

Synthesis of Chalcone Derivatives: Inducing Apoptosis of HepG2 Cells via Regulating ROS and Mitochondrial Pathway

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1. Compounds spectrum analysis

1.1. (*E*)-5-(3-(4-fluorophenyl)acryloyl)-2-hydroxybenzamide (**a1**)

White solid, m.p. 222.2-222.7 °C. Yield 83%; ¹H NMR (400 MHz, DMSO-*d*6): δ 13.98 (s, 1H), 8.80 (s, 1H), 8.71 (s, 1H), 8.23 (d, *J* = 8.8 Hz, 2H), 7.99 (t, 2H), 7.94 (d, *J* = 15.5 Hz, 1H), 7.77 (d, *J* = 15.5 Hz, 1H), 7.36 (t, 2H), 7.07 (d, *J* = 8.6 Hz, 1H); ¹³C NMR (100 MHz, DMSO-*d*6): δ 186.97, 172.28, 165.84, 162.43, 142.60, 134.82, 131.97, 131.52, 130.10, 128.98, 122.08, 118.64, 116.51, 116.29, 114.30; HRMS (ESI): found 286.0890 for [M+H]⁺ (calcd. for C₁₆H₁₂FNO₃, 285.0801). Purity: 99.0% (by HPLC).

1.2. (*E*)-5-(3-(3-chlorophenyl)acryloyl)-2-hydroxybenzamide (**a2**)

White solid, m.p. 212.3-212.8 °C. Yield 76%; ¹H NMR (400 MHz, DMSO-*d*6): δ 14.01 (s, 1H), 8.80 (s, 1H), 8.72 (s, 1H), 8.24 (d, *J* = 8.0 Hz, 2H), 8.07 (d, 2H), 7.83 (s, 1H), 7.74 (d, *J* = 15.6 Hz, 1H), 7.51 (t, 2H), 7.06 (d, *J* = 8.6 Hz, 1H); ¹³C NMR (100 MHz, DMSO-*d*6): δ 186.67, 171.78, 165.52, 151.26, 142.13, 137.55, 134.91, 134.28, 131.08, 130.51, 130.17, 128.79, 128.23, 123.71, 118.67, 114.32; HRMS (ESI): found 324.0410 for [M+Na]⁺ (calcd. for C₁₆H₁₂ClNO₃, 301.0506). Purity: 91.4% (by HPLC).

1.3. (*E*)-5-(3-(3-bromophenyl)acryloyl)-2-hydroxybenzamide (**a3**)

Pink solid, m.p. 211.4-211.5 °C. Yield 68%; ¹H NMR (400 MHz, DMSO-*d*6): δ 13.99 (s, 1H), 8.79 (s, 1H), 8.71 (s, 1H), 8.25 (d, *J* = 8.8 Hz, 2H), 8.16 (s, 1H), 8.05 (d, *J* = 15.6 Hz, 1H), 7.87 (d, *J* = 8.0 Hz, 1H), 7.74 (d, *J* = 15.2 Hz, 1H), 7.66 (d, *J* = 8.0 Hz, 1H), 7.46 (t, 1H), 7.06 (d, *J* = 8.7 Hz, 1H); ¹³C NMR (100 MHz, DMSO-*d*6): δ 190.71, 172.46, 170.94, 166.24, 144.33, 136.24, 134.54, 131.90, 130.87, 130.13, 127.02, 126.52, 123.24, 122.63, 117.84, 116.89; HRMS (ESI): found 369.9879 for [M+Na]⁺ (calcd. for C₁₆H₁₂BrNO₃, 345.0001). Purity: 95.1% (by HPLC).

1.4. (*E*)-5-(3-(4-bromophenyl)acryloyl)-2-hydroxybenzamide (**a4**)

Yellow solid, m.p. 226.9-227.5 °C. Yield 66%; ¹H NMR (400 MHz, DMSO-*d*6): δ 13.99 (s, 1H), 8.79 (s, 1H), 8.70 (s, 1H), 8.22 (d, *J* = 8.8 Hz, 2H), 8.01 (d, *J* = 15.6 Hz, 1H), 7.86 (d, *J* = 8.4 Hz, 2H), 7.73 (d, *J* = 15.8 Hz, 1H), 7.69 (d, *J* = 8.5 Hz, 2H), 7.06 (d, *J* = 8.7 Hz, 1H); ¹³C NMR (100 MHz, DMSO-*d*6): δ 187.09, 172.33, 165.90, 142.48, 134.87, 134.59, 132.37, 131.11, 130.16, 128.90, 124.32, 122.97, 118.69, 114.32; HRMS (ESI): found 346.0056 for [M+H]⁺ (calcd. for C₁₆H₁₂BrNO₃, 345.0001). Purity: 94.2% (by HPLC).

1.5. (*E*)-5-cinnamoyl-2-hydroxybenzamide (**a5**)

Yellow solid, m.p. 182.7-184.5 °C. Yield 74%; ¹H NMR (400 MHz, DMSO-*d*6): δ 14.01 (s, 1H), 8.82 (s, 1H), 8.72 (s, 1H), 8.23 (d, *J* = 8.8 Hz, 2H), 8.00 (d, *J* = 15.5 Hz, 1H), 7.89 (m, 2H), 7.78 (d, *J* = 15.5 Hz, 1H), 7.50 (m, 3H), 7.08 (d, *J* = 8.7 Hz, 1H); ¹³C NMR (100 MHz, DMSO-*d*6): δ 187.11, 172.06, 165.95, 135.29, 134.84, 130.99, 130.14, 129.31, 129.00, 122.19, 118.70, 118.19, 114.27; HRMS (ESI): found 268.0983 for [M+H]⁺ (calcd. for C₁₆H₁₃NO₃, 267.0895). Purity: 95.9% (by HPLC).

1.6. (*E*)-2-hydroxy-5-(3-(*o*-tolyl)acryloyl)benzamide (**a6**)

Yellow solid, m.p. 163.1-164.9 °C. Yield 68%; ^1H NMR (400 MHz, DMSO-*d*6): δ 14.02 (s, 1H), 8.83 (s, 1H), 8.72 (s, 1H), 8.21 (d, J = 9.0 Hz, 2H), 8.04 (d, J = 15.6 Hz, 1H), 7.99 (d, J = 8.2 Hz, 1H), 7.89 (d, J = 15.4 Hz, 1H), 7.41 – 7.23 (m, 3H), 7.07 (d, J = 8.7 Hz, 1H), 2.44 (d, J = 12.4 Hz, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 187.14, 172.18, 165.86, 140.99, 138.39, 134.82, 133.97, 131.27, 130.64, 130.09, 129.00, 127.20, 126.69, 123.12, 118.76, 114.20, 38.70 – 38.06; HRMS (ESI): found 282.1193 for [M+H] $^+$ (calcd. for C₁₇H₁₅NO₃, 281.1052). Purity: 99.1% (by HPLC).

1.7. (*E*)-2-hydroxy-5-(3-(*m*-tolyl)acryloyl)benzamide (**a7**)

White solid, m.p. 171.4-171.5 °C. Yield 62%; ^1H NMR (400 MHz, DMSO-*d*6): δ 14.00 (s, 1H), 8.83 (s, 1H), 8.71 (s, 1H), 8.20 (d, J = 8.9 Hz, 2H), 7.95 (d, J = 15.5 Hz, 1H), 7.72 (d, J = 15.5 Hz, 1H), 7.68 – 7.59 (m, 2H), 7.34 (t, J = 7.5 Hz, 1H), 7.25 (d, J = 7.4 Hz, 1H), 7.04 (d, J = 8.7 Hz, 1H), 2.36 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 187.16, 165.93, 143.90, 138.46, 135.22, 134.81, 131.69, 130.06, 129.55, 129.16, 128.90, 126.60, 121.89, 118.66, 114.24, 21.31; HRMS (ESI): found 282.1191 for [M+H] $^+$ (calcd. for C₁₇H₁₅NO₃, 281.1052). Purity: 99.8% (by HPLC).

1.8. (*E*)-2-hydroxy-5-(3-(*p*-tolyl)acryloyl)benzamide (**a8**)

Yellow solid, m.p. 190.5-191.7 °C. Yield 71%; ^1H NMR (400 MHz, DMSO-*d*6): δ 13.63 (s, 1H), 8.77 (s, 1H), 8.69 (s, 1H), 8.32 – 8.02 (m, 2H), 8.01 (d, J = 15.5 Hz, 1H), 7.85 – 7.65 (m, 2H), 7.49 – 7.12 (m, 3H), 7.03 (d, J = 8.7 Hz, 1H), 2.33 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 187.05, 172.33, 165.94, 143.86, 140.93, 134.73, 132.51, 130.02, 129.10, 127.01, 121.01, 118.66, 114.15, 20.99; HRMS (ESI): found 282.1228 for [M+H] $^+$ (calcd. for C₁₇H₁₅NO₃, 281.1052). Purity: 91.5% (by HPLC).

1.9. (*E*)-2-hydroxy-5-(3-(2-methoxyphenyl)acryloyl)benzamide (**a9**)

Yellow solid, m.p. 237.2-238.9 °C. Yield 71%; ^1H NMR (400 MHz, DMSO-*d*6): δ 13.97 (s, 1H), 8.81 (s, 1H), 8.70 (s, 1H), 8.20 (d, J = 8.8 Hz, 2H), 8.11 (d, J = 15.6 Hz, 1H), 7.99 (d, J = 7.6 Hz, 1H), 7.93 (d, J = 15.7 Hz, 1H), 7.47 (t, J = 7.3 Hz, 1H), 7.13 (d, J = 8.2 Hz, 1H), 7.09 – 7.05 (t, J = 7.7 Hz, 2H), 3.91 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 190.71, 170.94, 166.24, 157.37, 140.86, 134.54, 130.87, 130.57, 129.59, 127.84, 126.52, 121.88, 120.84, 117.84, 116.89, 113.50, 56.79; HRMS (ESI): found 298.1220 for [M+H] $^+$ (calcd. for C₁₇H₁₅NO₄, 297.1001). Purity: 97.7% (by HPLC).

1.10. (*E*)-2-hydroxy-5-(3-(3-methoxyphenyl)acryloyl)benzamide (**a10**)

Yellow solid, m.p. 194.5-195.2 °C. Yield 77%; ^1H NMR (400 MHz, DMSO-*d*6): δ 13.98 (s, 1H), 8.80 (s, 1H), 8.71 (s, 1H), 8.23 (d, J = 8.7 Hz, 2H), 7.98 (d, J = 15.5 Hz, 1H), 7.74 (d, J = 15.5 Hz, 1H), 7.46 (s, 2H), 7.40 (t, J = 8.0 Hz, 1H), 7.06 (d, J = 8.7 Hz, 2H), 3.84 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 187.15, 172.20, 165.97, 160.04, 143.87, 136.49, 134.76, 130.27, 129.94, 128.83, 122.38, 121.68, 118.52, 116.50, 114.22, 55.60; HRMS (ESI): found 298.1179 for [M+H] $^+$ (calcd. for C₁₇H₁₅NO₄, 297.1001). Purity: 98.9% (by HPLC).

1.11. (*E*)-2-hydroxy-5-(3-(4-methoxyphenyl)acryloyl)benzamide (**a11**)

Yellow solid, m.p. 212.8-212.8 °C. Yield 59%; ^1H NMR (400 MHz, DMSO-*d*6): δ 13.95 (s, 1H), 8.80 (s, 1H), 8.70 (s, 1H), 8.21 (d, J = 8.7 Hz, 2H), 7.93 – 7.81 (m, 3H), 7.74 (d, J = 15.4 Hz, 1H), 7.05 (d, J = 8.5 Hz, 3H), 3.84 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 187.04, 172.21, 165.90, 161.50,

143.73, 134.56, 131.02, 129.83, 129.16 – 128.96, 127.70, 119.32, 118.45, 114.58, 114.14, 55.85; HRMS (ESI): found 320.0902 for $[M+Na]^+$ (calcd. for $C_{17}H_{15}NO_4$, 297.1001). Purity: 95.5% (by HPLC).

1.12. (*E*)-5-(3-(2,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (**a12**)

Yellow solid, m.p. 193.7-193.7 °C. Yield 76%; 1H NMR (400 MHz, DMSO-*d*6): δ 13.84 (s, 1H), 8.73 (s, 1H), 8.53 (s, 1H), 8.21 – 8.07 (m, 2H), 8.05 (d, J = 15.5 Hz, 1H), 7.98 (m, 1H), 7.79 (d, J = 15.5 Hz, 1H), 6.98 (d, J = 8.5 Hz, 2H), 6.65 (s, 1H), 3.88 (d, J = 23.9 Hz, 3H), 2.54 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 196.13, 187.13, 171.72, 165.66, 163.27, 160.01, 138.00, 134.09, 130.19, 129.61, 128.25 – 128.05, 119.03, 118.50, 117.96, 114.30, 106.85, 98.70, 56.28, 55.88, 26.62; HRMS (ESI): found 328.1198 for $[M+H]^+$ (calcd. for $C_{18}H_{17}NO_5$, 327.1107). Purity: 91.3% (by HPLC).

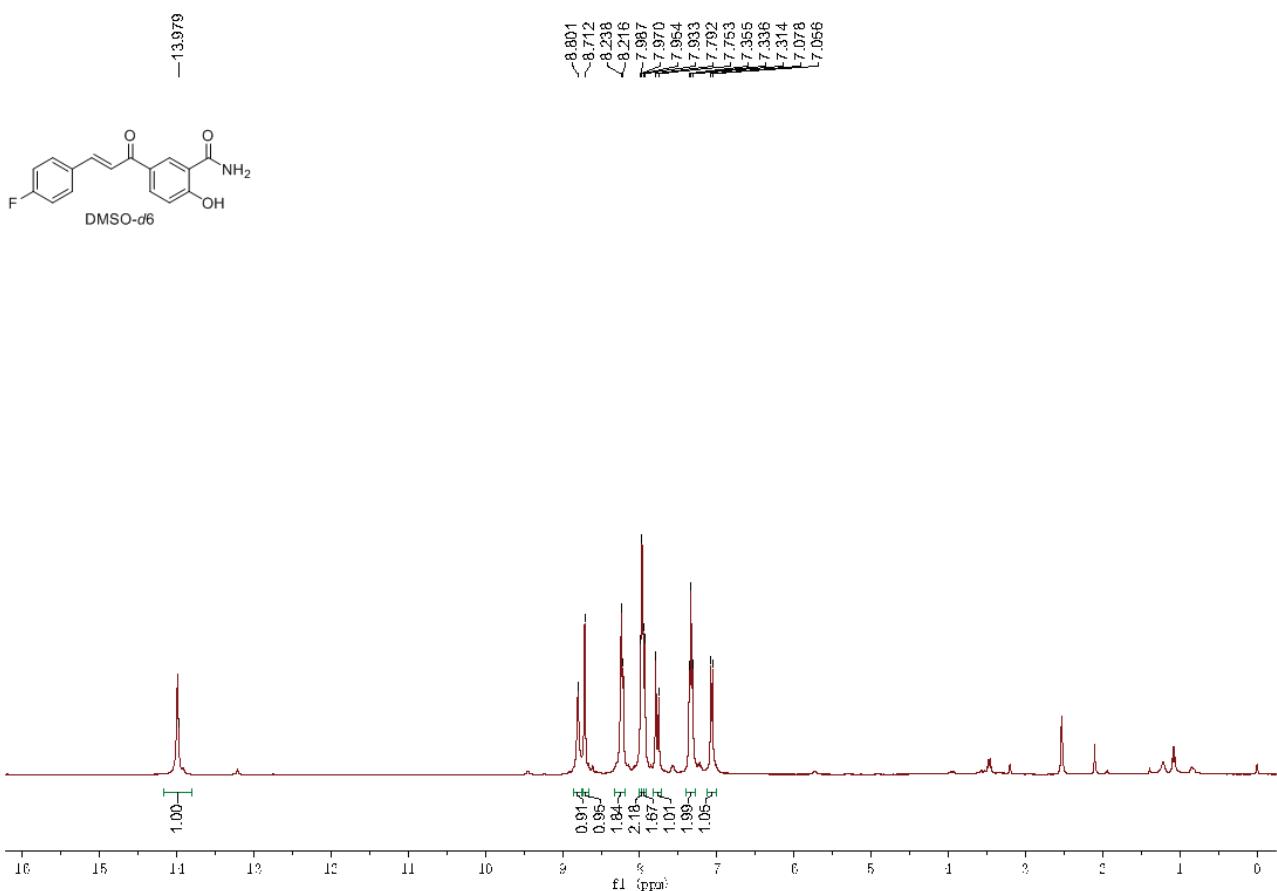
1.13. (*E*)-5-(3-(3,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (**a13**)

Yellow solid, m.p. 213.9-214.6 °C. Yield 85%; 1H NMR (400 MHz, DMSO-*d*6): δ 13.91 (s, 1H), 8.78 (s, 1H), 8.70 (s, 1H), 8.31 – 8.13 (m, 2H), 7.85 (d, J = 15.4 Hz, 1H), 7.73 (d, J = 15.4 Hz, 1H), 7.51 (s, 1H), 7.44 (d, J = 8.3 Hz, 1H), 7.06 (d, J = 8.7 Hz, 2H), 3.86 (d, J = 16.6 Hz, 6H); ^{13}C NMR (101 MHz, DMSO-*d*6): δ 187.22, 186.68, 172.24, 165.65, 151.78 – 151.54, 149.40, 144.56 – 144.18, 134.83 (s), 129.99, 129.32, 128.09, 123.83, 119.87, 118.42, 114.45, 112.11, 111.78, 56.14; HRMS (ESI): found 350.1010 for $[M+Na]^+$ (calcd. for $C_{18}H_{17}NO_5$, 327.1107). Purity: 94.8% (by HPLC).

