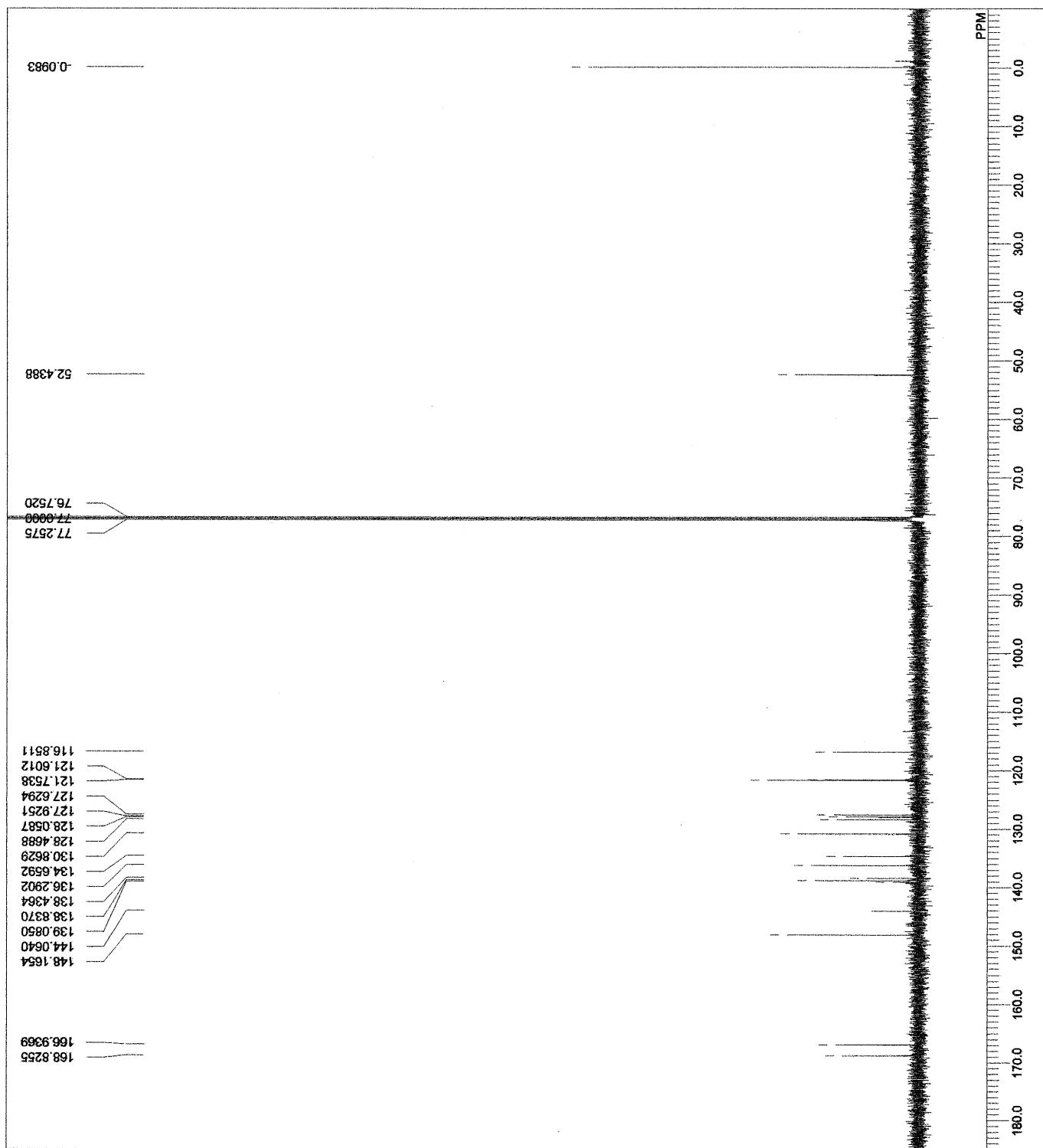
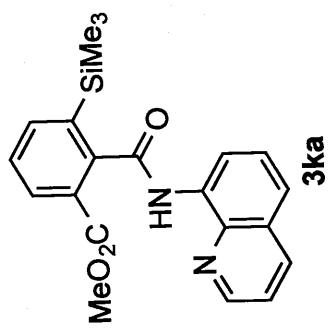


DFILE skk-p169-Ar(CF₃)-C-1-1.als
 COMINT skk-p169-Ar(CF₃)-C
 DATIM 2013-07-08 16:59:06
 OBNUC 13C
 EXMOD carbon,13C
 OBFHQ 125.77 MHz
 OBSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 26214
 FREQU 31446.54 Hz
 SCANS 387
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IRNUC 1H
 CTEMP 24.8 c
 CDCL3
 SLVNT 77.00 ppm
 EXREF 1.00 Hz
 BF 60
 RGAIN

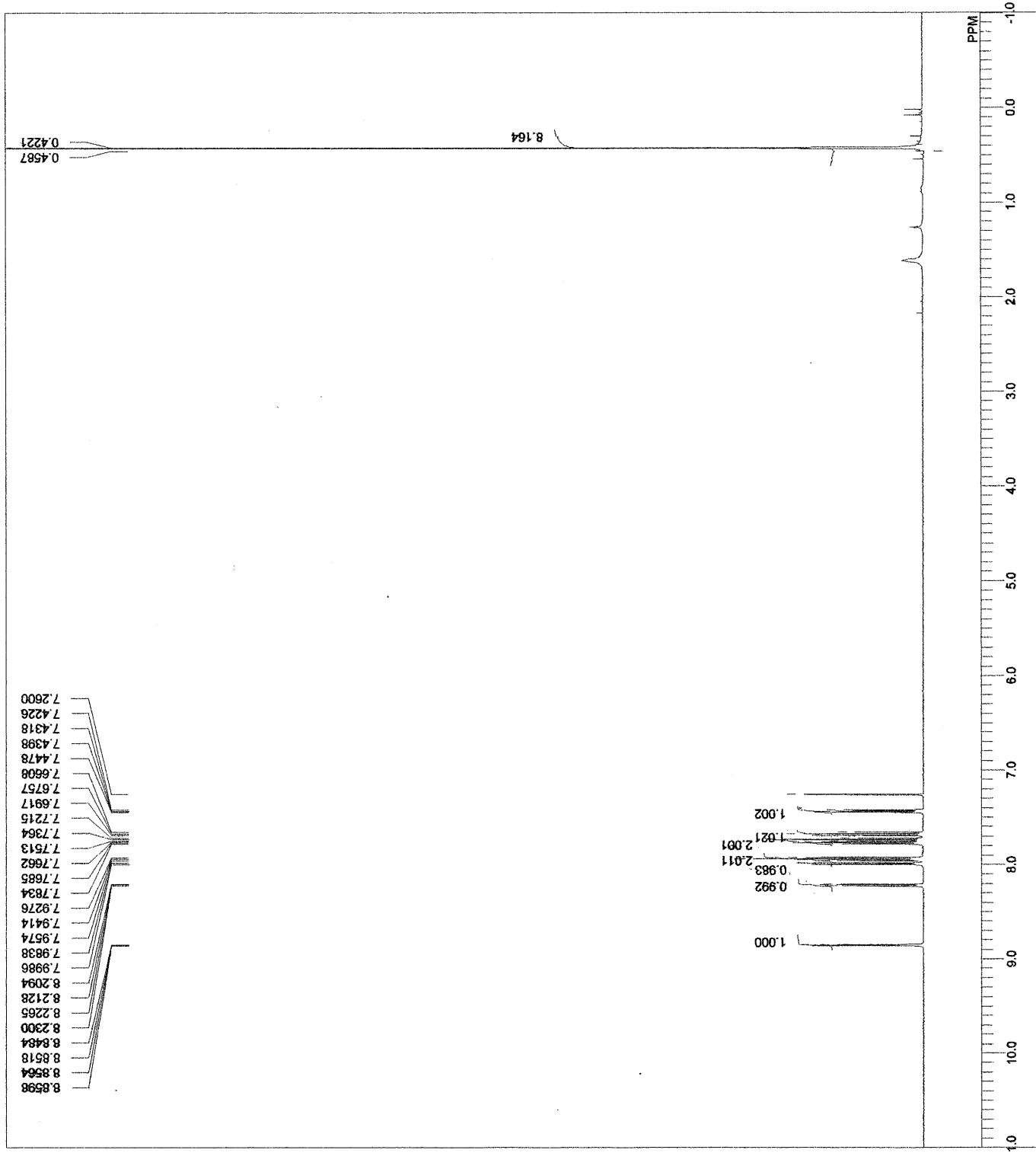
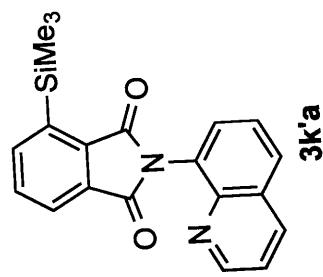


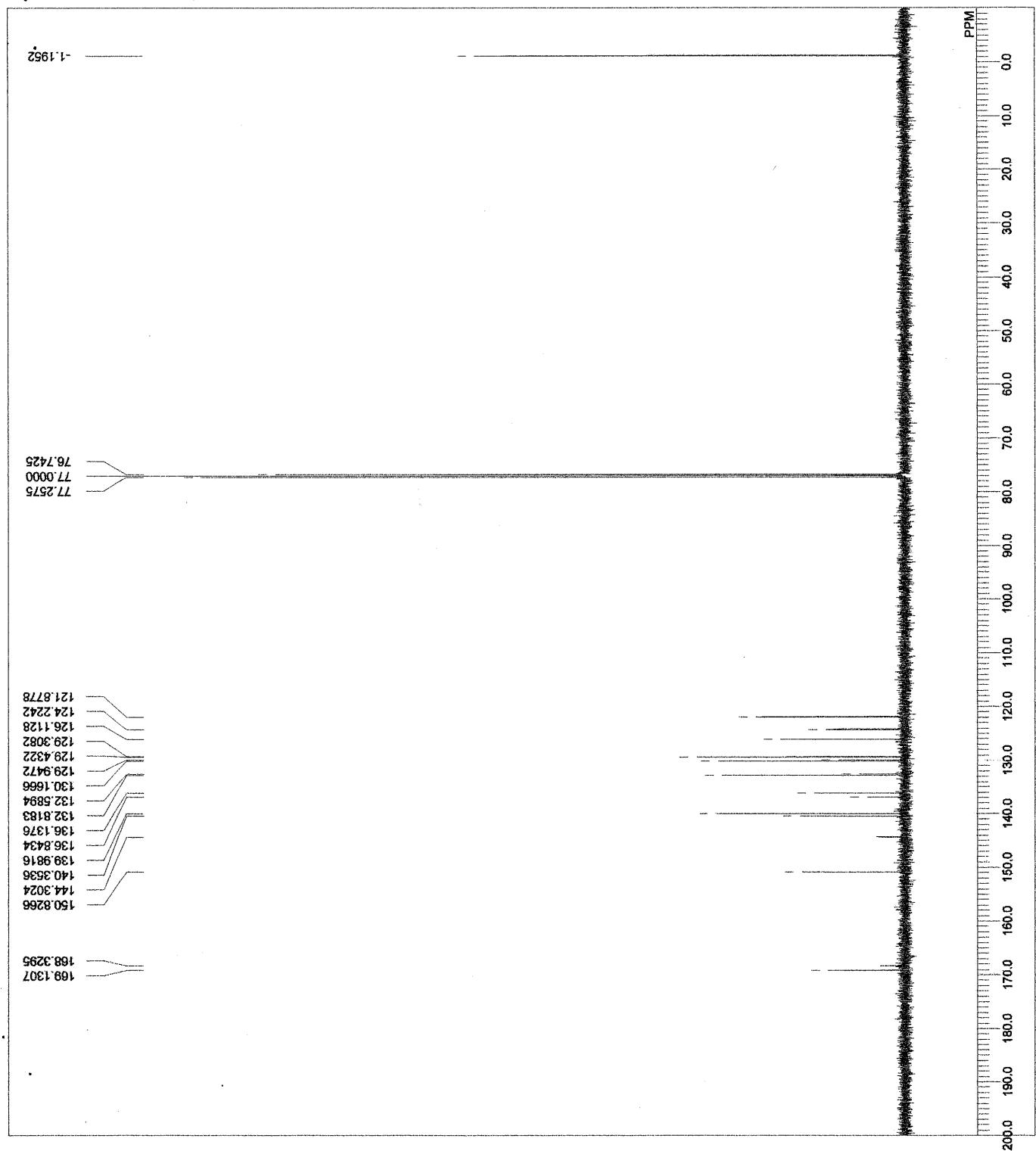


skk-p41-ArCO2Me-1-C-1-1.jdf
 skk-p41-ArCO2Me-1-C
 13C
 carbon.xpg
 DFILE
 COMNT
 DATIM
 OBNUC
 EXMOD
 OBFRQ
 OBSET
 OBFIN
 POINT
 FREQU
 SCANS
 ACQTM
 PD
 PW1
 IRNUC
 CTTEMP
 SLVNT
 EXREF
 BF
 RGAIN
 1002
 0.8336 sec
 3.0000 sec
 3.40 usec
 25.8 c
 CDCl₃
 77.00 ppm
 0.12 Hz
 60

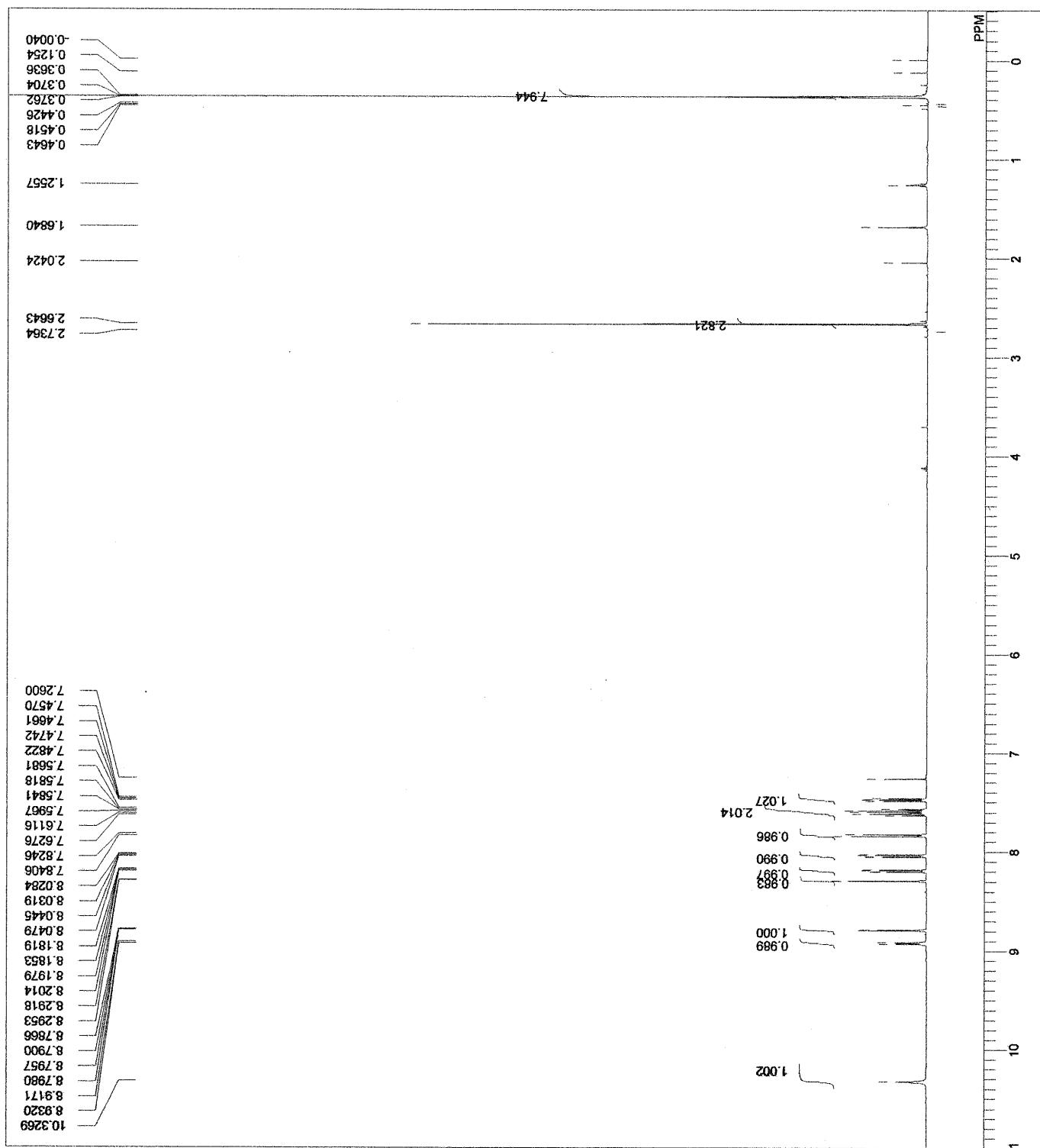
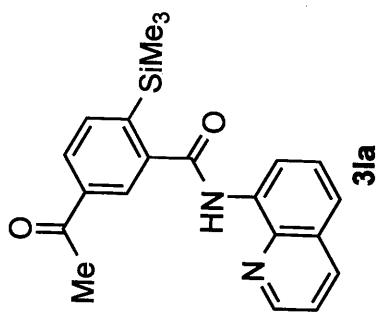


skk-p31-ArCO₂Me-Si-H $\frac{3}{2}k^{\prime \alpha}$



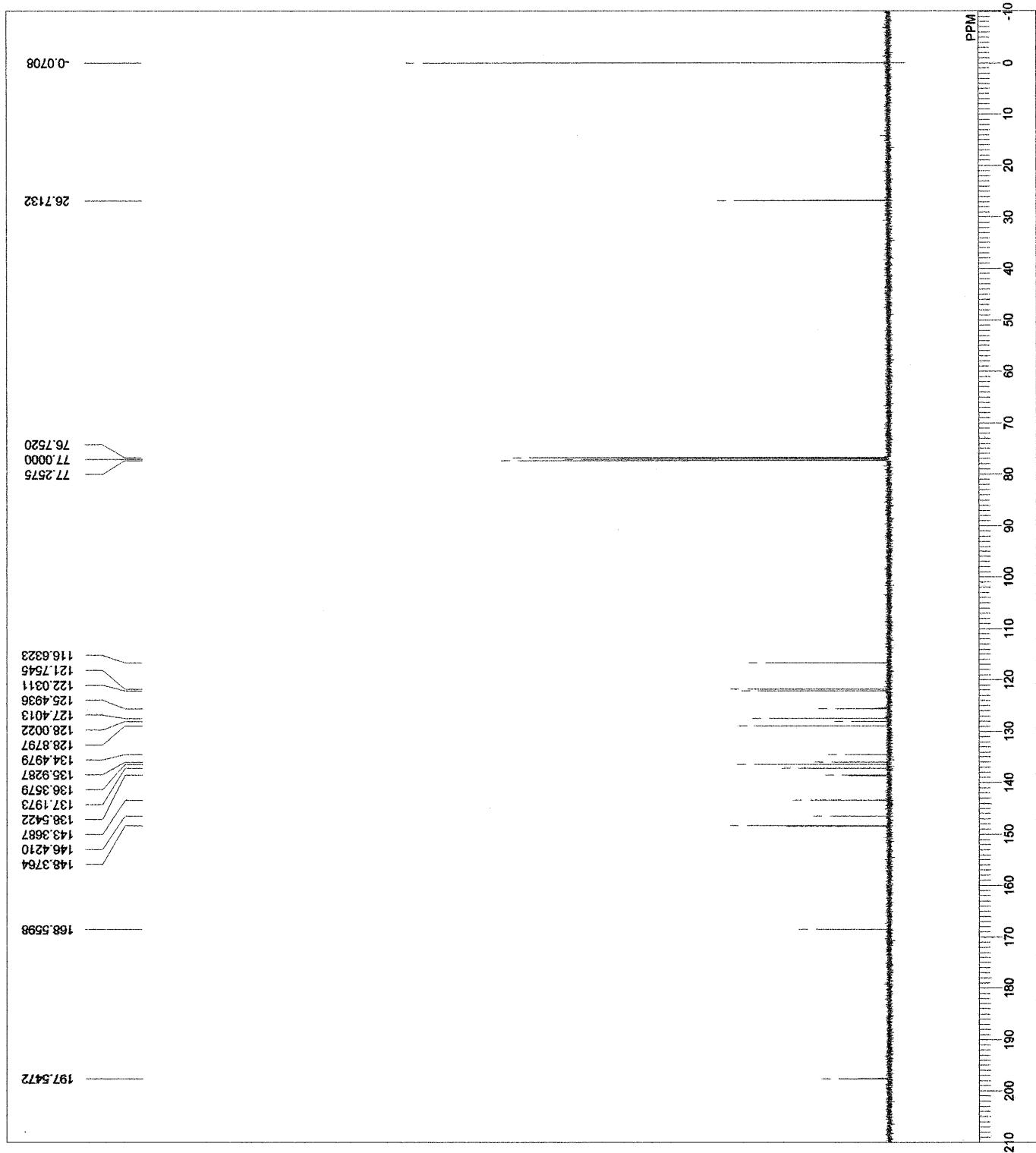


DFILE	skk-p185-ANMF-H-1.ais
COMIT	skk-p185-ANMF-H
DATIM	2013-07-29 14:23:07
OBNUC	1H
EXMOD	proton
OBFRQ	500.16 MHz
OSETB	2.41 kHz
OBFIN	6.01 Hz
POINT	13107
FREQU	7507.51 Hz
SCANS	16
ACQ1 TM	1.7459 sec
PD	5.0000 sec
PW1	5.55 usec
IRNUC	1H
CTEMP	22.8 °C
SLVNT	CDCl ₃
EXREF	7.26 ppm
BF	0.12 Hz
RGAIN	34



