

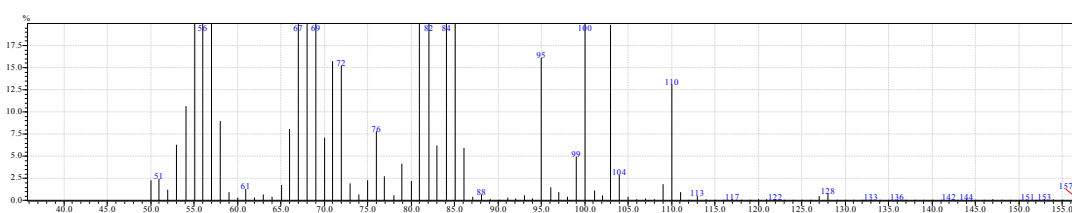
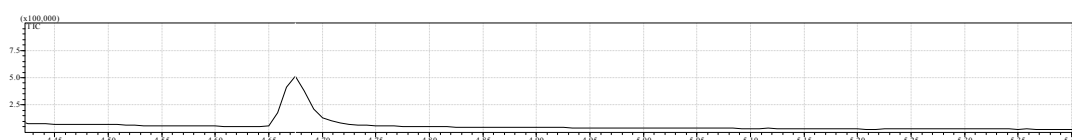
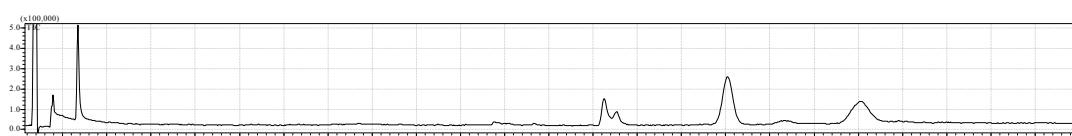
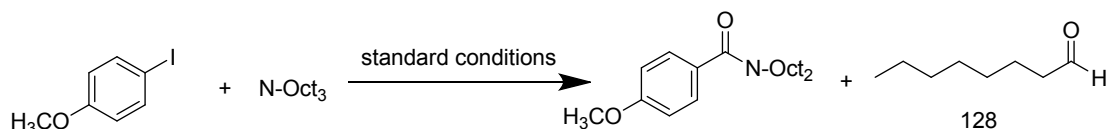
A Novel Pd-Catalyzed *N*-Dealkylative Carbonylation of Tertiary Amines for the Preparation of Amides

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General procedures : A Tube was charged with 4-iodoanisole(0.3 mmol, 70.2 mg), Pd(PhCN)₂Cl₂ (5 mol%, 5.7 mg), tributylamine(2.0 equiv. 111 mg),copper oxide(3.0 equiv, 72 mg) together with1mL PhCN . Once sealed,the tube was purged several times with CO(1atm,25mL,no balloon), then pressurized to 1 atm at room temperature and heated in an oil bath at 100 °C for 16 h. The tube was then cooled to room temperature and vented to discharge the excess CO. The reaction mixture was concentrated in *vacuo*, The crude residue was purified by silica gel chromatography.

Control Experiment.



[MS Spectrum]

of Peaks 547

Raw Spectrum 4.675 (scan : 82)

Background No Background Spectrum

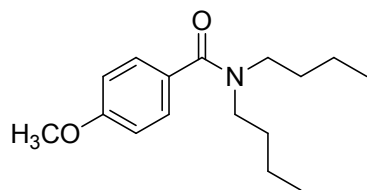
Base Peak m/z 84.05 (Inten : 59,993)

m/z Absolute Intensity Relative Intensity

50.00	1364	2.27	73.00	1146	1.91
51.00	1476	2.46	73.95	392	0.65
52.00	732	1.22	74.95	1345	2.24
53.00	3772	6.29	76.00	4655	7.76
54.05	6399	10.67	76.95	1628	2.71
55.05	43386	72.32	78.05	354	0.59
56.05	59555	99.27	78.95	2461	4.10
57.05	57408	95.69	80.05	1302	2.17
58.05	5377	8.96	81.00	26465	44.11
59.00	553	0.92	82.05	27777	46.30
60.00	210	0.35	83.00	3722	6.20

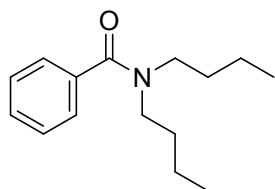
84.05	59993	100.00	105.00	230	0.38
85.05	21268	35.45	106.00	79	0.13
86.05	3549	5.92	107.00	142	0.24
87.10	257	0.43	108.00	65	0.11
88.10	404	0.67	109.05	1075	1.79
91.10	210	0.35	110.05	7808	13.01
92.10	116	0.19	111.05	562	0.94
93.10	335	0.56	112.00	49	0.08
94.00	159	0.27	113.00	294	0.49
95.00	9663	16.11	114.00	94	0.16
96.10	858	1.43	115.00	42	0.07
97.05	546	0.91	116.00	34	0.06
98.00	214	0.36	117.00	71	0.12
99.00	2985	4.98	122.00	74	0.12
100.05	12144	20.24	123.00	26	0.04
101.10	673	1.12	127.00	297	0.50
102.05	367	0.61	128.00	455	0.76
102.95	11866	19.78	129.00	52	0.09
104.00	1758	2.93	130.00	13	0.02

Characterization Data for Products



N,N-Dibutyl-4-methoxybenzamide⁽¹⁾ **3a**

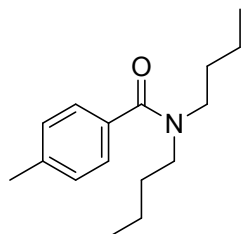
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ: 7.15 (d, *J*=8.5 Hz, 2H), 6.70 (d, *J*=8.5 Hz, 2H), 3.60 (s, 3H), 3.28 (s, 2H), 3.09 (s, 2H), 1.38 (m, 4H), 1.09 (m, 4H), 0.77 (s, 3H), 0.64 (s, 3H); ¹³C NMR (125 MHz, CDCl₃) δ: 171.0, 159.9, 129.3, 127.9, 112.9, 54.6, 48.4, 44.0, 30.2, 29.2, 19.3, 13.2; LRMS (EI 70 eV) *m/z* (%): 263 (M⁺, 7), 220 (6), 135 (100), 107 (7), 77 (9).



N,N-dibutylbenzamide⁽¹⁾ **3b**

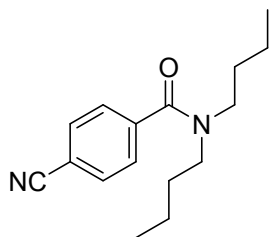
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ: 7.33–7.24 (m, 5H), 3.41 (s, 2H), 3.11 (s, 2H), 1.57 (s, 2H), 1.46–1.27 (m, 4H), 1.02 (m, 5H), 0.70 (s, 3H); ¹³C NMR (125

MHz,CDCl₃) δ : 171.6, 137.3, 128.9, 128.2, 126.4, 8.7, 44.4, 30.7, 29.6, 20.2, 19.6, 13.8, 13.5; LRMS(EI70ev)*m/z*(%): 233 (M⁺,7.6), 190 (12.8), 148 (5.2), 105 (100), 77 (24).



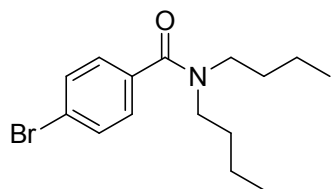
N,N-Dibutyl-4-methylbenzamide⁽¹⁾ **3c**

Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ : 7.21 (d, *J* = 7.5 Hz, 2H), 7.13 (d, *J* = 7.5 Hz, 2H), 3.43 (s, 2H), 3.16 (s, 2H), 2.34 (s, 3H), 1.59 (s, 2H), 1.44 (s, 2H), 1.35 (s, 2H), 1.10 (s,2H), 0.92 (s,3H), 0.75 (s,3H); ¹³CNMR (125MHz,CDCl₃) δ : 171.7, 138.8, 134.3, 128.8, 126.4, 48.7, 44.3, 30.7, 29.6, 21.2, 20.7, 19.6, 14.0, 13.8; LRMS (EI70ev)*m/z*(%): 247 (M⁺,9), 204 (9), 119 (100), 120 (9), 65 (7).



N,N-Dibutyl-4-cyanobenzamide⁽²⁾ **3d**

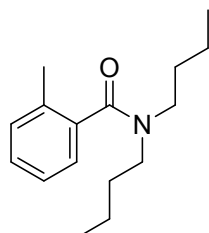
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ : 7.63 (d, *J* =8.0 Hz, 2H), 7.39 (d, *J* = 8.0 Hz, 2H), 3.43-3.40(m, 2H), 3.10–3.00 (m, 2H), 1.63–1.50 (m, 2H), 1.43-1.31(m, 4H),1.08-.04(m, 2H),0.92-0.89(m,3H),0.73-0.70(m,3H);¹³C NMR (125 MHz, CDCl₃) δ :169.6,142.0,132.3,127.1,118.3,112.9,48.7,44.5,30.7,29.5,20.2, 19.6, 13.8, 13.5; LRMS(EI70ev)*m/z*(%):258(M⁺,3.4),215(14.5),173(9.4),130(100),102(24.2).



4-bromo-N,N-Dibutylbenzamide **3e**

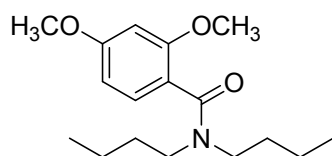
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ :7.54 (d, *J* =8.0 Hz, 2H), 7.25 (d, *J* = 8.5 Hz, 2H), 3.48 (s, 2H), 3.18 (s, 2H), 1.64 (s, 2H), 1.49 (s, 2H), 1.41 (s, 2H), 1.19-1.15

(m,2H), 0.98 (s,3H), 0.82 (s,3H); ^{13}C NMR (125MHz,CDCl₃) δ : 170.7, 136.3, 131.3, 127.8, 122.7, 48.8, 44.7, 30.80, 29.6, 20.3, 19.7, 13.9, 13.6; LRMS (EI70ev) m/z (%): 311(M⁺,8.7), 313 (8.9), 270 (18.8), 183 (100),155 (20),76 (20); IR(neat,cm⁻¹): 2954, 2868, 1633, 1421, 831, 572. HRMS m/z (ESI) for C₁₅H₂₃BrNO (M+H)⁺calcd : 312.0958, found: 312.0952.



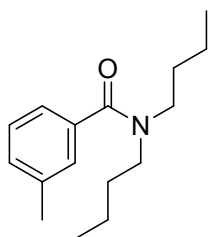
N,N-Dibutyl-2-methylbenzamide⁽¹⁾ **3f**

Yellow liquid; ^1H NMR(500MHz,CDCl₃) δ : 7.27–7.22 (m,1H), 7.20-7.17 (m,2H), 7.16–7.13 (m,1H), 3.70 (s,1H), 3.32 (s,1H), 3.09–3.01 (m, 2H), 2.29 (s, 3H), 1.71–1.61 (m, 2H), 1.44-1.38 (m, 4H), 1.16–1.06 (m, 2H), 0.99 (t, $J = 7.5$ Hz, 3H), 0.76 (t, $J = 7.5$ Hz, 3H); ^{13}C NMR (125 MHz, CDCl₃) δ : 170.8, 137.2, 133.8, 130.2, 128.5, 125.8, 125.7, 48.0, 44.0, 30.6, 29.6, 20.4, 19.7, 18.9, 13.9, 13.5; LRMS (EI70ev) m/z (%): 247 (M⁺,6.4), 202 (8.4), 145 (55.2), 119 (100),77 (30.3).



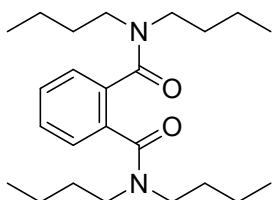
N,N-dibutyl-2,4-dimethoxybenzamide⁽³⁾ **3g**

Yellow liquid; ^1H NMR (500 MHz, CDCl₃) δ : 7.11 (d, $J = 8.0$ Hz, 1H), 6.53–6.41 (m, 2H), 3.81 (s, 3H), 3.78 (s, 3H), 3.69–3.27 (m, 2H), 3.08 (t, $J = 7.5$ Hz,2H),1.67–1.58 (m, 2H), 1.39-1.36 (m, 4H), 1.14-1.09 (m, 2H), 0.97 (t, $J = 7.5$ Hz, 3H), 0.76 (t, $J = 7.5\text{Hz}$,3H); ^{13}C NMR (125MHz, CDCl₃) δ : 169.2, 161.2, 156.6, 128.5, 119.9, 104.6, 98.5, 55.6 (2C), 48.3, 44.0, 30.6, 29.3, 20.1, 19.7, 13.9, 13.6; LRMS (EI70ev) m/z (%): 293 (M⁺,7.9), 250 (6.2),194 (6), 165(100), 122(5.7).



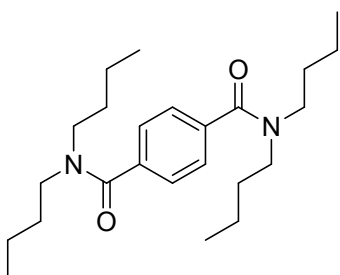
N,N-Dibutyl-3-methylbenzamide⁽⁴⁾ **3h**

Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ: 7.25 (t, *J* = 7.5 Hz, 1H), 7.19–7.14 (m, 2H), 7.12 (d, *J* = 7.5 Hz, 1H), 3.48 (s, 2H), 3.18 (s, 2H), 2.36 (s, 3H), 1.64 (s, 2H), 1.48 (s, 2H), 1.40 (s, 2H), 1.14 (s, 2H), 0.97 (s, 3H), 0.78 (s, 3H); ¹³C NMR (125 MHz, CDCl₃)δ: 171.7, 137.9, 137.2, 129.6, 128.0, 126.9, 123.3, 48.7, 44.3, 30.7, 29.5, 21.2, 20.2, 19.6, 13.8, 13.5; LRMS (EI70ev) *m/z*(%): 247(M⁺,11.3), 204(11.6), 162(4.7), 119(100), 91(22.7).



*N*¹,*N*¹,*N*²,*N*²-tetrabutylphthalamide⁽⁴⁾ **3i**

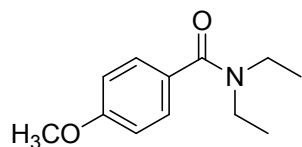
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ: 7.38-7.31 (m, 2H), 7.26-7.20 (m, 2H), 3.38 (s, 4H), 3.16–3.07 (m, 4H), 1.56-1.62 (m, 4H), 1.44-1.47 (m, 4H), 1.33-1.37 (m, 4H), 1.10 (m, 4H), 0.93 (t, *J* = 7.5 Hz, 6H), 0.76 (t, *J* = 7.5 Hz, 6H); ¹³C NMR (125 MHz, CDCl₃)δ: 169.6, 134.9, 128.3, 126.1, 48.8, 44.6, 30.5, 29.5, 20.3, 19.7, 13.9, 13.5; LRMS (EI70ev) *m/z*(%): 388(M⁺, 7.2), 260(17.3), 204(18.3), 128(100), 105(12.5).



*N*¹,*N*¹,*N*⁴,*N*⁴-tetrabutylterephthalamide **3j**

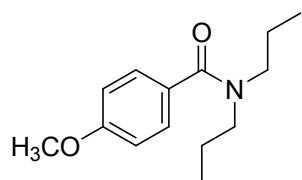
Yellow solid, mp 72.1~74.0°C; ¹H NMR (500 MHz, CDCl₃) δ 7.38 (s, 4H), 3.53–3.43 (m, 4H), 3.22–3.12 (m, 4H), 1.64 (d, *J*=7.0 Hz, 4H), 1.46 (d, *J*=6.5, 4H), 1.43-1.39 (m, 4H), 1.15-1.11 (m, 4H), 0.98 (t, *J* = 7.0 Hz, 6H), 0.80 (t, *J* = 7.0 Hz, 6H); ¹³C

NMR (125 MHz, CDCl₃) δ 171.0, 138.0, 126.5, 48.7, 44.4, 30.8, 29.6, 20.2, 19.7, 13.8, 13.6; IR(KBr, cm⁻¹): 2948, 2921, 2862, 1620, 1504, 1454, 1424, 1106, 736;
LRMS(EI70ev) m/z (%): 388(M⁺, 6.5), 281(12.9), 260(100), 105(22.6), 73 (16);
HRMS m/z (ESI) for C₂₄H₄₁N₂O₂ (M+H)⁺ calcd : 389.3163, found: 389.3150.



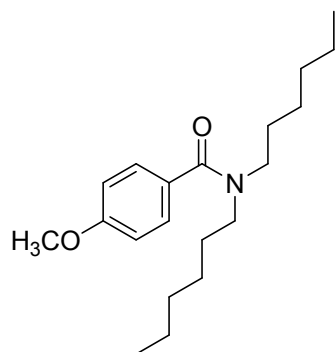
N,N-Diethyl-4-methoxybenzamide⁽⁵⁾ **3k**

Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ : 7.27 (d, *J* = 8.5 Hz, 2H), 6.82 (d, *J* = 8.5 Hz, 2H), 3.75 (s, 3H), 3.38 (m, 2H), 3.30 (m, 2H), 1.10 (s, 6H); ¹³C NMR (125 MHz, CDCl₃) δ : 171.4, 160.1, 129.4, 128.0, 113.7, 55.3, 43.4, 39.5, 14.0, 13.3;
LRMS (EI 70 ev) m/z (%): 207(M⁺, 9.6), 207(10.3), 135(100), 92(8.9), 77(14.8).



N,N-Dipropyl-4-methoxybenzamide⁽⁶⁾ **3l**

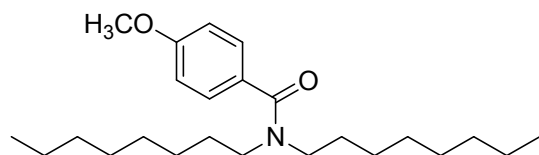
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ : 7.23 (d, *J* = 8.5 Hz, 2H), 6.80 (d, *J* = 8.5 Hz, 2H), 3.73 (s, 3H), 3.31 (s, 2H), 3.14 (s, 2H), 1.53 (s, 4H), 0.85 (s, 3H), 0.69 (s, 3H); ¹³C NMR (125 MHz, CDCl₃) δ : 171.7, 160.2, 129.6, 128.3, 113.6, 55.2, 51.3, 21.3, 11.2; LRMS(EI70ev) m/z (%): 235(M⁺, 8.4), 164(6.1), 135(100), 107(6.4), 77(11.3).



N,N-dihexyl-4-methoxybenzamide **3m**

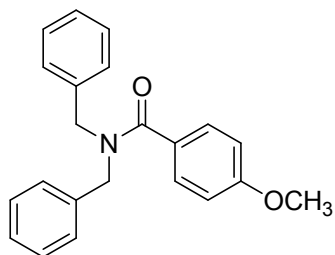
Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ : 7.32 (d, *J* = 8.5 Hz, 2H), 6.89 (d, *J* = 8.5 Hz, 2H), 3.82 (s, 3H), 3.44 (s, 2H), 3.23 (s, 2H), 1.61-1.52 (m, 4H), 1.29-1.16 (m,

12H), 0.87(s, 6H); ^{13}C NMR (125MHz, CDCl_3) δ : 171.5, 160.2, 129.7, 128.3, 113.6, 55.2, 49.1, 44.9, 31.3, 26.6, 22.5, 13.9; LRMS (EI70ev) m/z (%): 319 (M^+ , 5.0), 248(6.4), 192 (4.8), 135 (100), 77 (6.6); IR(neat, cm^{-1}): 2931, 2857, 1633, 1297, 1028, 843. HRMS m/z (ESI) for $\text{C}_{20}\text{H}_{34}\text{NO}_2$ ($\text{M}+\text{H}$) $^+$ calcd : 320.2584, found: 320.2578.



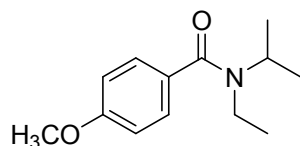
4-methoxy-N,N-dioctylbenzamide 3n

Yellow liquid; ^1H NMR (500 MHz, CDCl_3) δ : 7.31 (d, $J = 8.5$ Hz, 2H), 6.89 (d, $J = 9.0$ Hz, 2H), 3.82 (s, 3H), 3.43 (s, 2H), 3.29 (s, 2H), 1.59 (s, 4H), 1.27 (s, 20H), 0.88 (t, $J = 10.0$, 6H); ^{13}C NMR(125MHz, CDCl_3) δ : 171.7, 160.4, 129.7, 128.3, 113.7, 55.2, 47.4, 42.1, 31.7, 29.1, 28.6, 26.9, 22.7, 14.3; LRMS (EI70ev) m/z (%): 375 (M^+ , 5.8), 276 (8.1), 220 (4.7), 135 (100), 156 (6.8); IR(neat, cm^{-1}): 2921, 2842, 1633, 1461, 839, 765; HRMS m/z (ESI) for $\text{C}_{24}\text{H}_{42}\text{NO}_2$ ($\text{M}+\text{H}$) $^+$ calcd : 376.3210, found: 376.3205.



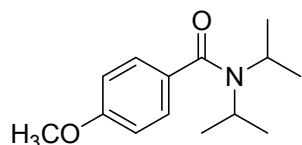
***N,N*-Dibenzyl-4-methoxybenzamide⁽⁷⁾ 3o**

White solid; ^1H NMR (500 MHz, CDCl_3) δ : 7.47 (d, $J = 8.0$ Hz, 2H), 7.31-7.17 (m, 10H), 6.84 (d, $J = 8.0$ Hz, 2H), 4.66 (s, 2H), 4.48 (s, 2H), 3.71 (s, 3H); ^{13}C NMR (125MHz, CDCl_3) δ : 172.0, 160.7, 136.6, 128.7, 127.7, 127.3, 127.1, 113.3, 54.9, 51.4, 47.0; Melting point : 114-116 $^\circ\text{C}$; LRMS (EI70ev) m/z (%): 331 (M^+ , 2.3) , 240(28.9), 135(100), 107(8.8), 77(12.4).



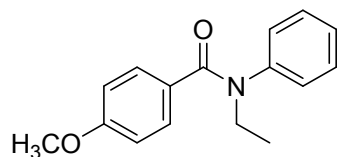
***N*-ethyl-*N*-isopropyl-4-methoxybenzamide 3p**

Yellow liquid; ^1H NMR (500 MHz, CDCl_3) δ : 7.16 (d, $J = 8.5$ Hz, 2H), 6.75 (d, $J = 8.5$ Hz, 2H), 3.66 (s, 4H), 3.21 (s, 2H), 1.03 (s, 12H); ^{13}C NMR (125 MHz, CDCl_3) δ : 171.2, 159.9, 128.6, 127.7, 113.5, 54.8, 50.2, 41.5, 22.5, 20.8; IR(neat, cm^{-1}): 2973, 1609, 1250, 1024, 840, 591; LRMS (EI70ev) $m/z(\%)$: 221 (M^+ , 9.4), 192 (5.9), 135 (100), 107 (6.7), 77 (11.5). HRMS m/z (ESI) for $\text{C}_{13}\text{H}_{20}\text{NO}_2$ ($\text{M}+\text{H}$) $^+$ calcd : 222.1489, found: 222.1483



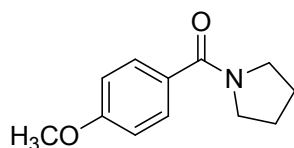
N,N-diisopropyl-4-methoxybenzamide⁽⁸⁾ **3q**

Yellow liquid; ^1H NMR (500 MHz, CDCl_3) δ : 7.19 (d, $J = 8.5$ Hz, 2H), 6.80 (d, $J = 8.5$ Hz, 2H), 3.73 (s, 5H), 1.25 (s, 12H). ^{13}C NMR (125 MHz, CDCl_3) δ : 170.9, 159.9, 128.6, 127.5, 113.5, 54.8, 20.8; LRMS (EI70ev) $m/z(\%)$: 235 (M^+ , 6.5), 192 (18.1), 135 (100), 107 (5.7), 77 (9.3).