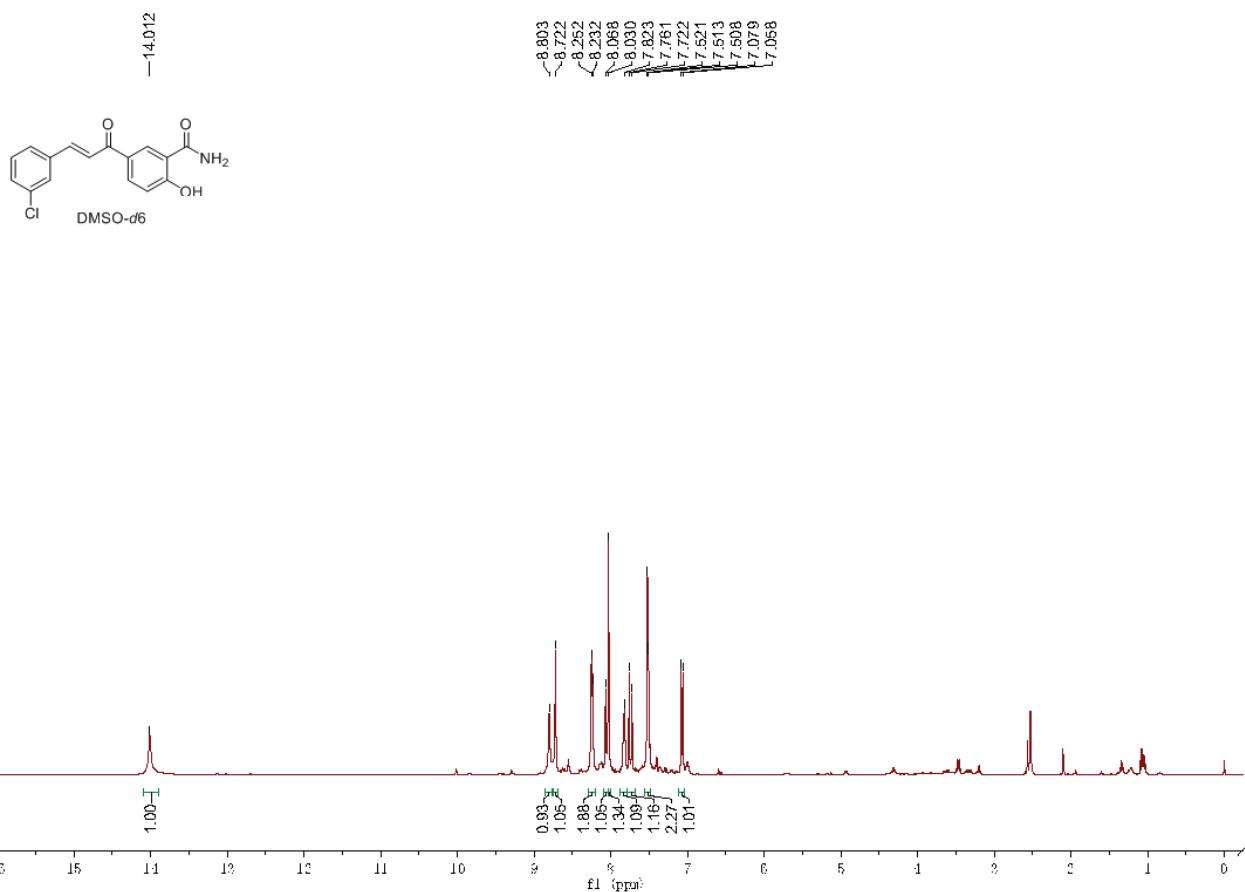
1.14. (*E*)-2-hydroxy-5-(3-(3,4,5-trimethoxyphenyl)acryloyl)benzamide (**a14**)

Yellow solid, m.p. 259.9-260.8 °C. Yield 82%; 1H NMR (400 MHz, DMSO-*d*6): δ 13.90 (s, 1H), 8.79 (s, 1H), 8.70 (s, 1H), 8.24 (d, J = 8.6 Hz, 1H), 8.17 (s, 1H), 7.89 (d, J = 15.5 Hz, 1H), 7.70 (d, J = 15.5 Hz, 1H), 7.22 (s, 2H), 7.05 (d, J = 8.6 Hz, 1H), 3.87 (s, 6H), 3.72 (s, 3H); ^{13}C NMR (100 MHz, DMSO-*d*6): δ 187.36, 172.20, 165.75, 153.59, 144.35, 140.23, 134.93, 130.82, 130.17, 129.19, 121.59, 118.36, 114.62, 107.02, 60.63, 56.62; HRMS (ESI): found 358.1356 for $[M+H]^+$ (calcd. for $C_{19}H_{19}NO_6$, 357.1212). Purity: 95.0% (by HPLC).

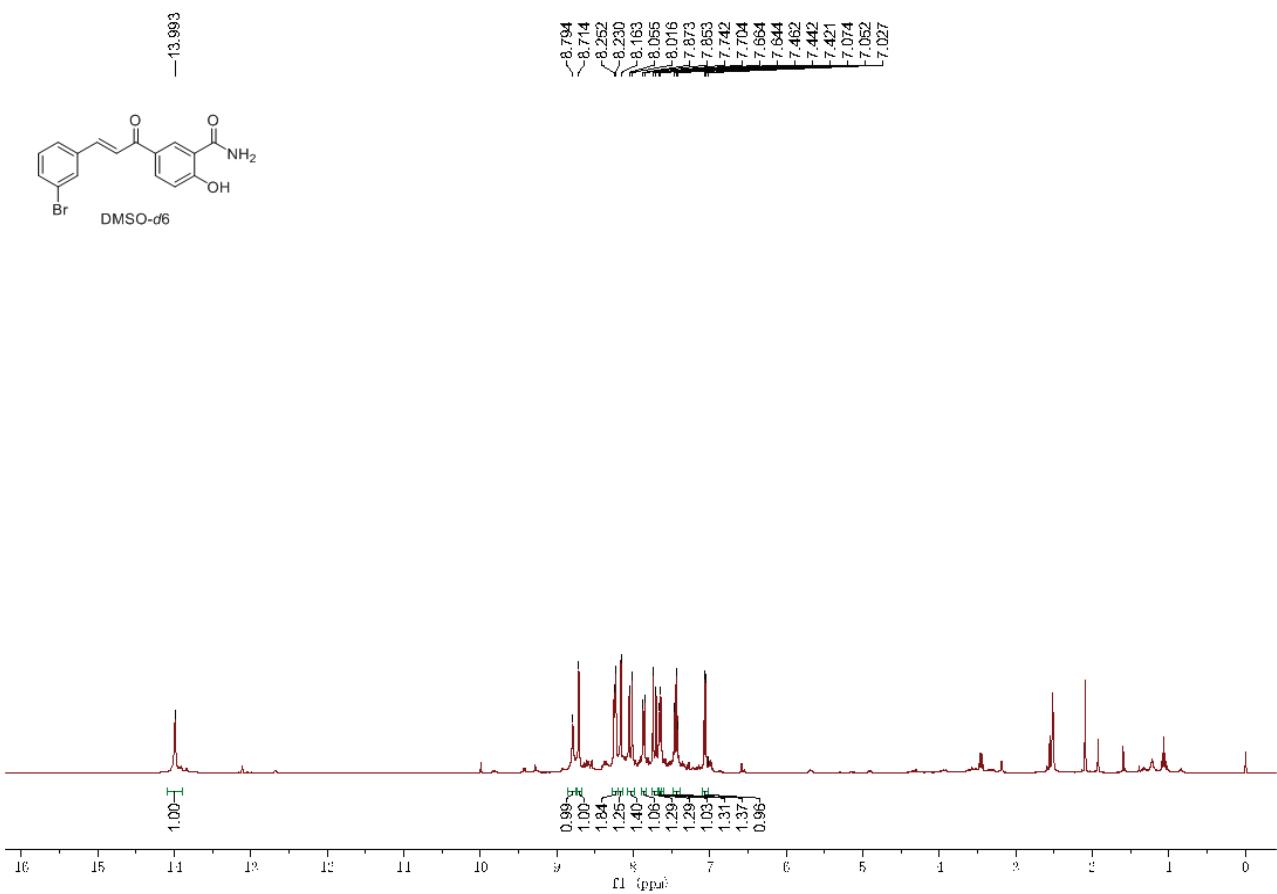
2. H-NMR information of the synthesized compounds

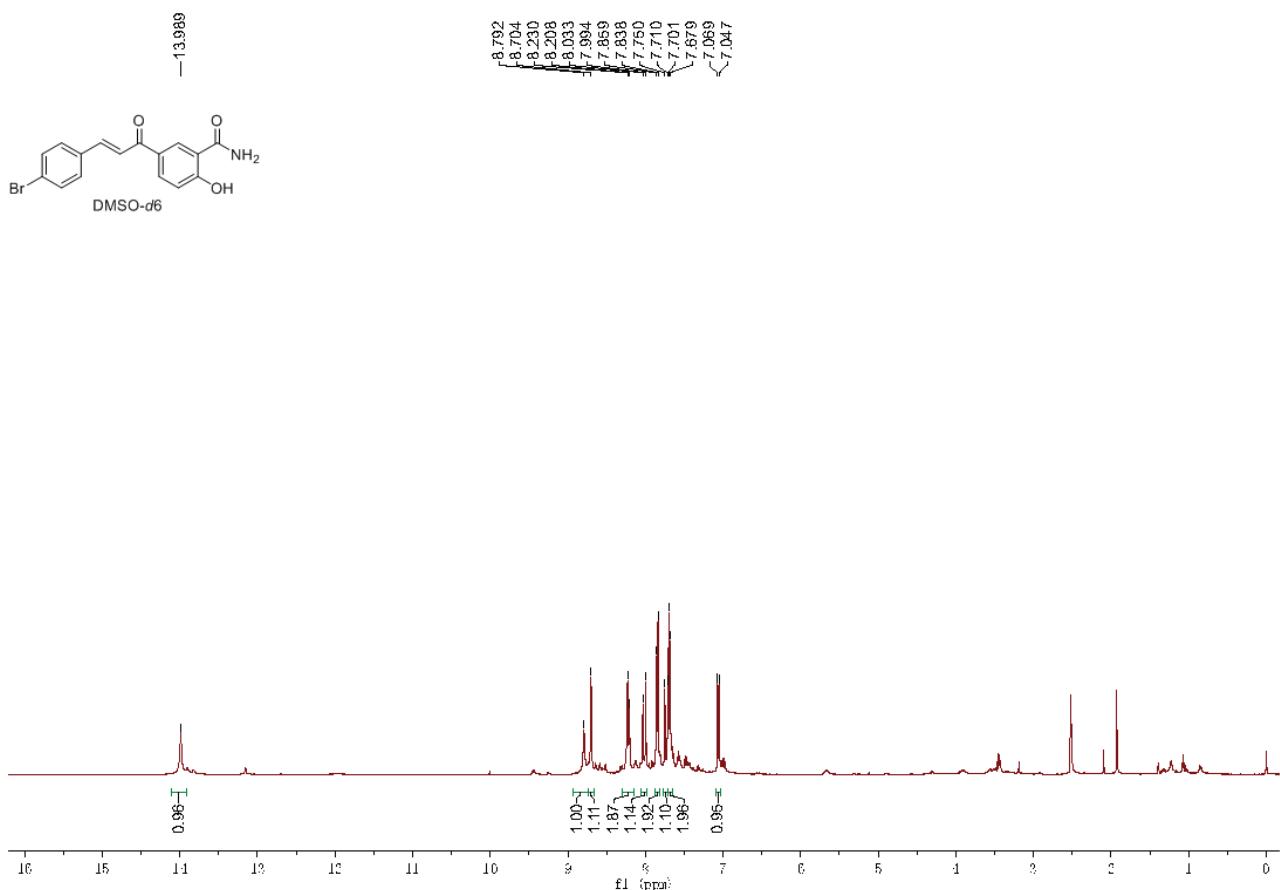


(E)-5-(3-(4-fluorophenyl)acryloyl)-2-hydroxybenzamide (**a1**)

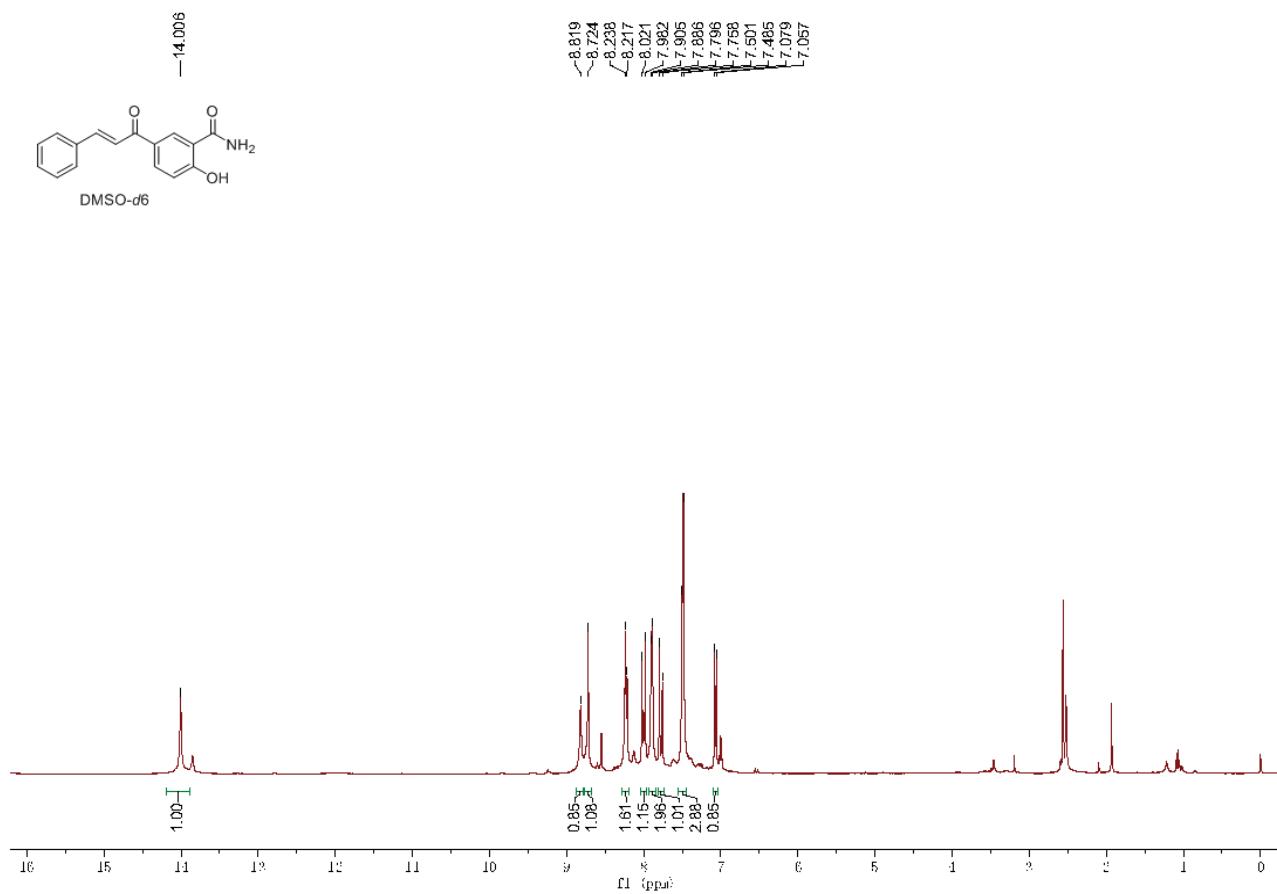


(*E*)-5-(3-(3-chlorophenyl)acryloyl)-2-hydroxybenzamide (**a2**)