Aromatic-sines

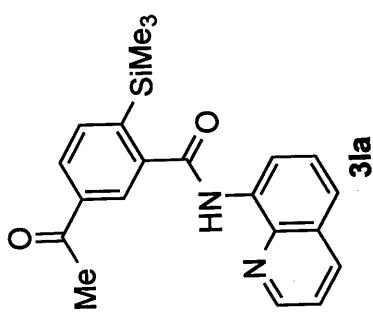
3la



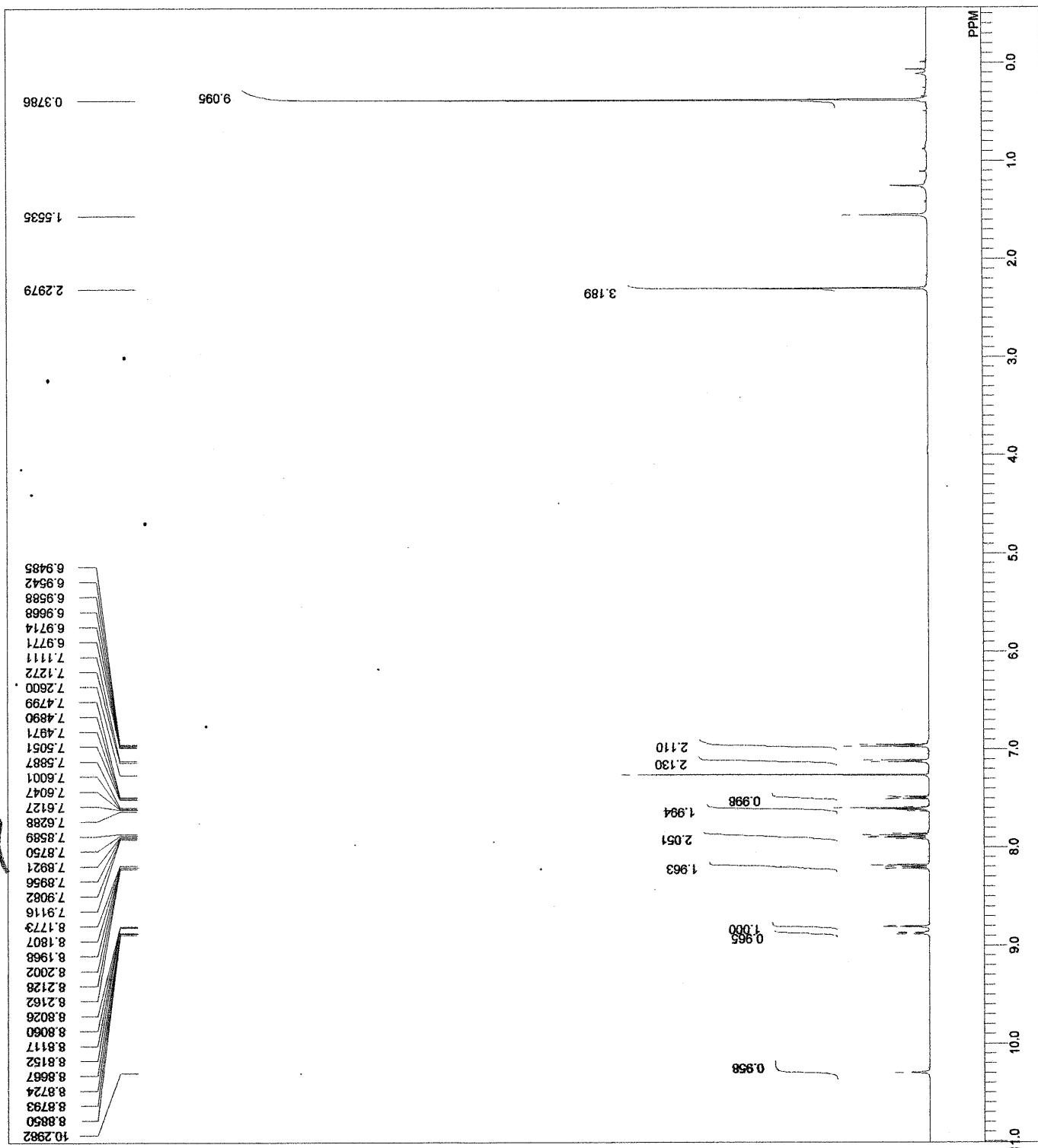
```

DFILE      skk-p185-ArCOMe-C-1.als
COMNT      skk-p185-ArMeF-C
DATIM      2013-07-29 14:25:49
OBNUC      13C
EXMOD      carbonJkp
OBFRQ      125.77 MHz
OFSET      7.87 kHz
OBFIN      4.21 Hz
POINT      26214
FREQU      31446.54 Hz
SCANS      386
ACQTM      0.8336 sec
PD         3.0000 sec
PW1        3.40 usec
IRNUC      1H
CTEMP      23.6 c
SLVNT      CDCl3
EXREF      77.00 ppm
BF         0.12 Hz
RGAIN      60