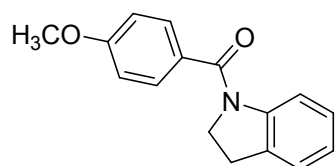
N-ethyl-*N*-phenyl-4-methoxybenzamide **3r**

Yellow liquid; ^1H NMR (500 MHz, CDCl_3) δ : 7.15 (dd, $J = 9.0, 8.0$ Hz, 4H), 7.05 (t, $J = 7.5$ Hz, 1H), 6.94 (d, $J = 7.5$ Hz, 2H), 6.55 (d, $J = 9.0$ Hz, 2H), 3.88 (q, $J = 7.0$ Hz, 2H), 3.62 (s, 3H), 1.12 (t, $J = 7.0$ Hz, 3H); ^{13}C NMR (125 MHz, CDCl_3) δ : 169.6, 160.9, 143.6, 130.7, 129.0, 128.3, 127.7, 126.3, 112.6, 55.3, 45.3, 12.6. LRMS (EI70ev) $m/z(\%)$: 255 (M^+ , 14.2), 135 (100), 107 (9.4), 92 (8.7), 77 (17.7). IR (neat, cm^{-1}): 2970, 2935, 1633, 1299, 840, 697. HRMS m/z (ESI) for $\text{C}_{16}\text{H}_{18}\text{NO}_2$ ($\text{M}+\text{H}$) $^+$ calcd : 256.1332, found: 256.1340.



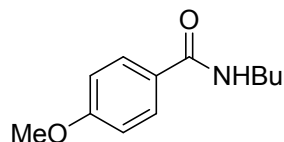
(4-methoxyphenyl)(pyrrolidin-1-yl)methanone⁽⁹⁾ **3s**

White solid, mp 78-79 °C; ¹H NMR (500 MHz, CDCl₃) δ 7.52 (d, *J* = 9.0 Hz, 2H), 6.90 (d, *J* = 8.5 Hz, 2H), 3.84 (s, 3H), 3.64 (t, *J* = 6.9 Hz, 2H), 3.48 (t, *J* = 6.6 Hz, 2H), 1.99 – 1.92 (m, 2H), 1.91 – 1.83 (m, 2H); ¹³C NMR (125 MHz, CDCl₃) δ 169.5, 160.8, 129.5, 129.2, 113.4, 55.3, 49.8, 46.3, 26.5, 24.5; LRMS (EI 70ev) *m/z* (%) : 205 (M⁺, 23.5), 135 (100), 107(10.9), 92 (9.8), 77 (16.9).



indolin-1-yl(4-methoxyphenyl)methanone **3t**

White solid, mp 102-103°C; ¹H NMR (500 MHz, CDCl₃) δ: 7.53 (d, *J* = 8.5 Hz, 2H), 7.19 (d, *J* = 7.5 Hz, 1H), 7.10 (s, 1H), 6.99 (t, *J* = 7.1 Hz, 1H), 6.93 (d, *J* = 8.5 Hz, 2H), 4.10 (t, *J* = 8.0 Hz, 2H), 3.84 (s, 3H), 3.09 (t, *J* = 8.0 Hz, 2H); ¹³C NMR (125 MHz, CDCl₃) δ 168.9, 161.4, 142.9, 132.5, 129.4, 129.1, 127.2, 124.9, 123.7, 113.8, 55.4, 50.7, 28.2; IR(KBr, cm⁻¹): 2916, 1633, 1482, 1246, 1028, 845, 753; LRMS (EI 70ev) *m/z* (%) : 253(12.2), 135(100), 107(10.9), 92(11.5), 77(16.6); HRMS *m/z* (ESI) for C₁₆H₁₆NO₂ (M+H)⁺ calcd : 254.1176, found: 254.1174.



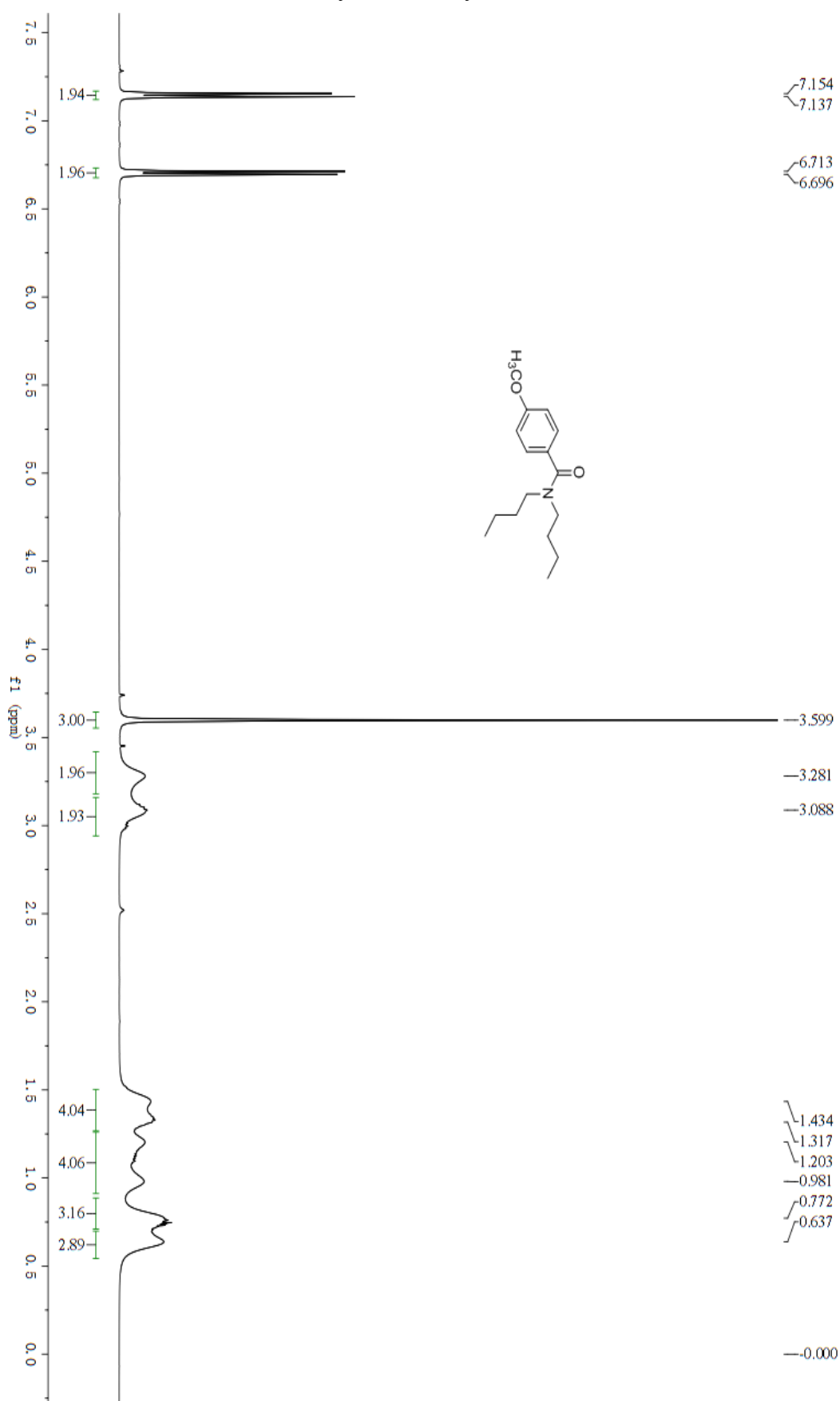
N-butyl-4-methoxybenzamide⁽¹⁰⁾ **3u**

Yellow liquid; ¹H NMR (500 MHz, CDCl₃) δ: 7.73 (d, *J* = 9.0 Hz, 2H), 6.91 (d, *J* = 9.0 Hz, 2H), 6.11 (s, 1H), 3.84 (s, 3H), 3.44 (dd, *J*₁ = 13.0, *J*₂ = 7.0 Hz, 2H), 1.64-1.54 (m, 2H), 1.41 (m, 2H), 0.95 (t, *J* = 7.5 Hz, 3H); ¹³C NMR (125 MHz, CDCl₃) δ 167.0, 162.1, 128.6, 127.2, 113.7, 55.4, 39.7, 31.8, 20.1, 13.7.; LRMS (EI70ev) *m/z*(%): 207 (M⁺, 13.1), 164 (13.3), 135(100), 92 (9.3), 77 (11.0).

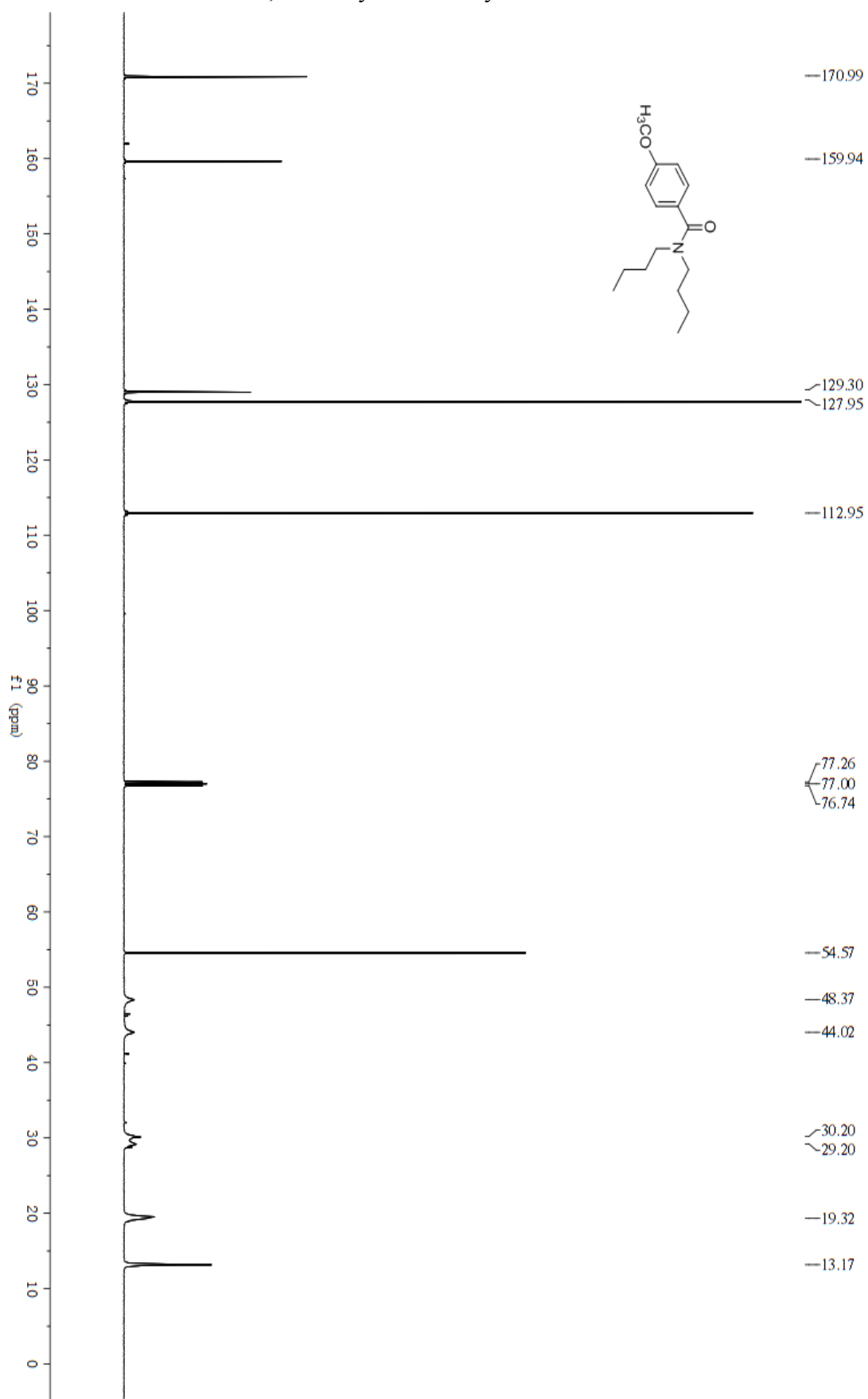
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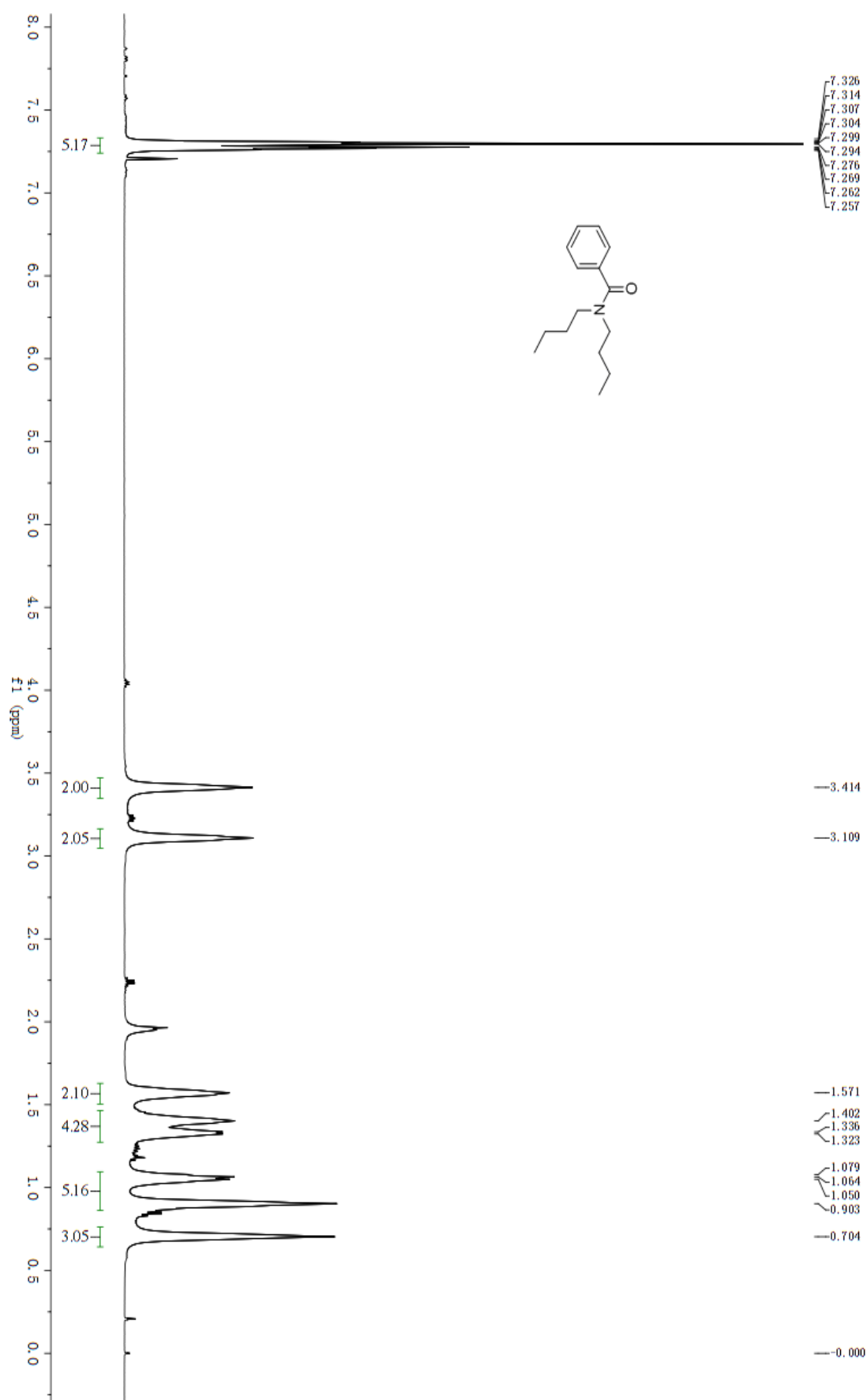
N,N-Dibutyl-4-methoxybenzamide **3a**