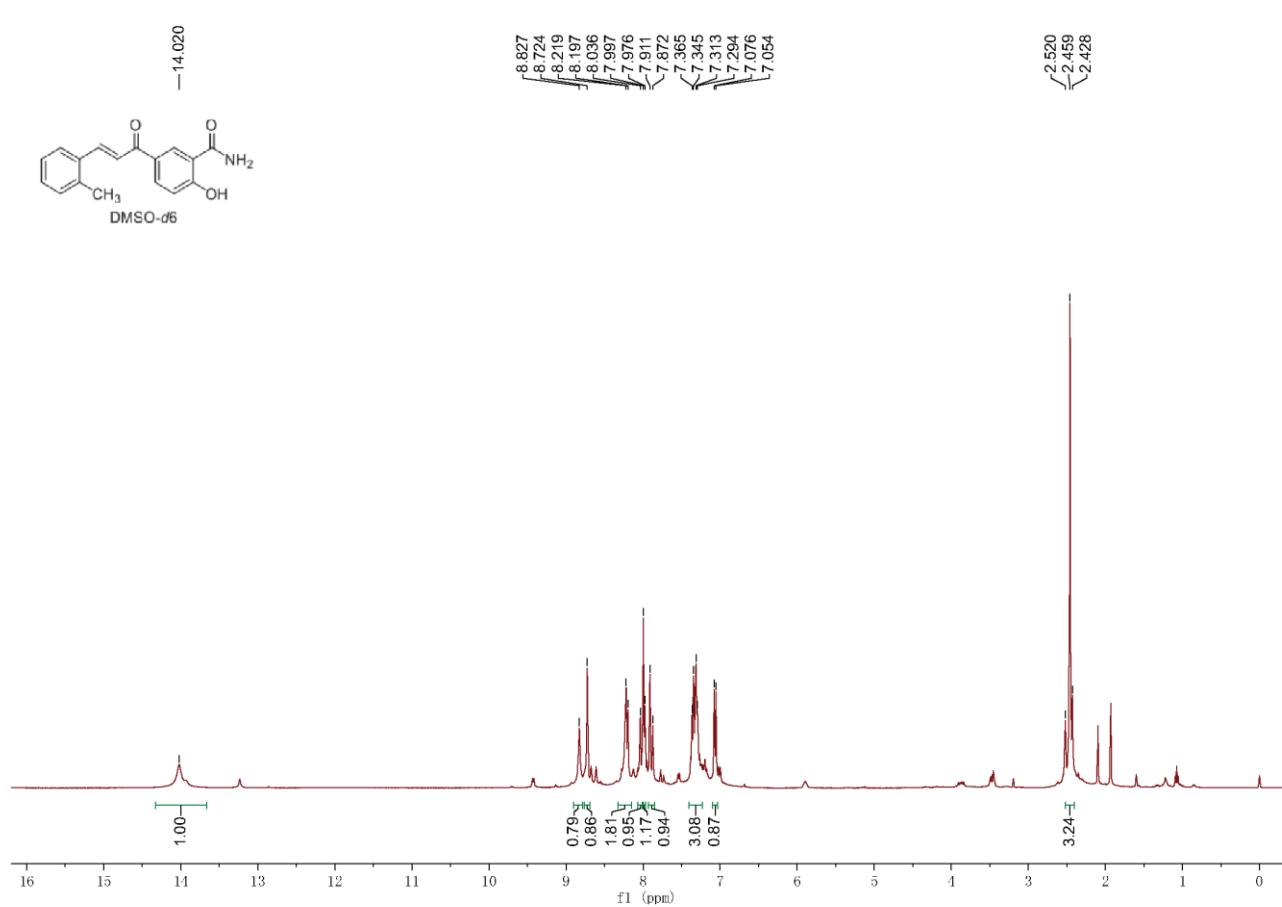




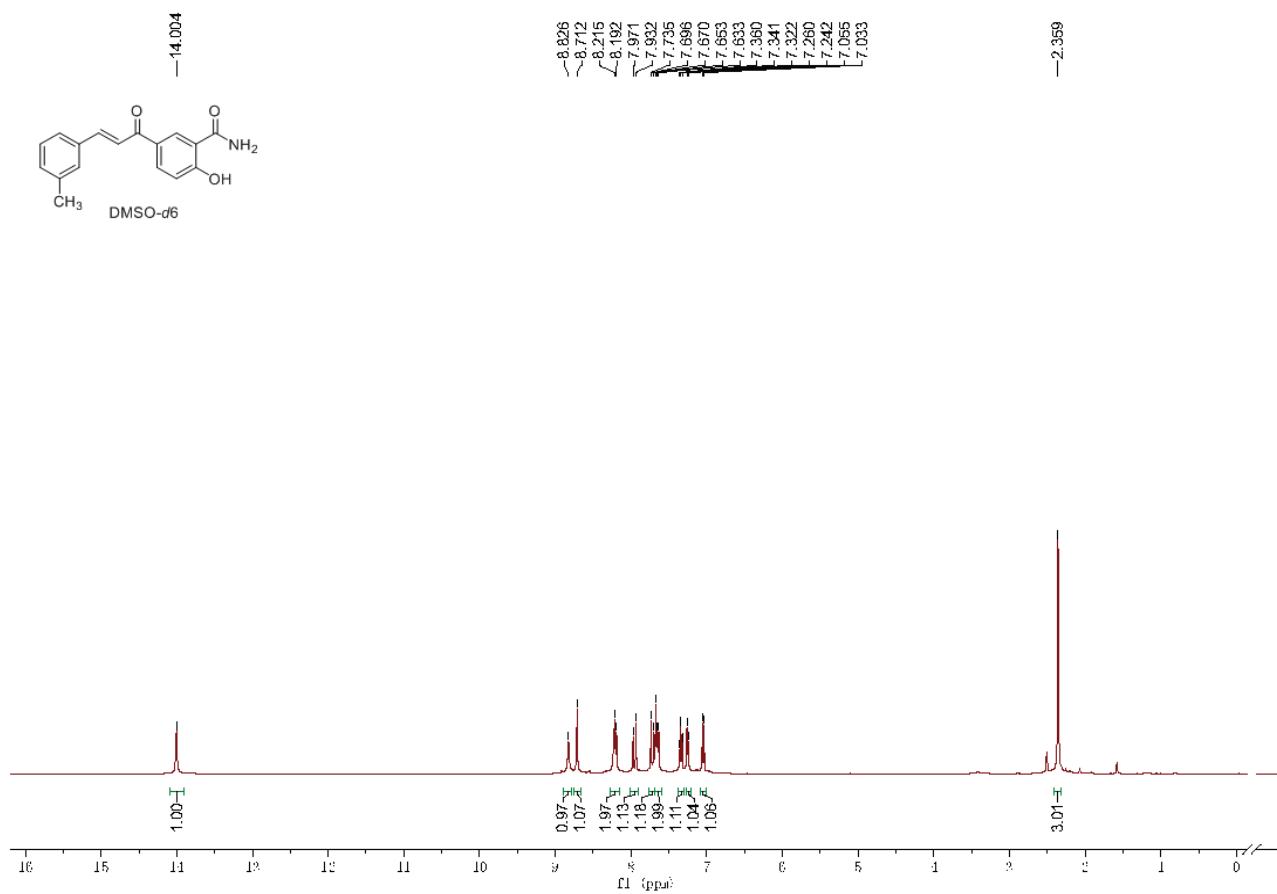
(E)-5-(3-(4-bromophenyl)acryloyl)-2-hydroxybenzamide (a4)



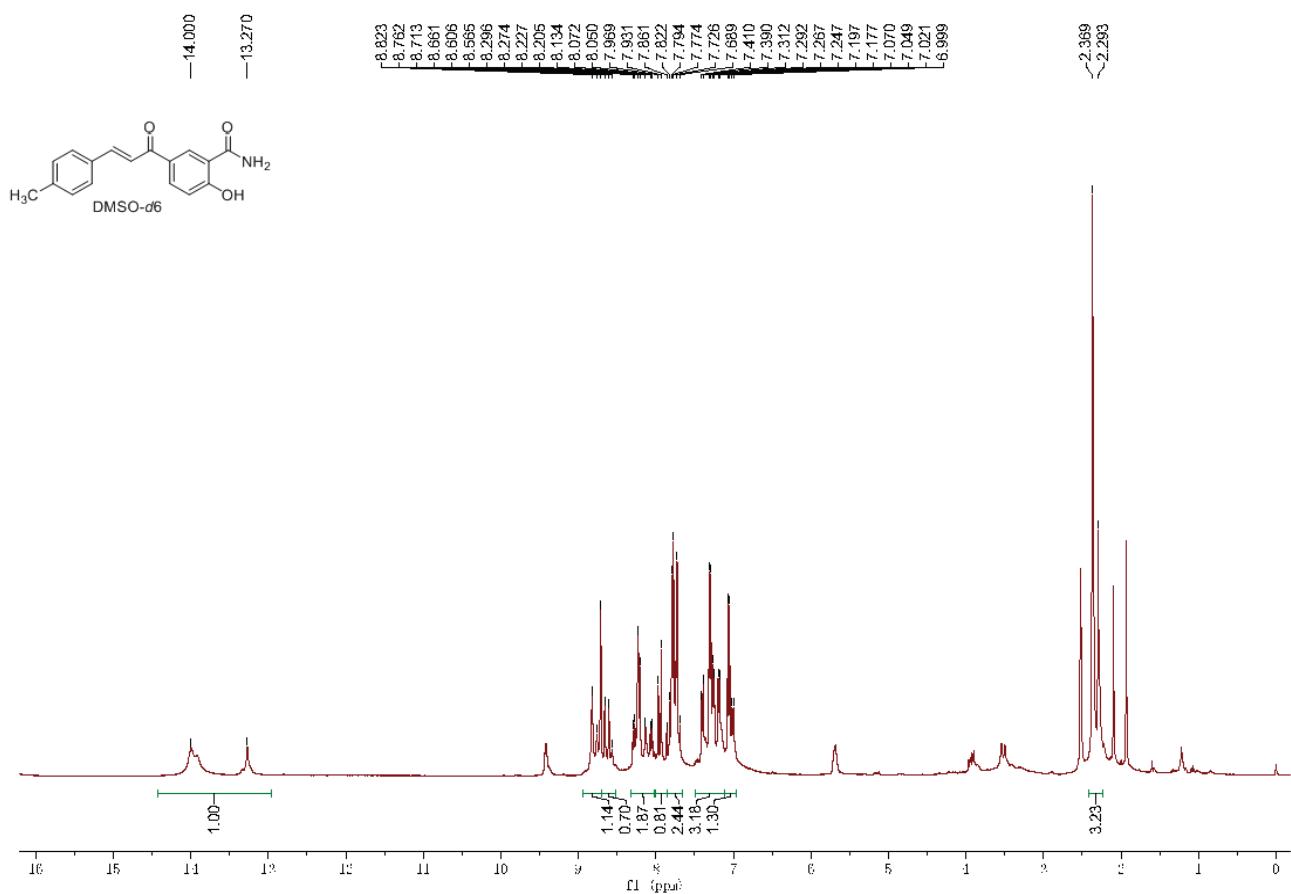
(E)-5-cinnamoyl-2-hydroxybenzamide (**a5**)



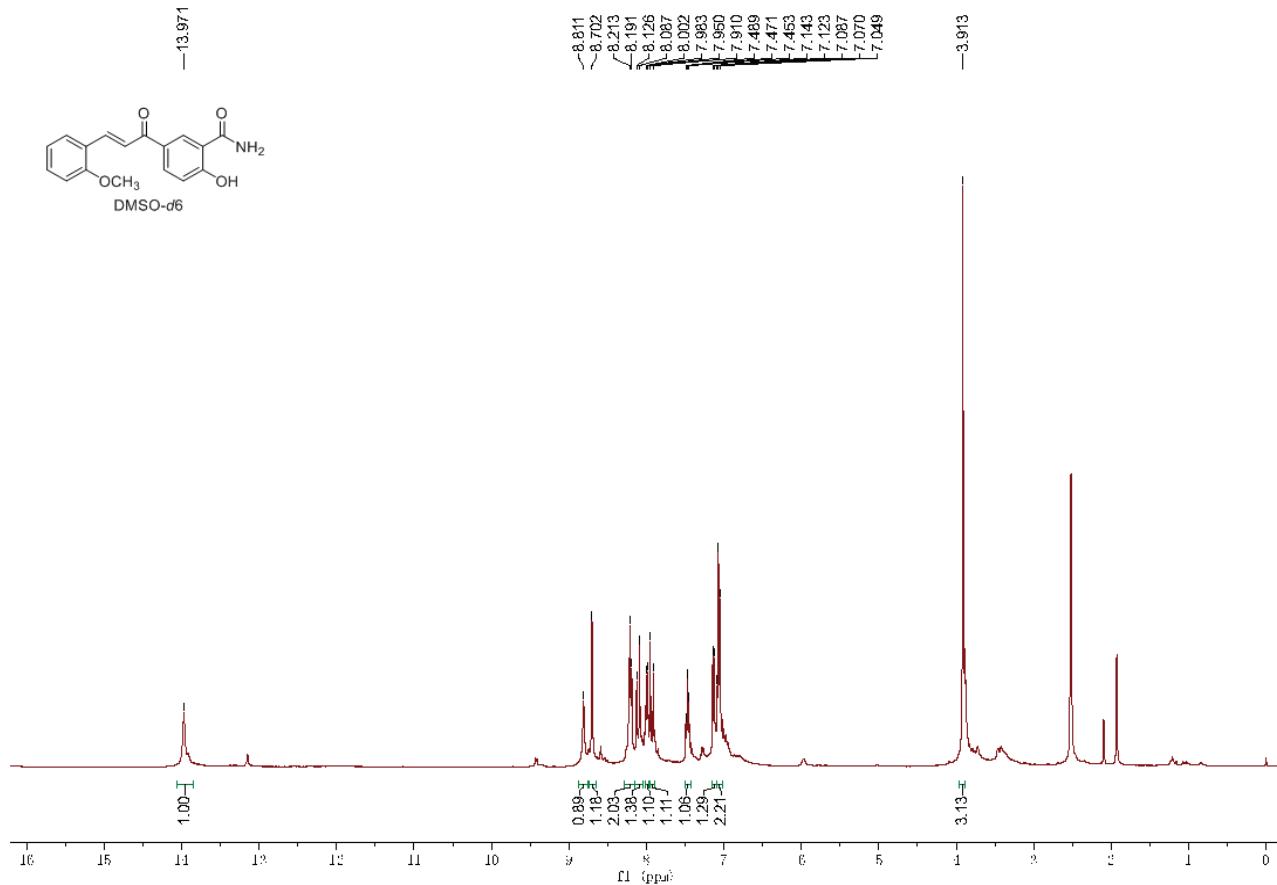
(E)-2-hydroxy-5-(3-(o-tolyl)acryloyl)benzamide (a6)