```



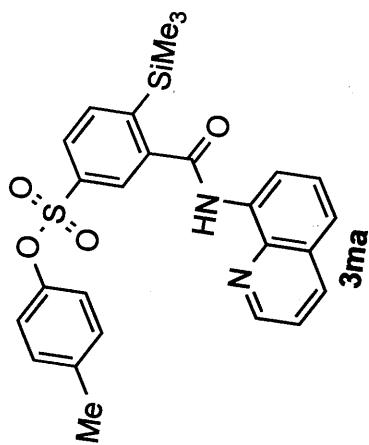
skk-p41-ArSO₃Ar-SiMe₃-H **3ma**



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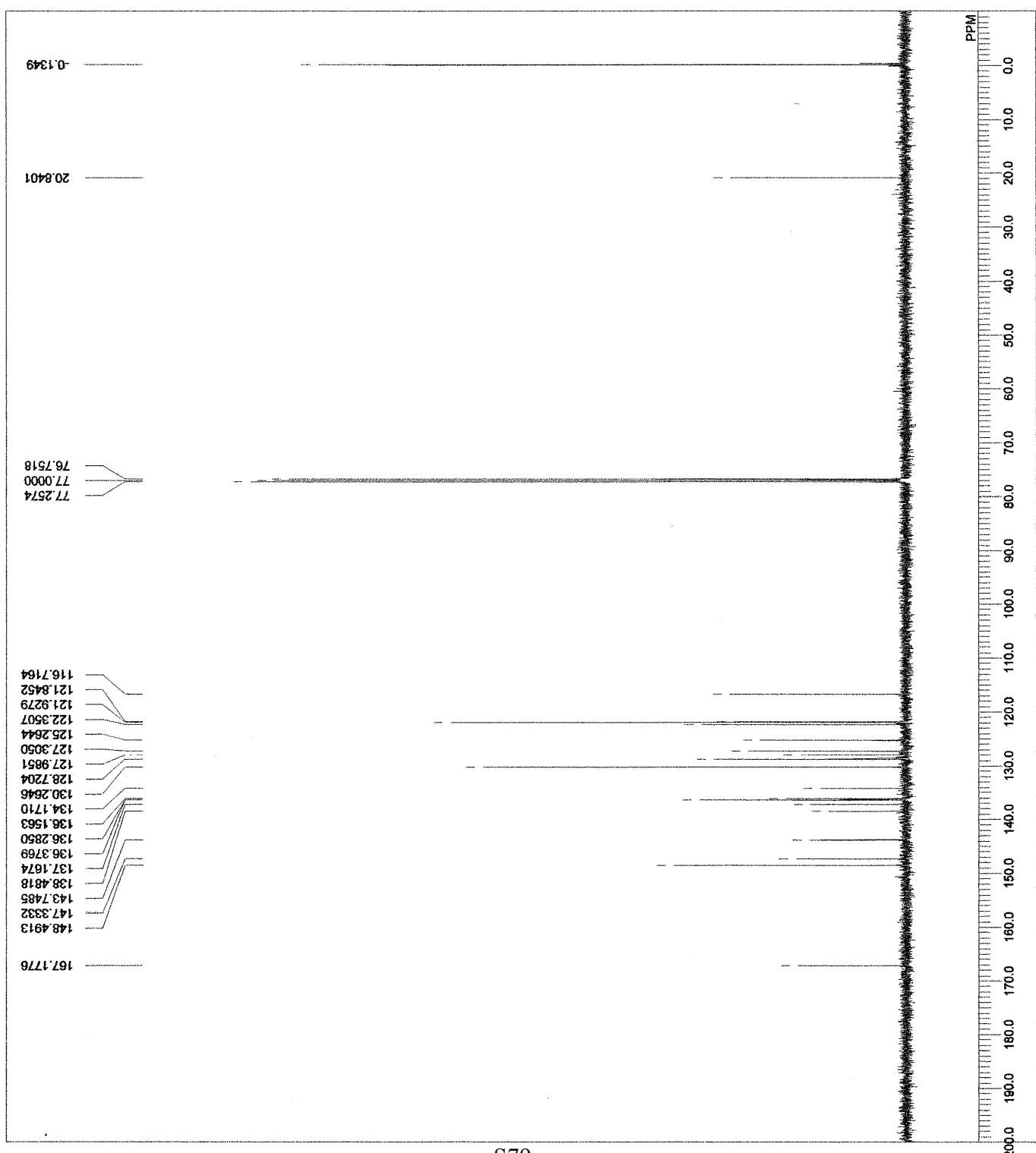
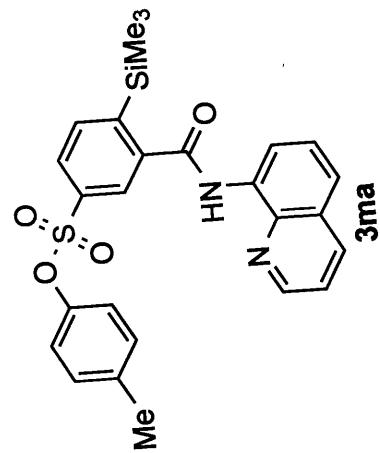
skk-p41-ArSO3Ar-SiMe3-H-3-1.dff
skk-p41-ArSO3Ar-SiMe3-H
2013-11-06 14:30:30
1H
protein
500.16 MHz
2.41 kHz
6.01 Hz
16384
9384.38 Hz
8
1.7459 sec
5.0000 sec
5.55 usec
1H
24.1 c
CDCl3
7.26 ppm
0.12 Hz
32

```

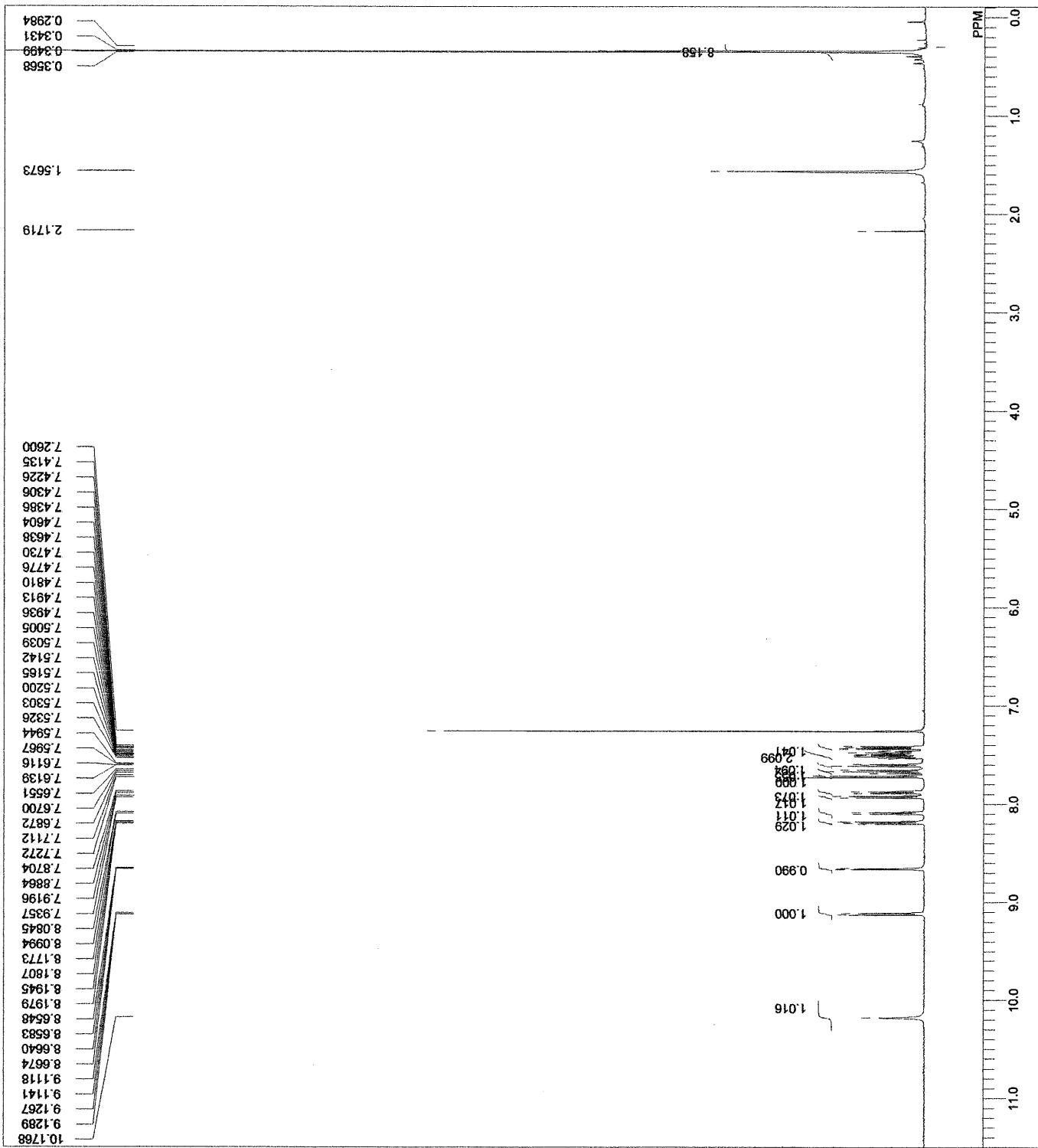
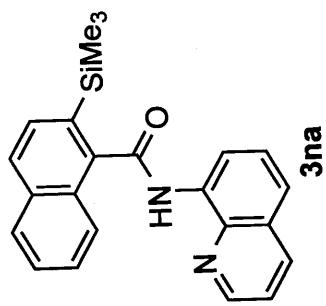


skk-p41-ArSO3Ar-SiMe3-C_copy4-1-1.prf
 skk-p41-ArSO3Ar-SiMe3-C
 2013-10-11 17:24:00
 13C
 carbon-13D
 EXMOD
 OBFRQ
 OBSET
 OBFIN
 POINT
 FREQU
 SCANS
 ACQTM
 PD
 PW1
 INUC
 CTEMP
 SLVNT
 EXREF
 BF
 RGAIN

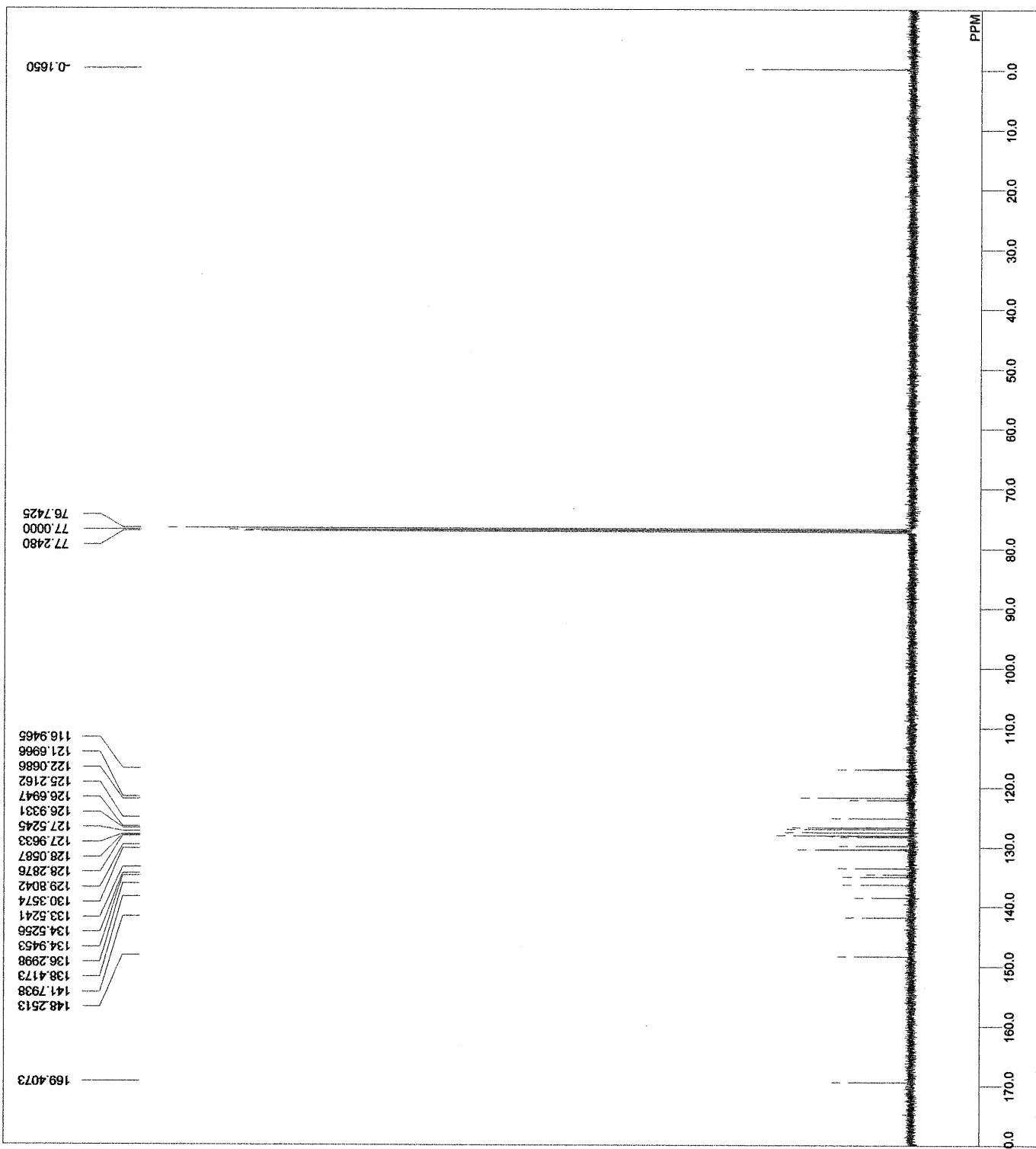
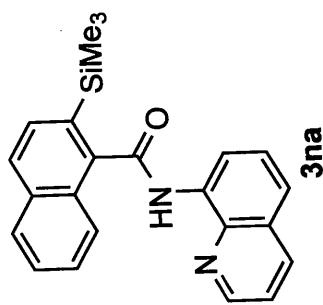
125.77 MHz
 7.87 kHz
 4.21 Hz
 32767
 31878.79 Hz
 151
 0.0000 sec
 3.0000 sec
 3.40 Usec
 1H
 26.0 c
 CDCl₃
 77.00 ppm
 1.00 Hz
 60



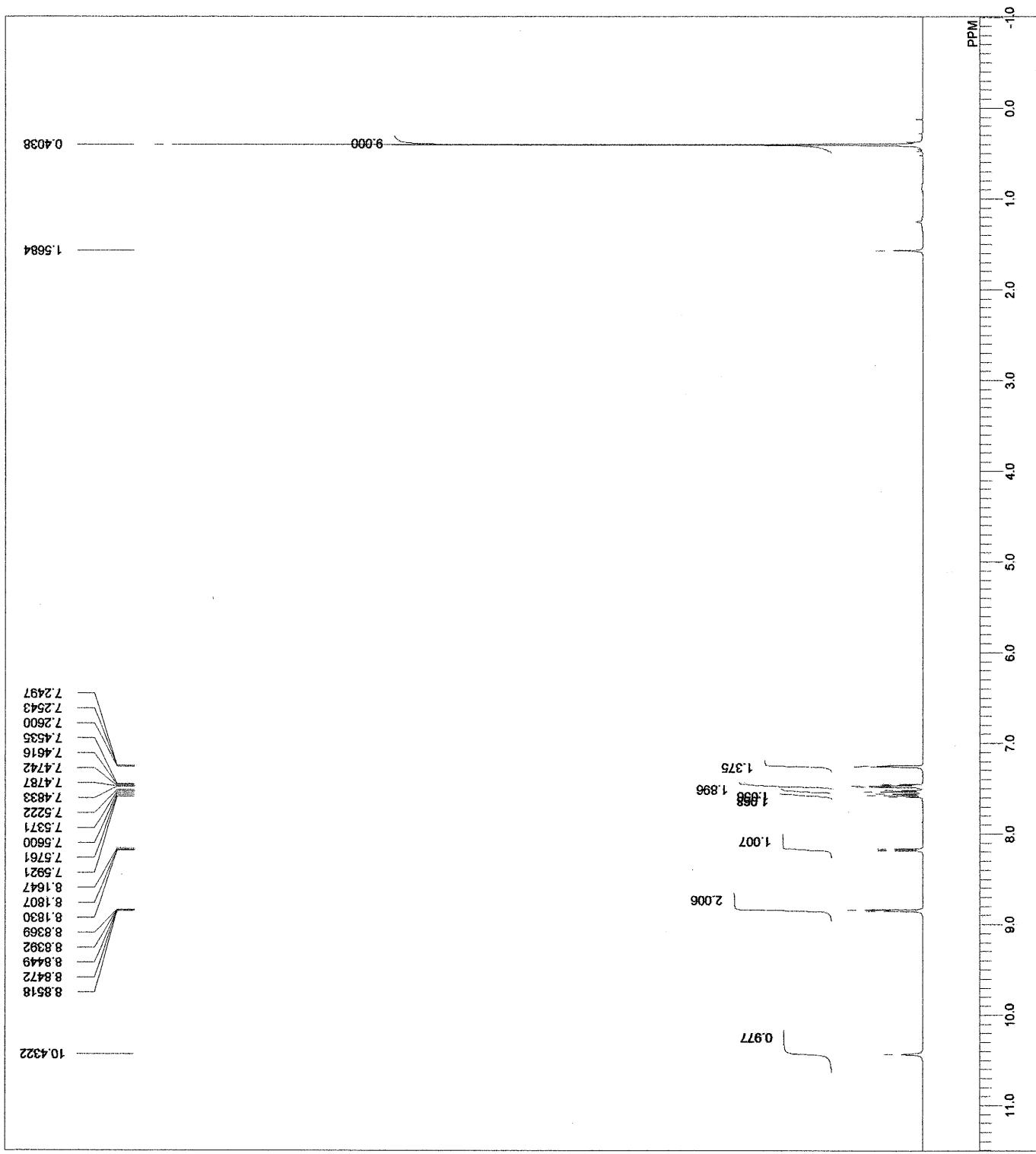
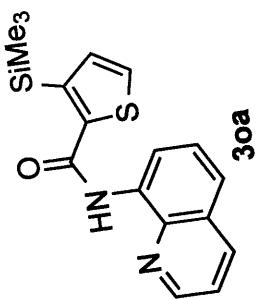
3na



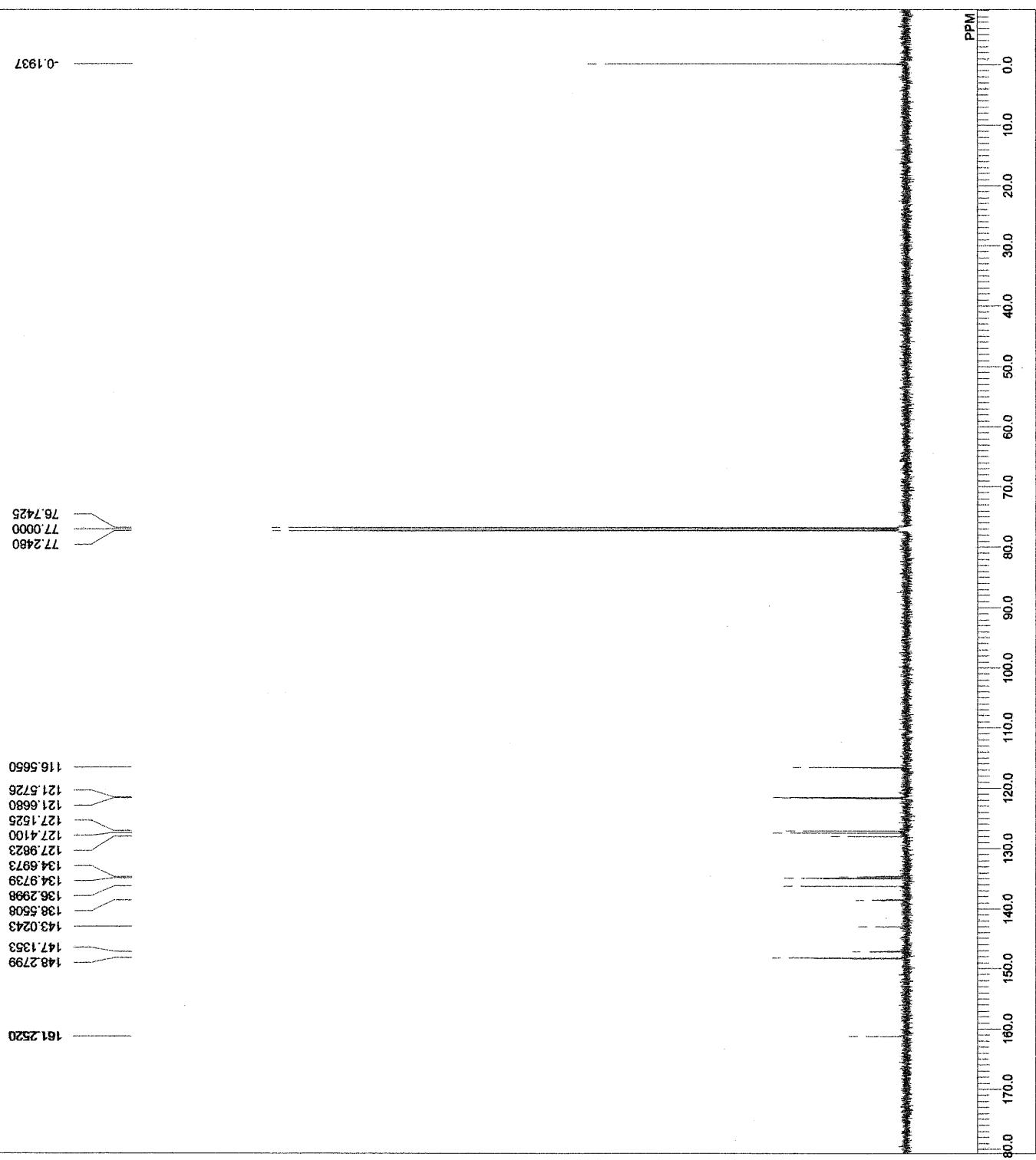
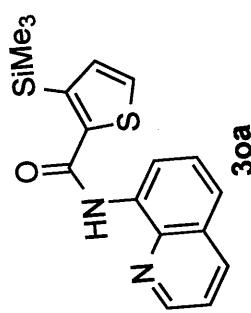
DFILE skk-p192-ArNaphth-C-1-1.jdf
 COMNT skk-p192-ArNaphth-C
 DATIM 2013-08-05 14:11:56
 EXNUC 13C
 EXMOD carbon.ixp
 OBFREQ 125.77 MHz
 OFFSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 32767
 FREQU 39308.18 Hz
 SCANS 551
 ACQTIM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IRNUC 1H
 CTEMP CDCl₃
 SLVNT 77.00 ppm
 EXREF 0.12 Hz
 BF 60
 RGAIN



skk-bk2-p135-8AqCOthiophene-H-1-1.xls
 skk-bk2-p135-8AqCOthiophene-H
 1H
 proton-1D
 500.16 MHz
 2.41 kHz
 6.01 Hz
 16384
 9334.38 Hz
 8
 SCANS 1,7459 sec
 ACQTM 5.0000 sec
 PD 5.55 usec
 PW1 1H
 INUC CTEMP 24.0 c
 SLVNT CDCl₃
 EXREF 7.26 ppm
 BF 1.00 Hz
 RGAIN 30

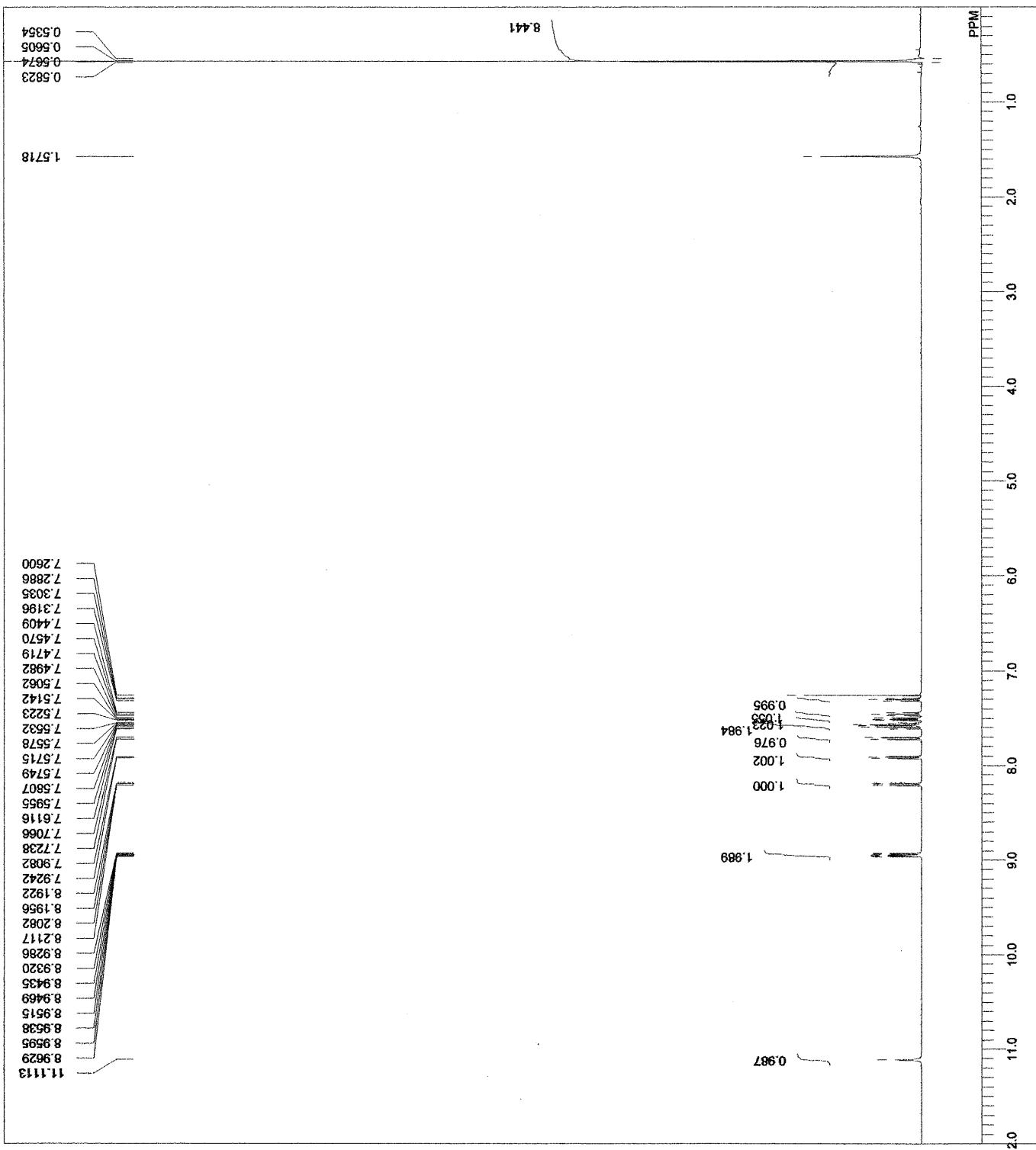
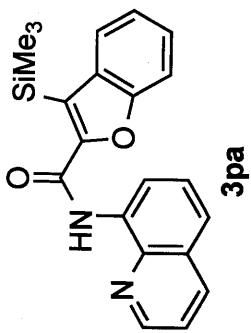


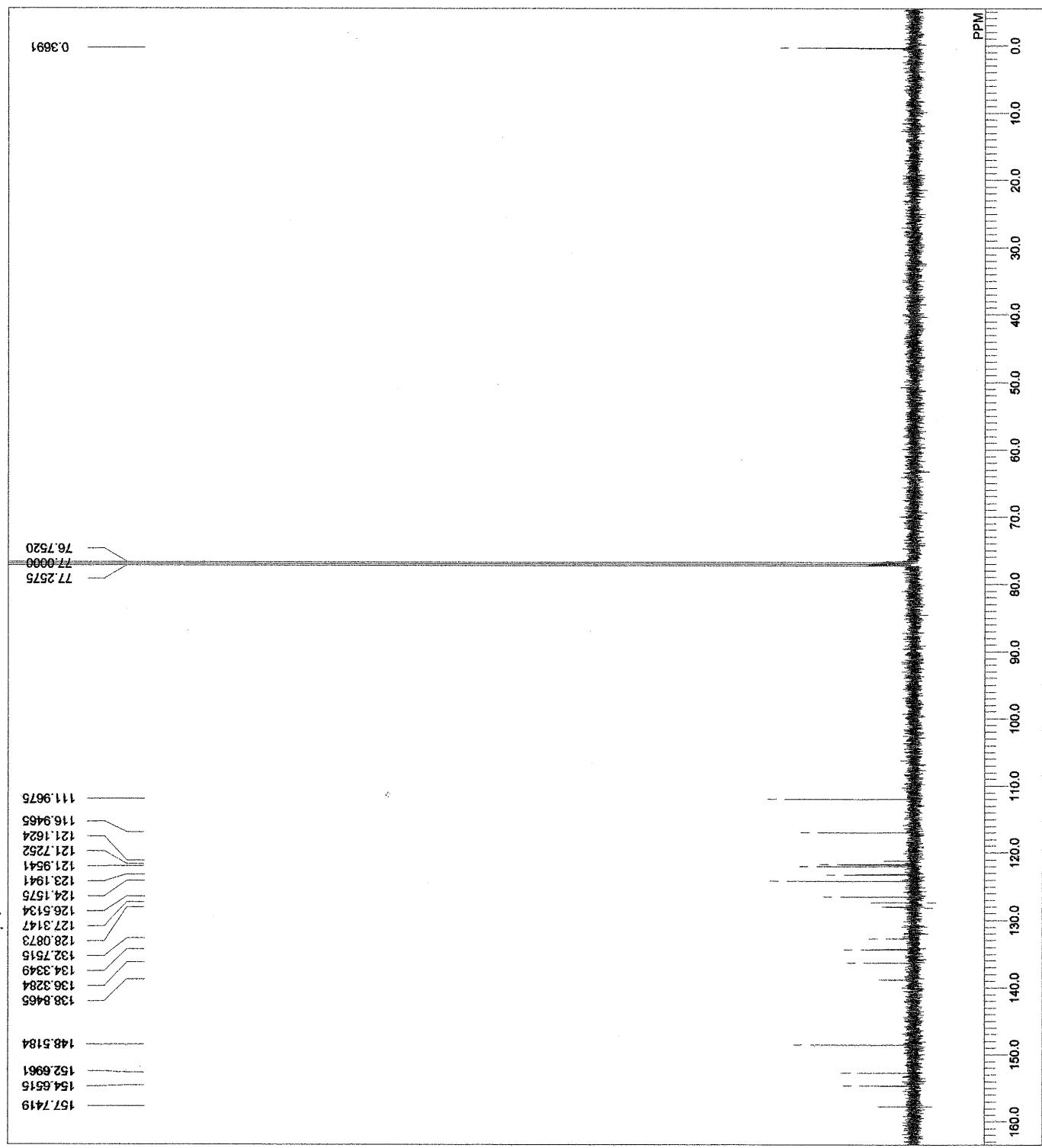
DFILE
 COMINT
 DATIM
 2013-11-01 12:55:60
 13C
 EXMOD
 carbon.kdp
 OBFRQ 125.77 MHz
 OBSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 32/167
 FREQU 39008.18 Hz
 SCANS 194
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 INUC 1H
 CTEMP 24.6 c
 CDCL3 77.00 ppm
 SVNT EXREF
 BF 1.00 Hz
 RGAIN 60



skk-p178-Benzofuran-(1)-H-f-1.ais
skk-p178-Benzofuran-(1)-H
2013-07-31 10:33:55

DFILE COMM DATIM OBNLU EXMO OBFRS OBSE OBFIN POINT FREQU SCANS ACQTTI PD PW1 IRNUUC CTEM SLVNT EXREF BF RGAIN

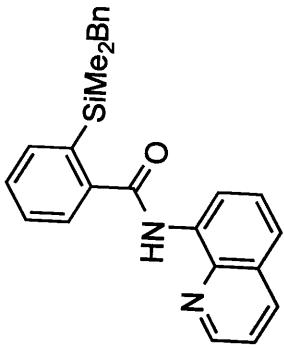


3pa

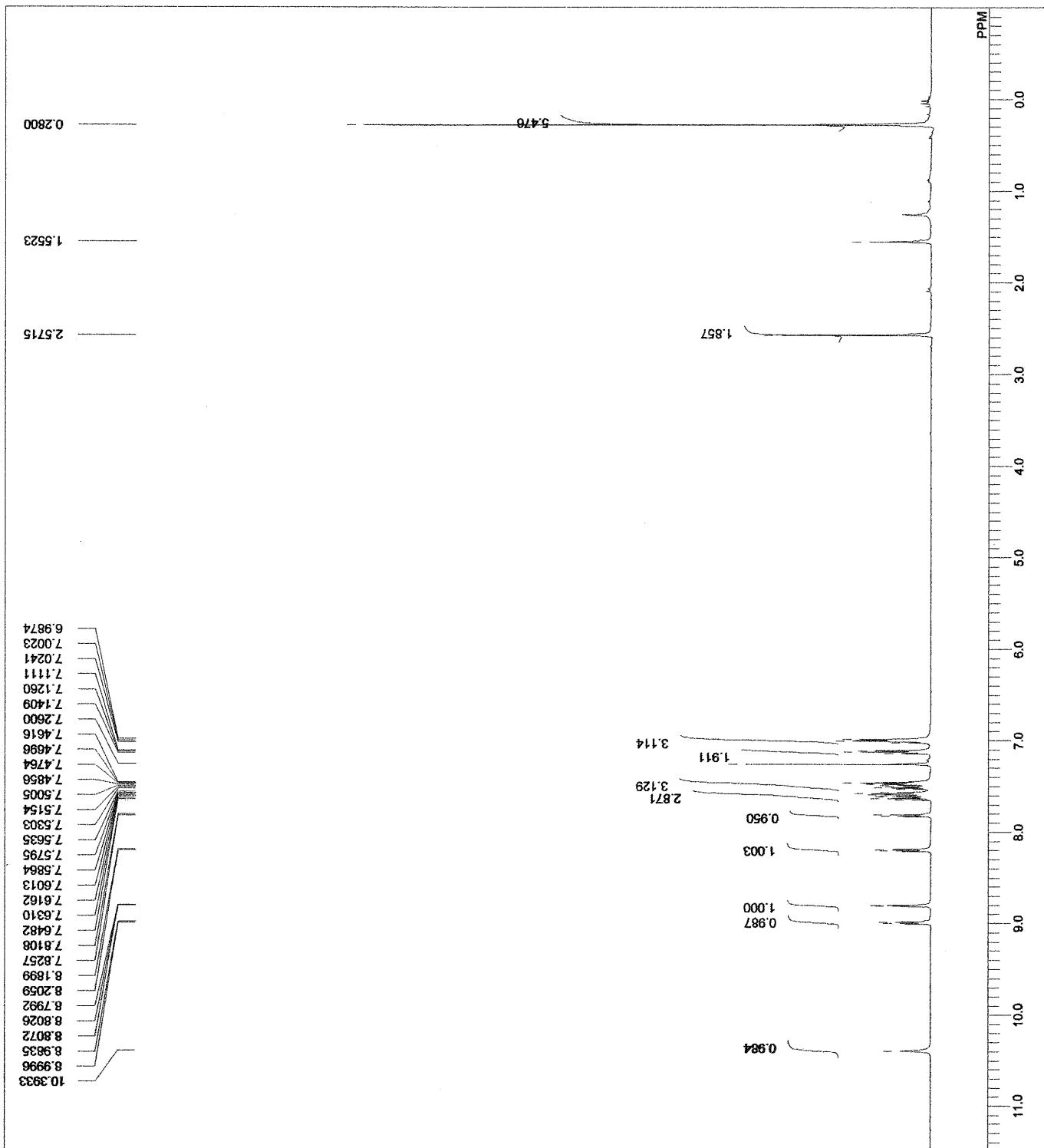
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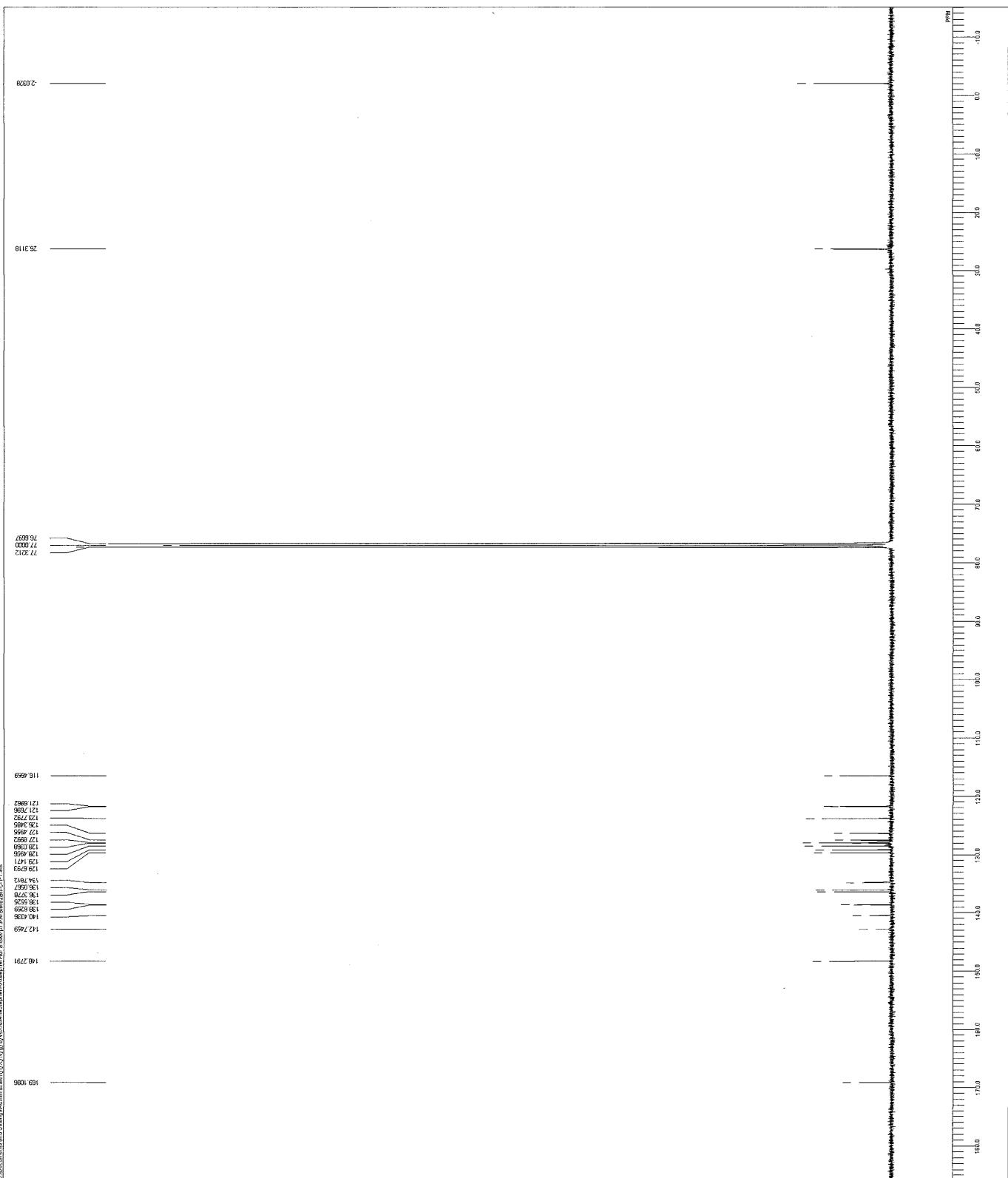
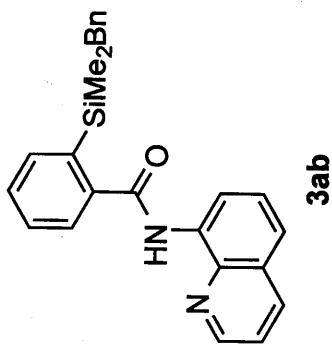
DFILE      skk-p3-4fSIM2Br-H-1-1.als
COMNT      skk-p3-4fSIM2Br-H
DATIM      2013-12-11 16:43:32
OBNUC      1H
EXMOD      proton,jxp
OFQRO      500.16 MHz
OFFSET     2.41 kHz
OFBIN      6.01 Hz
POINT      13107
FREQU      7507.51 Hz
SCANS      16
ACQTM      1.745s sec
PD         5.0000 sec
PW1        5.55 usec
IRNUC      1H
CTEMP      21.7 c
SLVNT      CDCL3
EXREF      7.26 ppm
BF         0.12 Hz
RGAIN      40

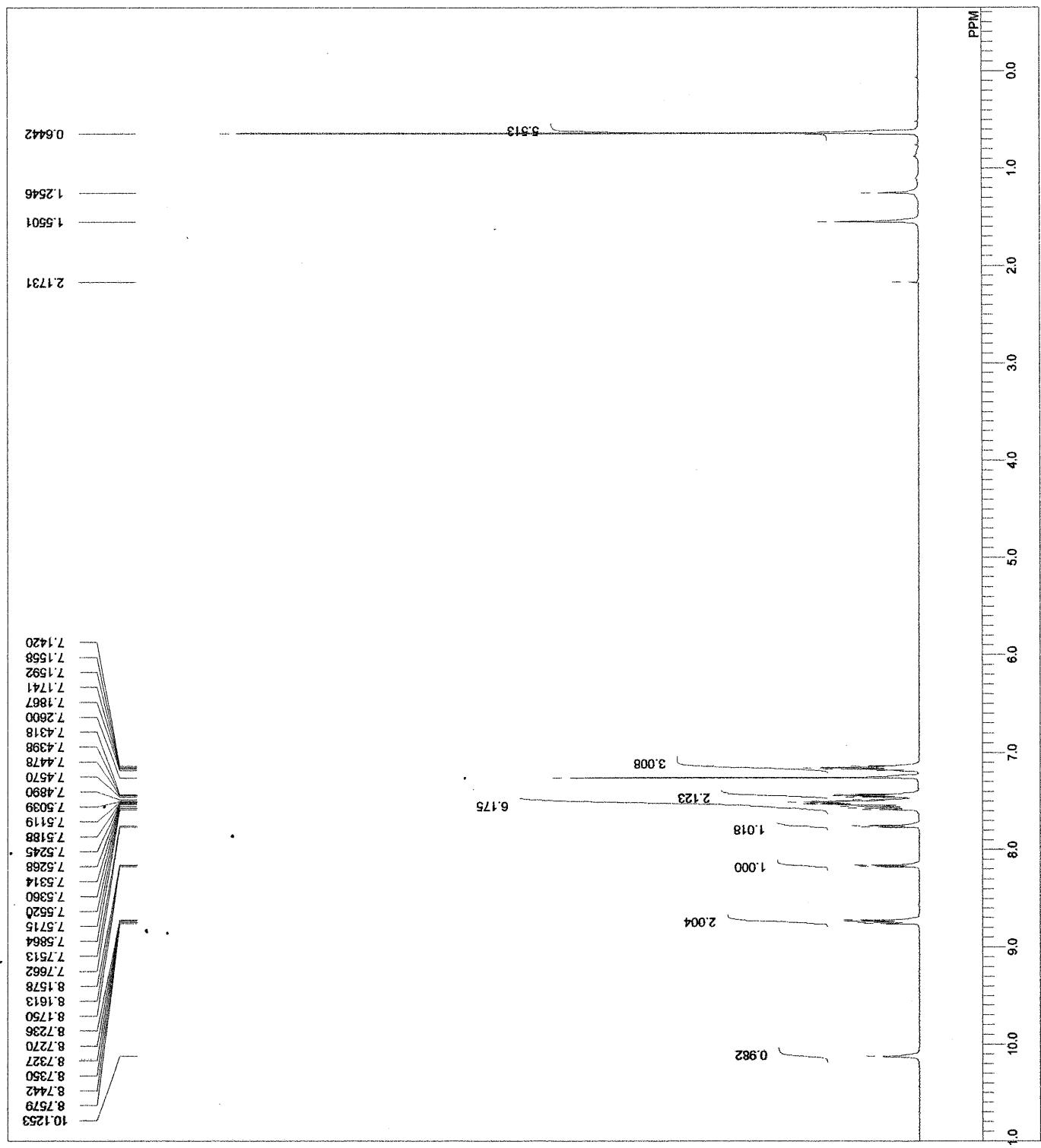
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3ab





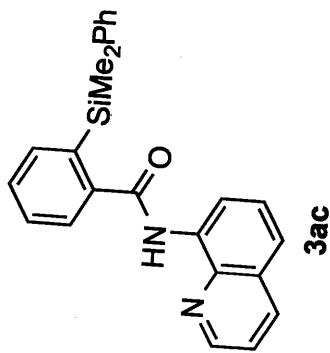