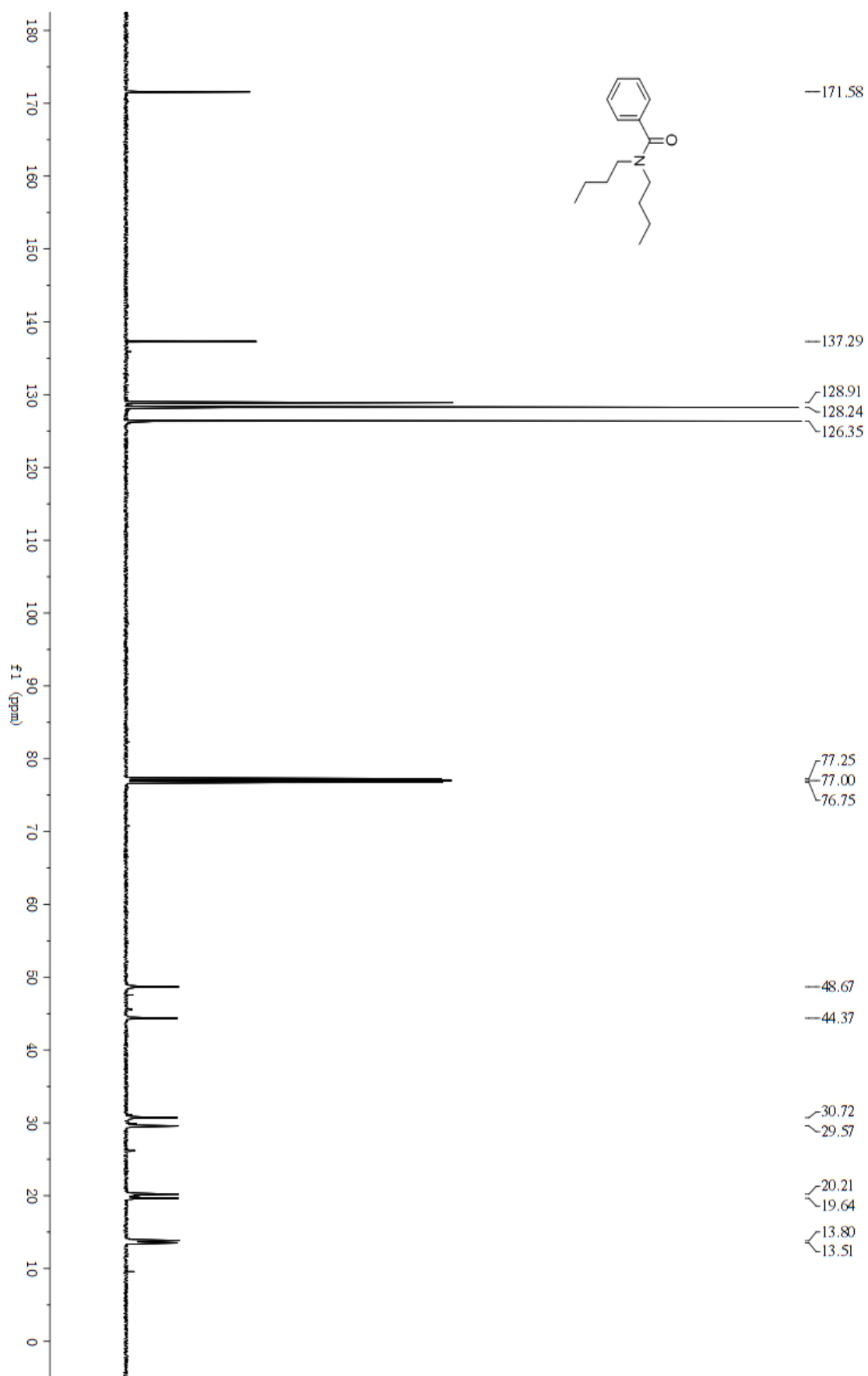
N,N-Dibutyl-4-methoxybenzamide **3a**



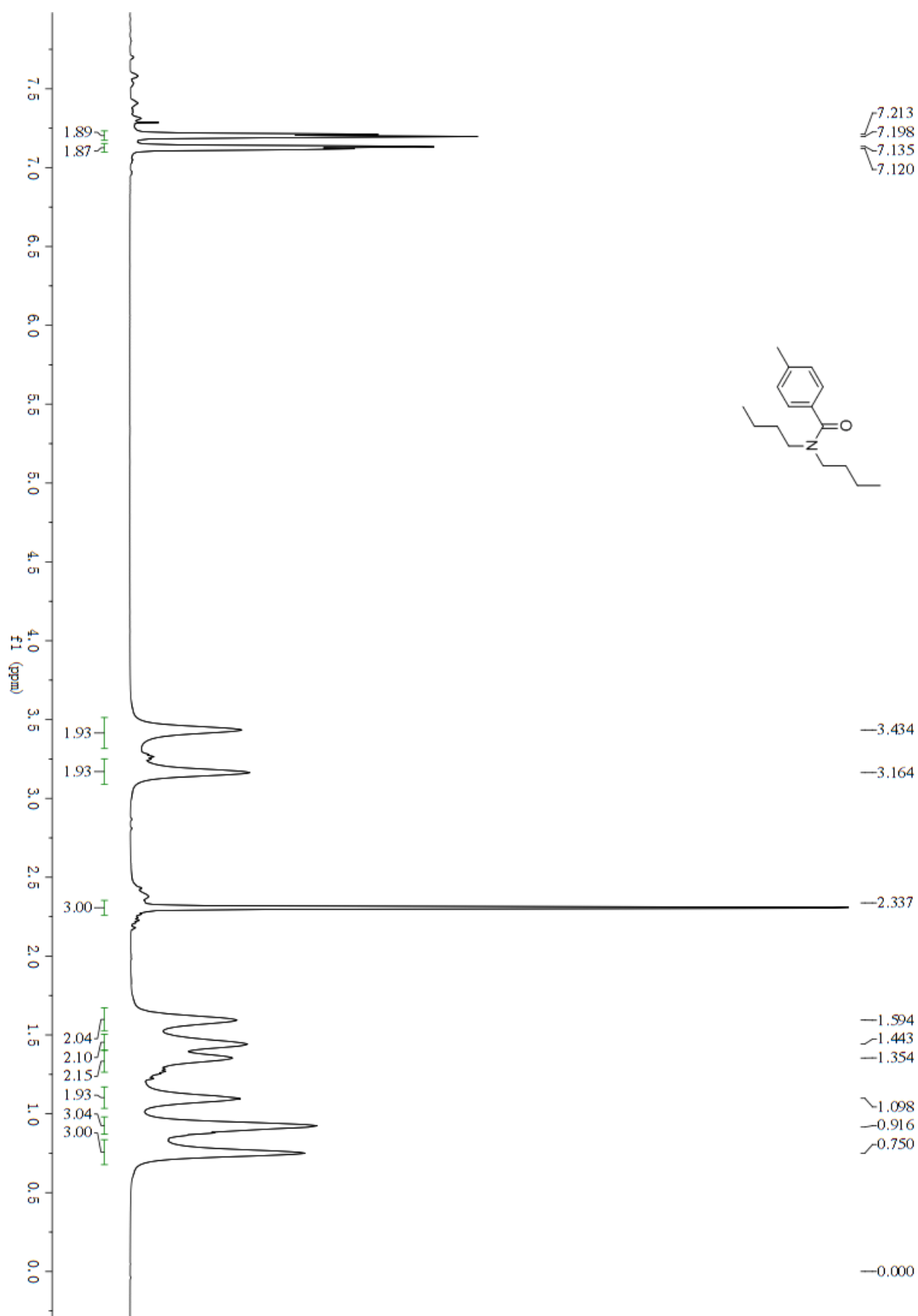
N,N-dibutylbenzamide **3b**



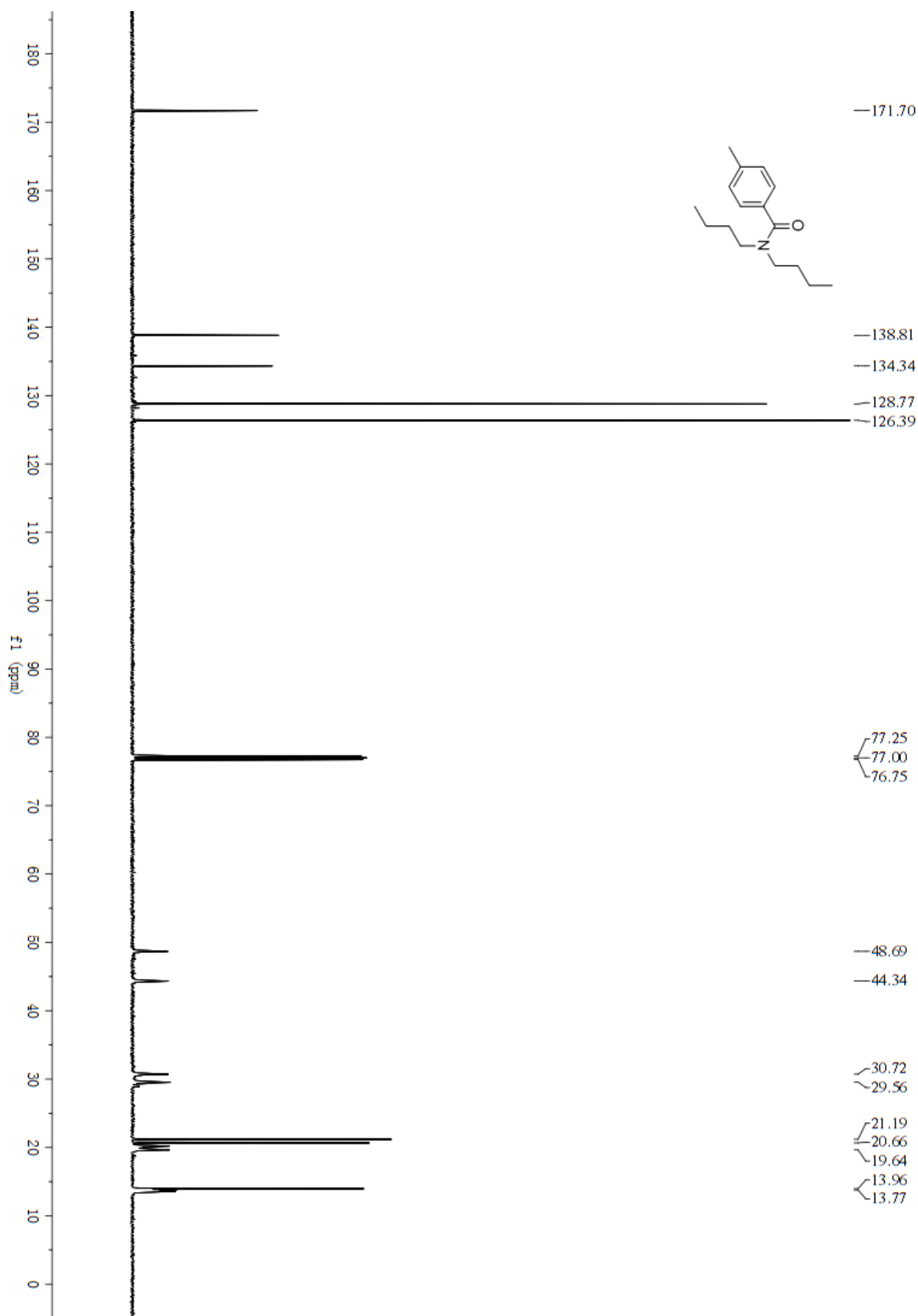
N,N-dibutylbenzamide **3b**



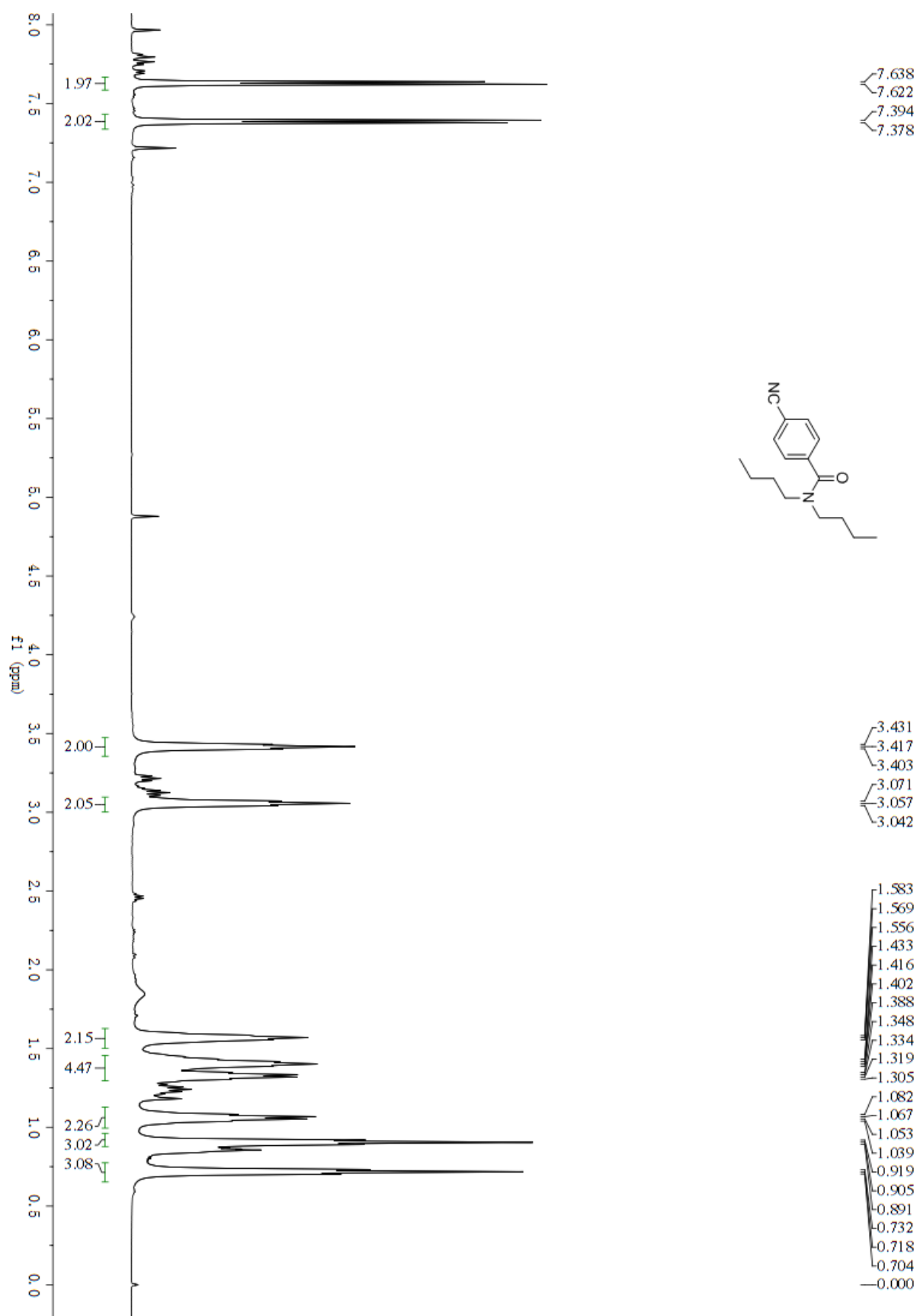
N,N-Dibutyl-4-methylbenzamide **3c**



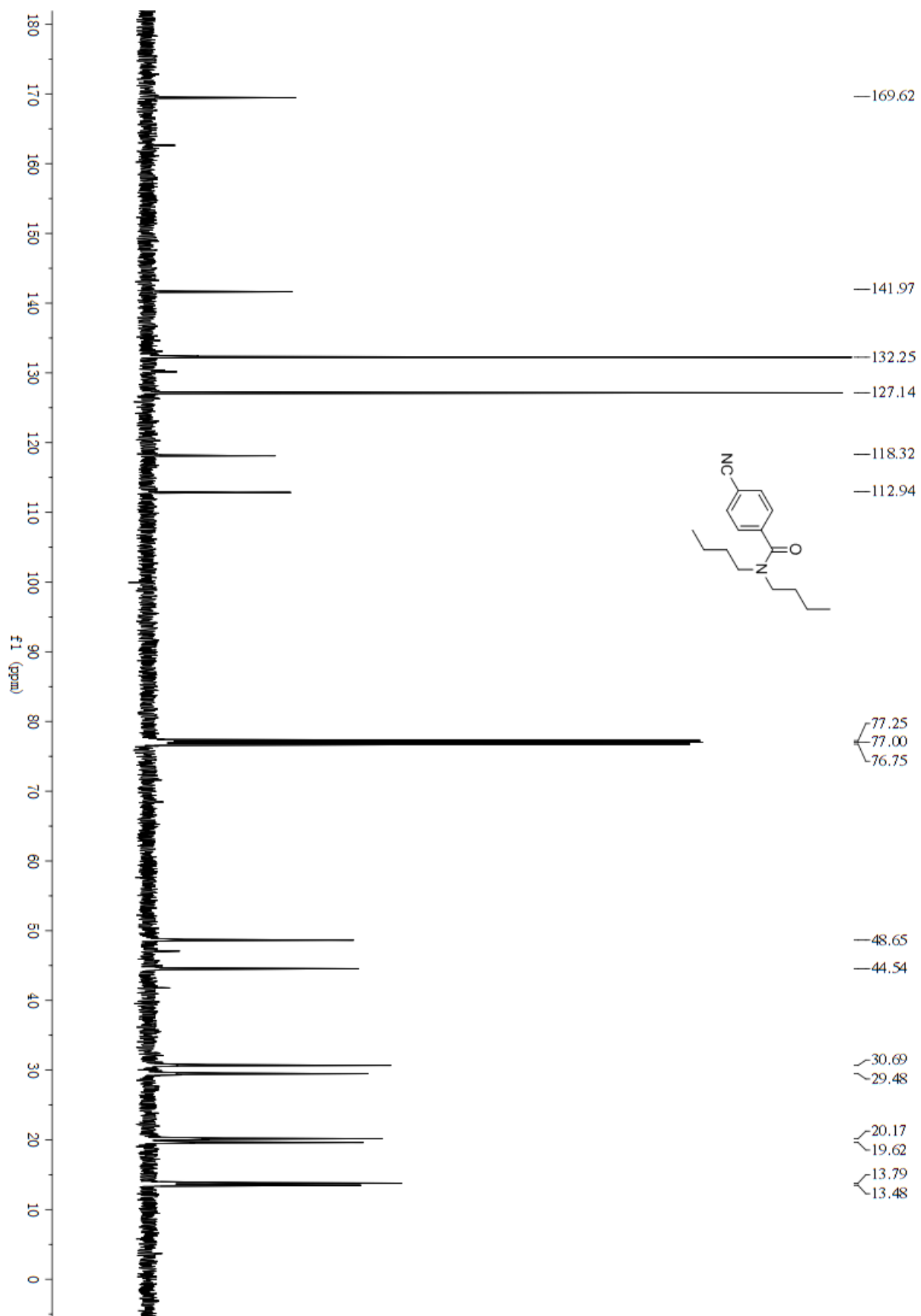
N,N-Dibutyl-4-methylbenzamide **3c**



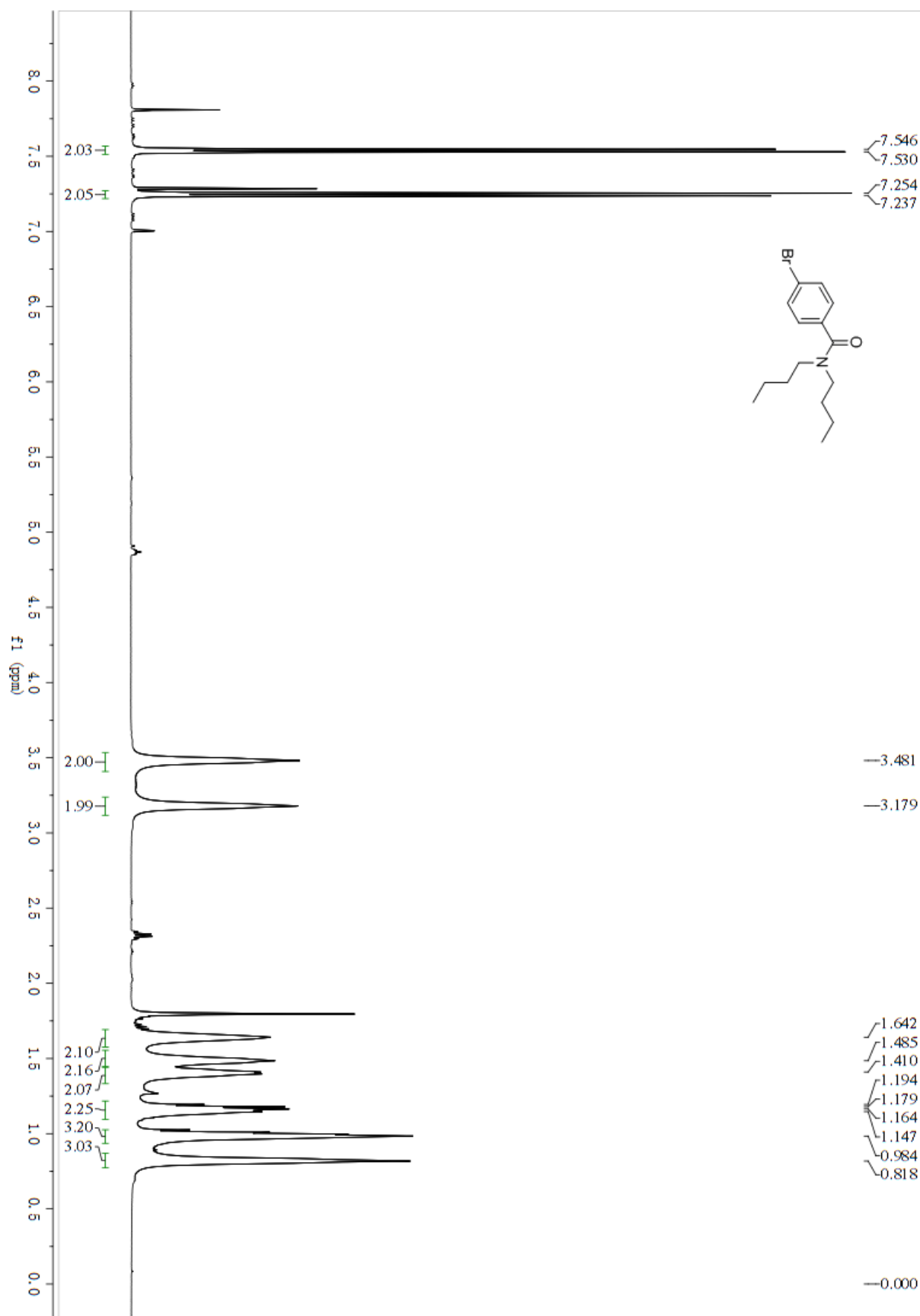
N,N-Dibutyl-4-cyanobenzamide **3d**



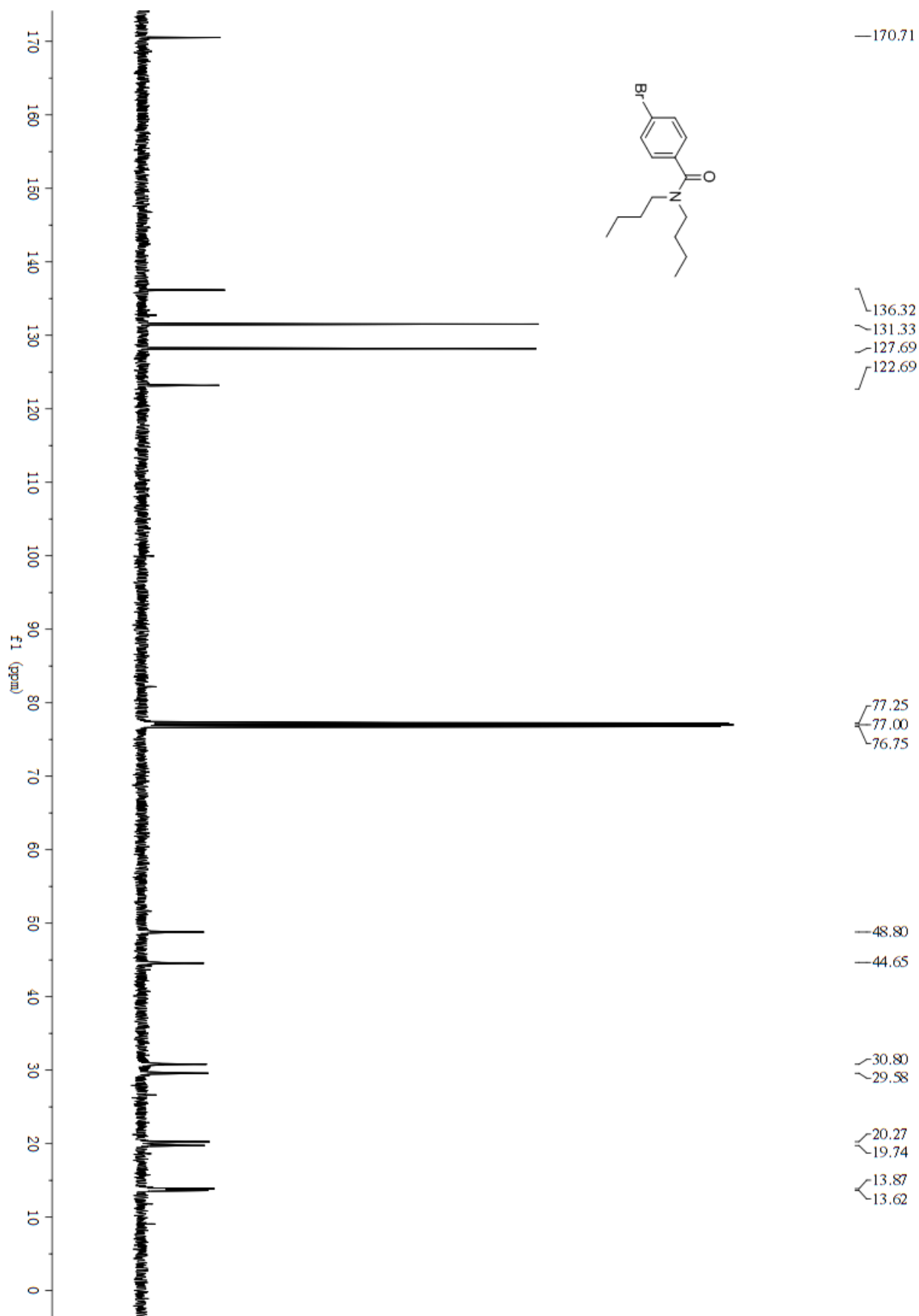