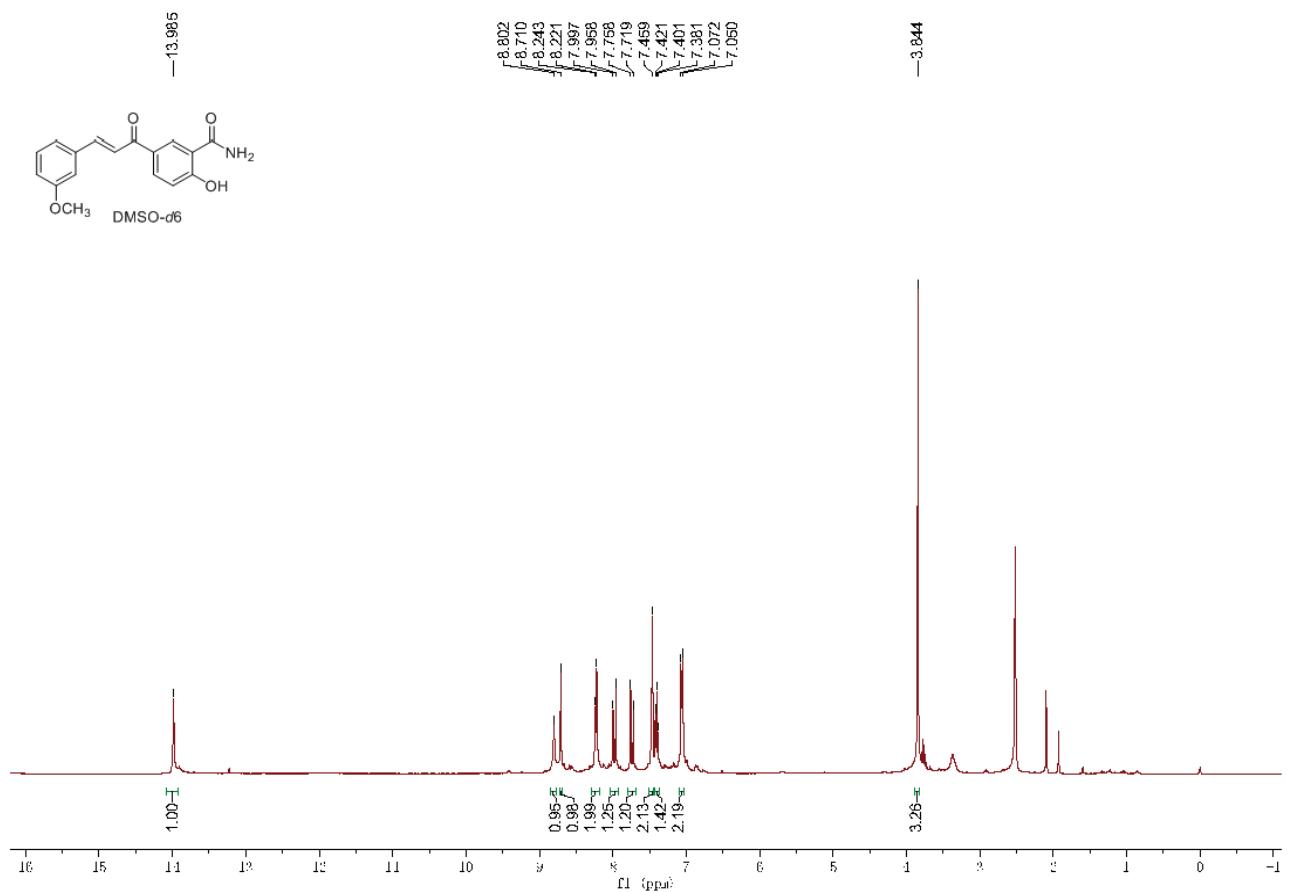
*(E)-2-hydroxy-5-(3-(*m*-tolyl)acryloyl)benzamide (a7)*



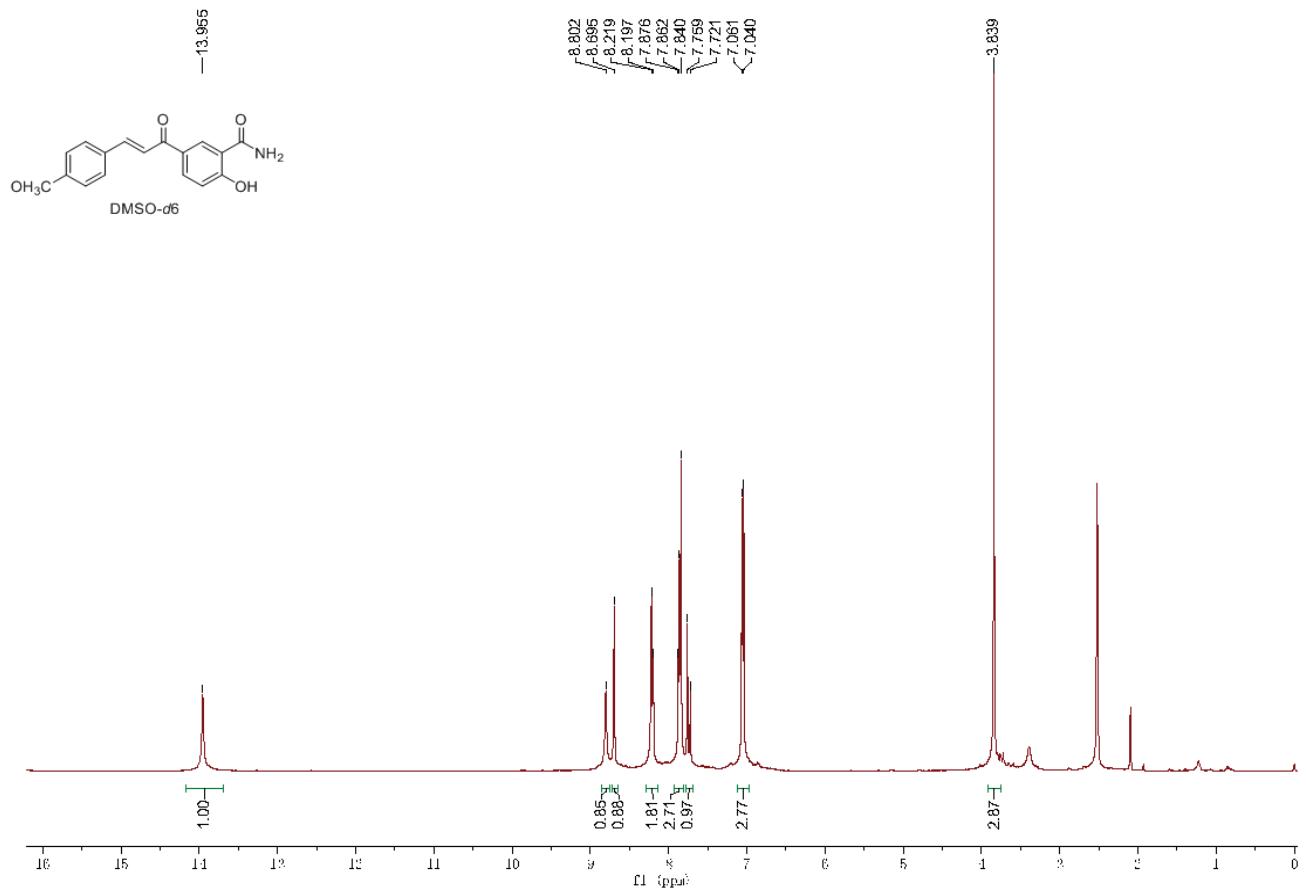
(*E*)-2-hydroxy-5-(3-(*p*-tolyl)acryloyl)benzamide (**a8**)



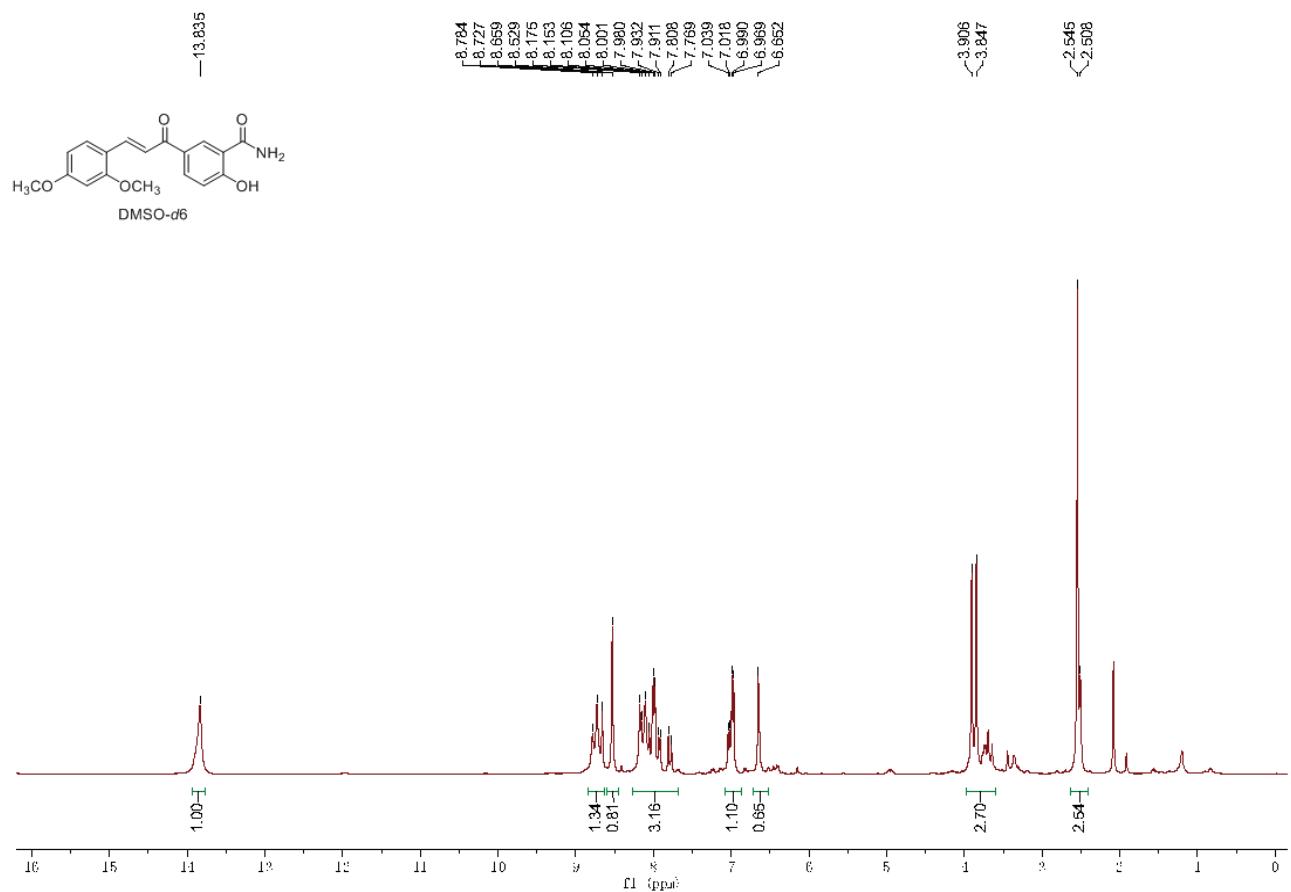
(E)-2-hydroxy-5-(3-(2-methoxyphenyl)acryloyl)benzamide (a9)



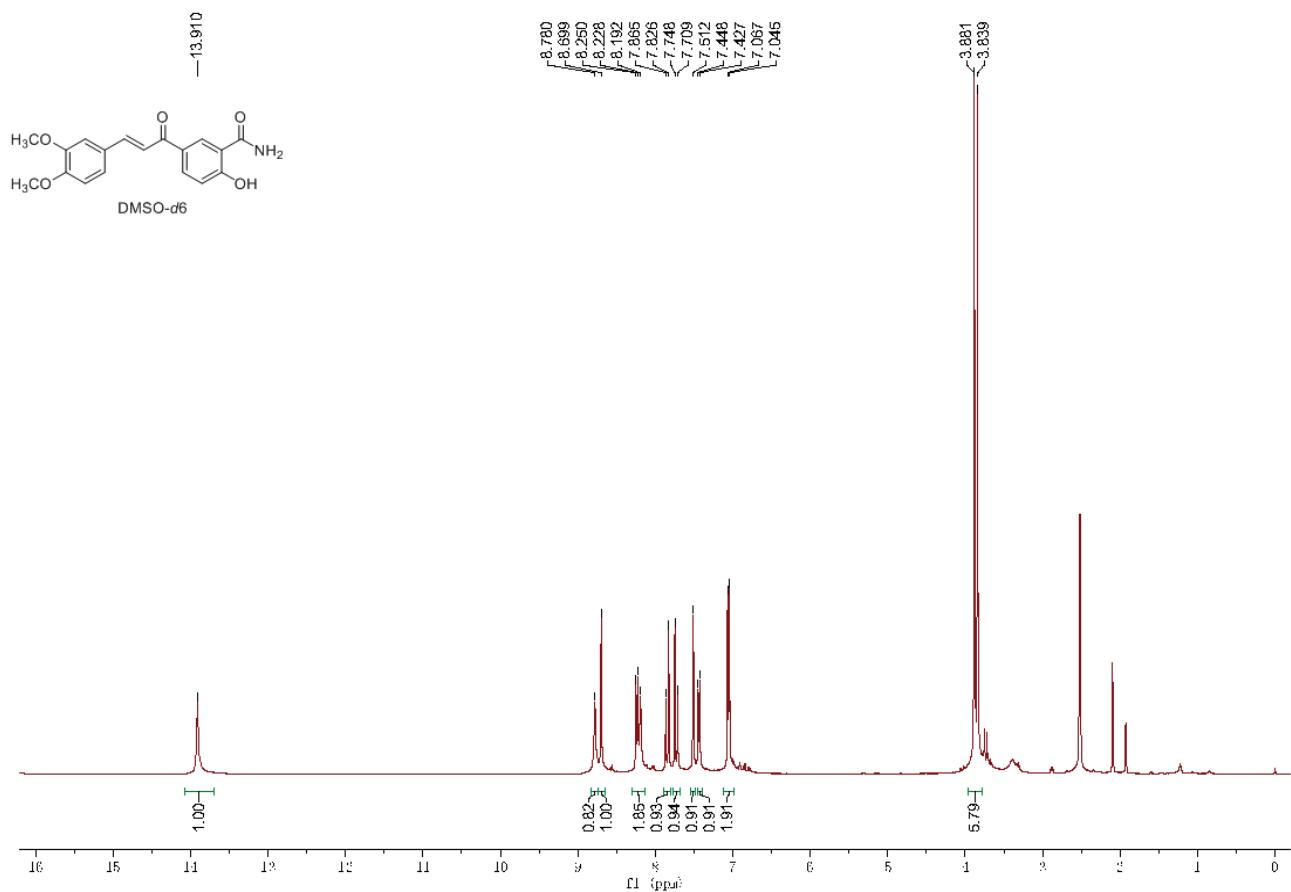
(E)-2-hydroxy-5-(3-(3-methoxyphenyl)acryloyl)benzamide (a10)



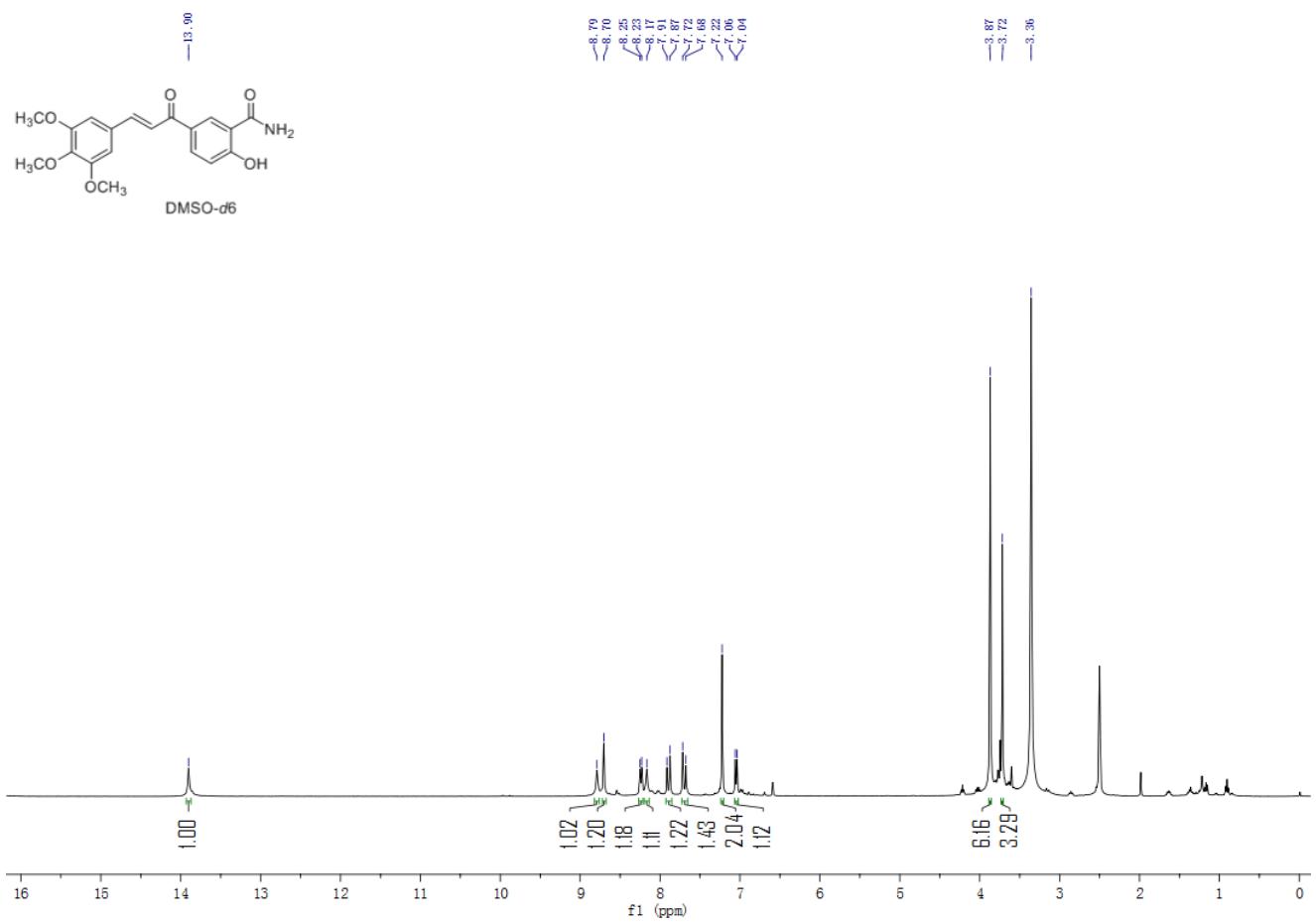
(E)-2-hydroxy-5-(3-(4-methoxyphenyl)acryloyl)benzamide (**a11**)