```

skk-p3-ArSiMe2Ph-H-1-1.dif
skk-p3-ArSiMe2Ph-H
1H
proton,j,xp
500.16 MHz
2.41 kHz
6.01 Hz
18384
9384.38 Hz
16
1.7459 sec
5.0000 sec
5.55 usec
CDCl3
21.5 c
7.26 ppm
0.12 Hz
40

```

DFILE
COMNT
DATIM
OBNUC
EXMOD
OBFRQ
OBSET
OBFIN
POINT
FREQU
SCANS
ACQTM
PD
PW1
IRNUC
CTEMP
SLVNT
EXREF
BF
RGAIN



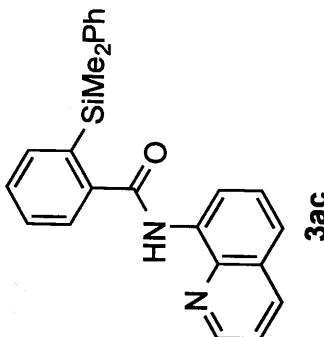
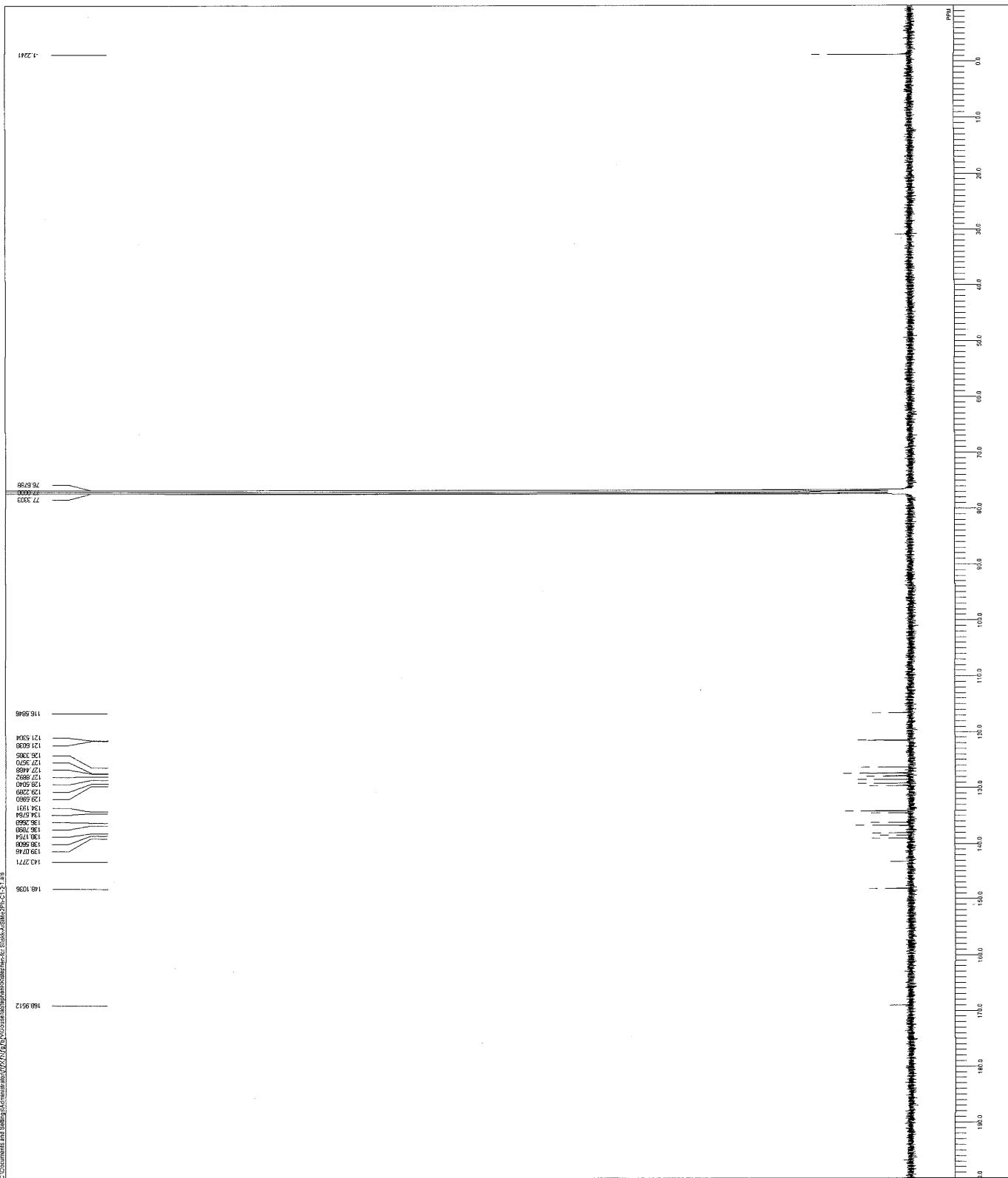
363

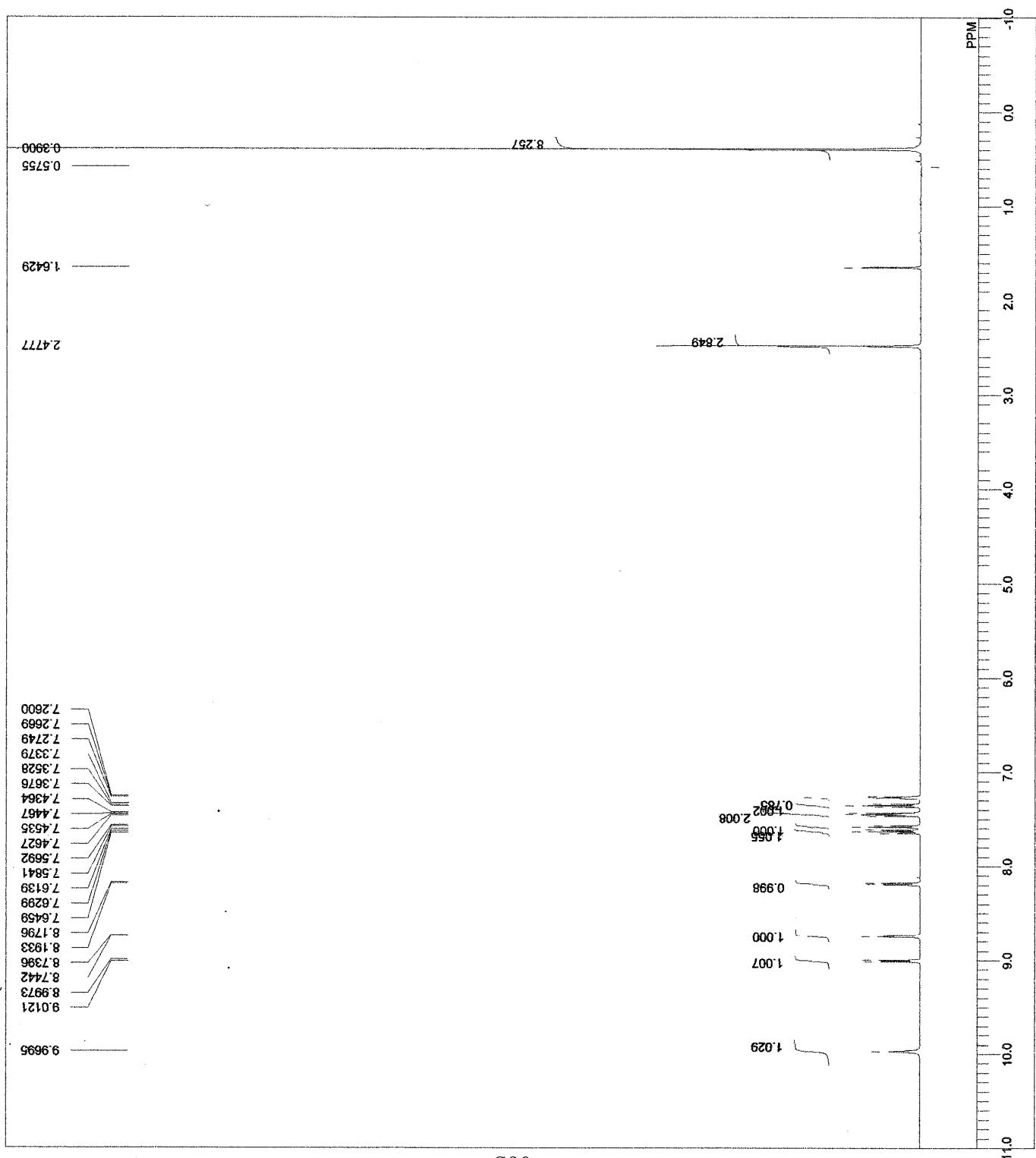
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1

DFILE skk-ArSiM@2Ph-C1-2-1.als

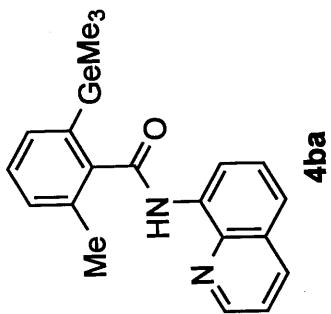




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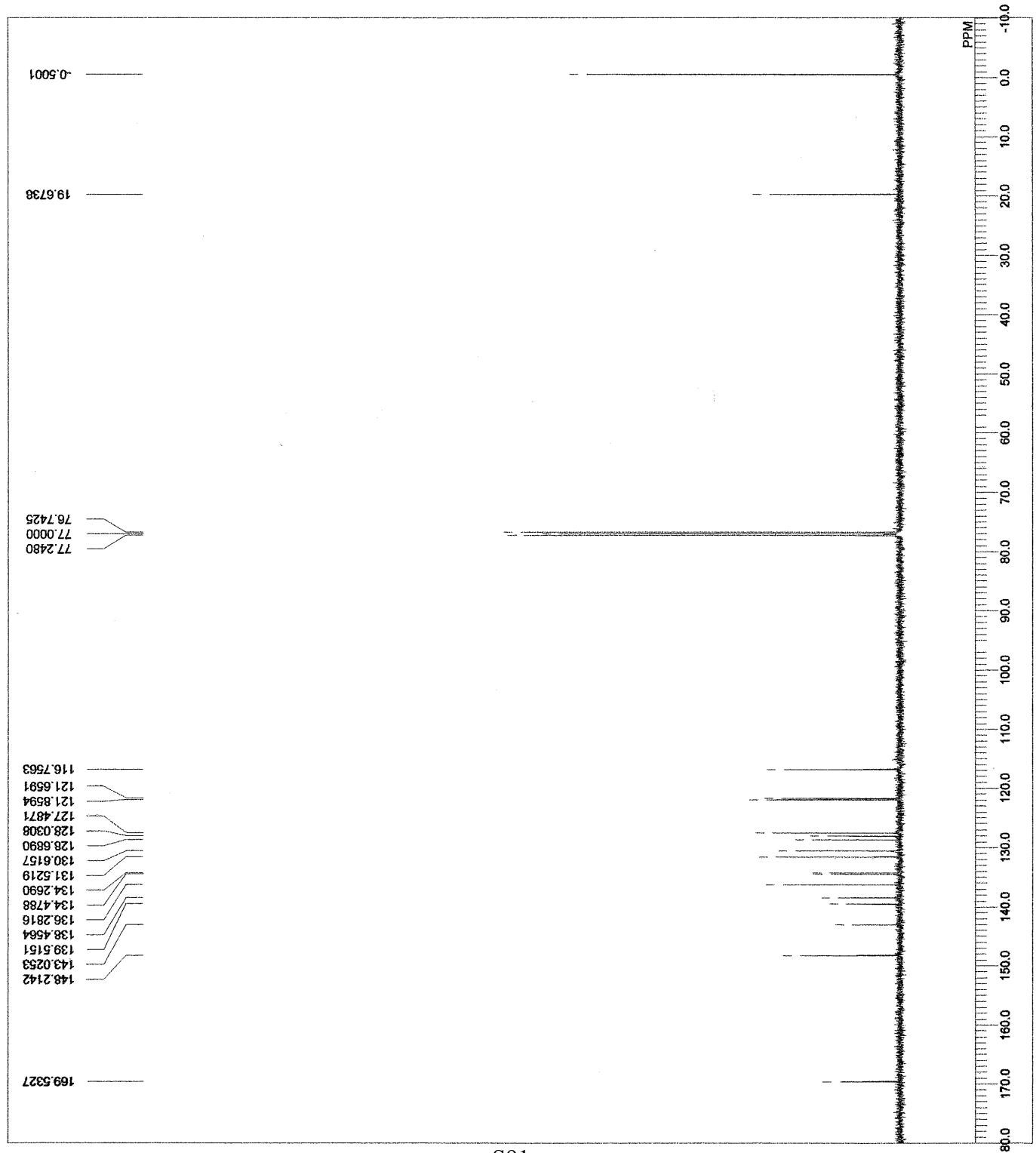
DFILE      skk-p202-ArMe-Ge-H-1-1.idf
COMNT
DATIM      2013-08-12 18:08:06
1H
EXMOD      proton-1D
OBFRQ      500.16 MHz
OSET       2.41 kHz
OBFIN      6.01 Hz
POINT      16384
FREQU     9334.38 Hz
SCANS      16
ACQTM      1.7459 sec
PD         5.0000 sec
PW1        5.55 usec
IRNUC      1H
CTEMP      26.3 c
SLVNT      CDCl3
EXREF      7.26 ppm
BF         1.00 Hz
RGAIN      34