N,N-Dibutyl-4-cyanobenzamide **3d**



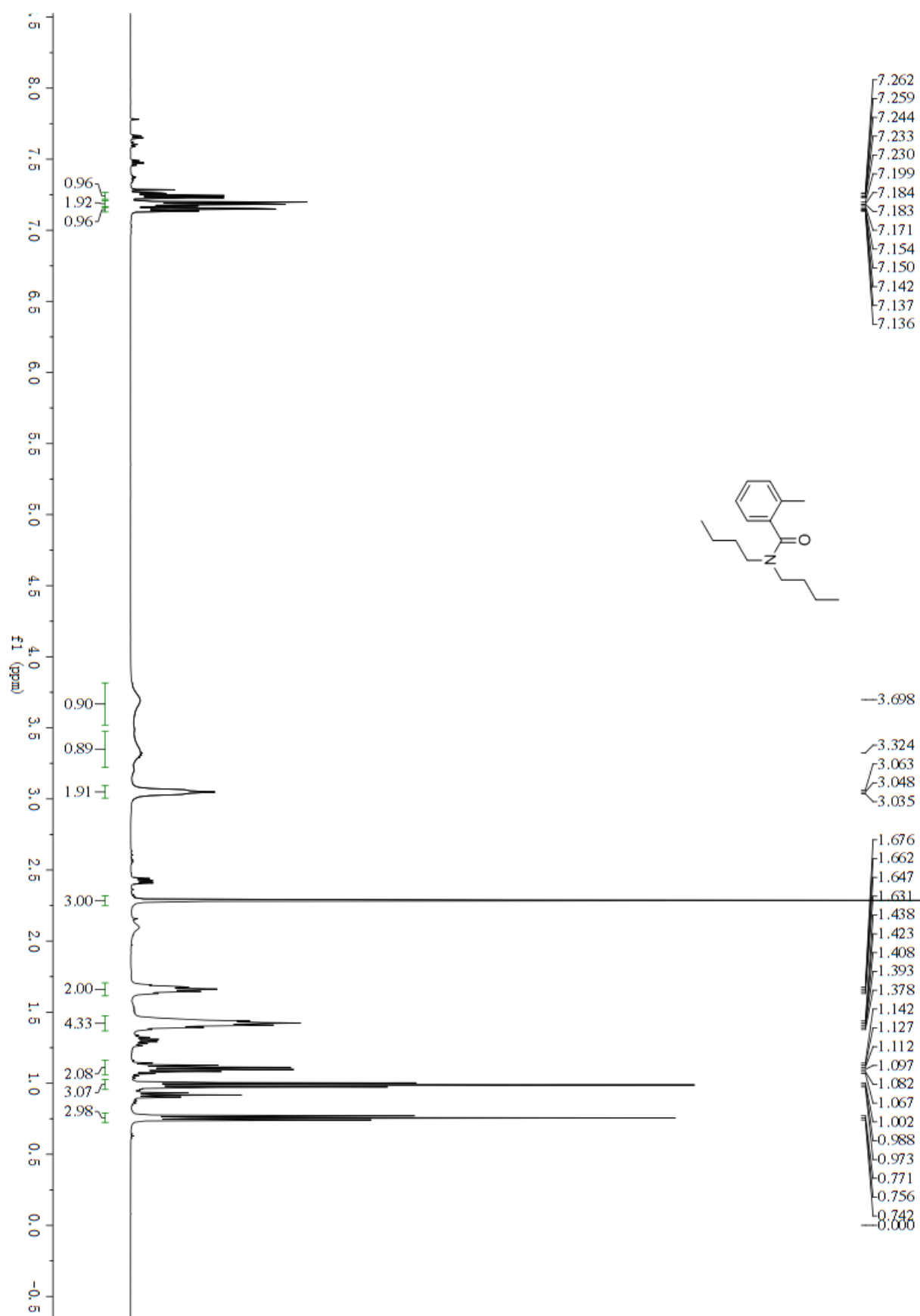
4-bromo-N,N-Dibutylbenzamide 3e



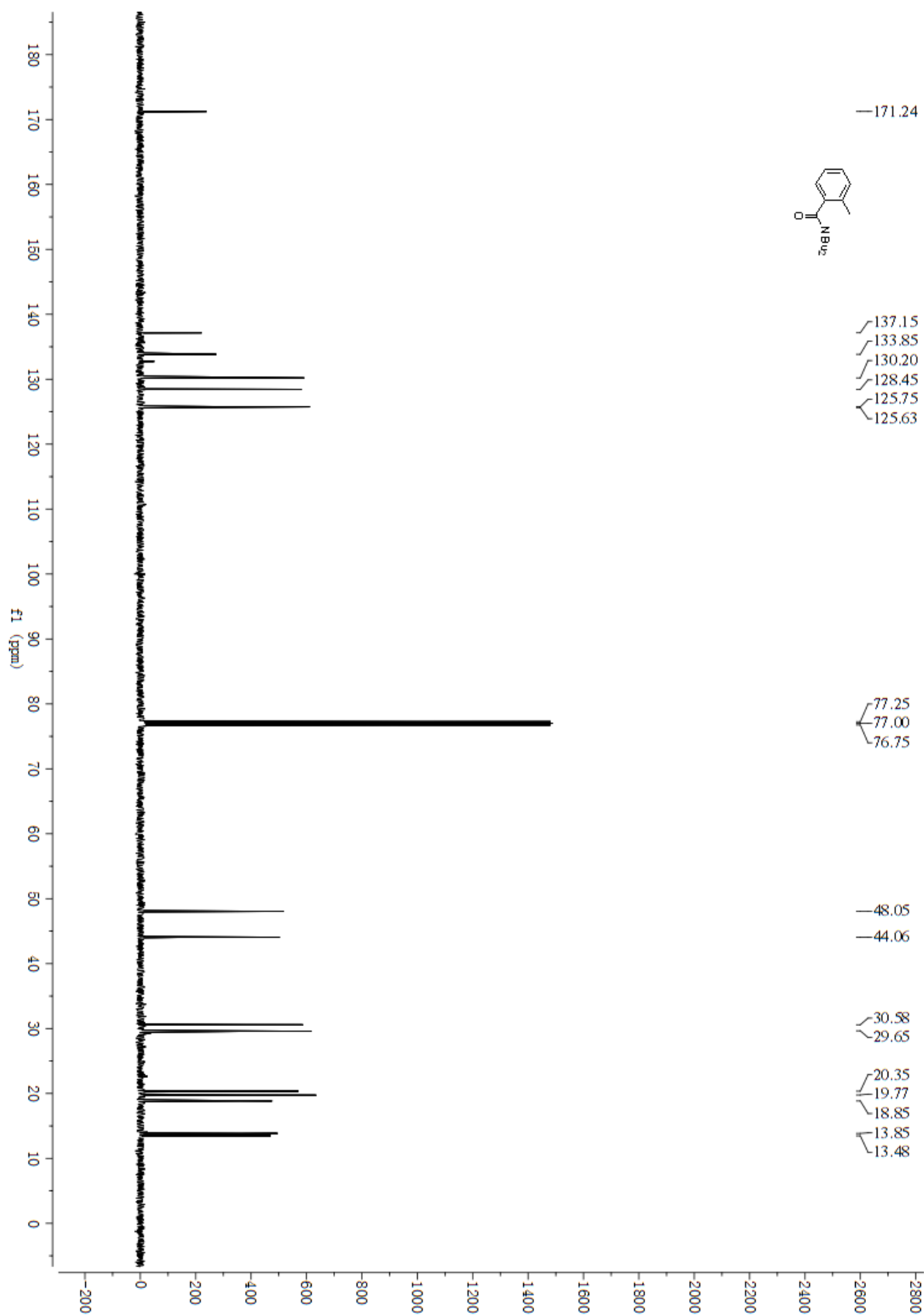
4-bromo-N,N-Dibutylbenzamide 3e



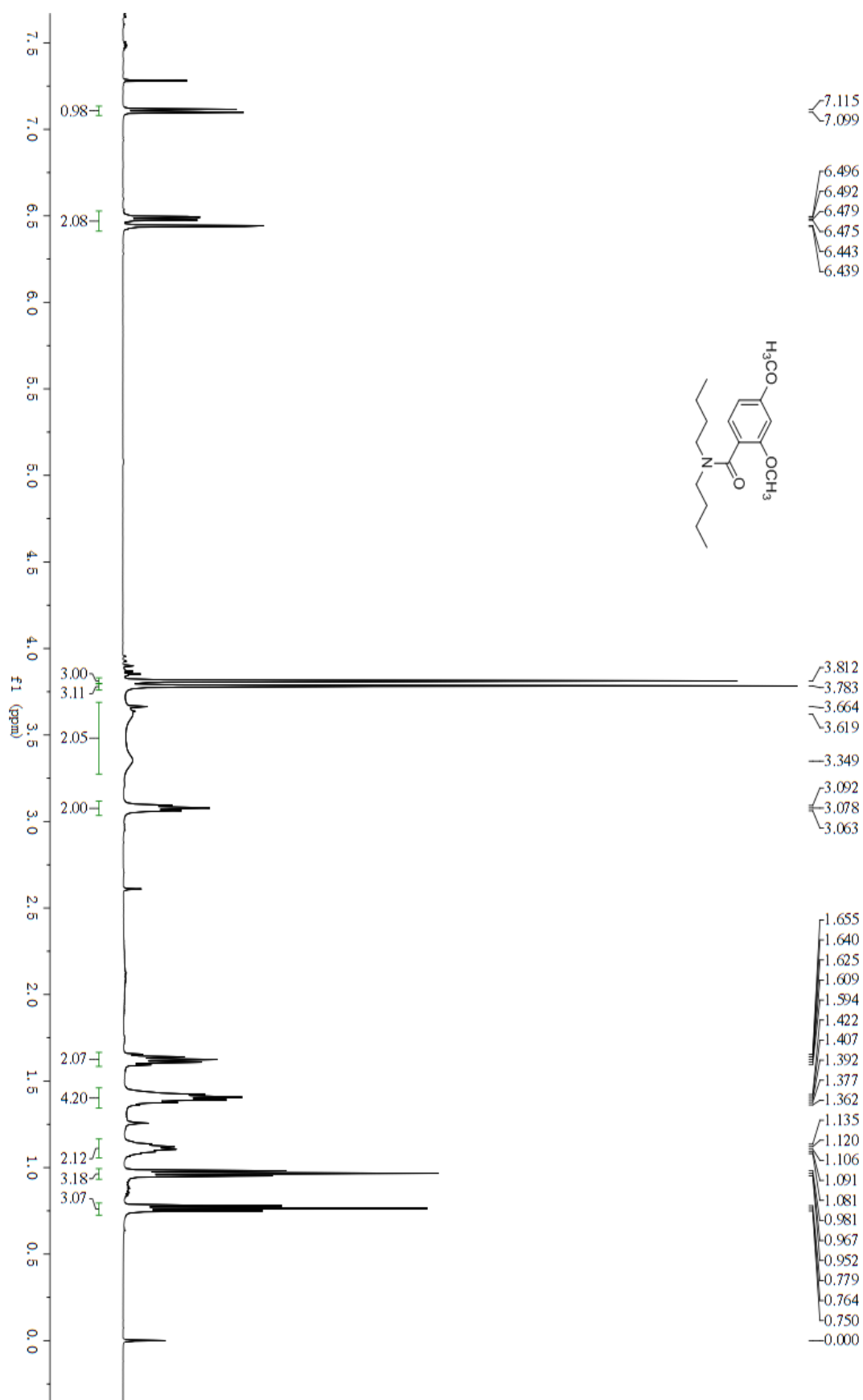
N,N-Dibutyl-2-methylbenzamide **3f**



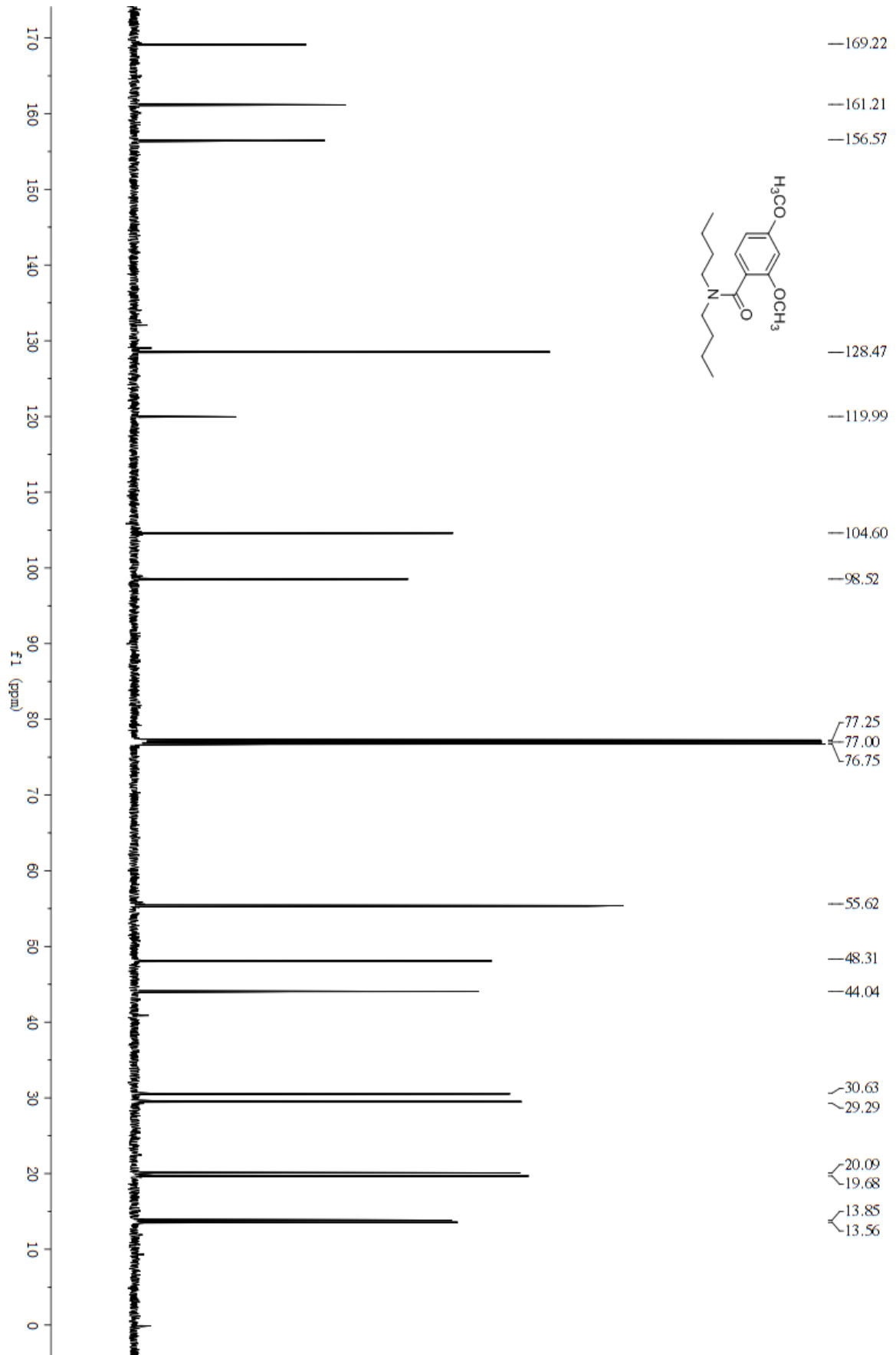
N,N-Dibutyl-2-methylbenzamide **3f**



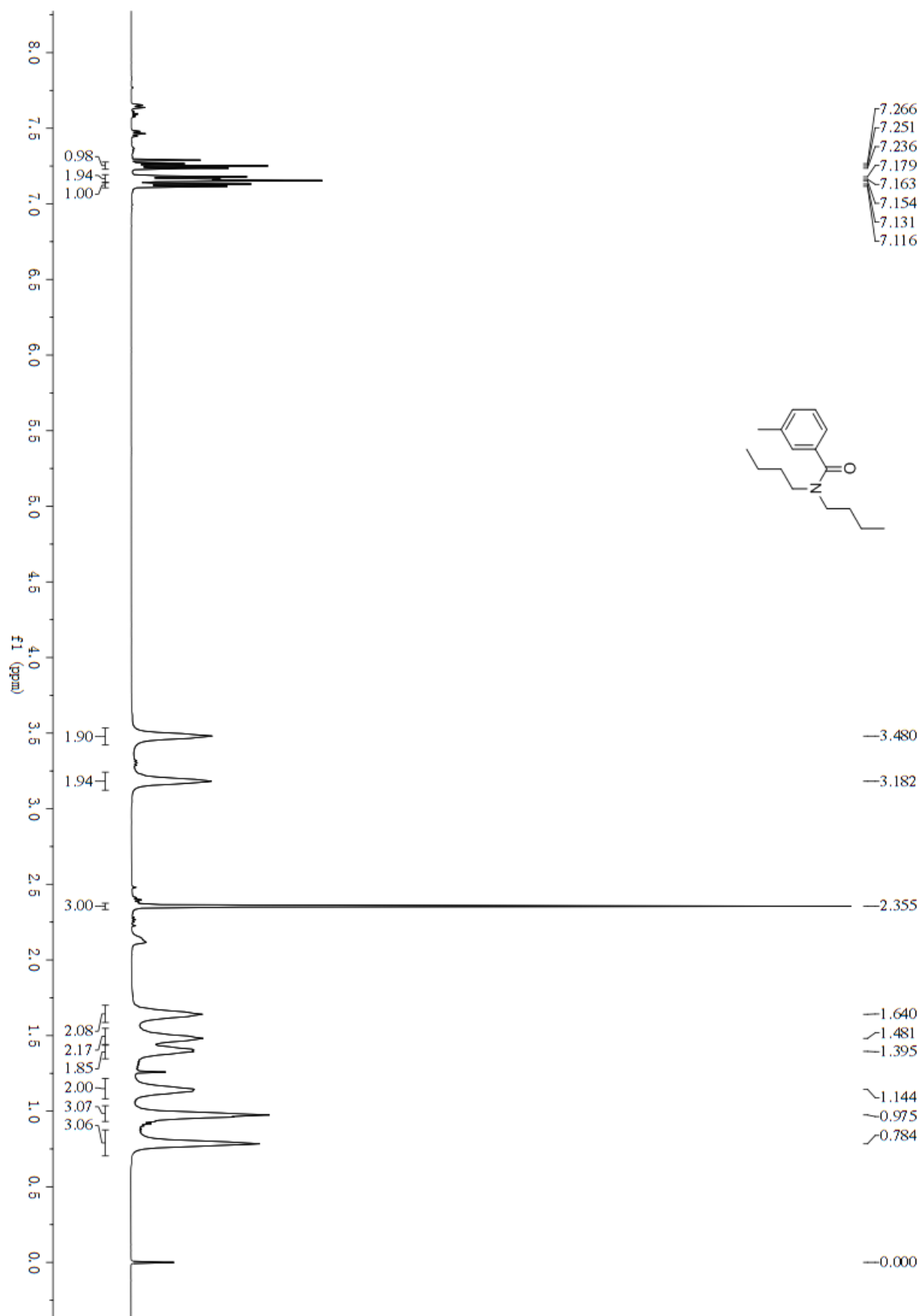
N,N-dibutyl-2,4-dimethoxybenzamide **3g**



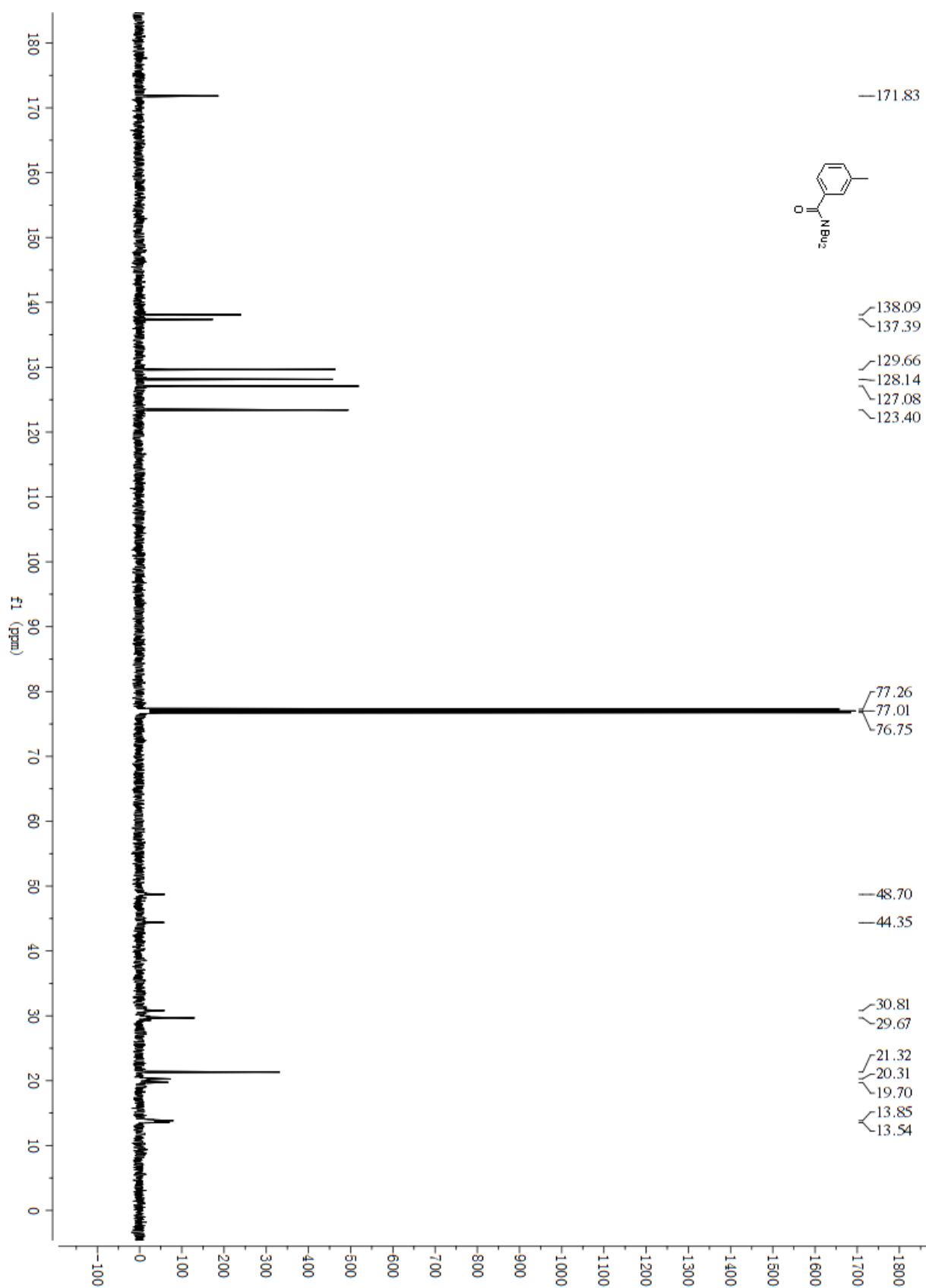
N,N-dibutyl-2,4-dimethoxybenzamide **3g**