(E)-5-(3-(2,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (**a12**)

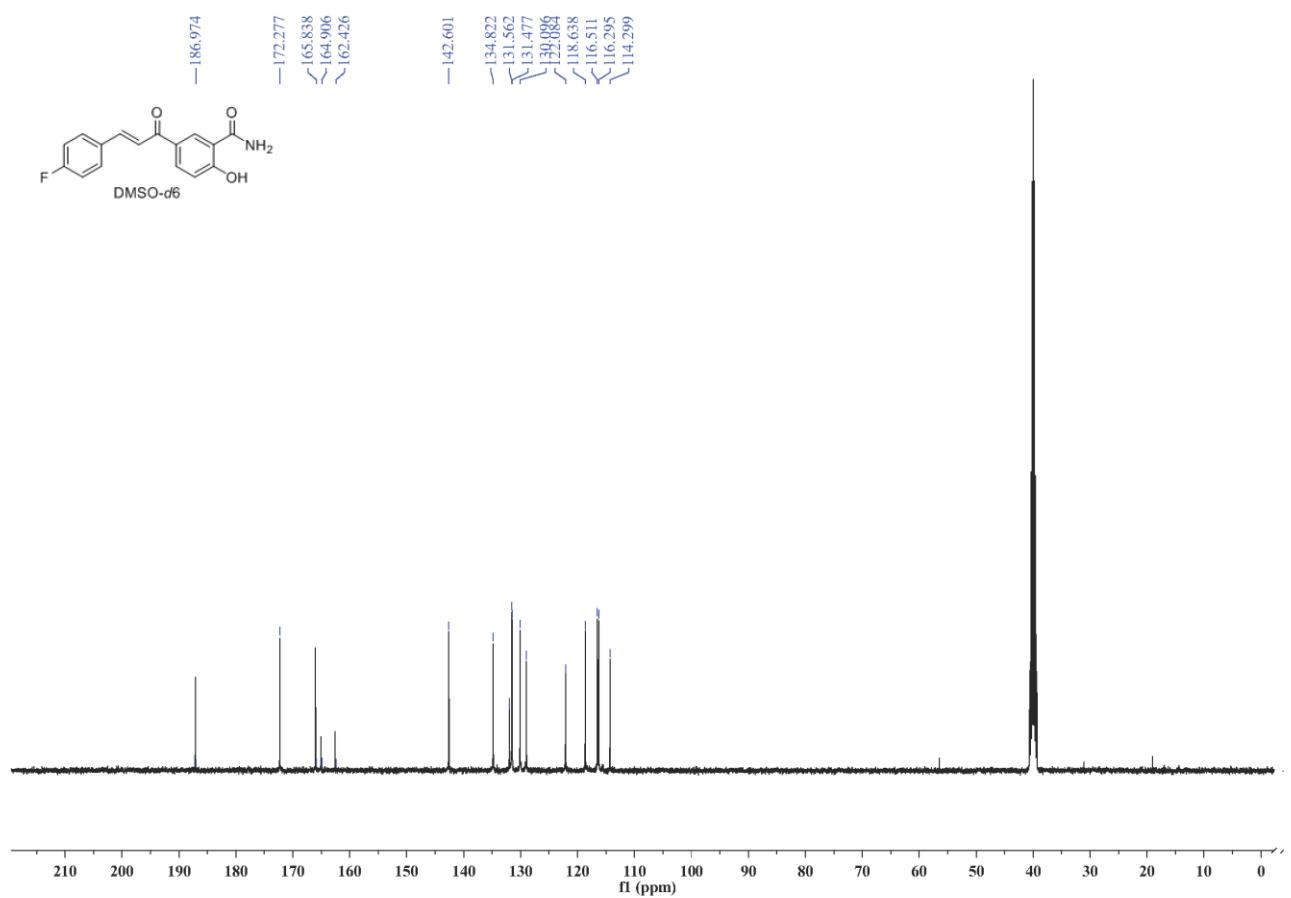


(E)-5-(3,4-dimethoxyphenyl)acryloyl-2-hydroxybenzamide (a13)

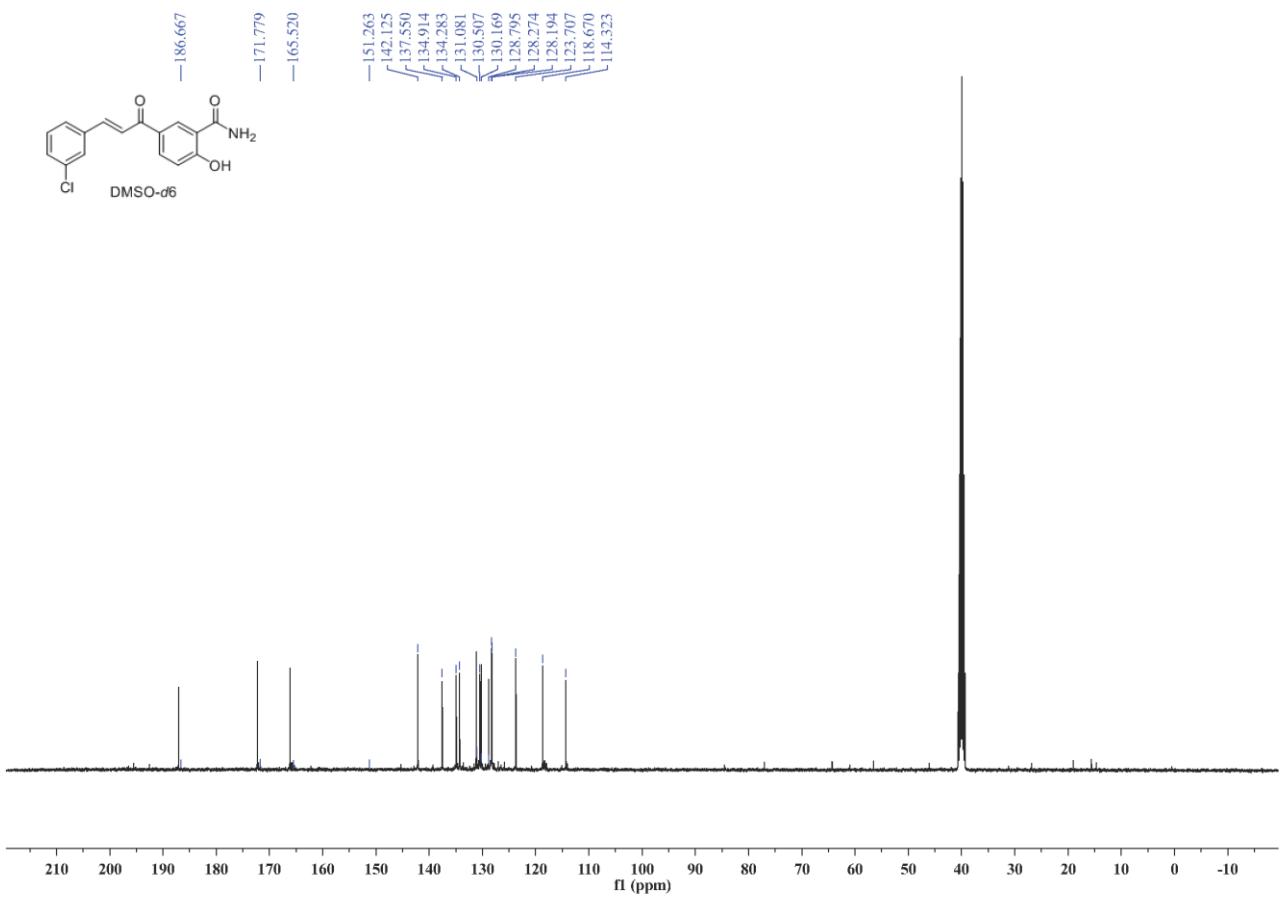


(E)-2-hydroxy-5-(3,4,5-trimethoxyphenyl)acryloylbenzamide (a14)

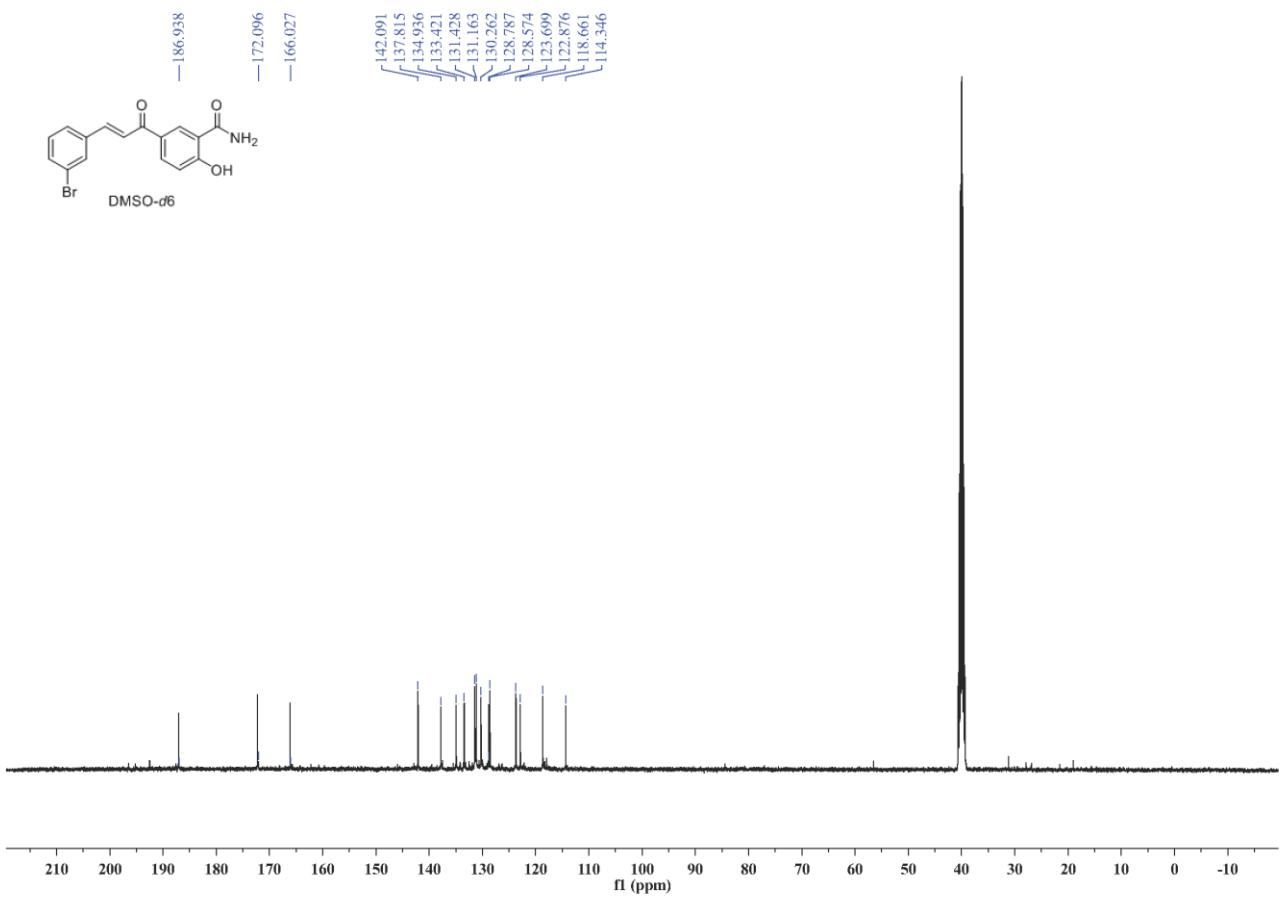
3. C-NMR information of the synthesized compounds



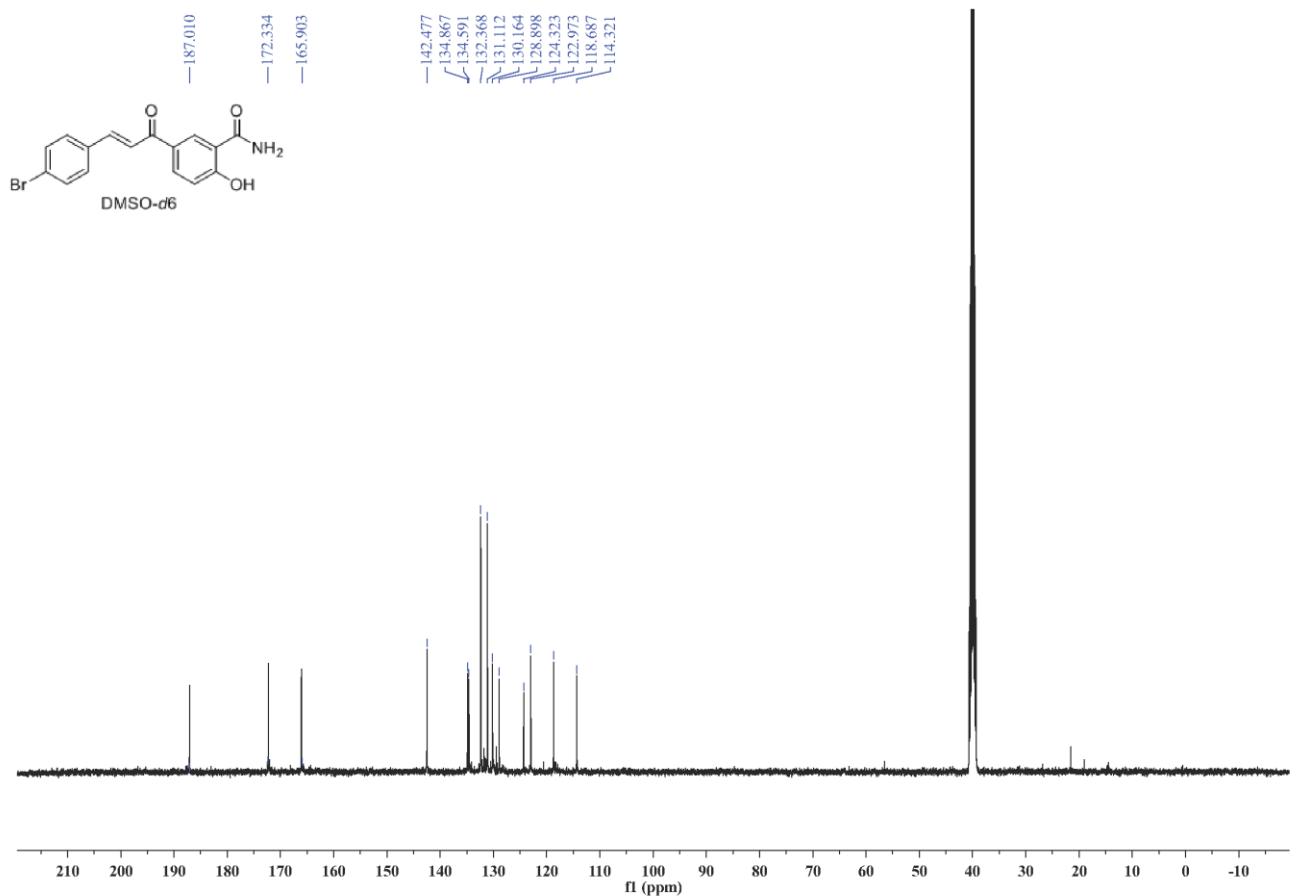
(*E*)-5-(3-(4-fluorophenyl)acryloyl)-2-hydroxybenzamide (**aI**)