```

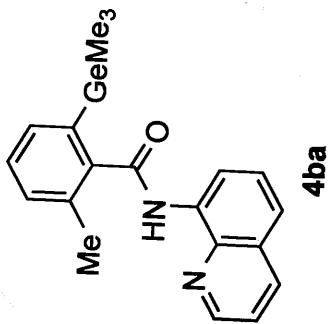


skk-p202-ArMe-Ge-C

4ba



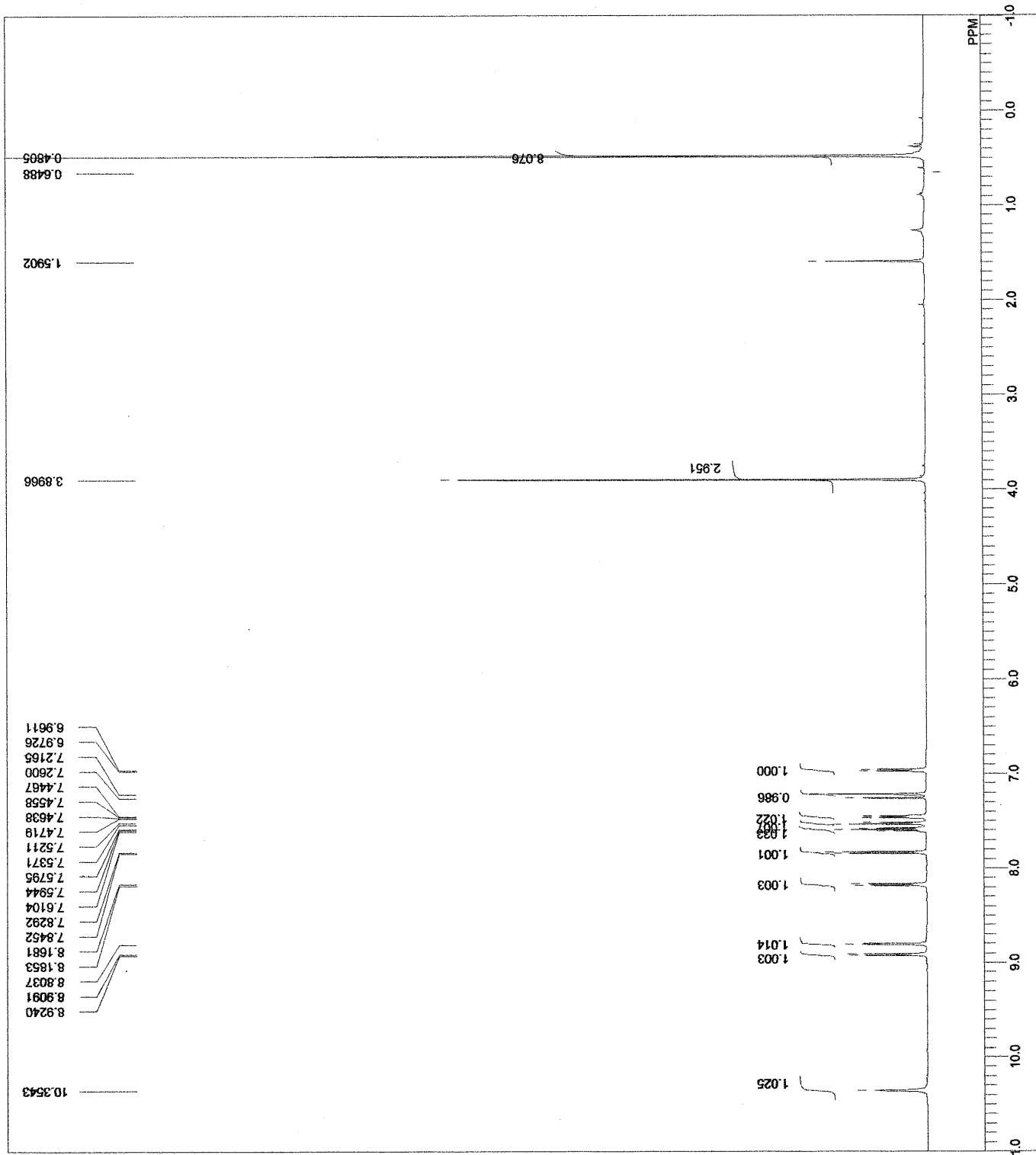
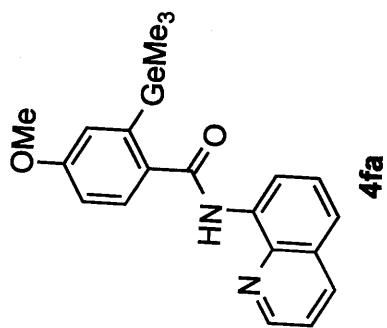
skk-p202-ArMe-Ge-C-1-1.als
skk-p202-ArMe-Ge-C-1
2013-08-12 18:11:17
13C
carbon.jpc
OBFRQ 125.77 MHz
OBFSET 7.87 kHz
OBFIN 4.21 Hz
POINT 26214
FREQU 31446.54 Hz
SCANS 122
ACQTM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usec
IRNUC 1H
CTEMP 26.6 c
SLVNT CDCl₃
EXREF 77.00 ppm
BF 1.00 Hz
RGAIN 60

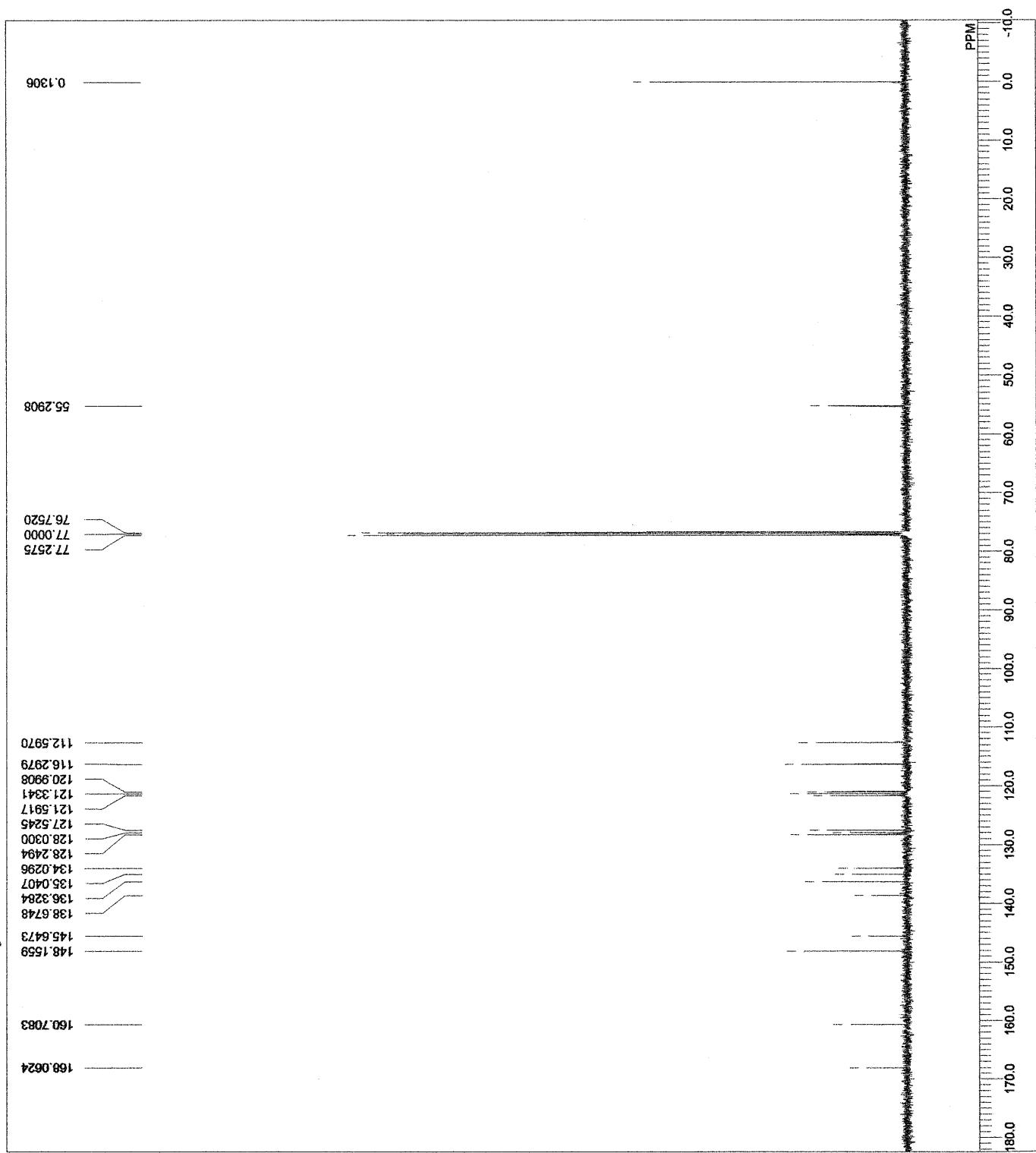


skk-p20-Ge-ArOME-H

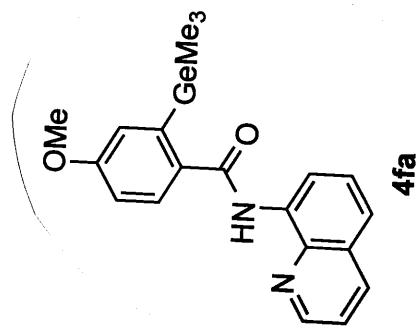
4fa

DIFILE skk-p20-Ge-AOMe-H-1-1-1df
COMNT skk-p20-Ge-AOMe-H
DATIM 2013-09-17 16:54:53
OBNUC 1H
EXMOD proton.jxp
OBFRQ 500.16 MHz
OBSET 2.41 kHz
OBFIN 6.01 Hz
POINT 16384
FREQU 9384.38 Hz
SCANS 16
ACQTIM 1.7439 sec
PD 5.0000 sec
PW1 5.55 usec
IRNUC 1H
CTEMP 25.0 c
SLVNT CDCL₃
EXREF 7.26 ppm
BF 1.20 Hz
RGAIN 38

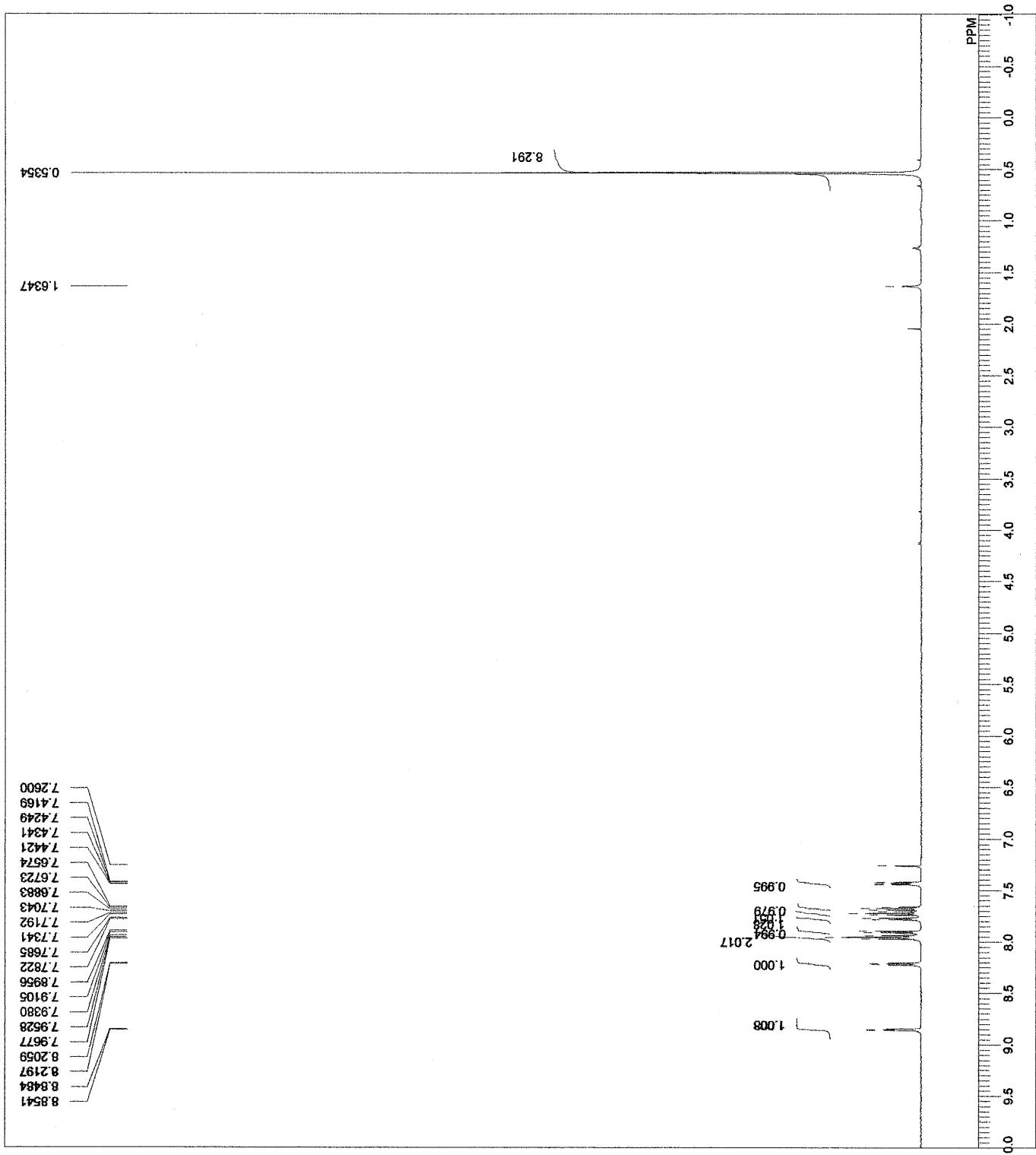
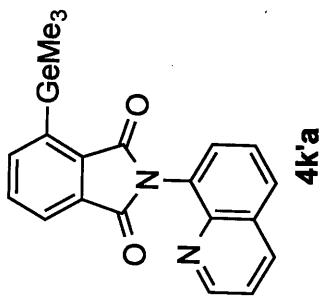