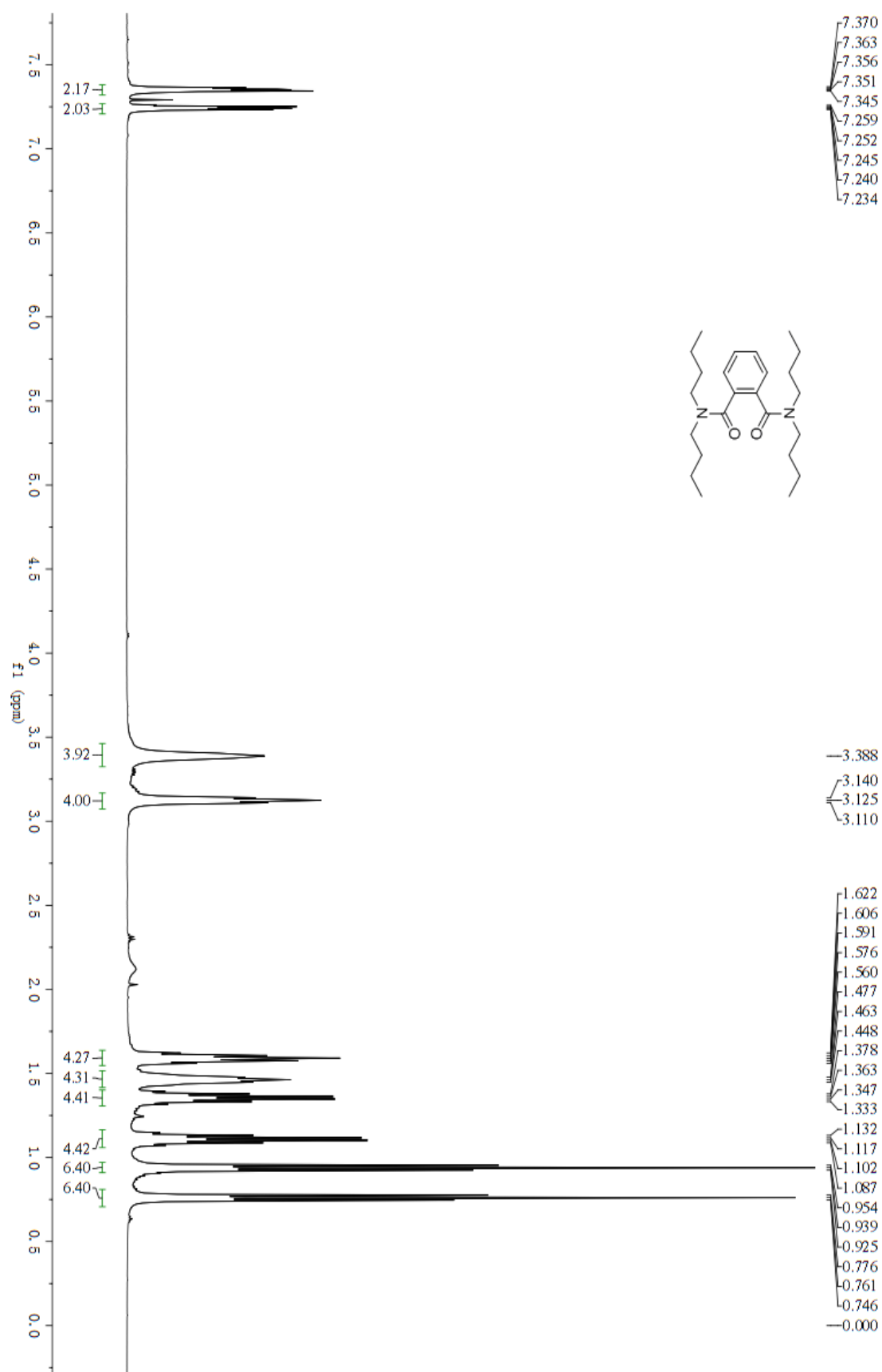
N,N-Dibutyl-3-methylbenzamide **3h**



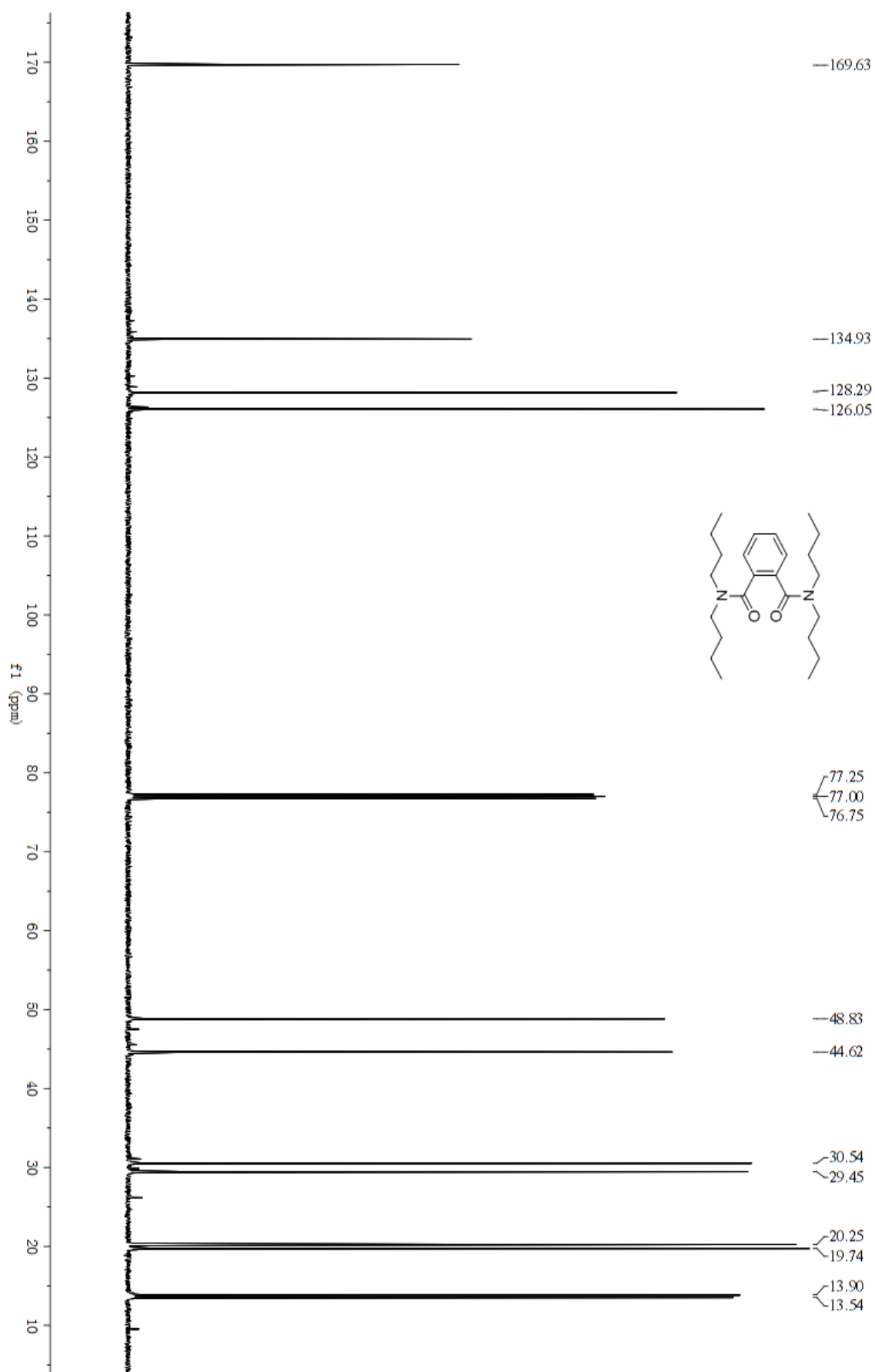
N,N-Dibutyl-3-methylbenzamide **3h**



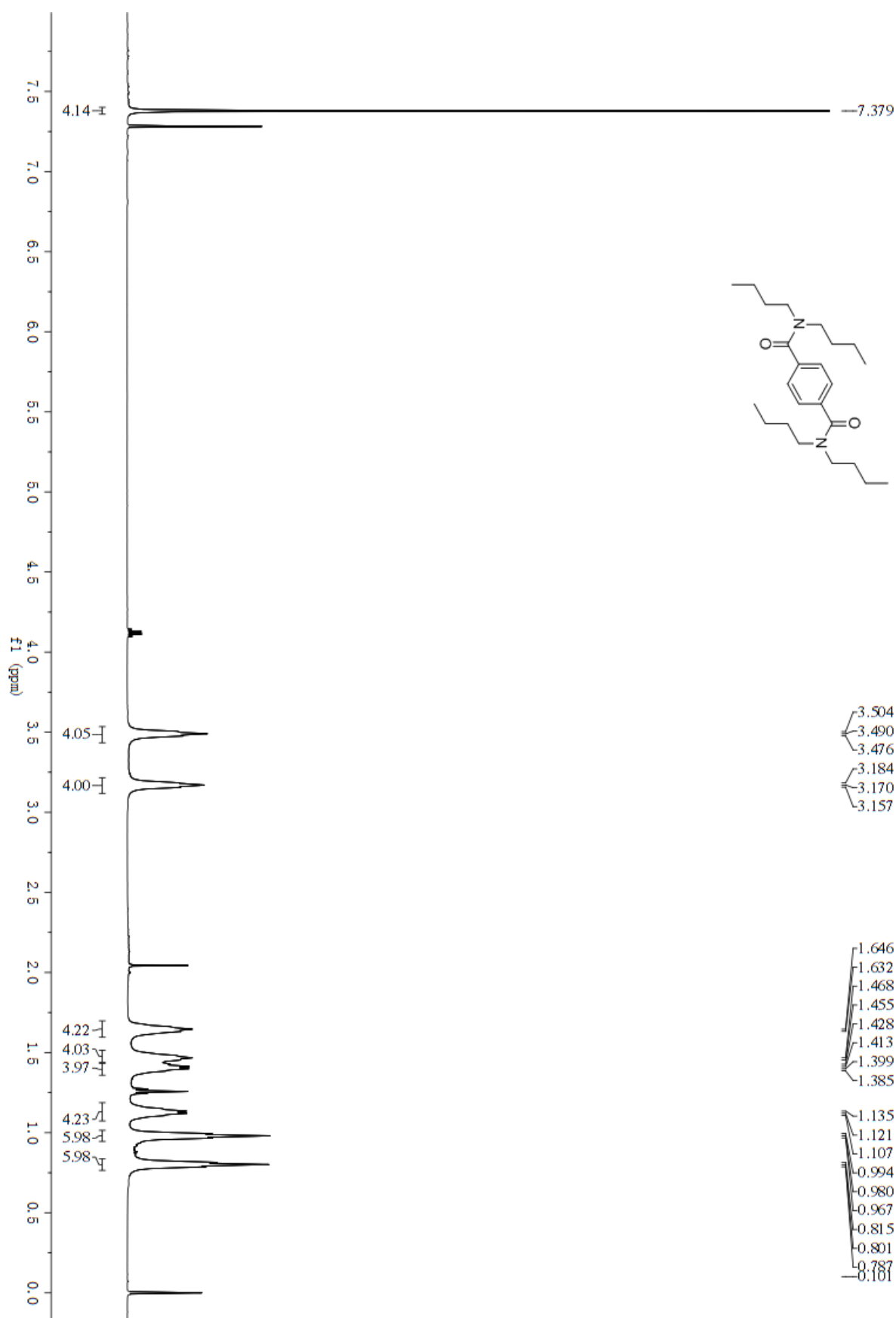
N¹,N¹,N²,N²-tetrabutylphthalamide **3i**



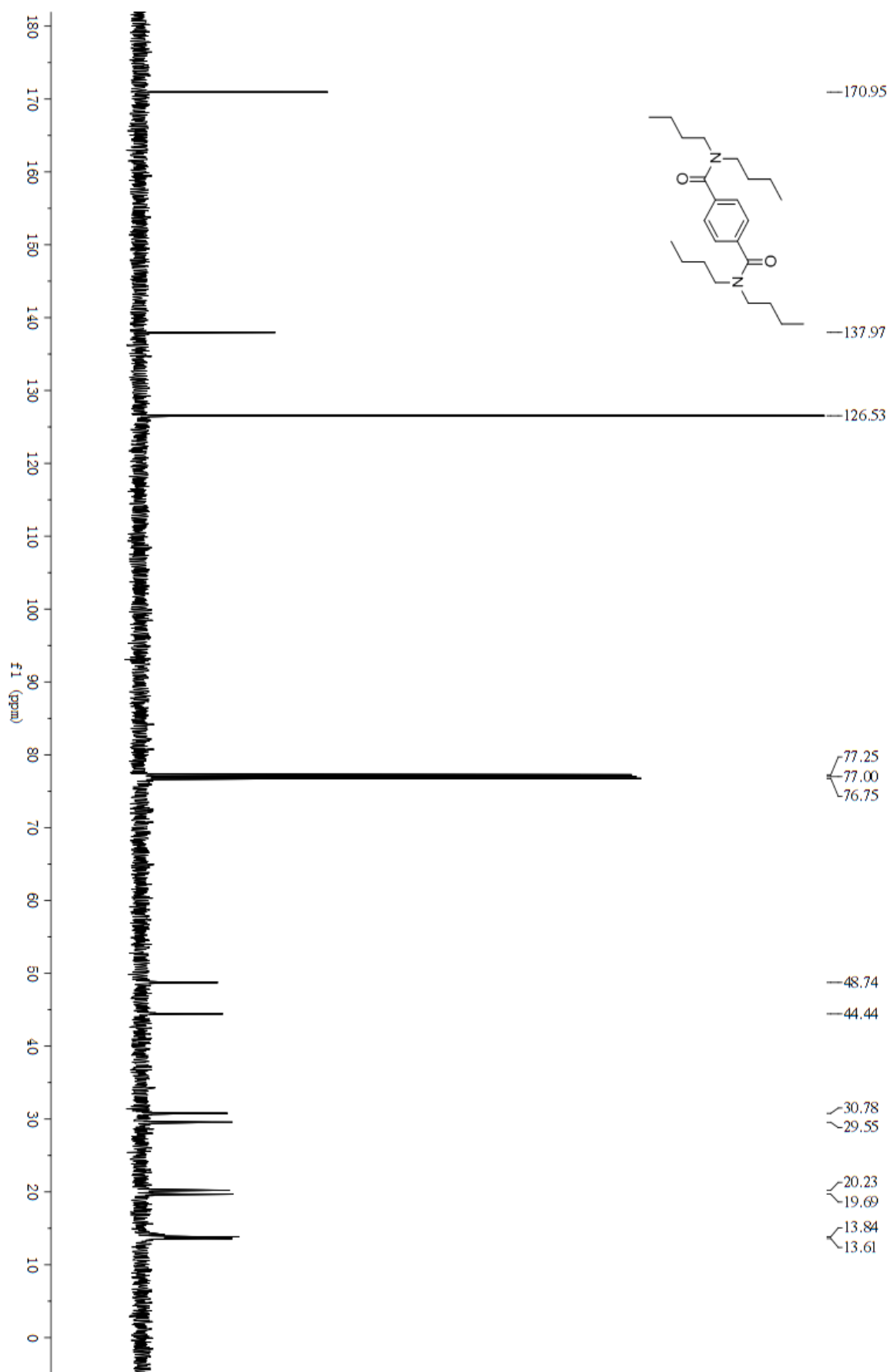
N¹,N¹,N²,N²-tetrabutylphthalamide **3i**



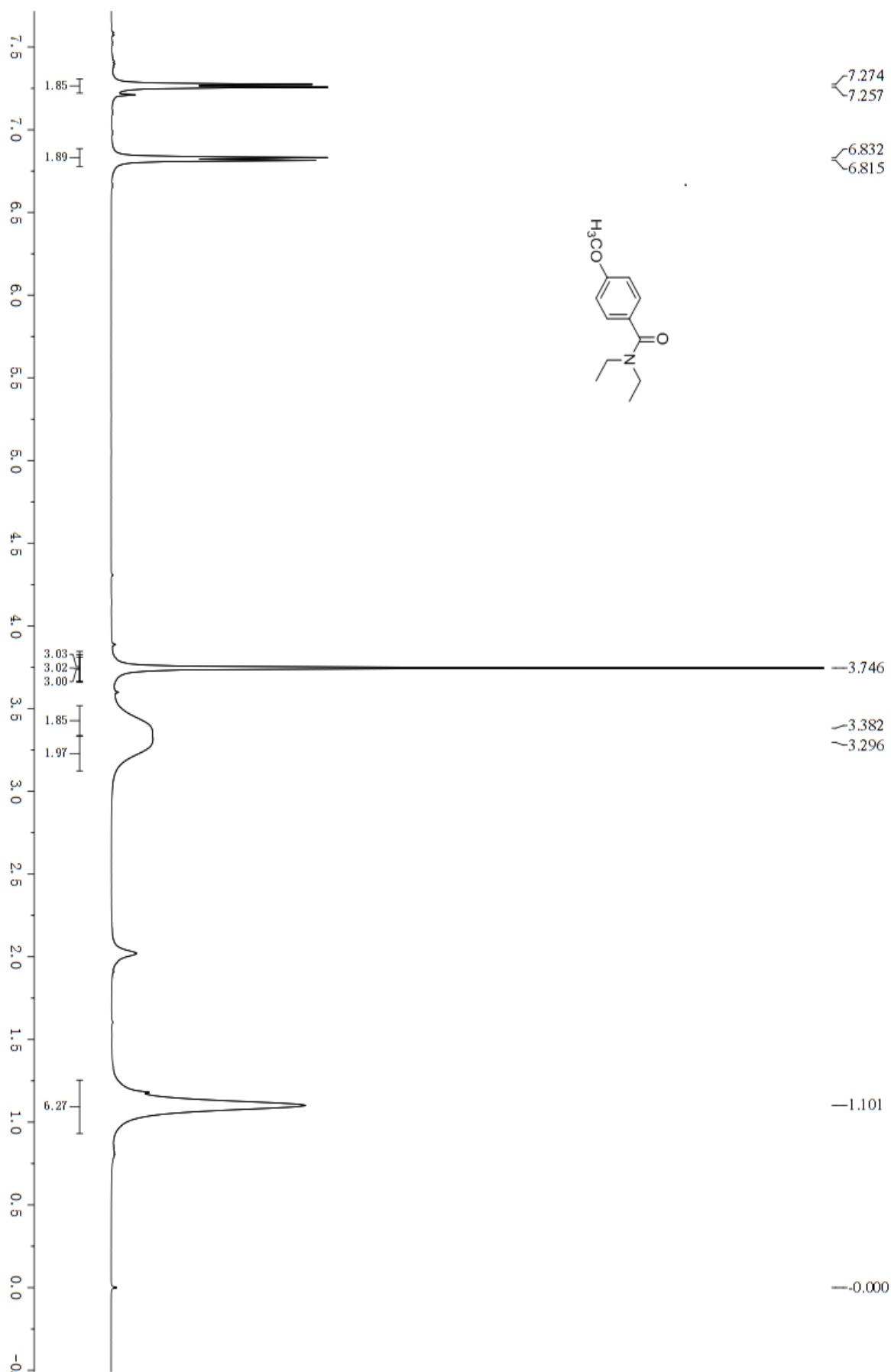
N¹,N¹,N⁴,N⁴-tetrabutylterephthalamide **3j**



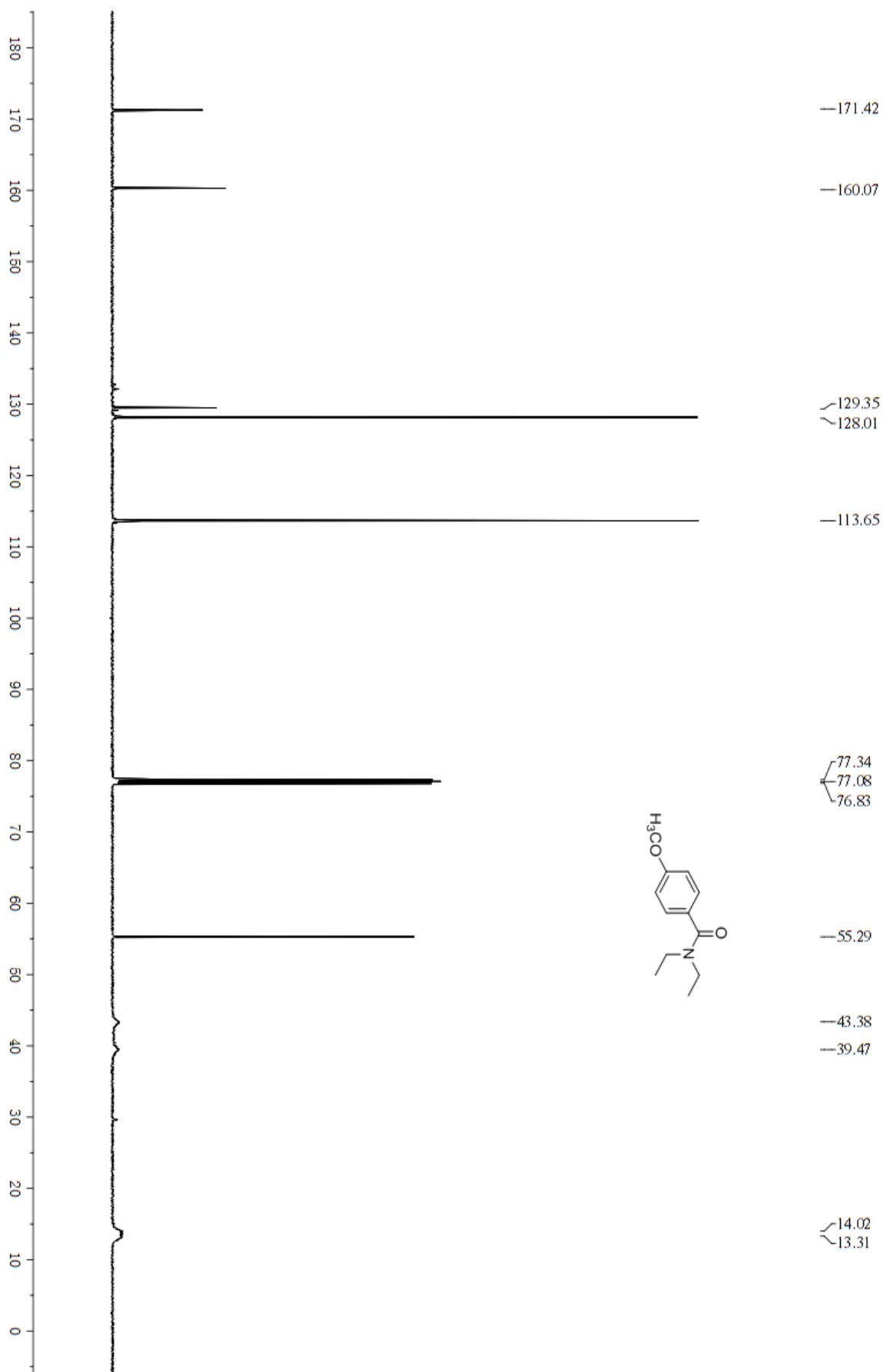
N^1,N^1,N^4,N^4 -tetrabutylterephthalamide **3j**