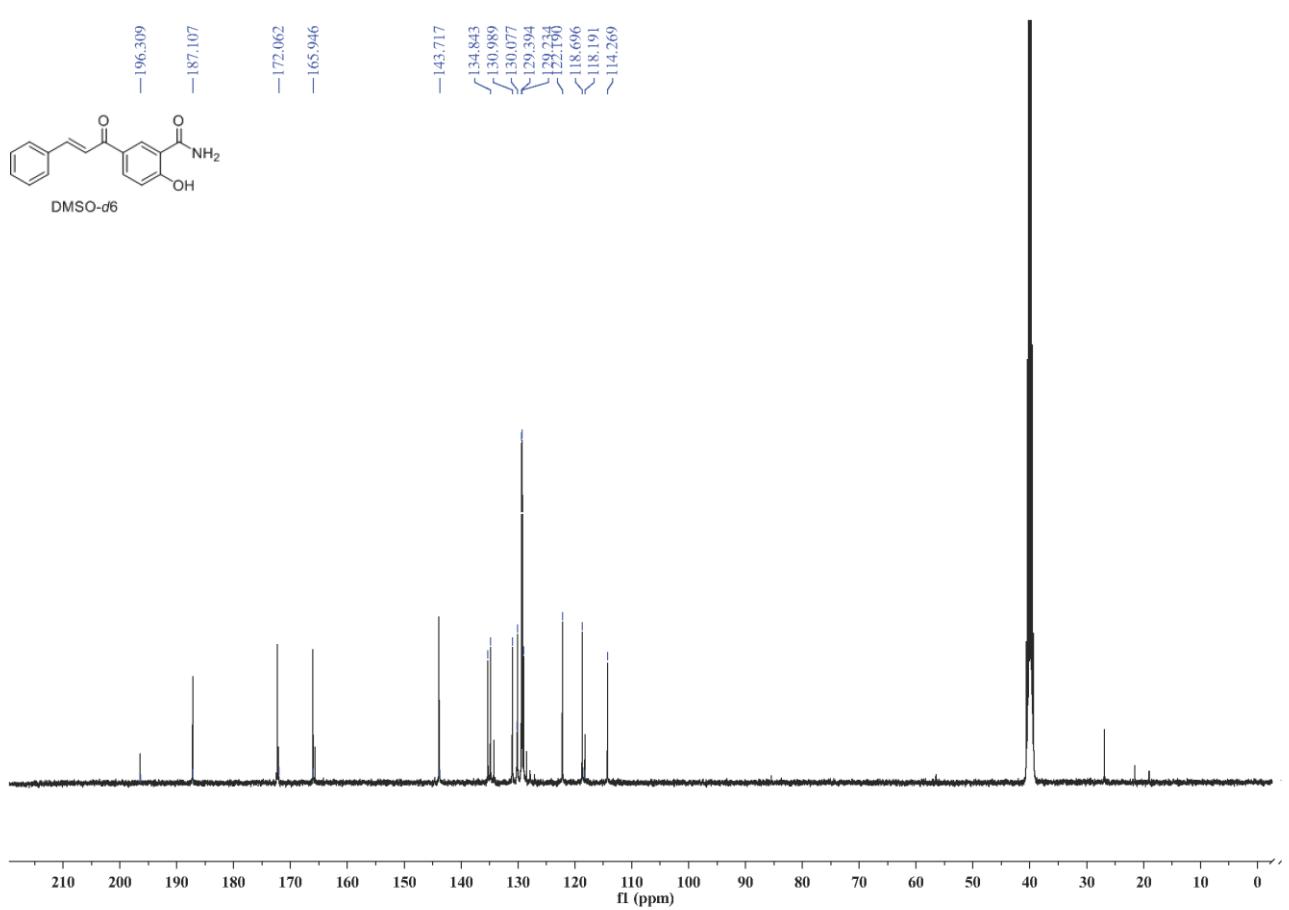
(E)-5-(3-(3-chlorophenyl)acryloyl)-2-hydroxybenzamide (a2)



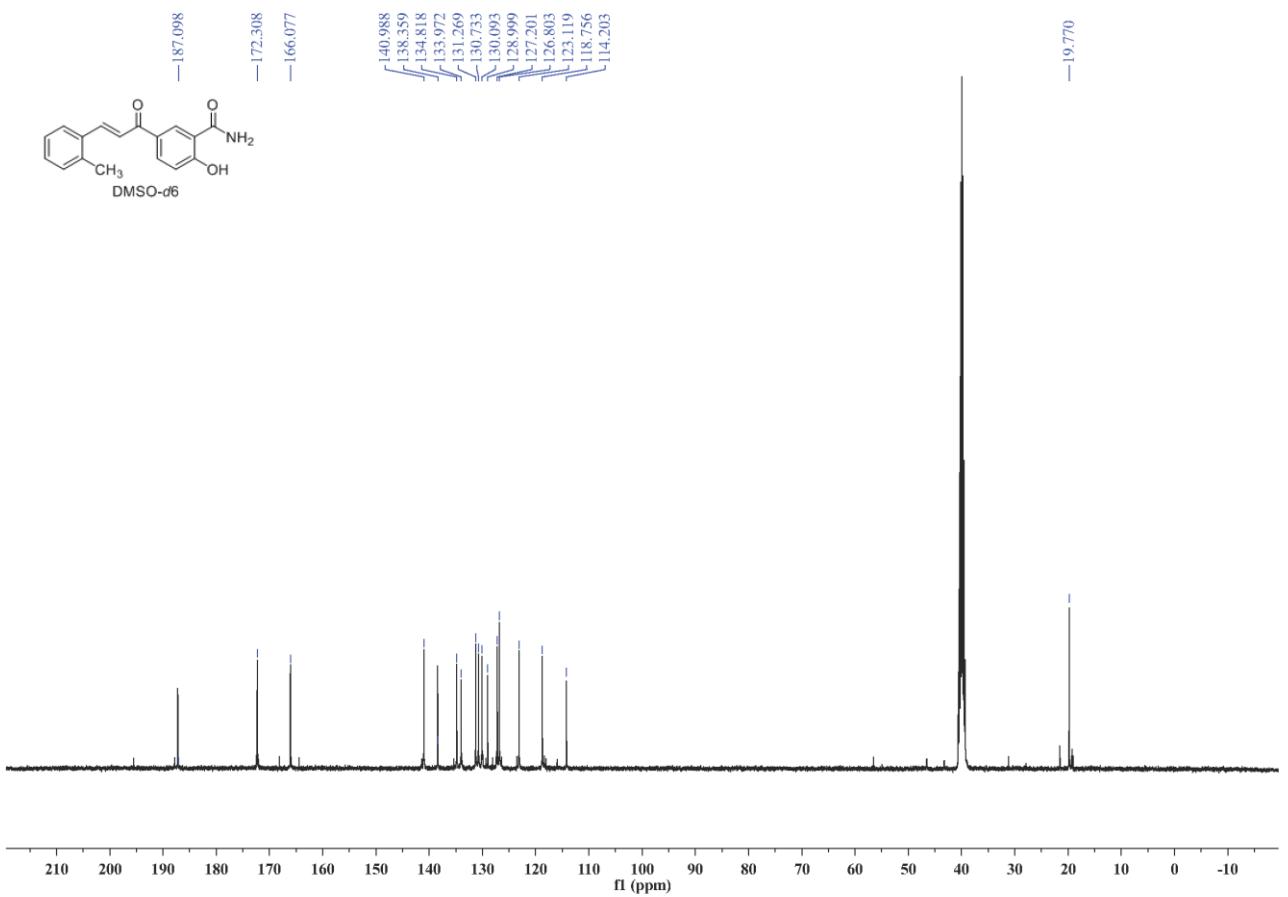
(E)-5-(3-(3-bromophenyl)acryloyl)-2-hydroxybenzamide (a3)



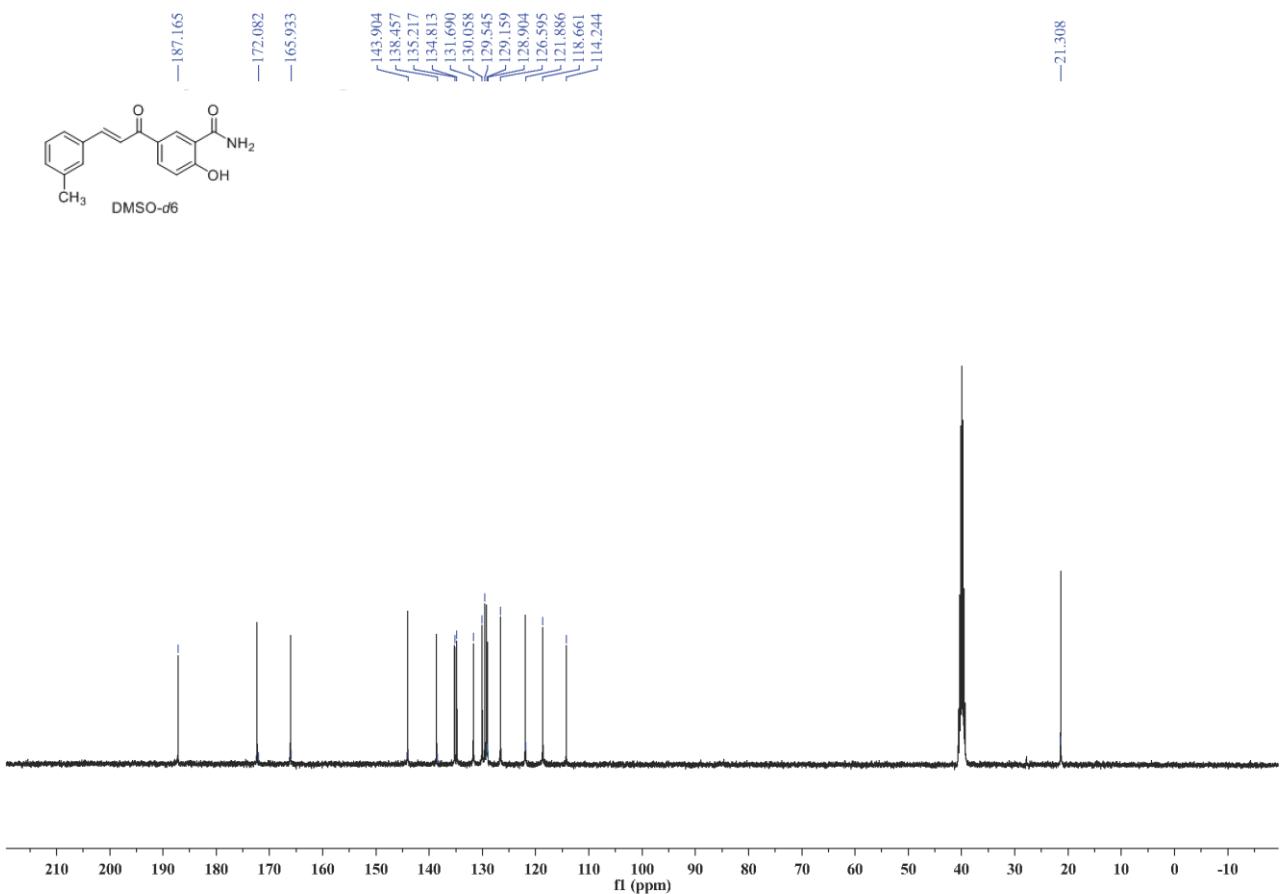
(E)-5-(3-(4-bromophenyl)acryloyl)-2-hydroxybenzamide (a4)



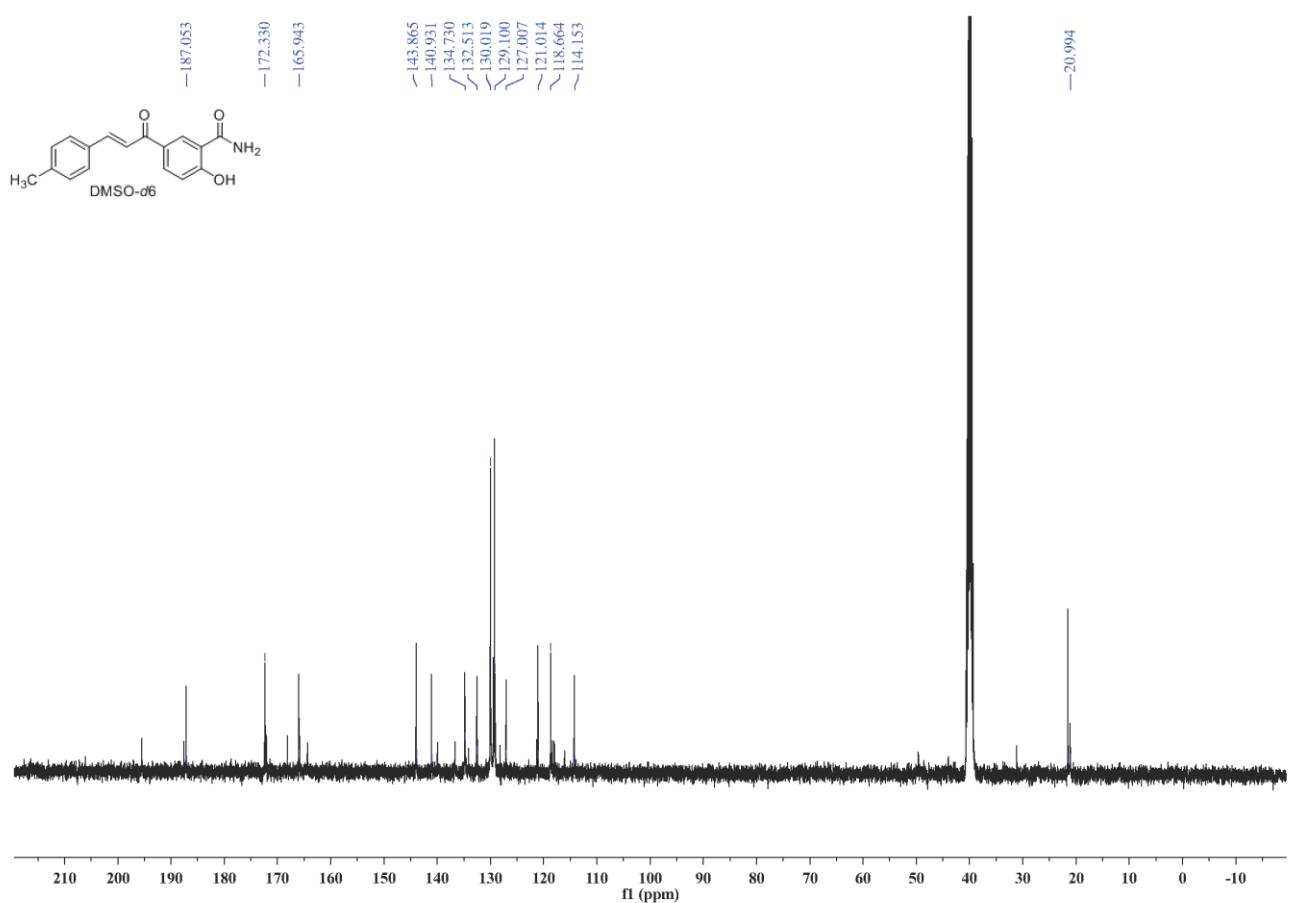
(E)-5-cinnamoyl-2-hydroxybenzamide (a5)