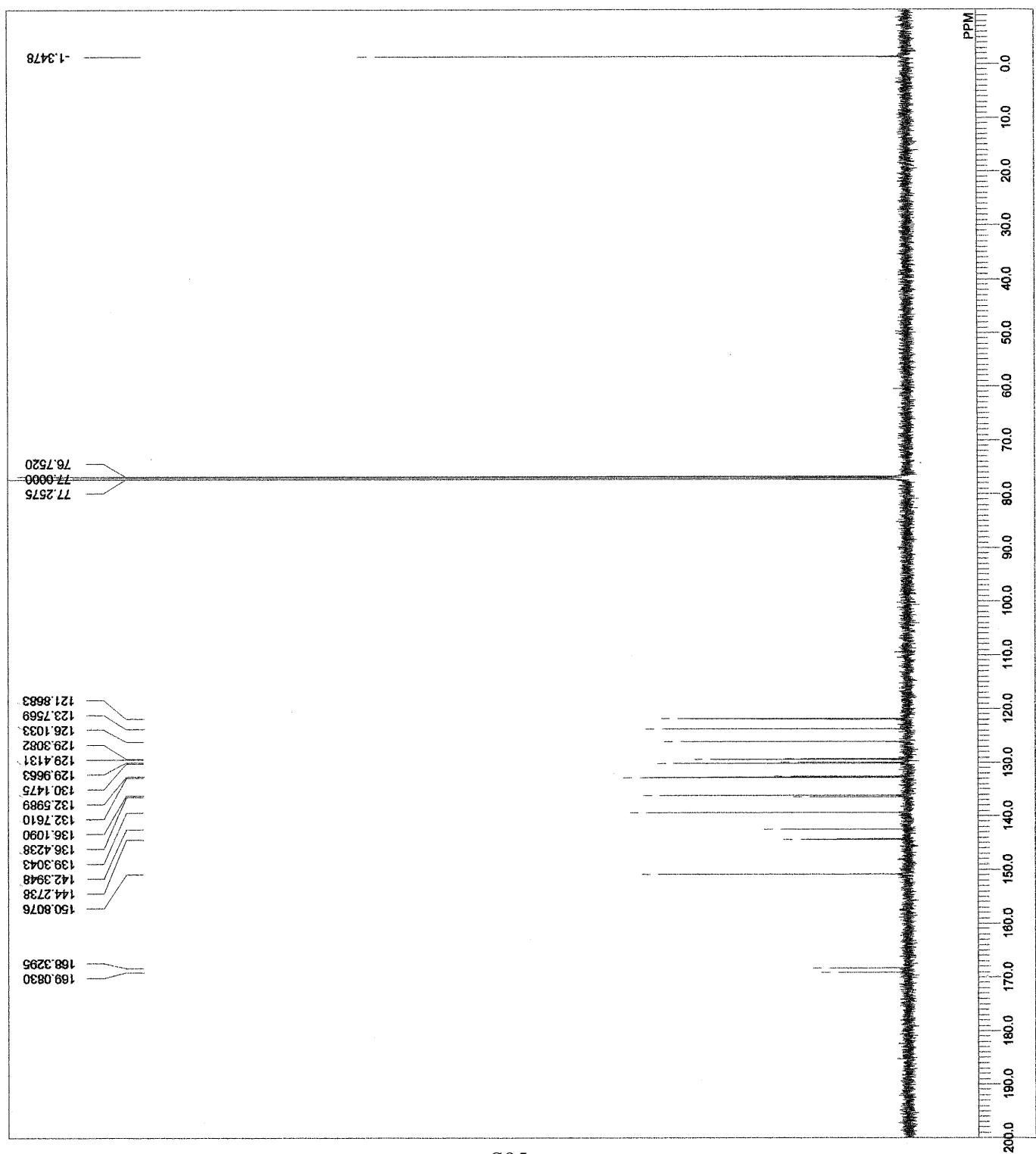
4fa

skk-p20-Ge-ArOMe-C-1-1.als
skk-p20-Ge-ArOMe-C
2013-09-17 16:57:35
13C
carbon-13Q
OBFRQ 125.77 MHz
OBSET 7.87 kHz
OBFIN 4.21 Hz
POINT 32767
FREQU 39308.18 Hz
SCANS 159
ACQTM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usec
IRNUC 1H
CTEMP 25.7 c
CDCL3 77.00 ppm
SLVNT EXREF
BF 1.20 Hz
RGAIN 60

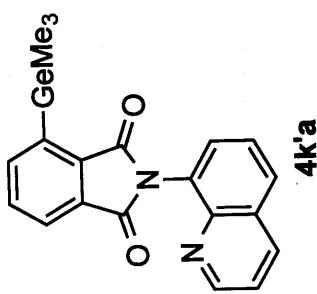


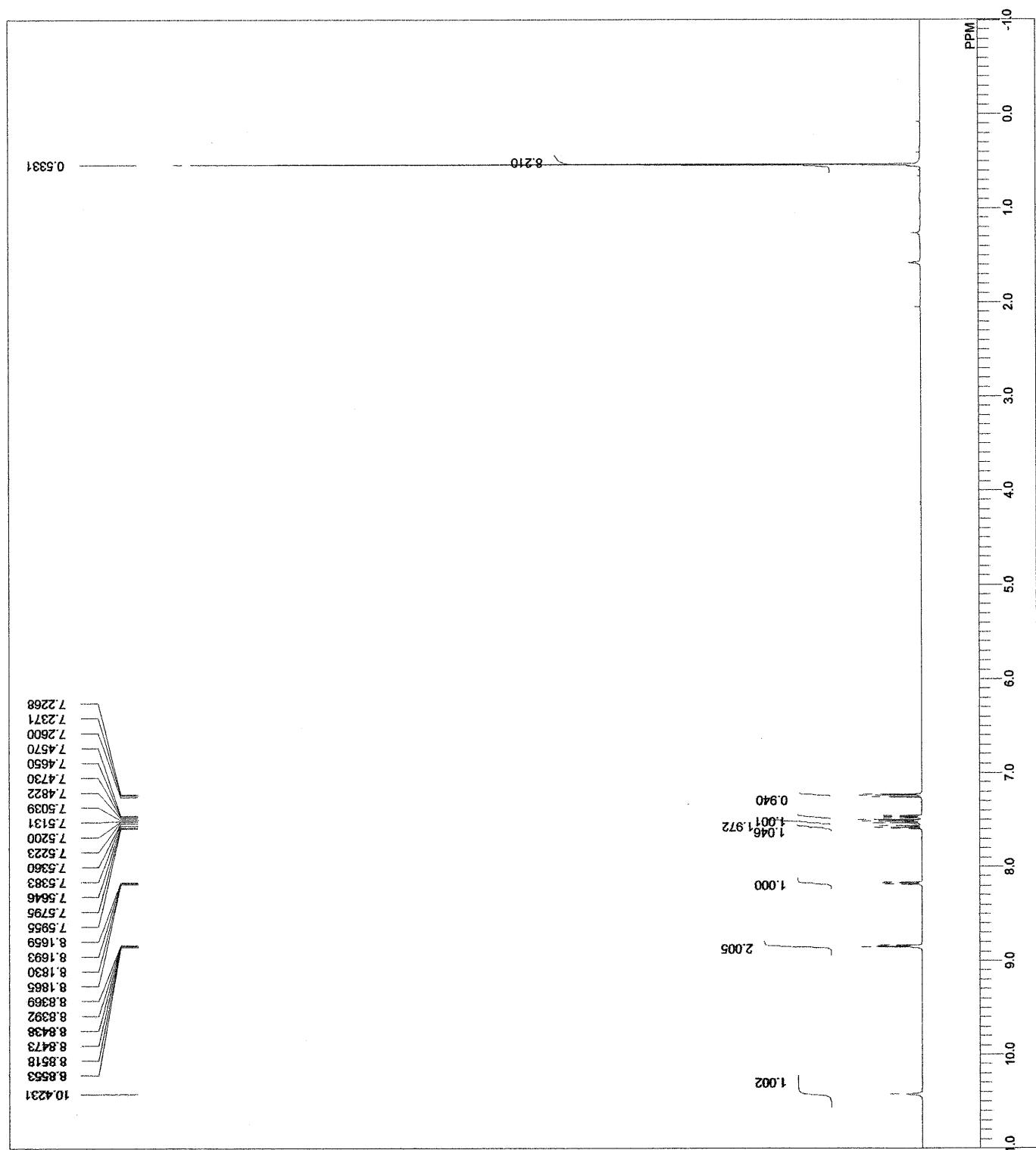
DFILE skk-p20-Ge-ArCO2Me-H-1-1.ais
 COMNT skk-p20-Ge-ArCO2Me-H
 DATIM 2013-09-17 16:33:03
 OBNUC 1H
 EXMOD proton JKO
 OBRQ 500.16 MHz
 OBSET 2.41 kHz
 OBEIN 6.01 Hz
 POINT 13107
 FREQU 7507.51 Hz
 SCANS 16
 ACQTM 1.7459 sec
 PD 5.0000 sec
 PW1 5.55 usec
 INUC 1H
 CTEMP 24.5 c
 SILVNT CDCl₃
 EXREF 7.26 ppm
 BF 1.20 Hz
 RGAIN 36

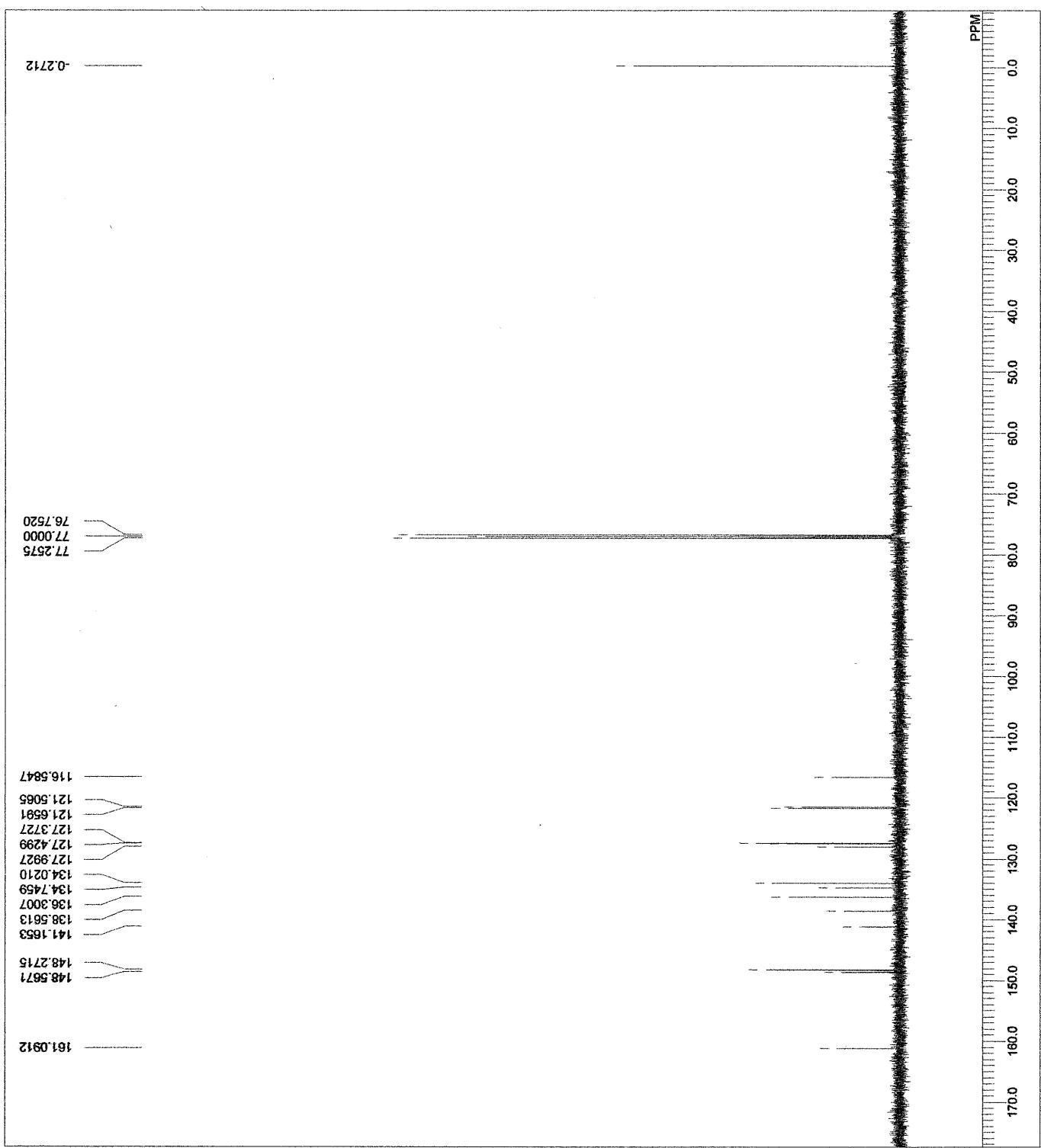


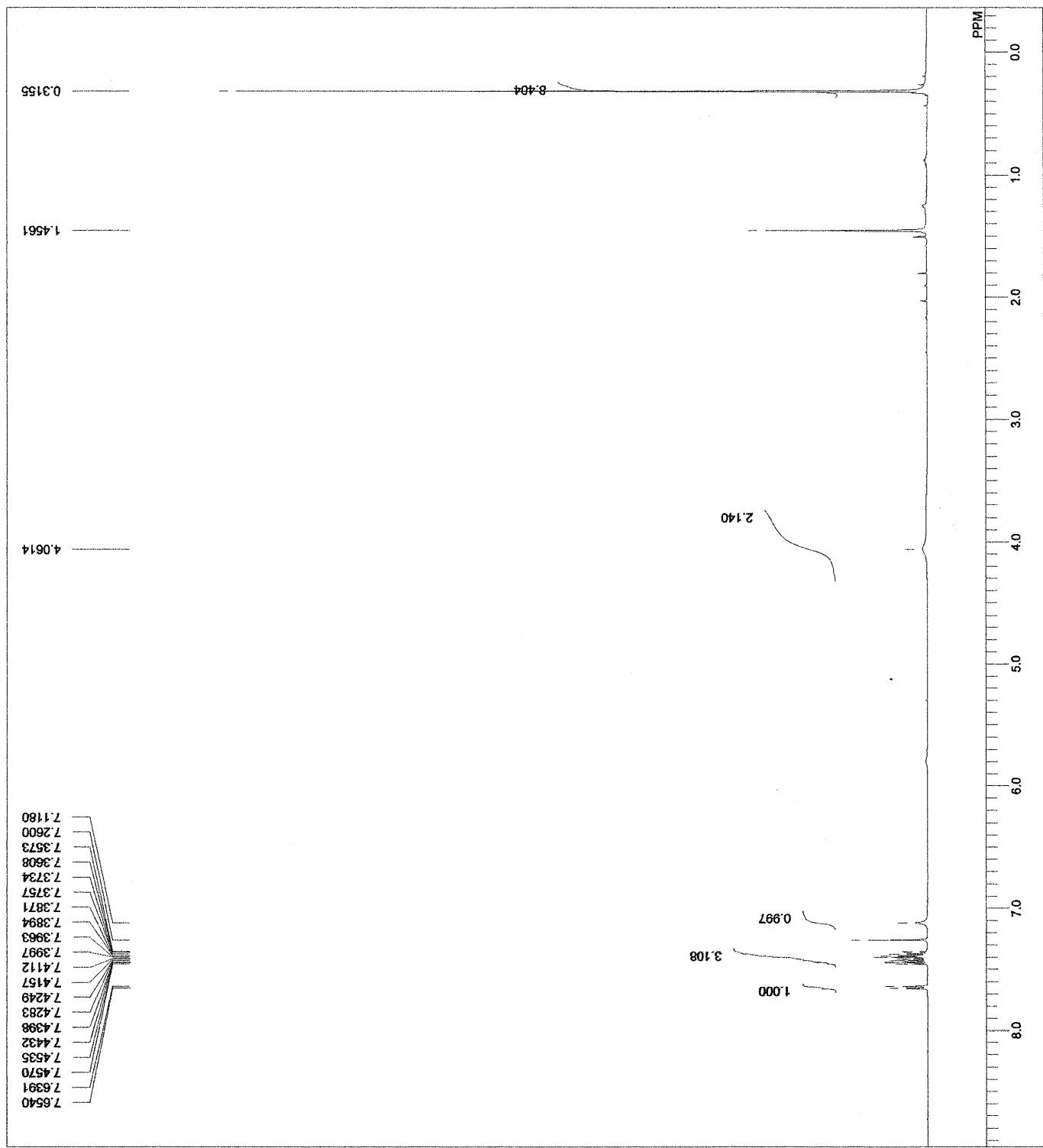


DFILE skk-p20-Ge-ArCO2Me-C-1-1.jdf
COMMENT skk-p20-Ge-ArCO2Me-C
DATIM 2013-09-17 16:38:03
OBNUC 13C
EXMOD carbon1.jxp
OBFREQ 125.77 MHz
OBSEIT 7.87 kHz
OBFIN 4.21 Hz
POINT 32767
FREQU 39308.18 Hz
SCANS 216
ACQTIM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usec
IRNUC 1H
CTEMP 25.3 c
SLVNT CDCl₃
EXREF 77.00 ppm
BF 1.20 Hz
RGAIN 60

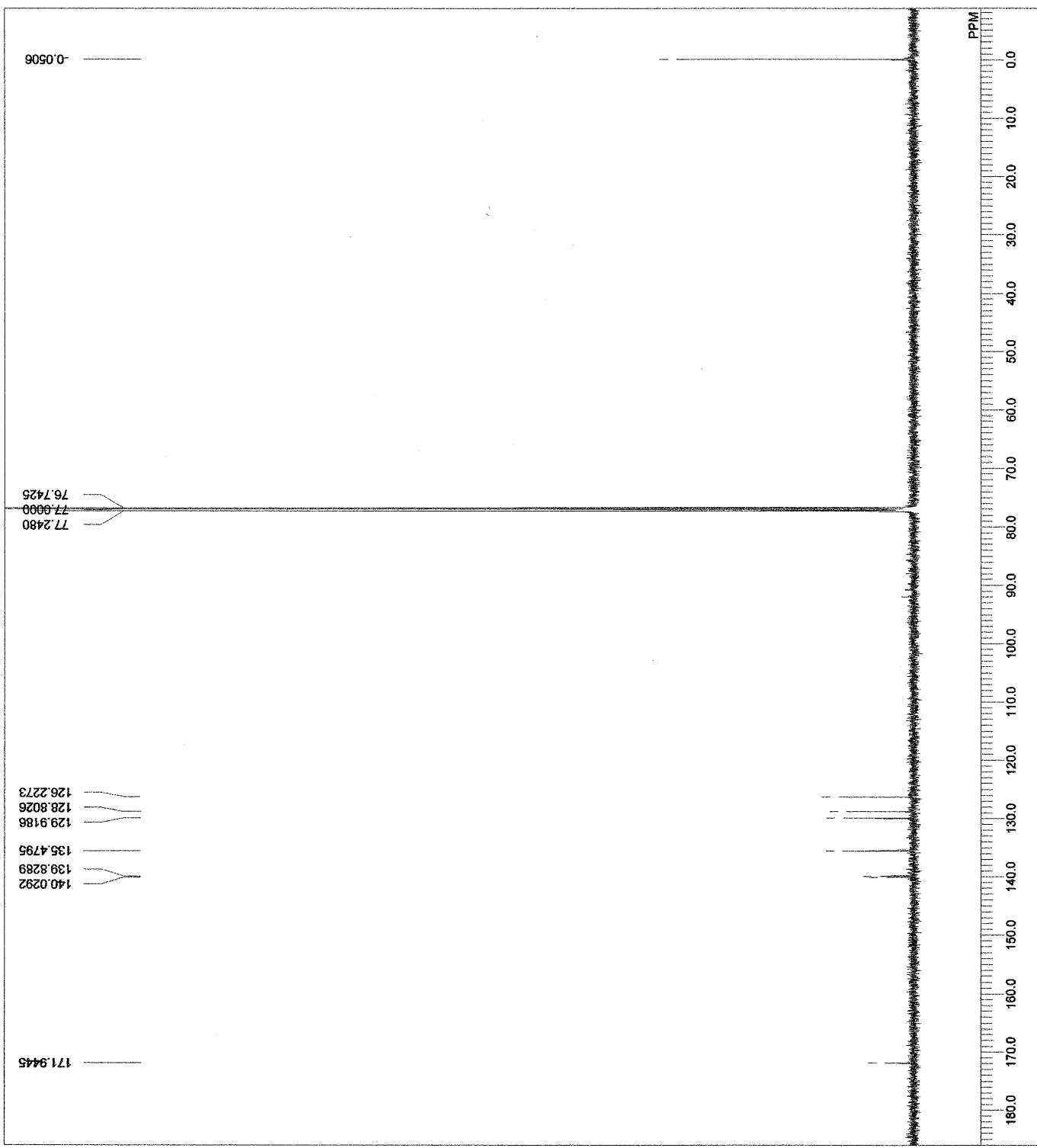
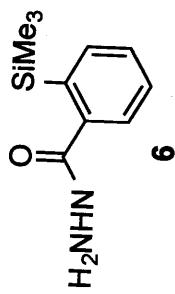