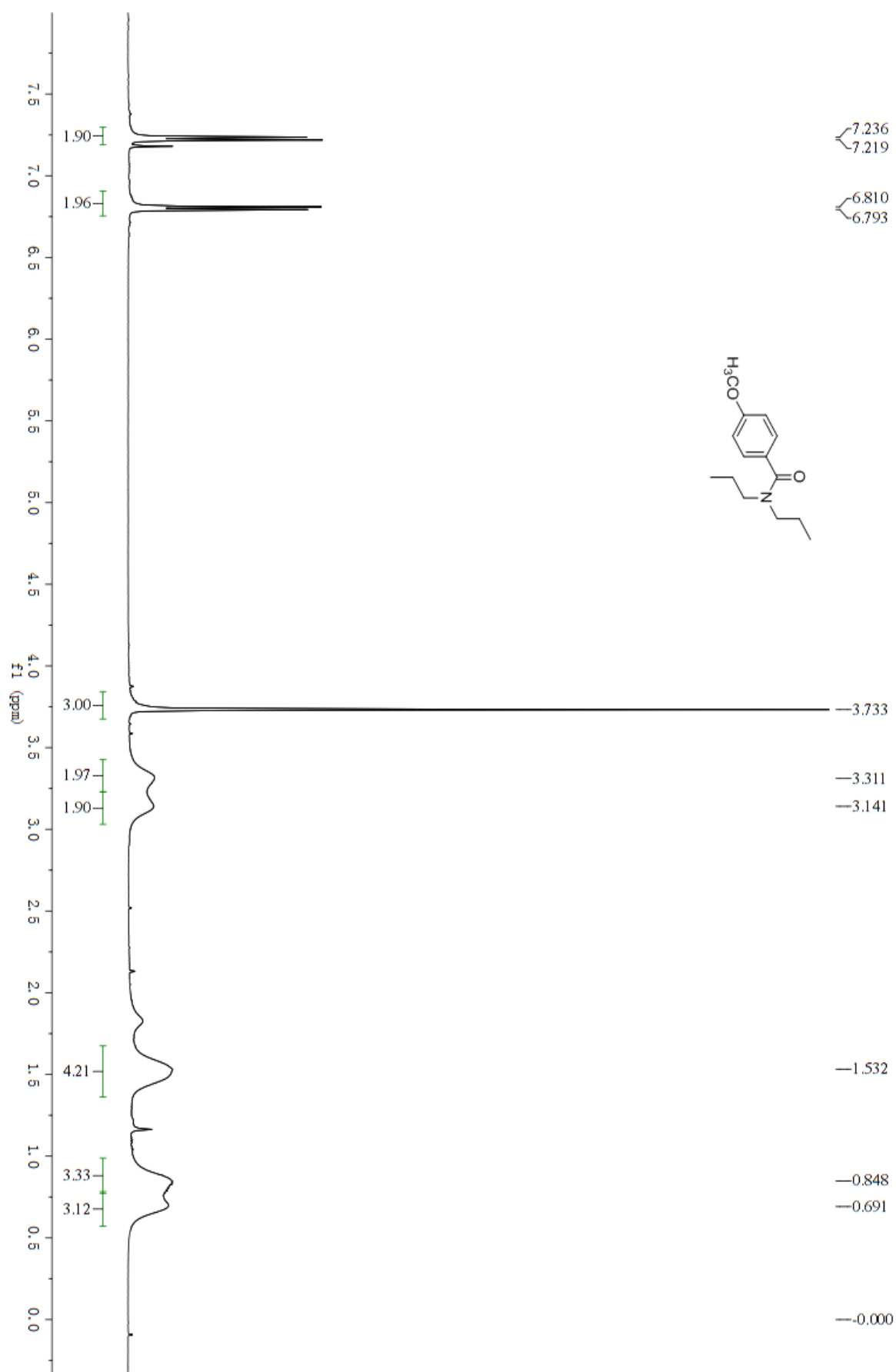
N,N-Diethyl-4-methoxybenzamide **3k**



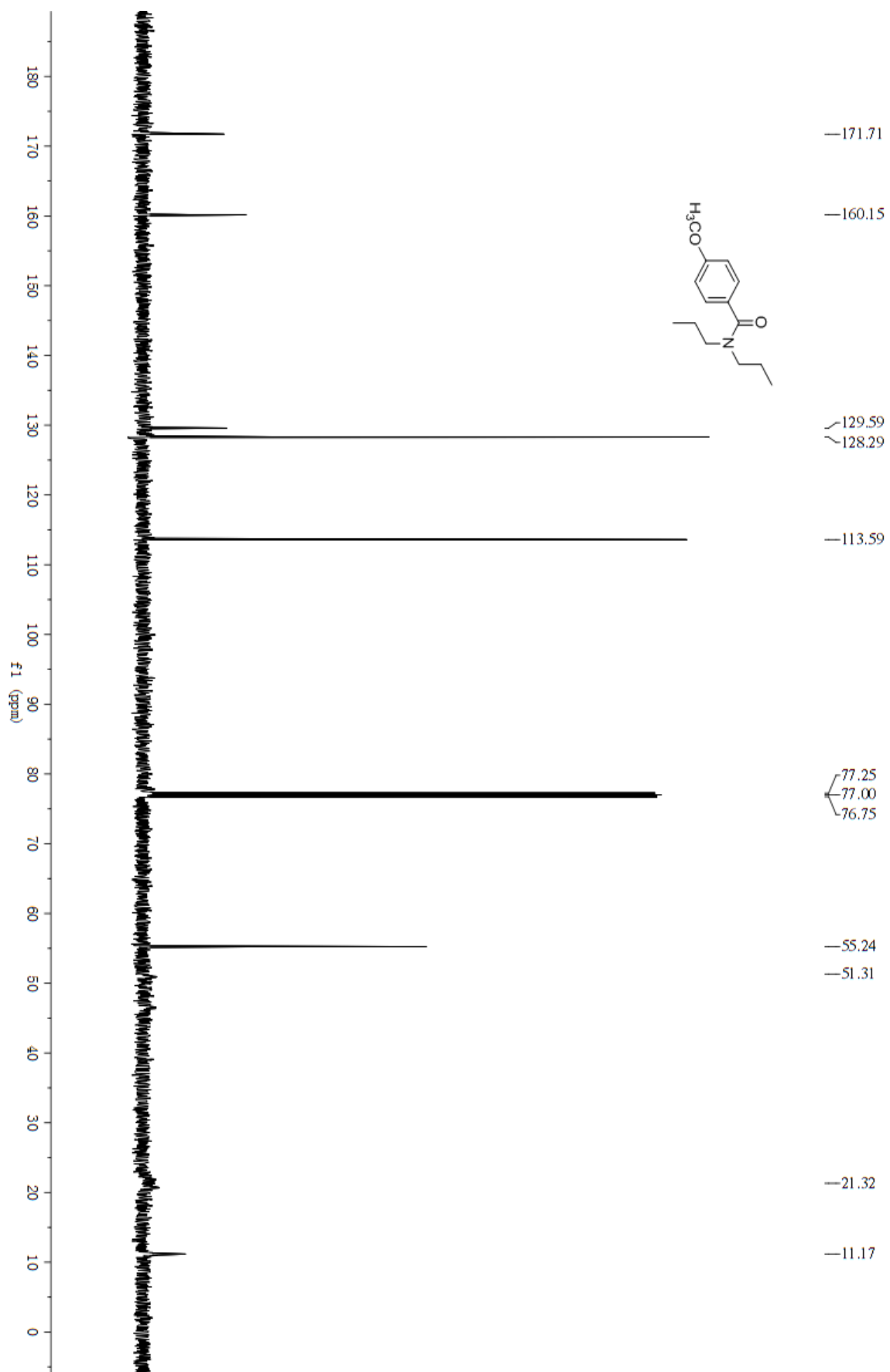
N,N-Diethyl-4-methoxybenzamide **3k**



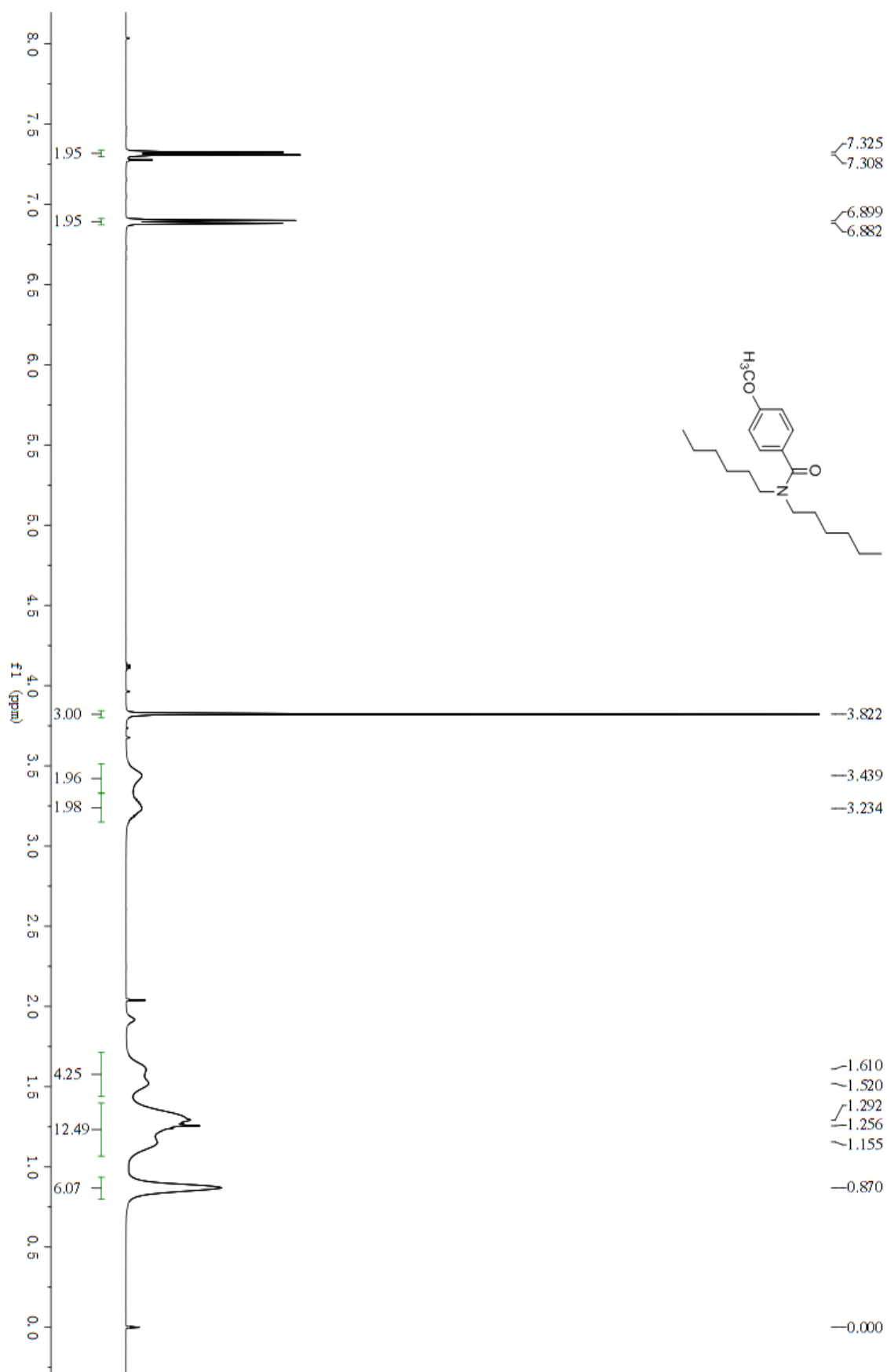
N,N-Dipropyl-4-methoxybenzamide **31**



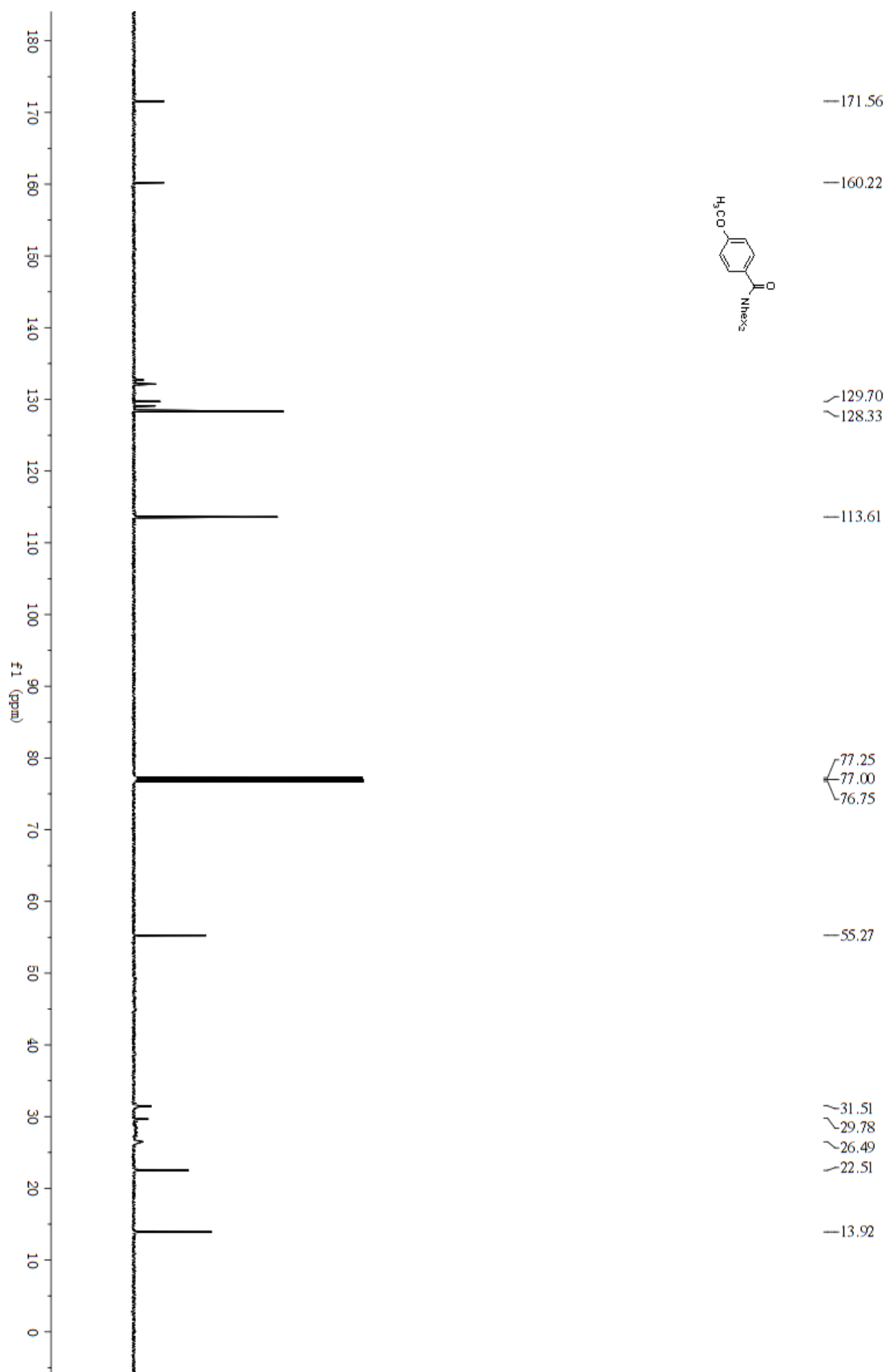
N,N-Dipropyl-4-methoxybenzamide **31**



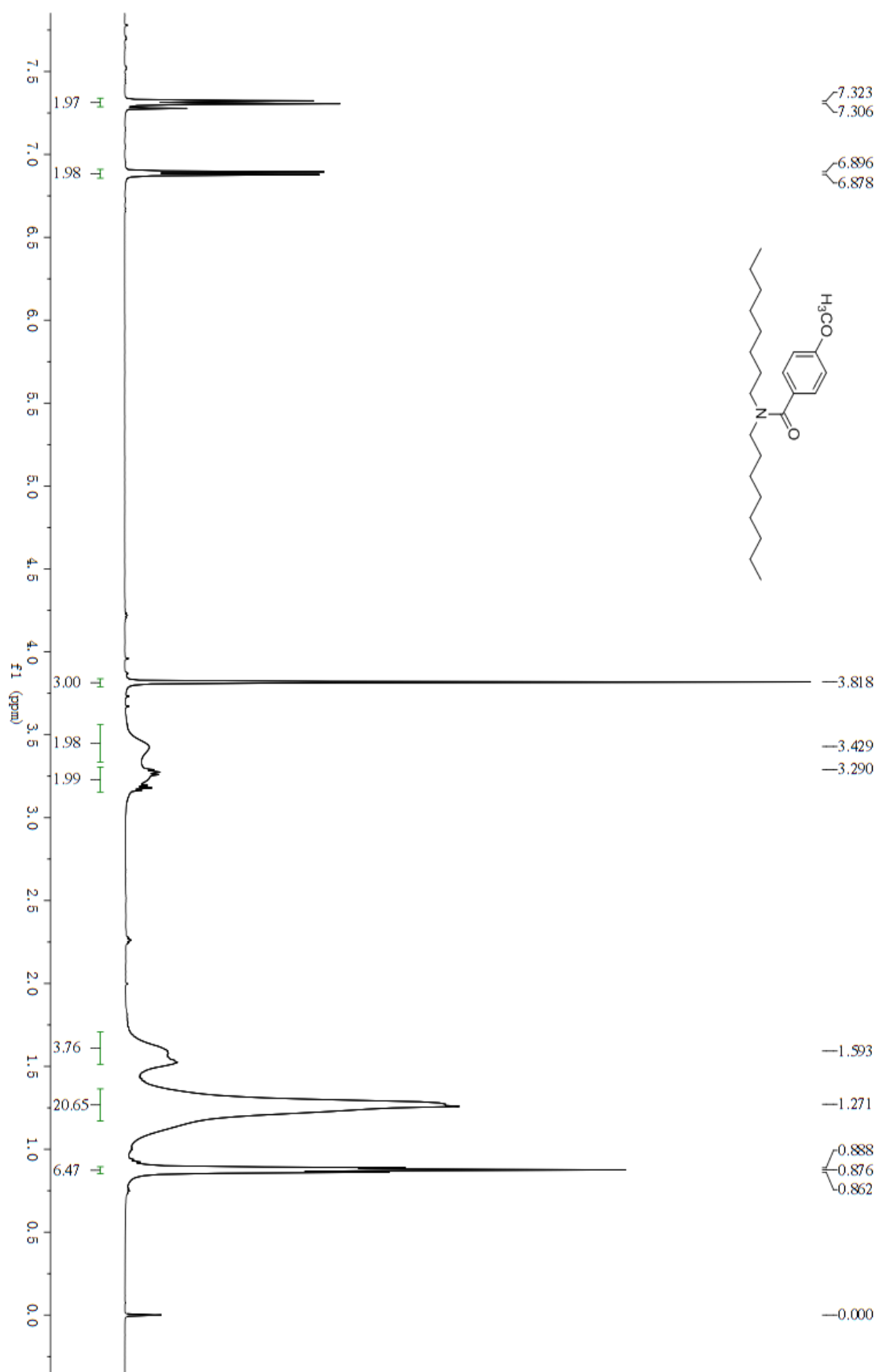
N,N-dihexyl-4-methoxybenzamide **3m**



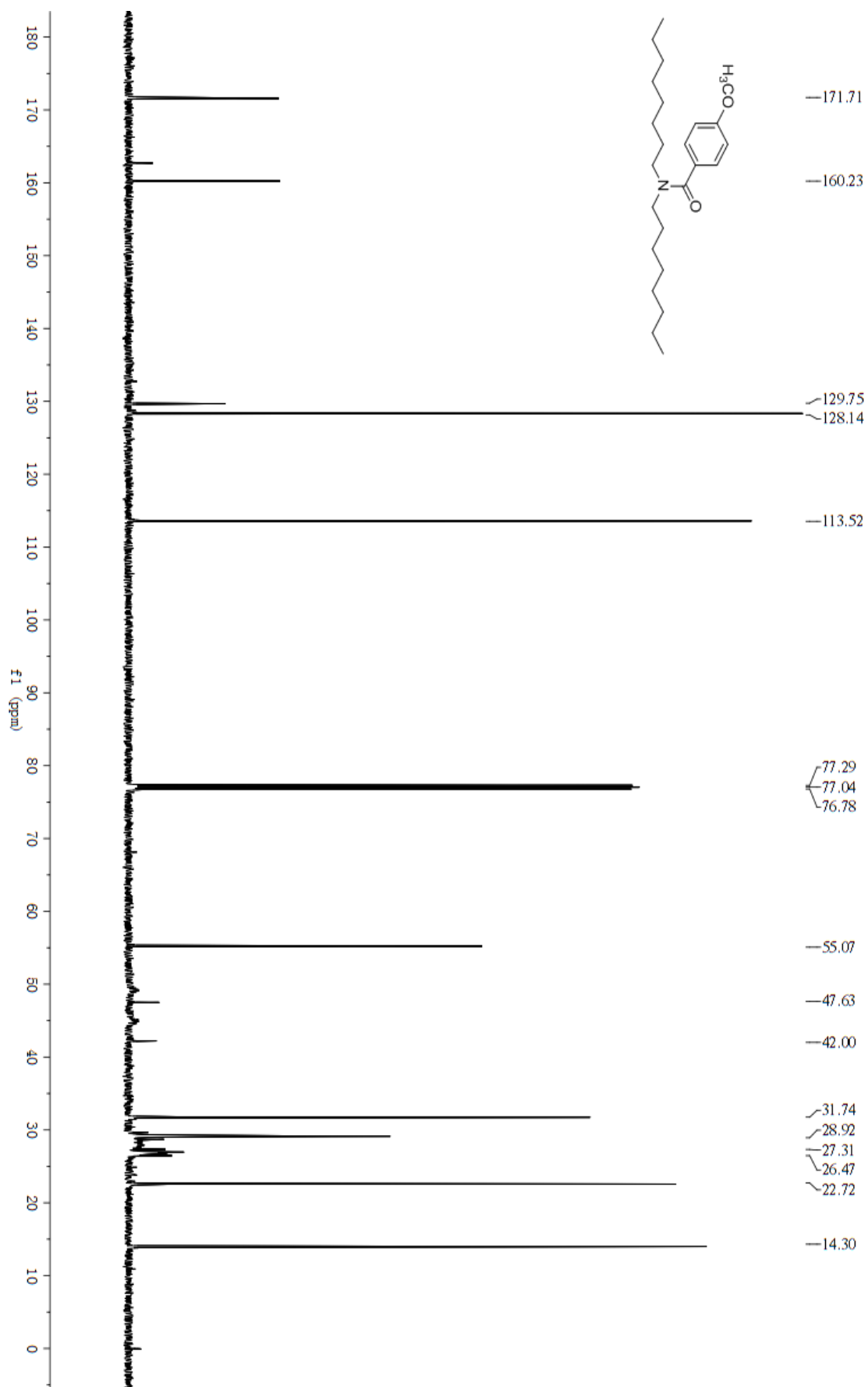
N,N-dihexyl-4-methoxybenzamide **3m**



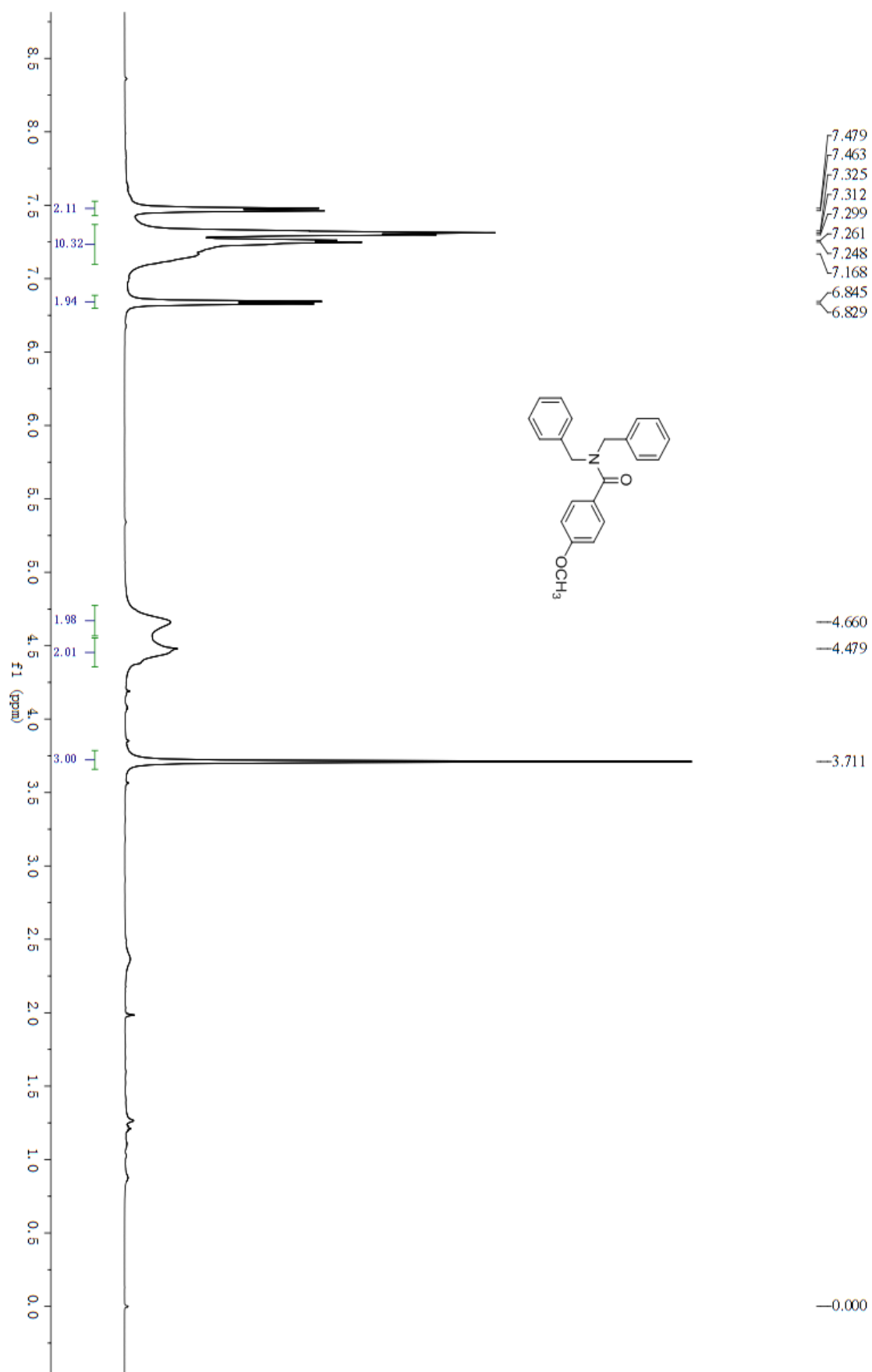