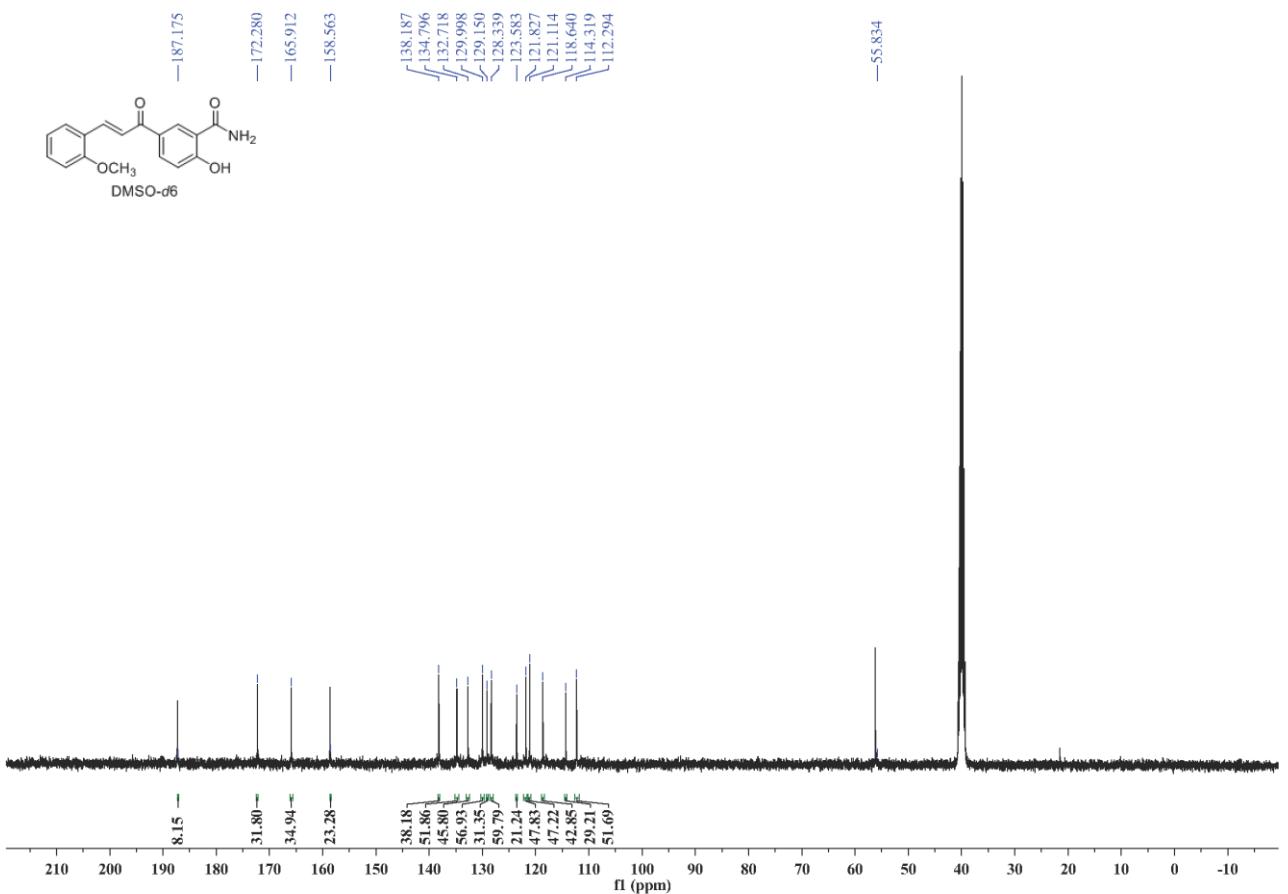
(E)-2-hydroxy-5-(3-(o-tolyl)acryloyl)benzamide(a6)



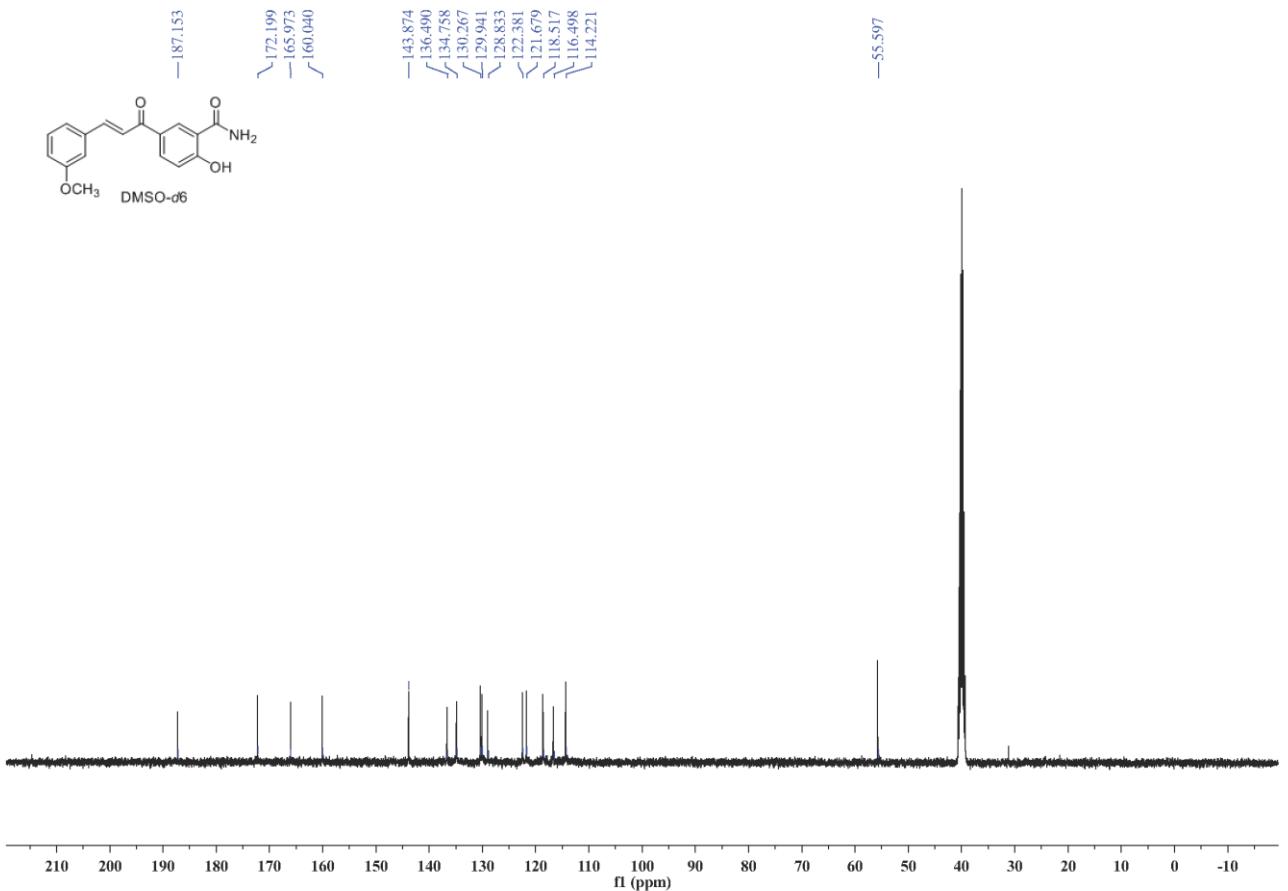
(E)-2-hydroxy-5-(3-(m-tolyl)acryloyl)benzamide (a7)



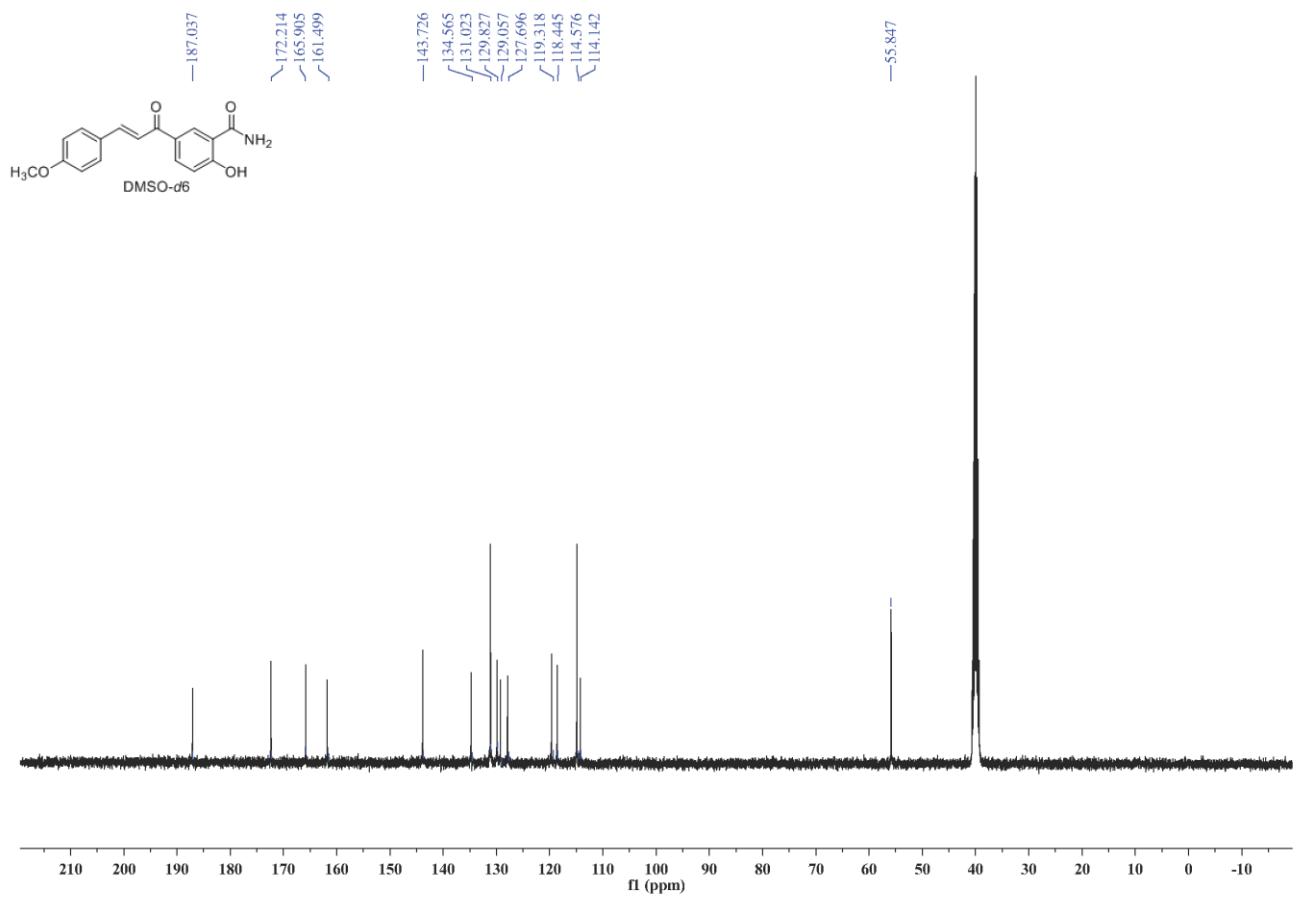
(E)-2-hydroxy-5-(3-(p-tolyl)acryloyl)benzamide (**a8**)



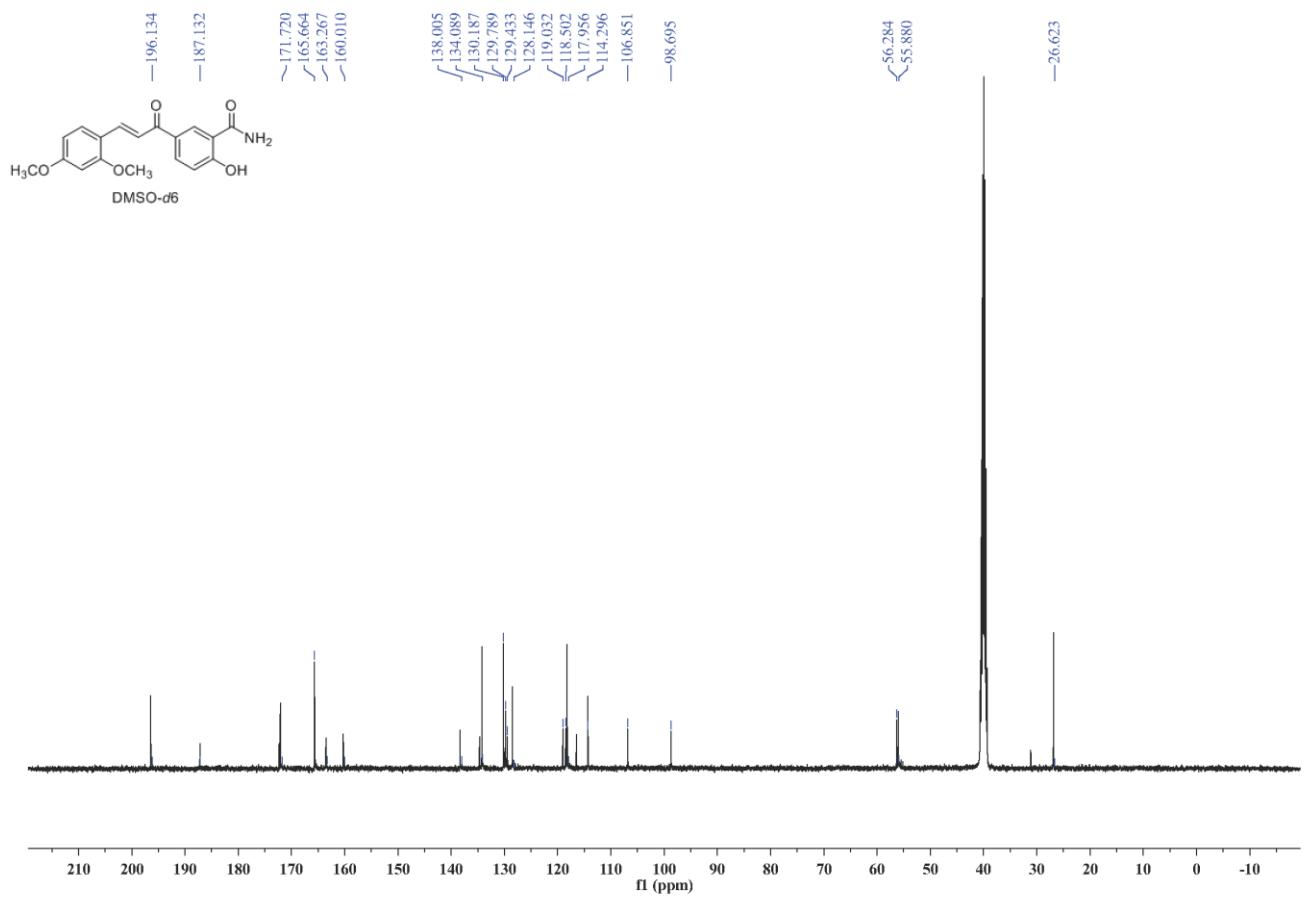
(E)-2-hydroxy-5-(3-(2-methoxyphenyl)acryloyl)benzamide(a9)



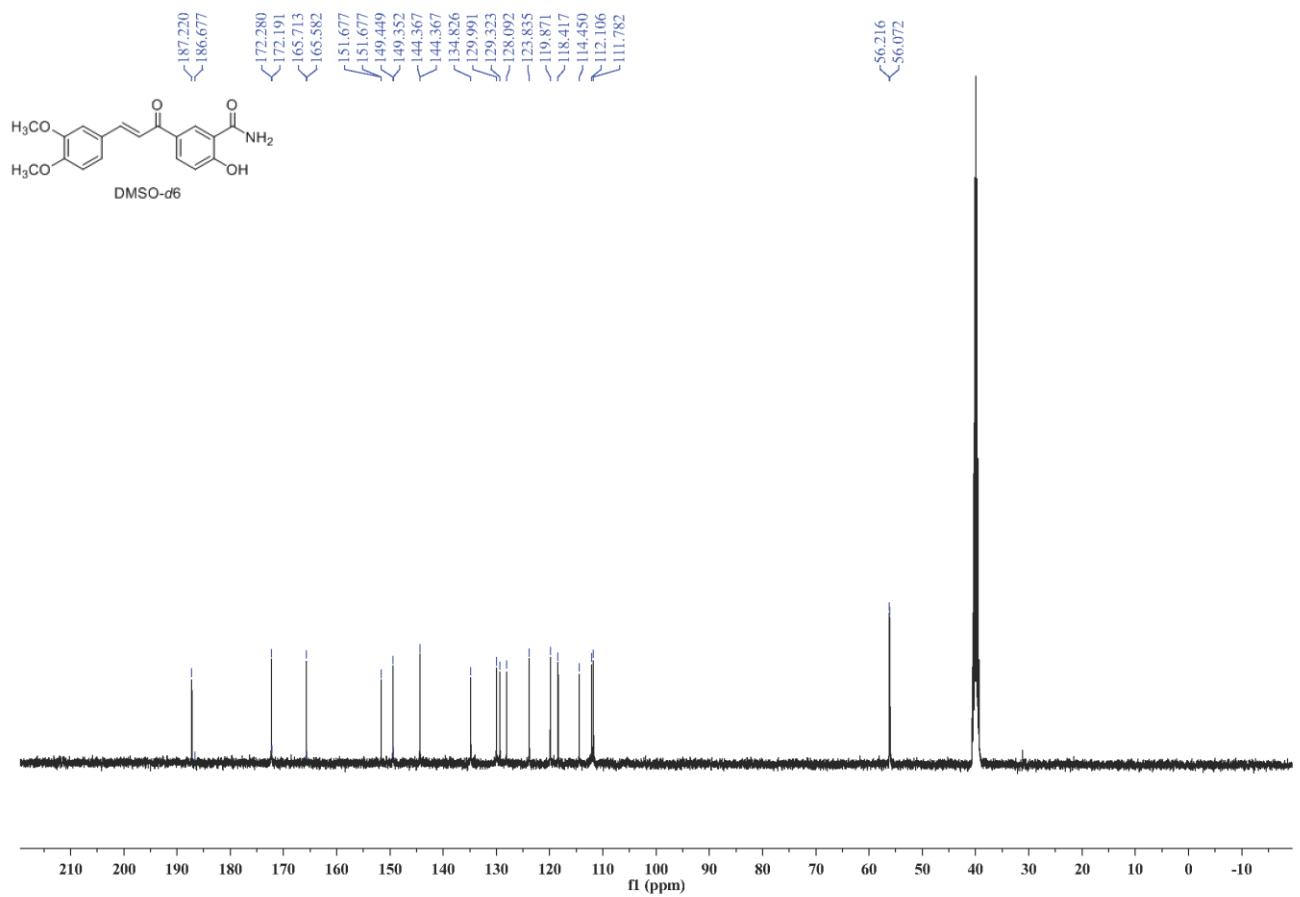
(E)-2-hydroxy-5-(3-(3-methoxyphenyl)acryloyl)benzamide (a10)