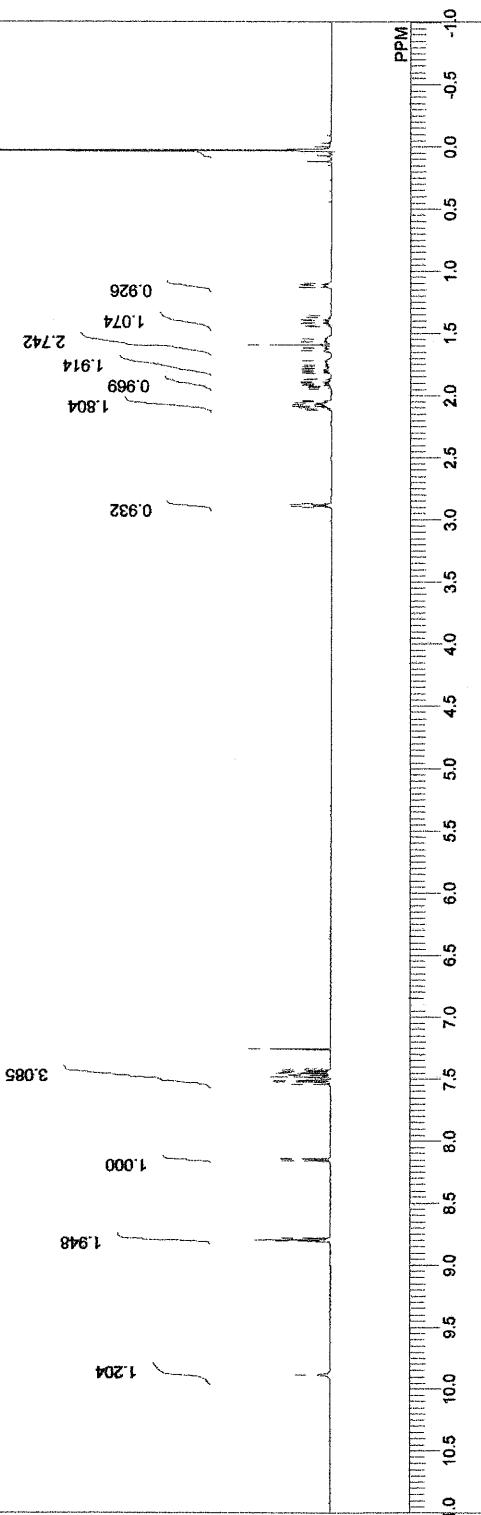
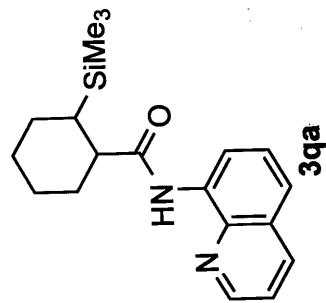
DFILE skk-hydrazone-1-1.dff
 COMNT skk-hydrazone-
 DATIM 2013-12-27 13:39:18
 OBNUC 13C
 EXMOD carbon-j,xp
 OBFRQ 125.77 MHz
 OFFSET 7.87 kHz
 OBFIN 4.21 Hz
 POINT 32767
 FREQU 39308.18 Hz
 SCANS 623
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IIRNUC 1H
 CTEMP 21.9 c
 SLVNT CDCl₃
 EXREF 77.00 ppm
 BF 1.00 Hz
 RGAIN 60



skk-p172-Cy-H-NMR

39a

DFILE	skk-p172-Cp-H-NMR-1.1.ais
COMNT	skk-p172-Cp-H-NMR
DATIM	2013-07-08 18:28:44
OBNUC	1H
EXMOD	protox1p9
OBFRQ	500.16 MHz
OBSET	2.41 kHz
OBFIN	6.01 Hz
POINT	13107
FREQU	7507.51 Hz
SCANS	16
ACQTM	1.7459 sec
PD [*]	5.0000 sec
PW1	5.55 usbc
IRNUC	1H
CTEMP	24.3 °C
SLVNT	CDCl ₃
EXREF	7.26 ppm
RFGAIN	0.12 Hz
	38



3qa

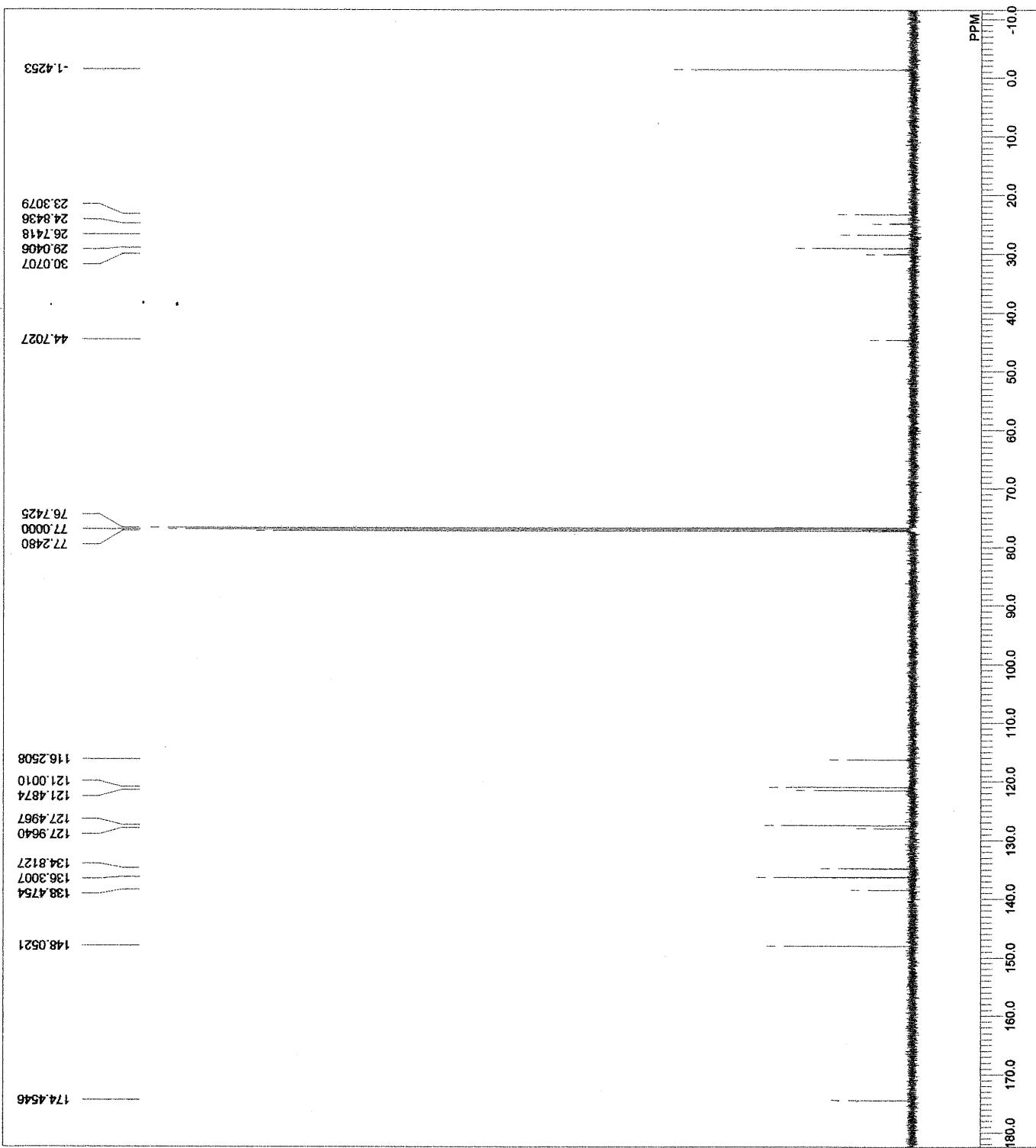
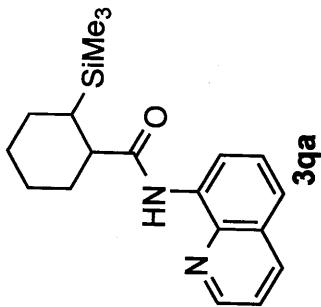
skk-p172-CyH-CNMR

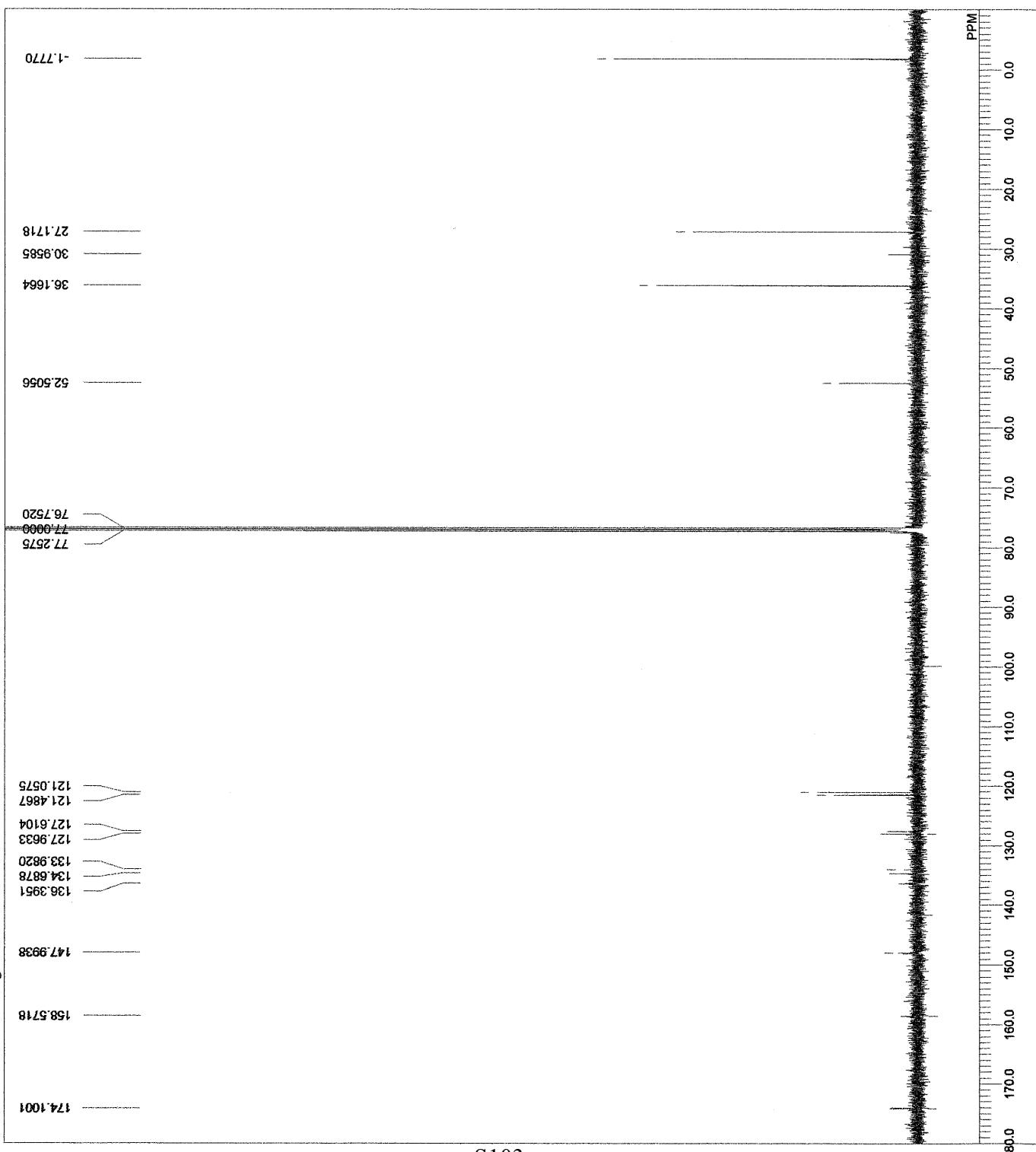
skk-p172-CyH-CNMR-1-1.als
skk-p172-CyH-CNMR
2013-07-08 18:31:45

13C
carbon,xp
OBFRQ 125.77 MHz
OBSET 7.87 kHz
OBFIN 4.21 Hz
POINT 26214

EXMOD 31446.54 Hz
SCANS 612
ACQTM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usec

FREQU 31446.54 Hz
IIRNUC 1H
CTEMP 24.9 C
SLVNT CDCl₃
EXREF 77.00 ppm
BF 0.12 Hz
RGAIN 60

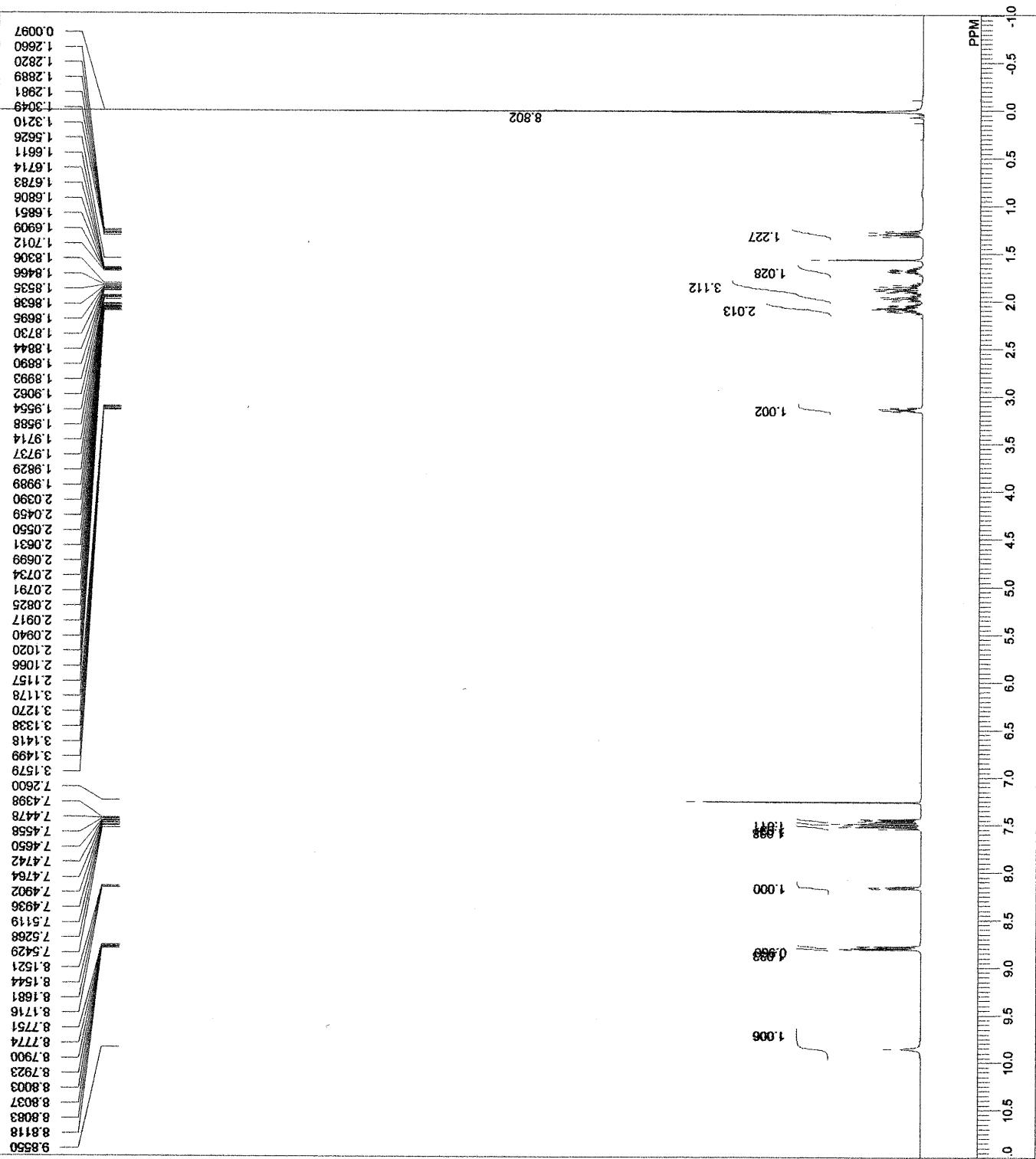
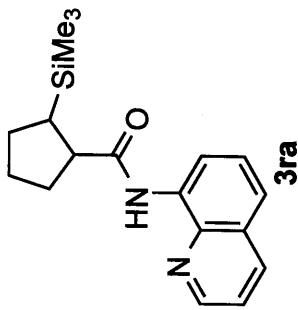