4-methoxy-N,N-dioctylbenzamide **3n**



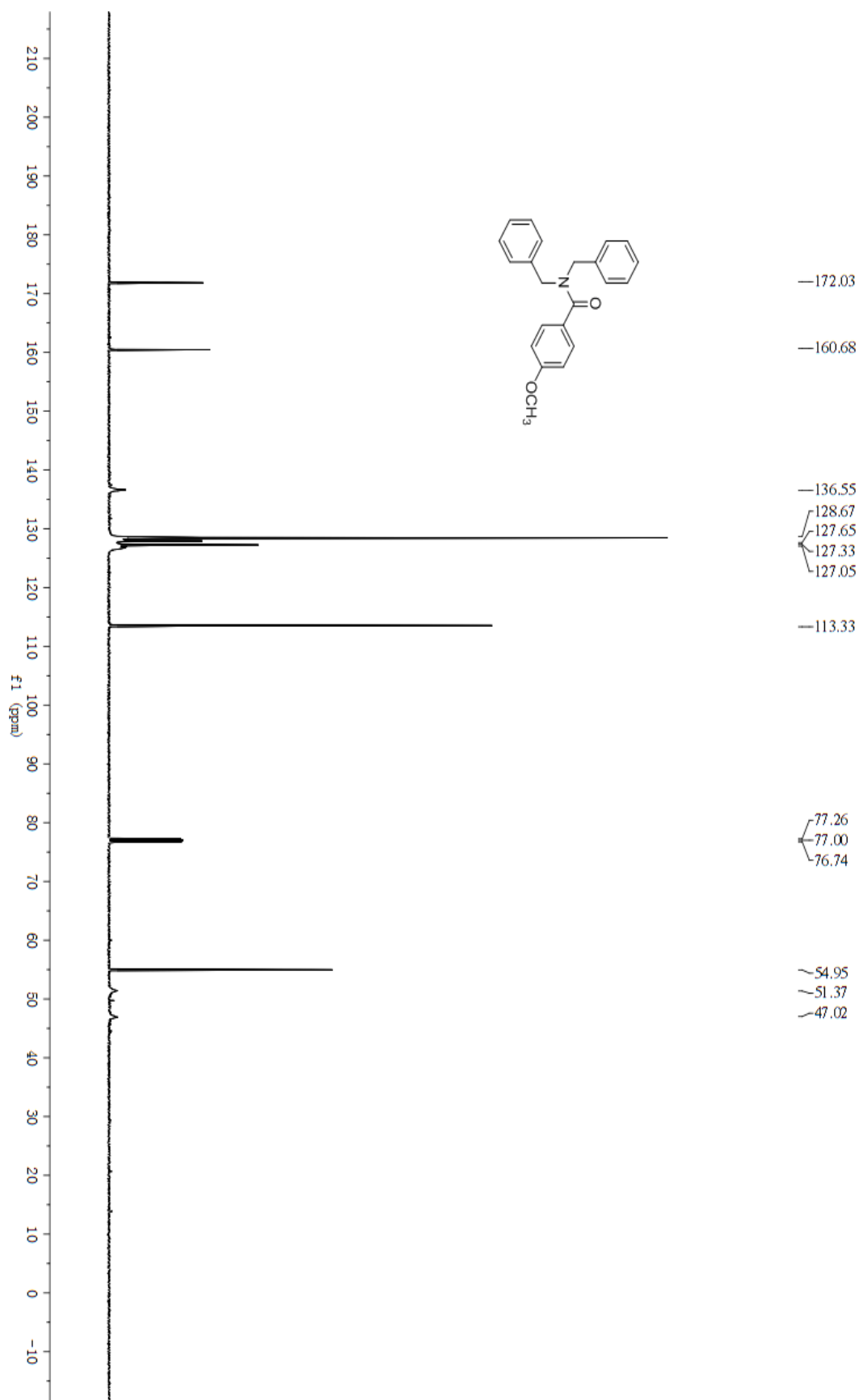
4-methoxy-N,N-dioctylbenzamide **3n**



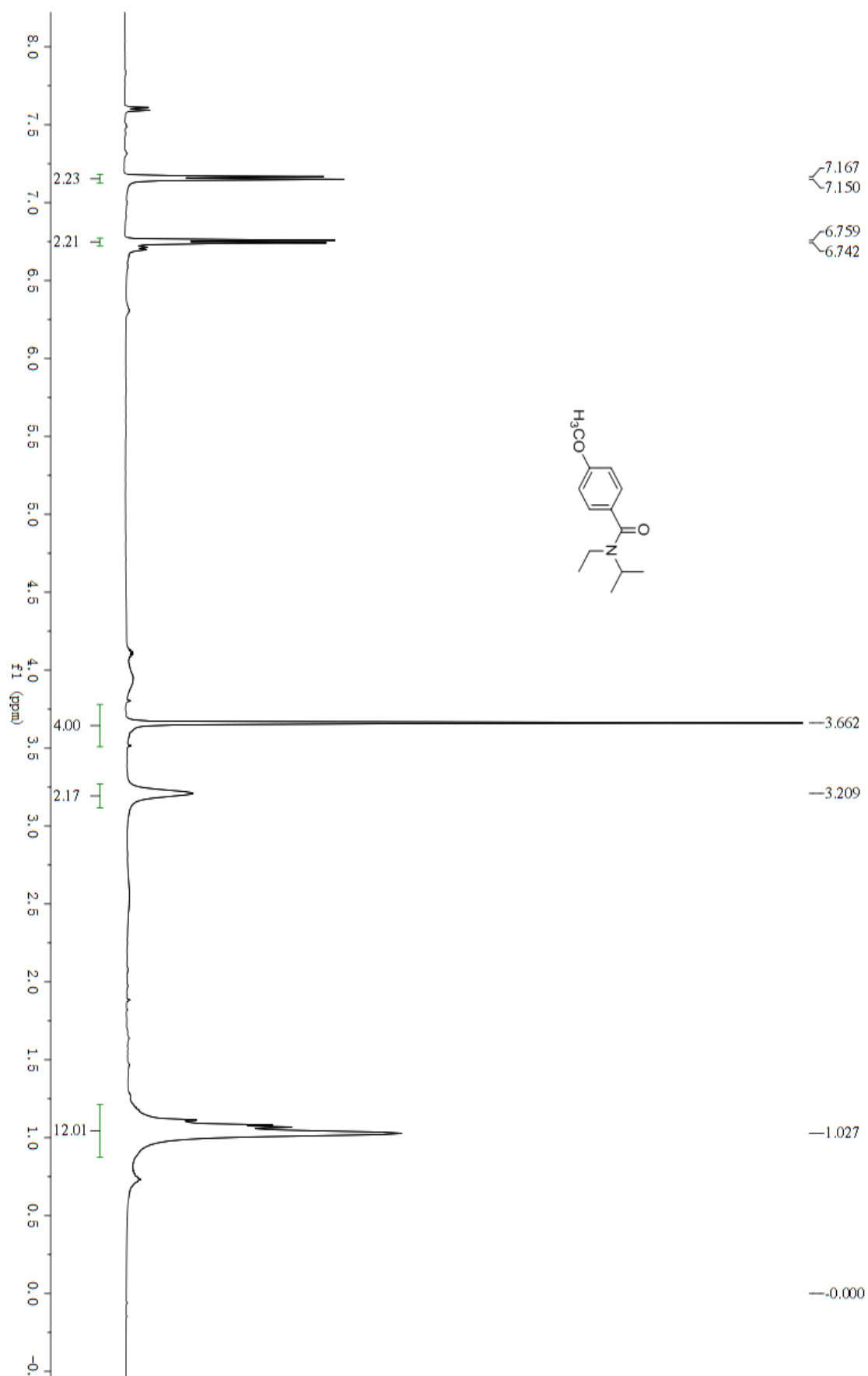
N,N-Dibenzyl-4-methoxybenzamide **3o**



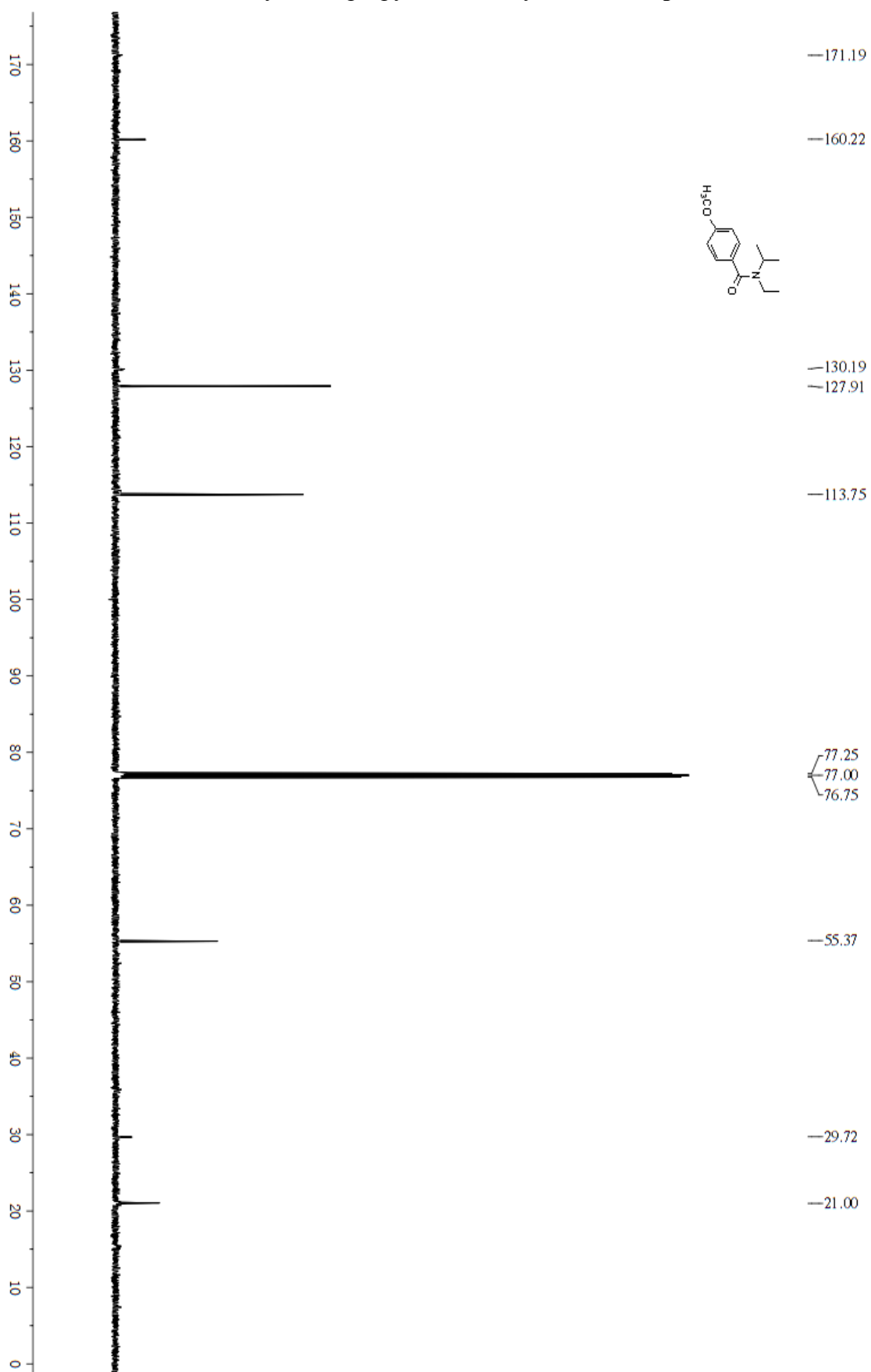
N,N-Dibenzyl-4-methoxybenzamide **3o**



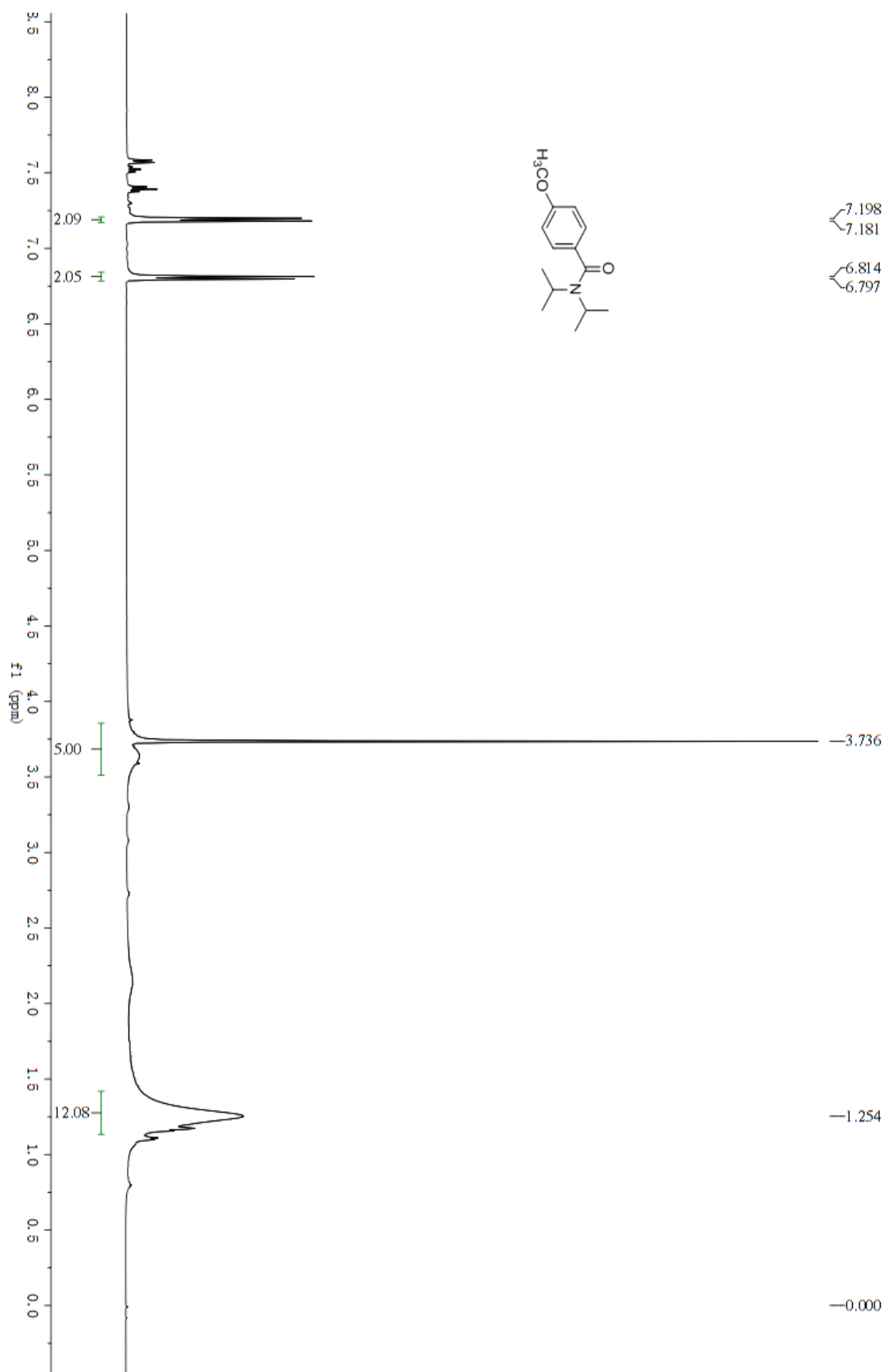
N-ethyl-*N*-isopropyl-4-methoxybenzamide **3p**



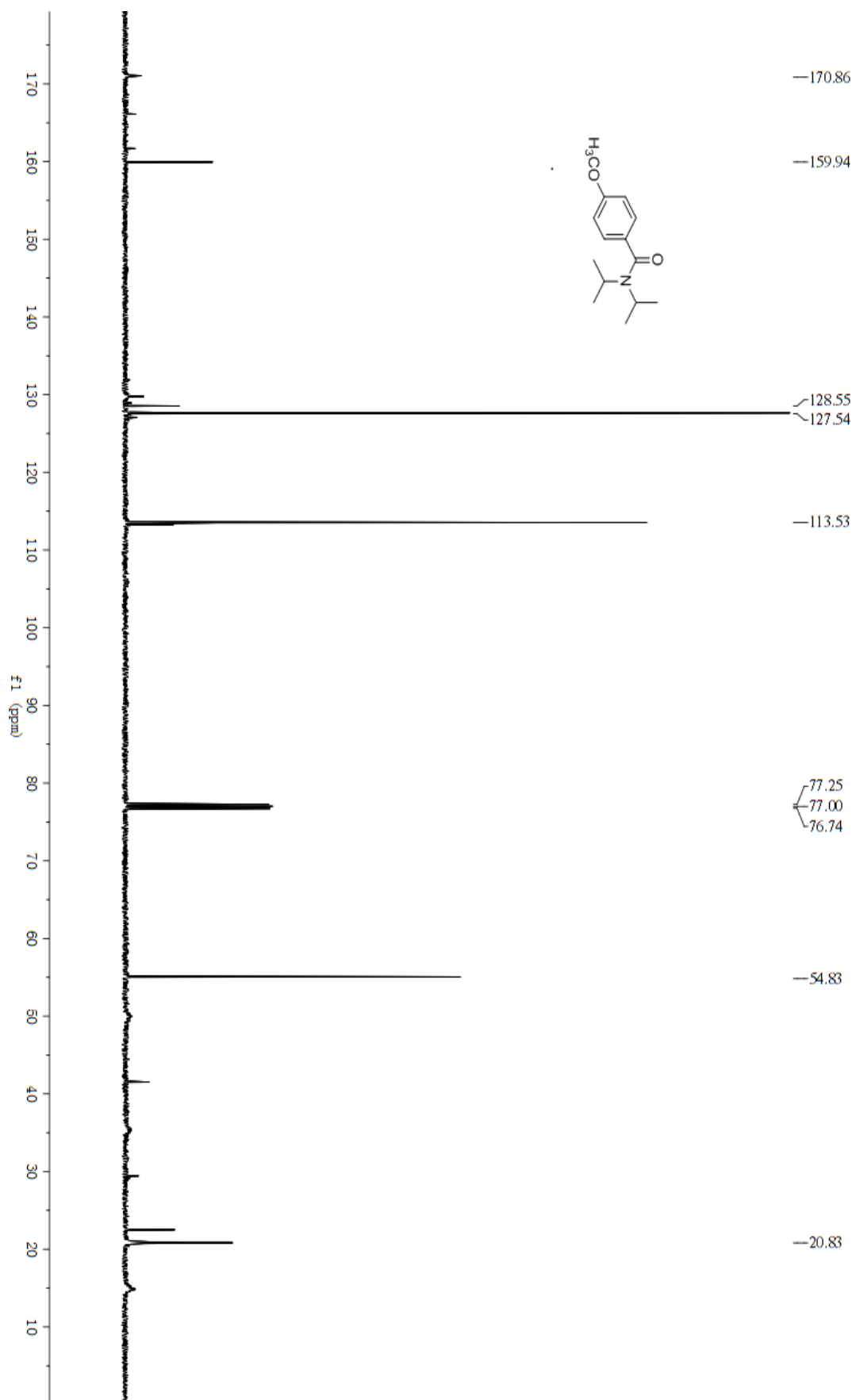
N-ethyl-*N*-isopropyl-4-methoxybenzamide **3p**



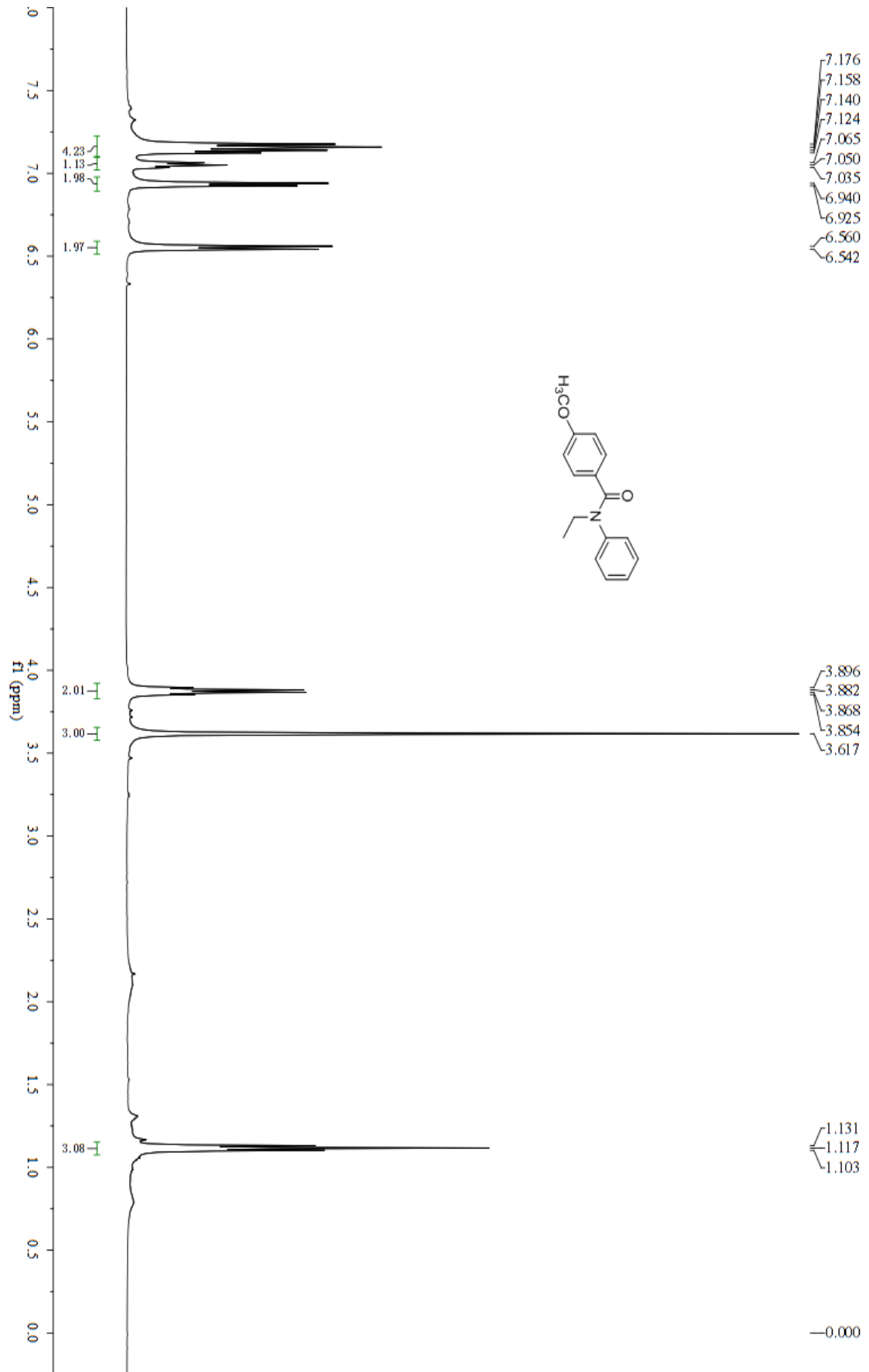
N,N-diisopropyl-4-methoxybenzamide **3q**



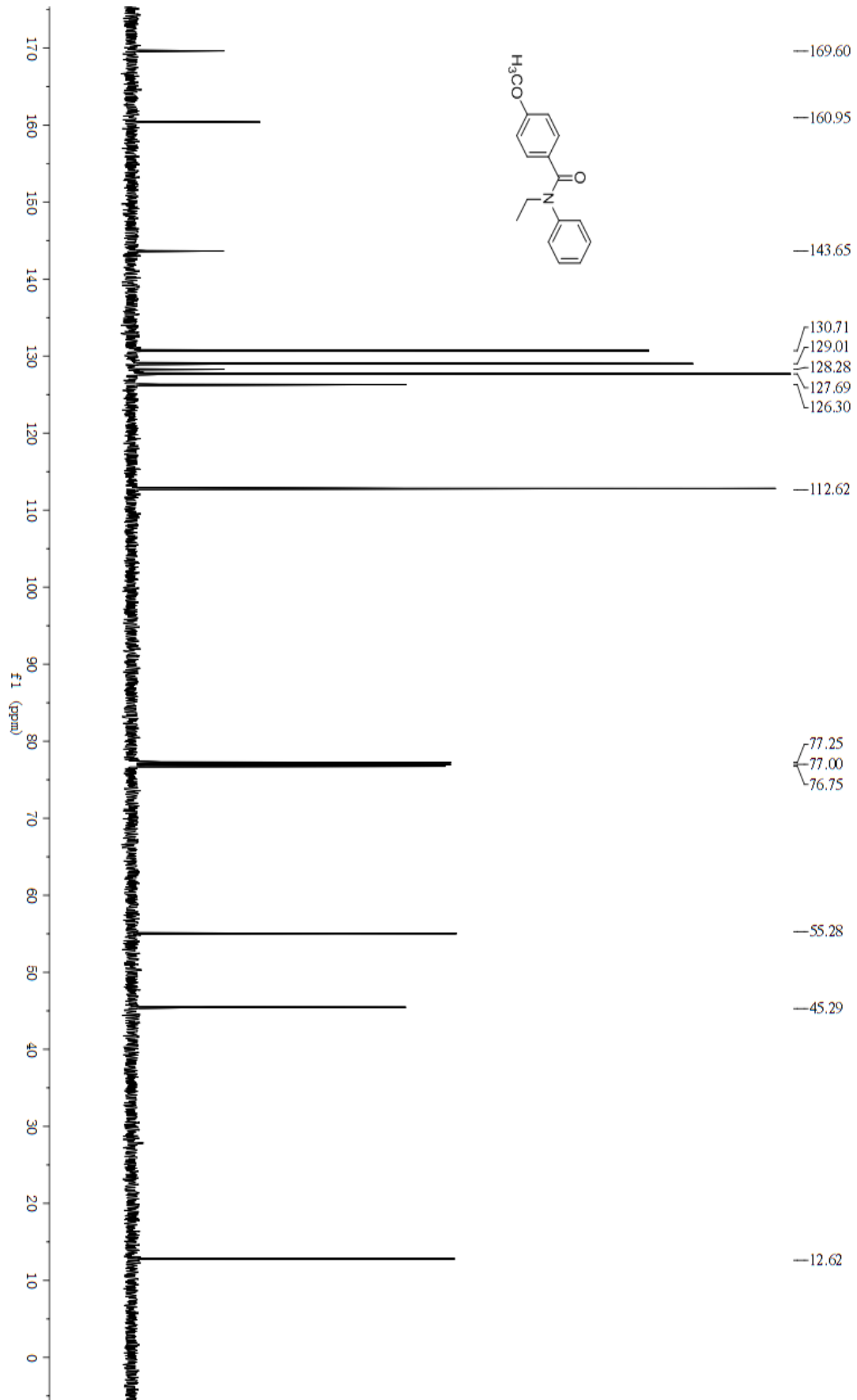
N,N-diisopropyl-4-methoxybenzamide **3q**



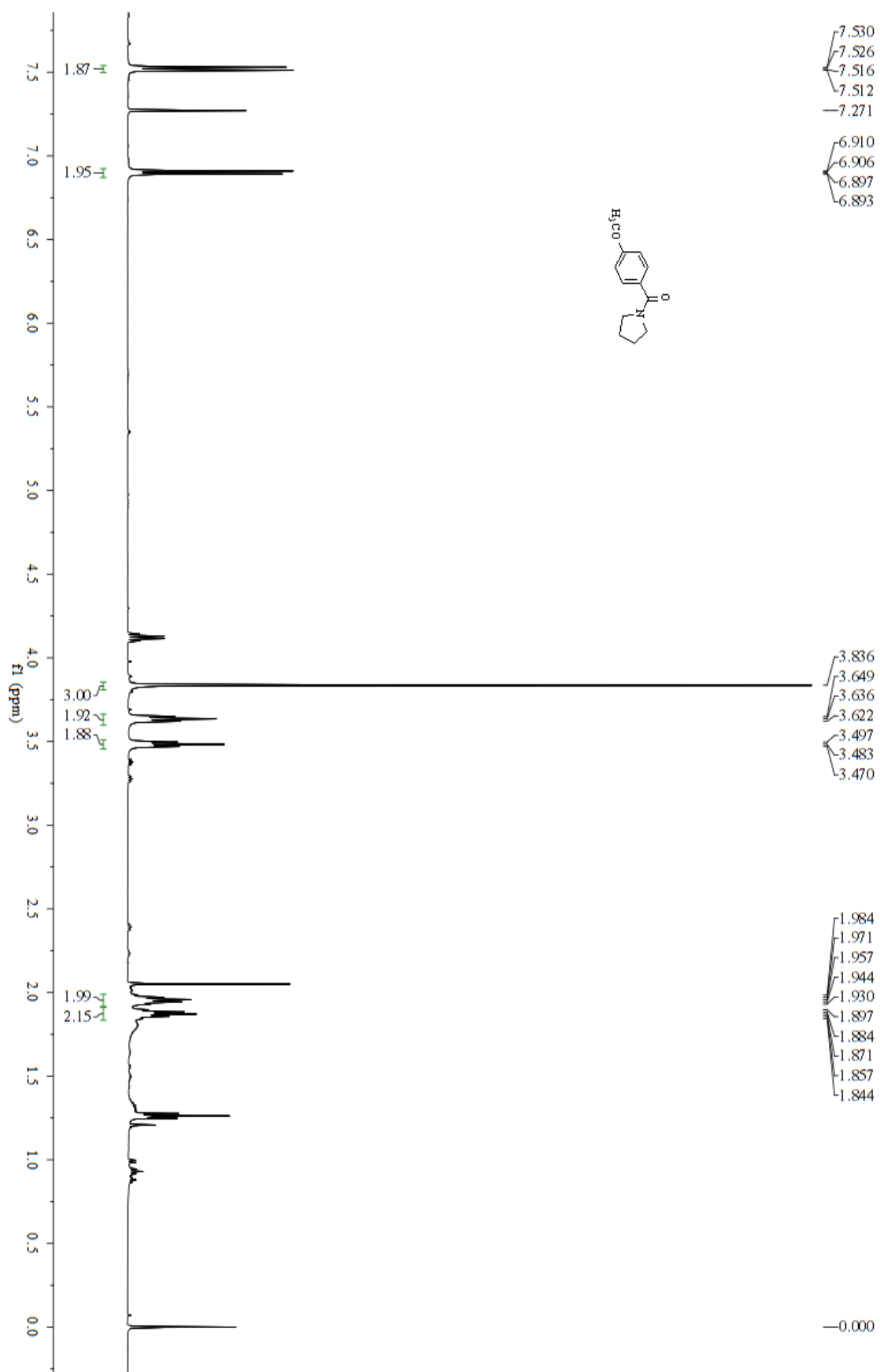
N-ethyl-*N*-phenyl-4-methoxybenzamide **3r**



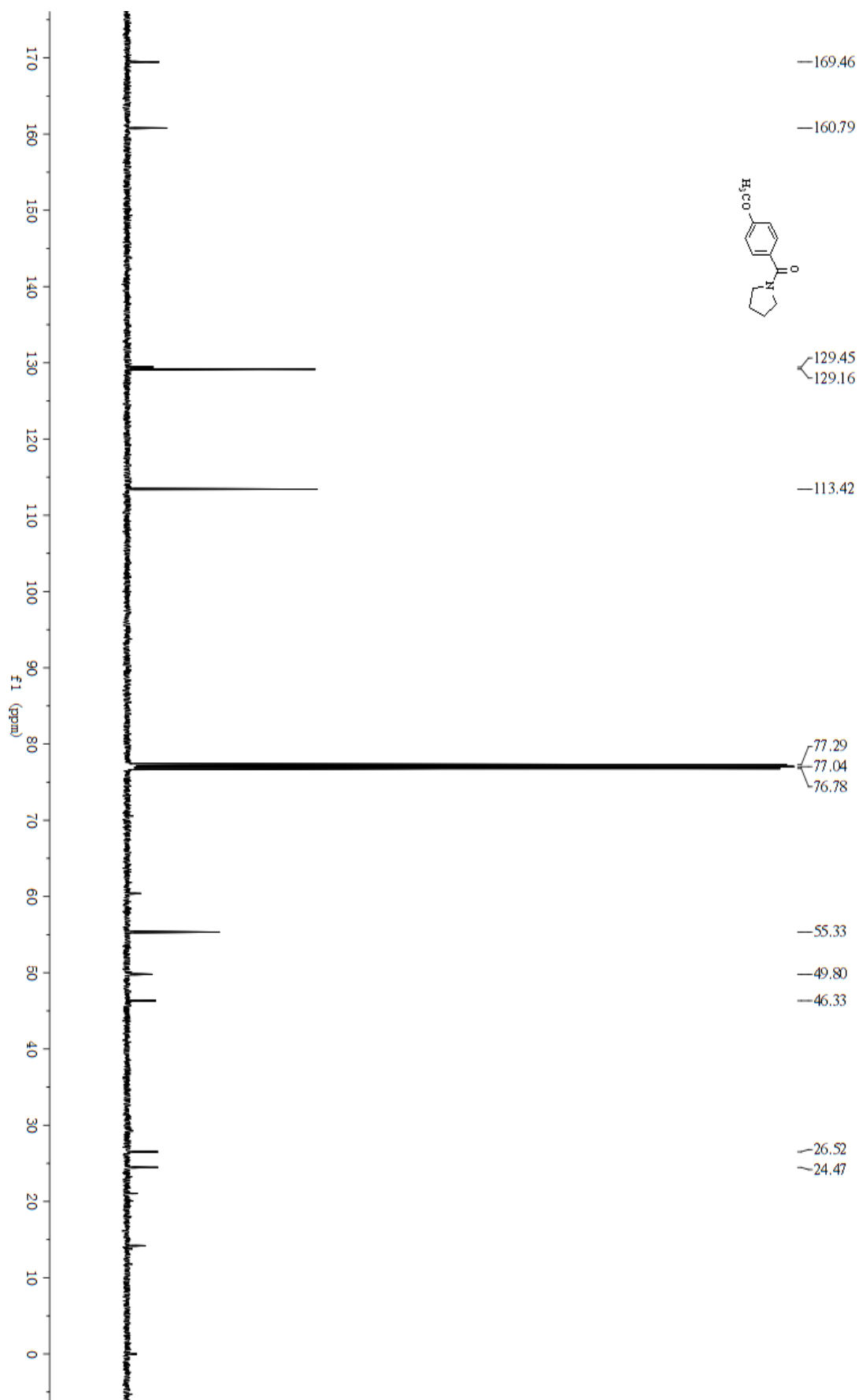
N-ethyl-*N*-phenyl-4-methoxybenzamide **3r**



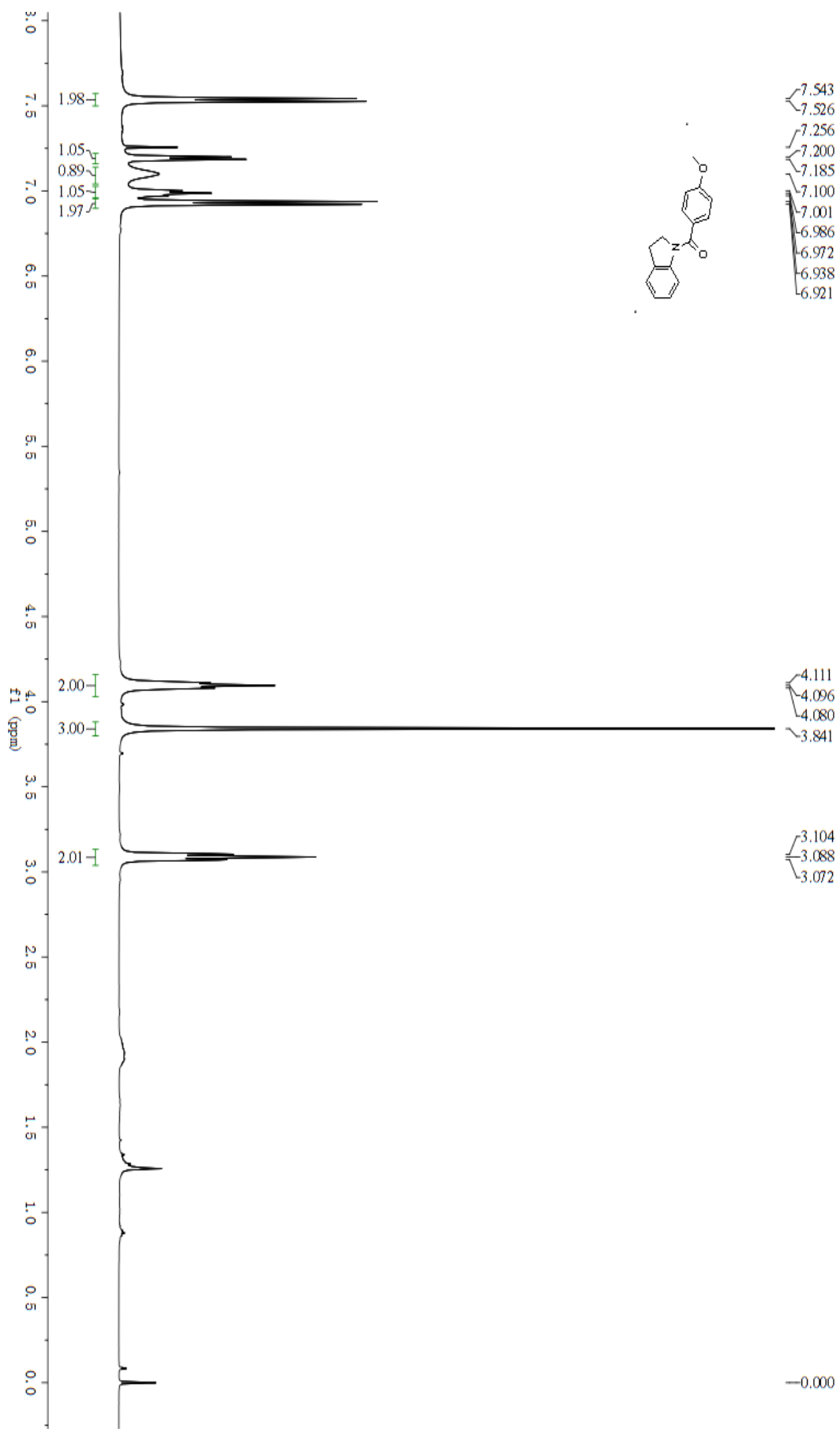
(4-methoxyphenyl)(pyrrolidin-1-yl)methanone **3s**



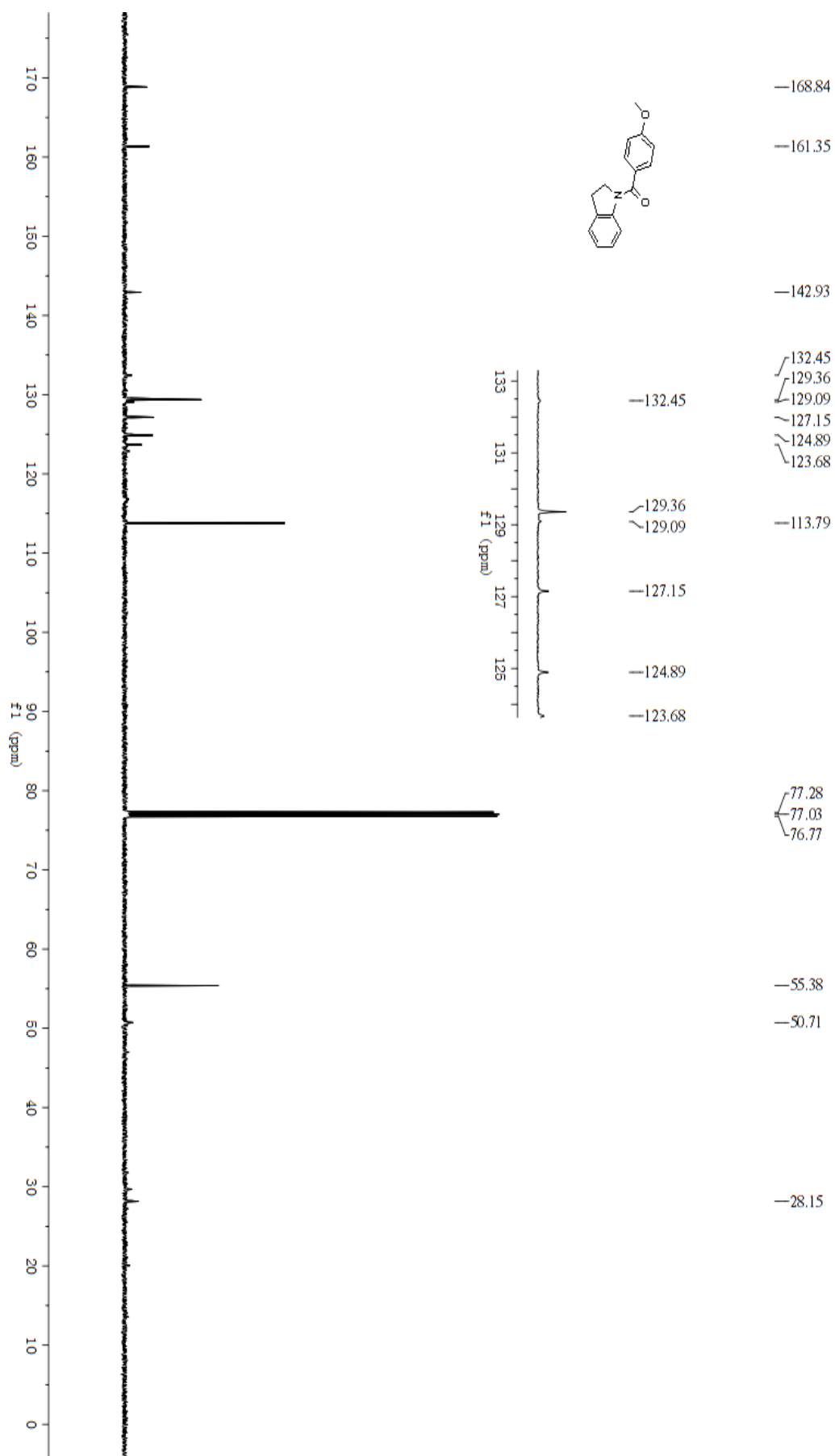
(4-methoxyphenyl)(pyrrolidin-1-yl)methanone **3s**



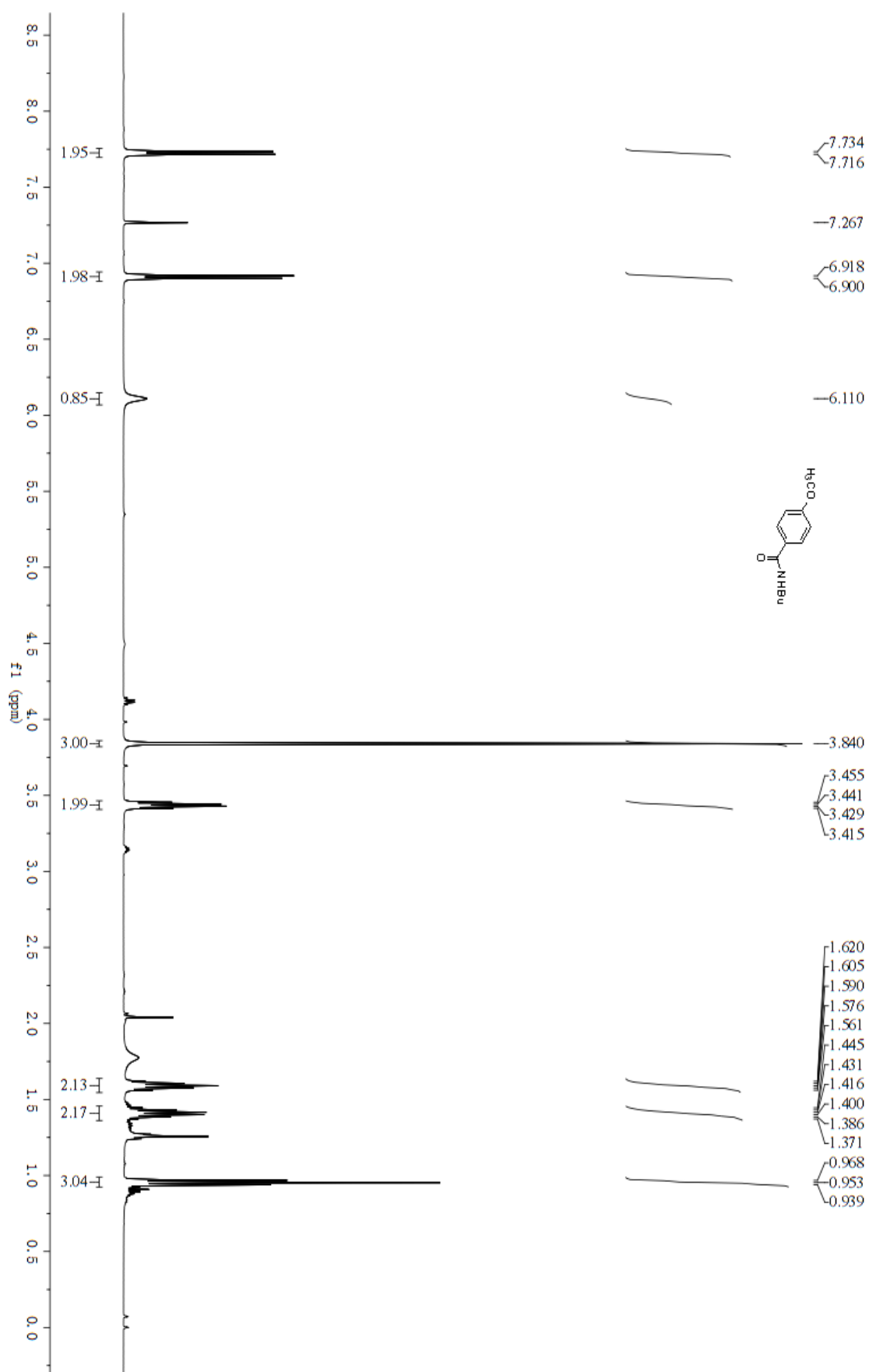
ndolin-1-yl(4-methoxyphenyl)methanone **3t**



olin-1-yl(4-methoxyphenyl)methanone **3t**



N-butyl-4-methoxybenzamide **3u**



N-butyl-4-methoxybenzamide **3u**