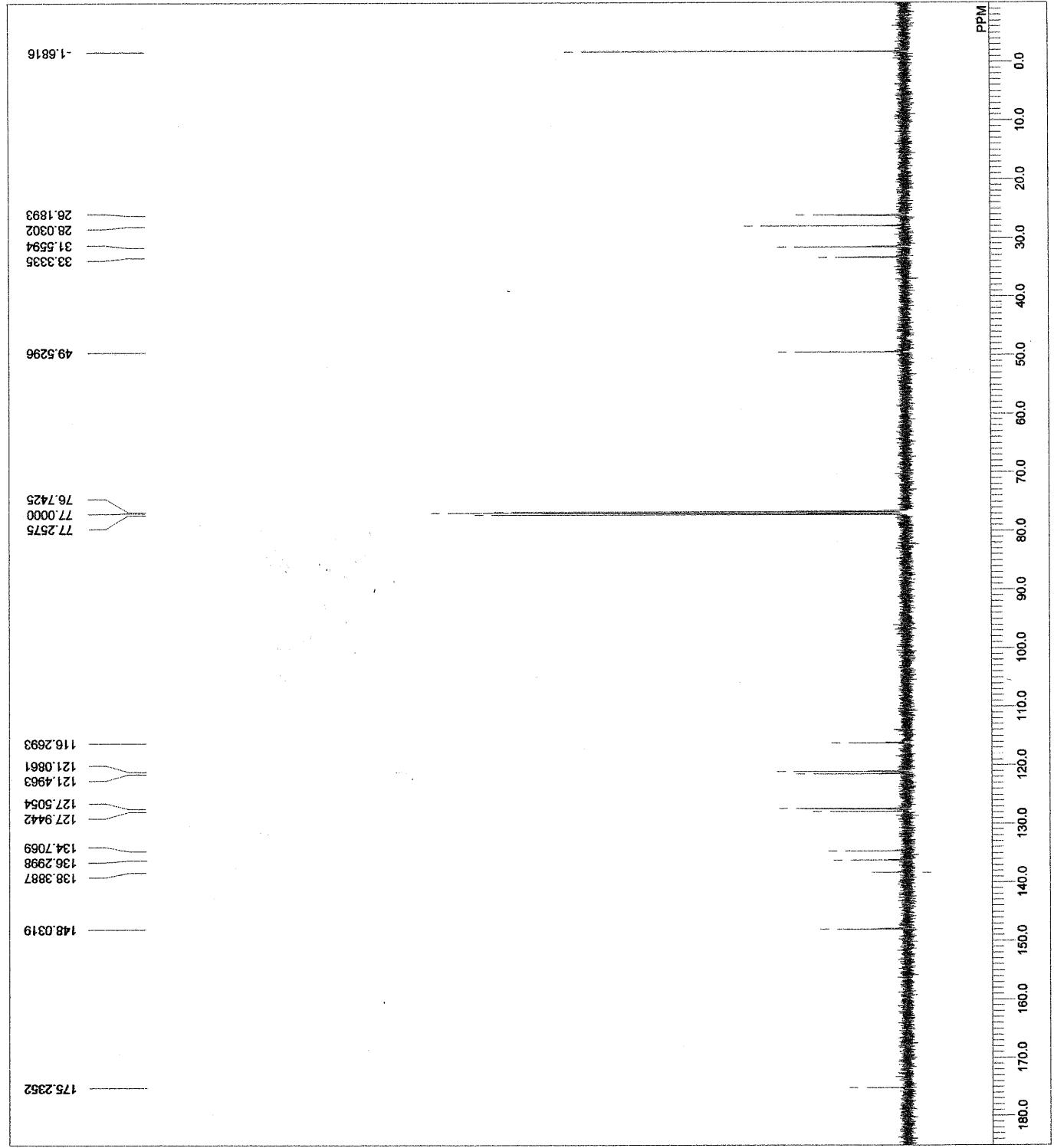


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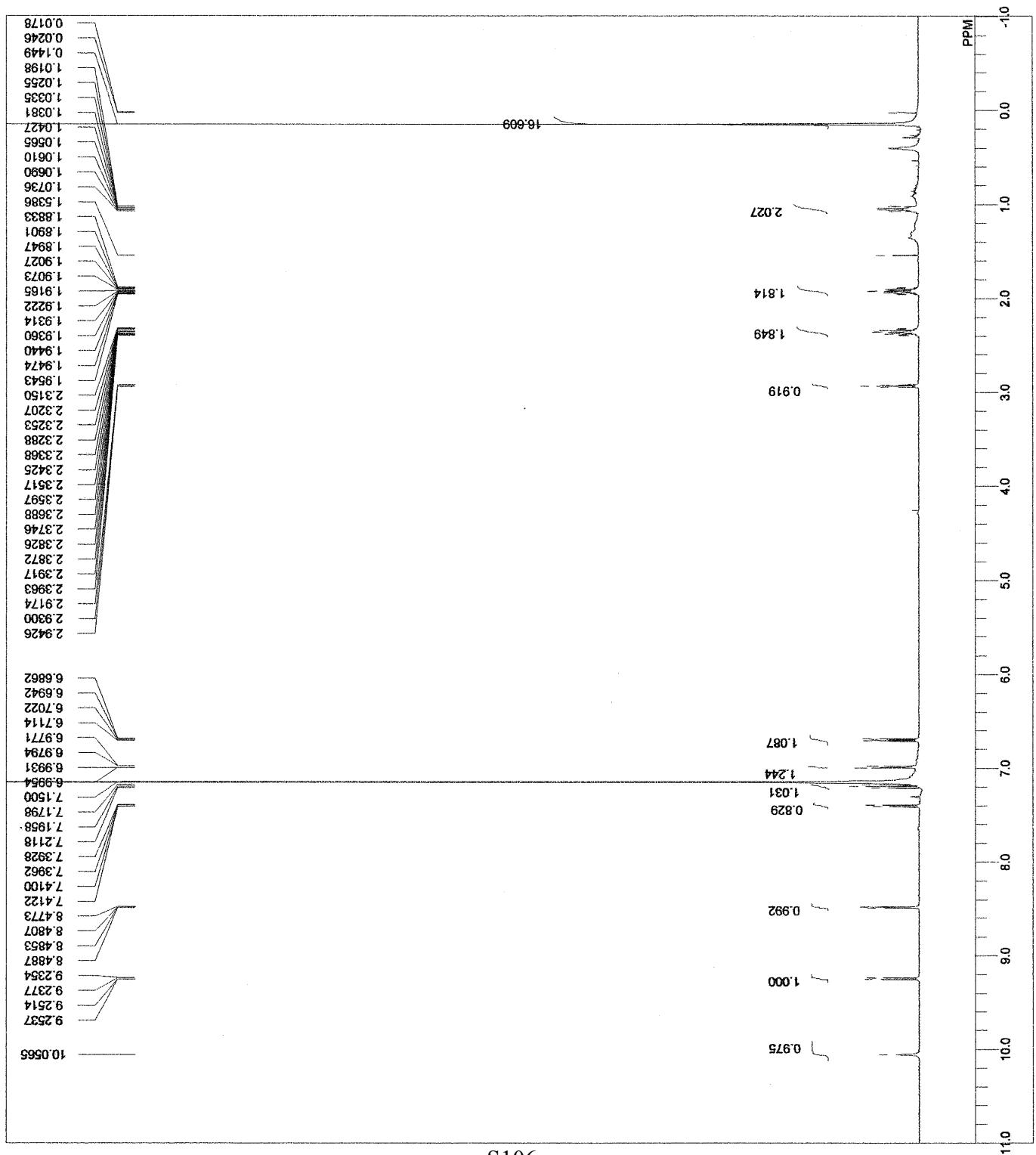
DFILE: skk-pg60-cycloPentane-SiMe3-H-3-1.ais
COMNT: skk-pg60-cycloPentane-SiMe3-H
DATIM: 2013-11-13 14:37:06
OBNUC: 1H
EXMOD: proton JRD
DPFRQ: 500.16 MHz
OBSET: 2.41 kHz
OBFIN: 6.01 Hz
POINT: 13107
FREQU: 7607.51 Hz
SCANS: 16
ACQTM: 1.7459 sec
PD: 5.0000 sec
PW1: 5.55 usec
IRNUC: 1H
CTEMP: 23.1 c
SLVNT: CDCl3
EXREF: 7.26 ppm
BF: 0.12 Hz
RGAIN: 32

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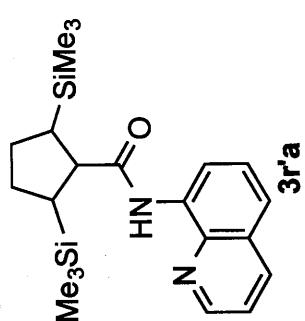




3r'a
skk-p73-inC6D6.g

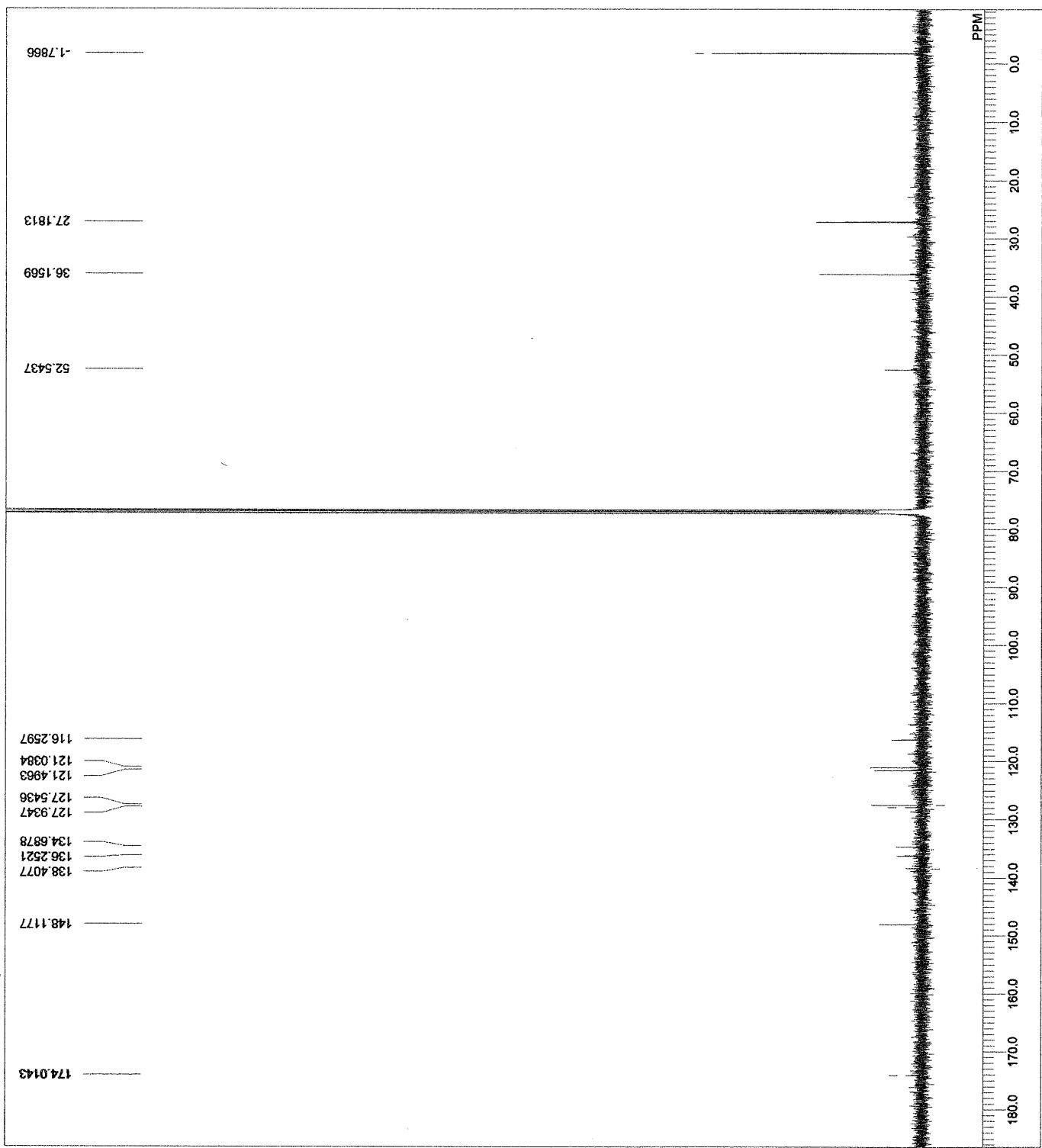
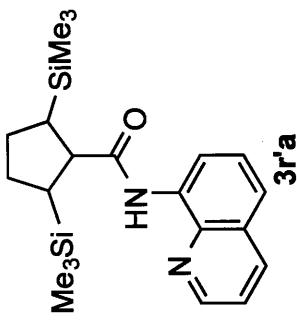


DFILE
COMNT
DATIM
OBNUC
EXMOD
OBFRQ
OBSET
OBFIN
POINT
FREQU
SCANS
8
ACQTM
PD
PW1
IRNUC
CTEMP
SLVNT
EXREF
BF
RGAIN
0.0178
0.0246
0.1449
0.1498
1.0255
1.0335
1.0381
1.0427
1.0565
1.0610
1.0690
1.0736
1.5386
1.8833
1.8901
1.8947
1.9027
1.9073
1.9165
1.9222
1.9314
1.9360
1.9440
1.9474
1.9543
2.3150
2.3207
2.3253
2.3288
2.3368
2.3425
2.3517
2.3597
2.3688
2.3746
2.3826
2.3872
2.3917
2.3963
2.9174
2.9300
2.9426
6.6862
6.6942
6.7022
6.7114
6.971
6.9794
6.9931
6.9954
7.1500
7.1798
7.1958
7.2118
7.3928
7.3962
7.4100
7.4122
8.4773
8.4807
8.4887
9.2377
9.2514
9.2537
10.0565
0.0178
0.0246
0.1449
0.1498
1.0255
1.0335
1.0381
1.0427
1.0565
1.0610
1.0690
1.0736
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1.8901
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1.9073
1.9165
1.9222
1.9314
1.9360
1.9440
1.9474
1.9543
2.3150
2.3207
2.3253
2.3288
2.3368
2.3425
2.3517
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2.3688
2.3746
2.3826
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2.3917
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2.9426
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6.6942
6.7022
6.7114
6.971
6.9794
6.9931
6.9954
7.1500
7.1798
7.1958
7.2118
7.3928
7.3962
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7.4122
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9.2537
10.0565

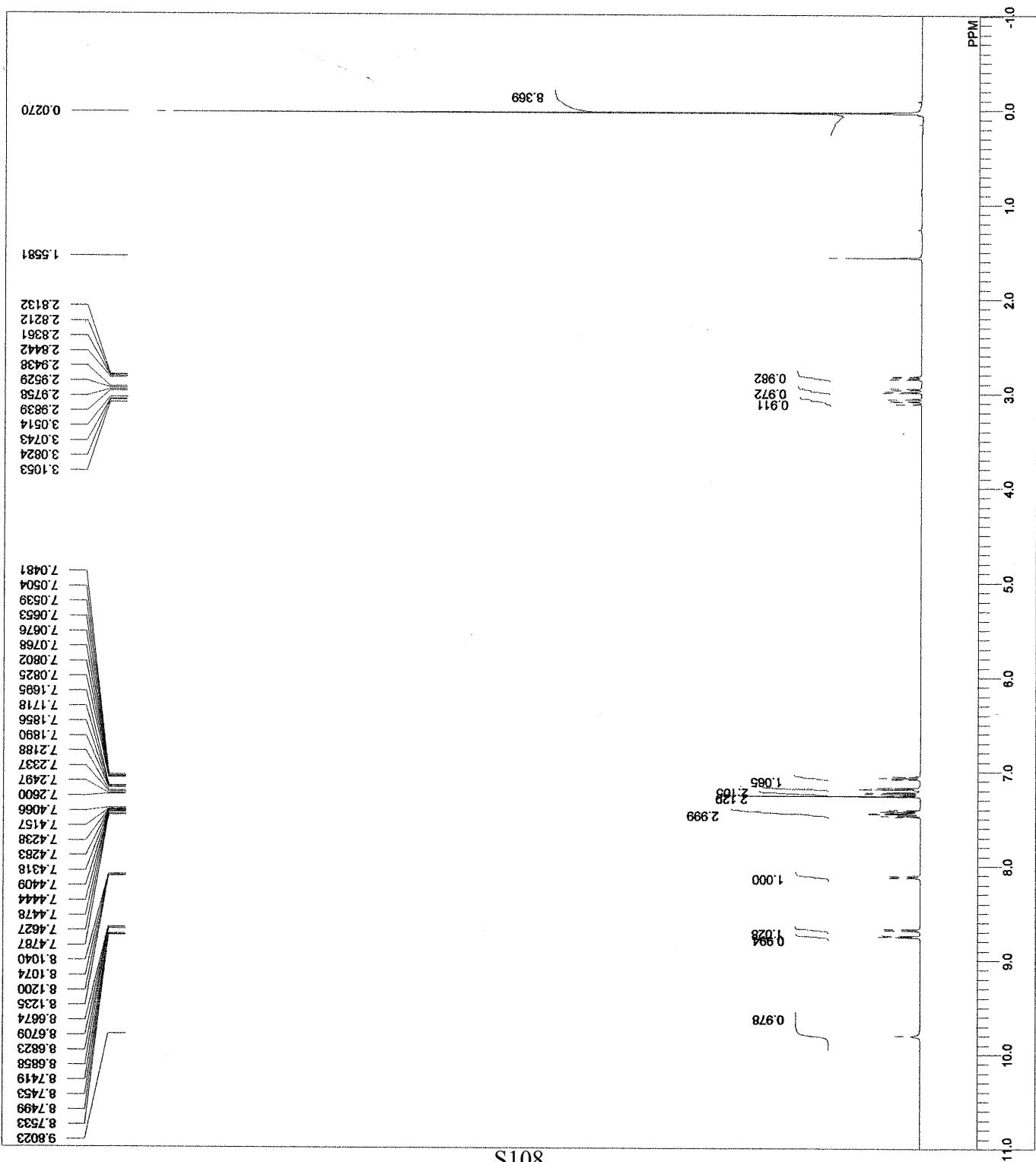


3r'a

DFILE skk-diSi-overnight-2-1.als
 skk-diSi-overnight
 2013-12-22 15:26:02
 13C
 carbon-13D
 EXMOD 125.77 MHz
 OBFRQ 7.87 kHz
 OBSET 4.21 Hz
 OBFIN 32767
 POINT 36308.18 Hz
 FREQU SCANS 17000
 ACQTM 0.8336 sec
 PD 3.0000 sec
 PW1 3.40 usec
 IRNUC 1H
 CTEMP 21.7 c
 SLYNT CDCL₃
 EXREF 77.00 ppm
 BF 0.12 Hz
 RGAIN 60



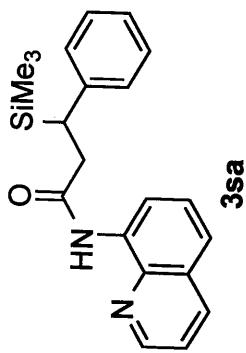
skk-p69-cinna-3sa

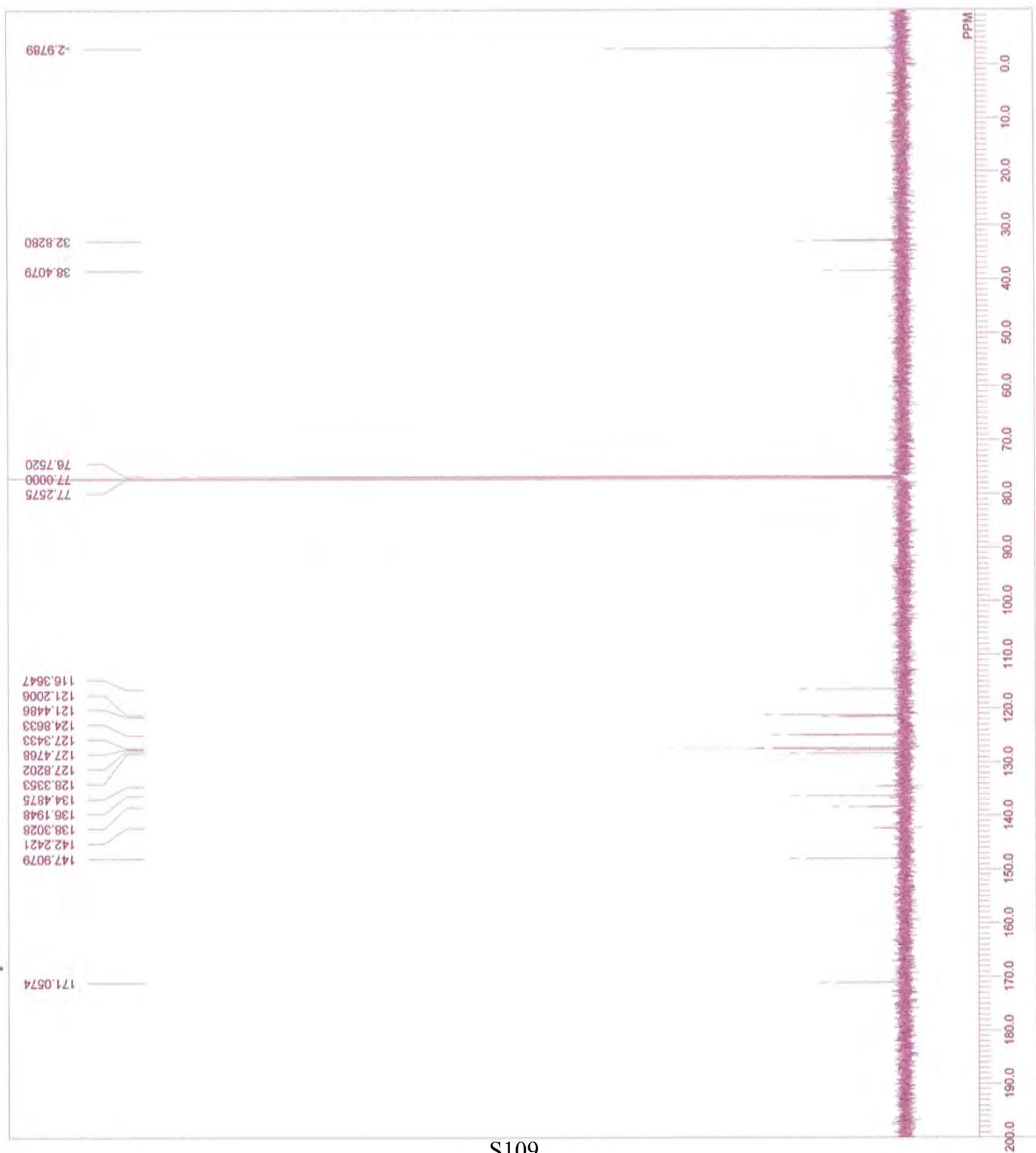


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skk-p69-cinna-1-1.jdp
skk-p69-cinna
2013-11-19 19:21:03
1H
proton.jdp
EXMOD
OBFRQ 500.16 MHz
OBSET 2.41 kHz
OBFIN 6.01 Hz
POINT 16384
FREQU 9384.38 Hz
SCANS 8
ACQTM PD
PW1 5.0000 sec
IRNUC 1H
CTEMP 22.8 c
SLVNT CDCl3
EAREF 7.26 ppm
BF 0.10 Hz
RGAIN 40

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skk-p174-TM-C-1-1.jdf
skk-p174-TM-C
2013-07-16 19:09:06
13C
carbon.jdp
DFILE
COMNT
DATTM
OBNUC
EXMOD
OBFRQ 125.77 MHz
OBSET 7.87 kHz
OBFIN 4.21 Hz
POINT 32767
FREQU 39308.18 Hz
SCANS 222
ACQTM 0.8336 sec
PD 3.0000 sec
PW1 3.40 usc
IRNUC 1H
CTEMP 24.4 c
SLVNT CDCl₃
EXREF 77.00 ppm
BF 0.12 Hz
RGAIN 60