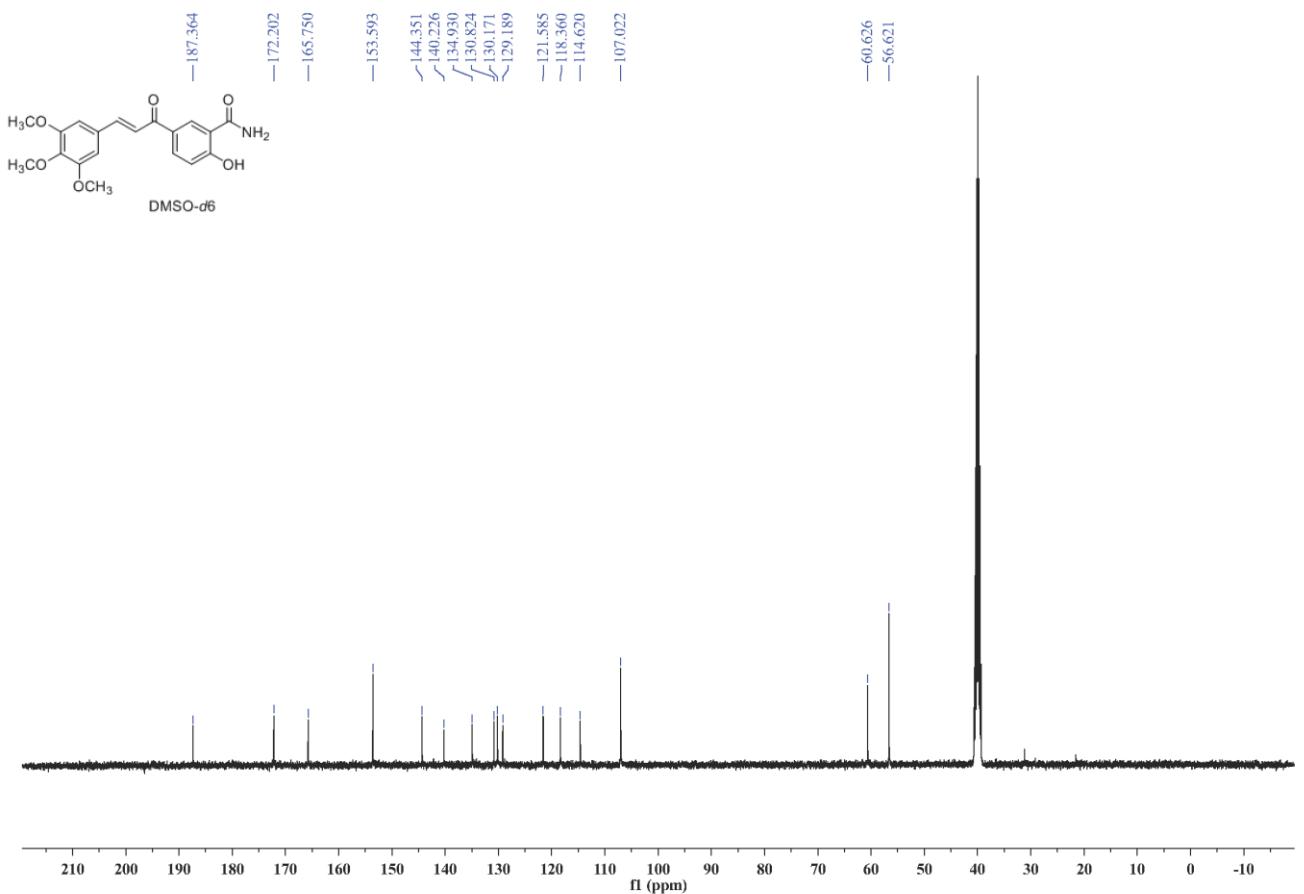
(E)-2-hydroxy-5-(3-(4-methoxyphenyl)acryloyl)benzamide (aII)



(E)-5-(3-(2,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (a12)

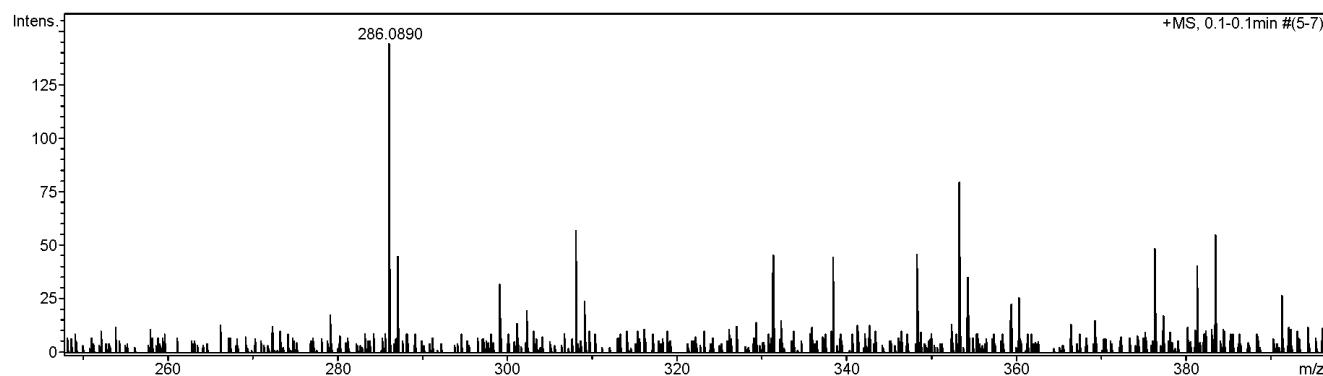
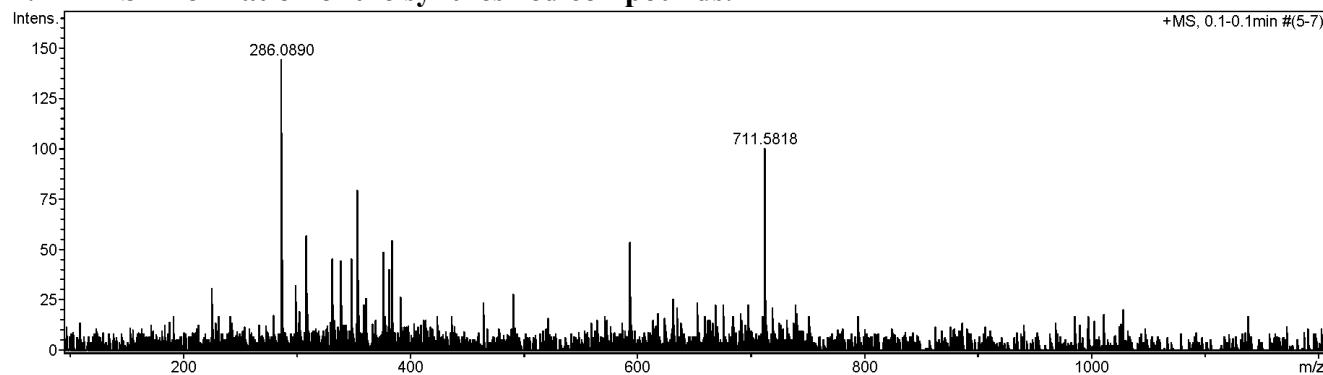


(*E*)-5-(3,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (**a13**)

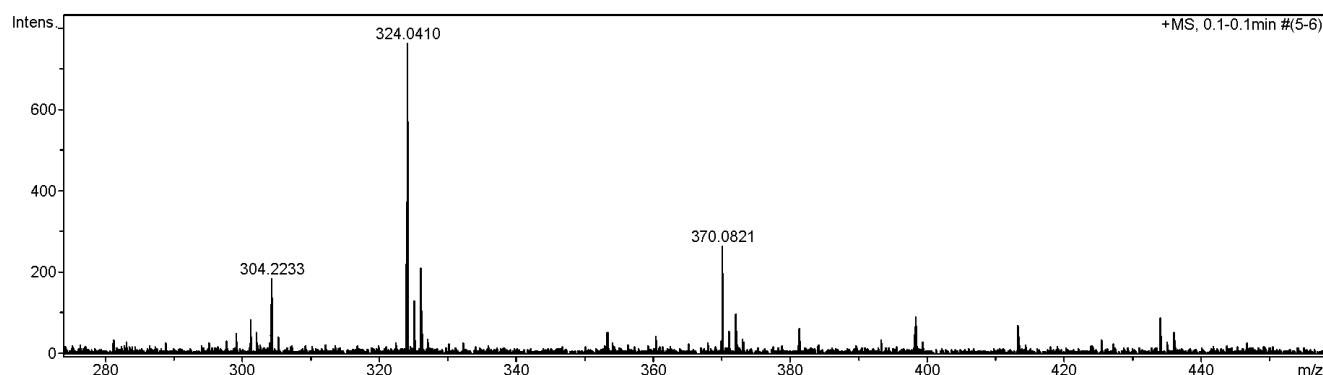
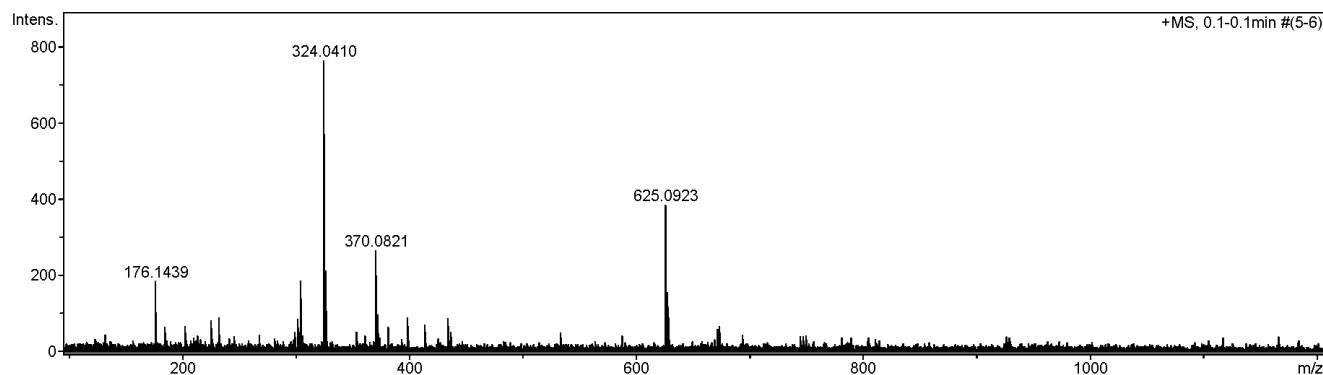


(E)-2-hydroxy-5-(3-(3,4,5-trimethoxyphenyl)acryloyl)benzamide (a14)

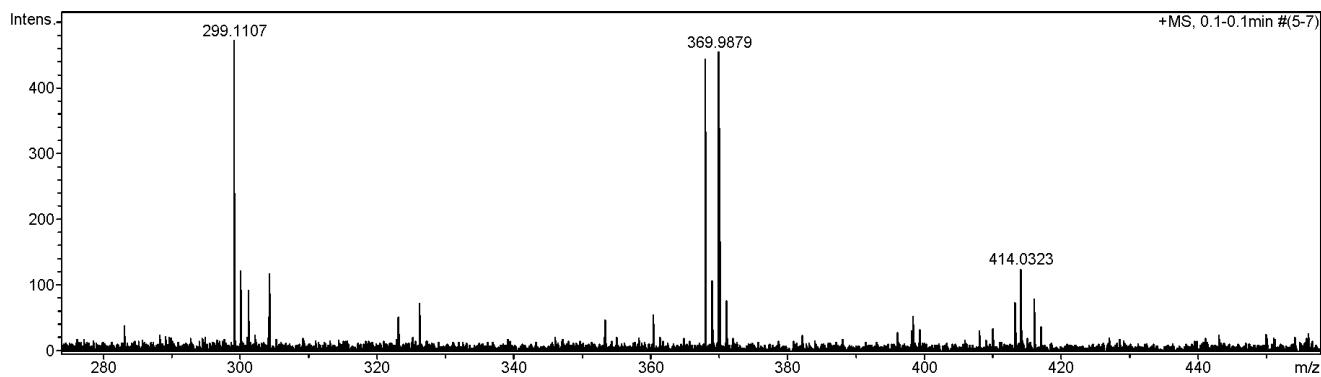
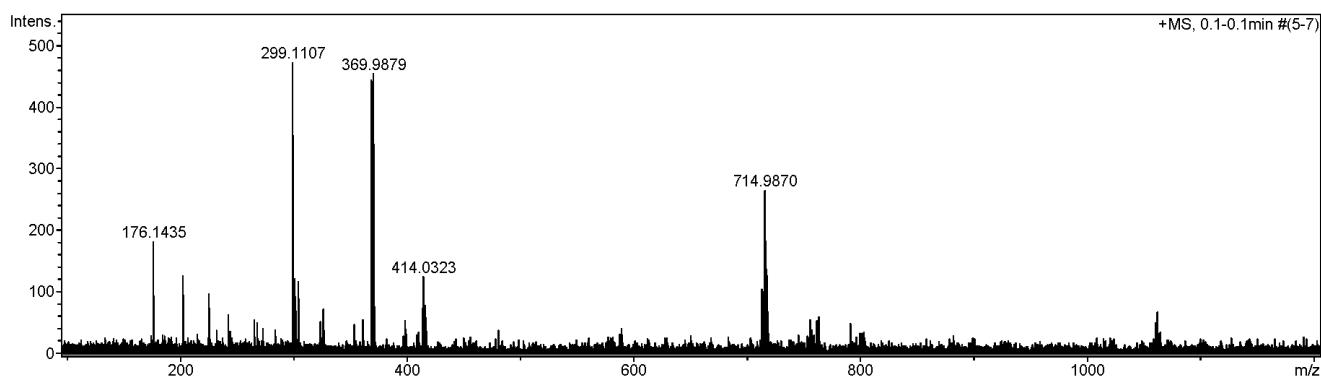
4. HRMS information of the synthesized compounds.



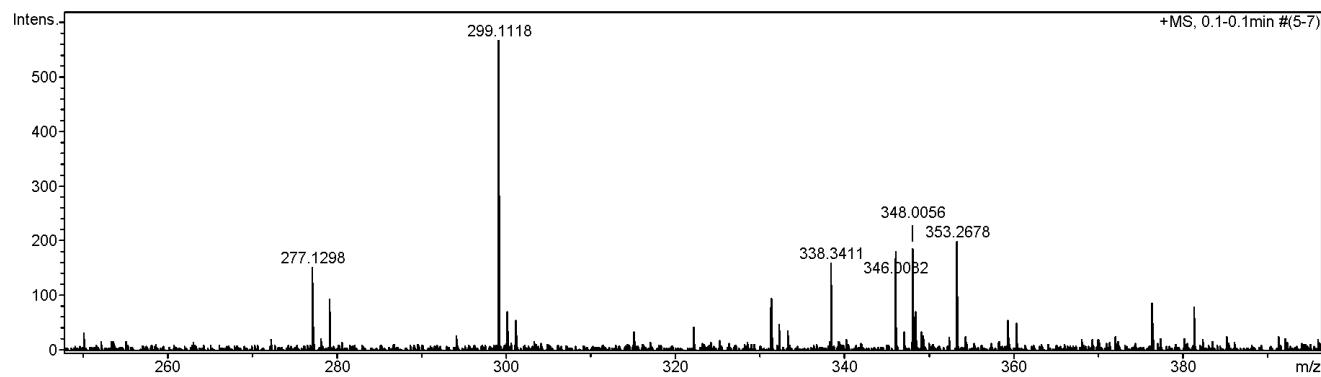
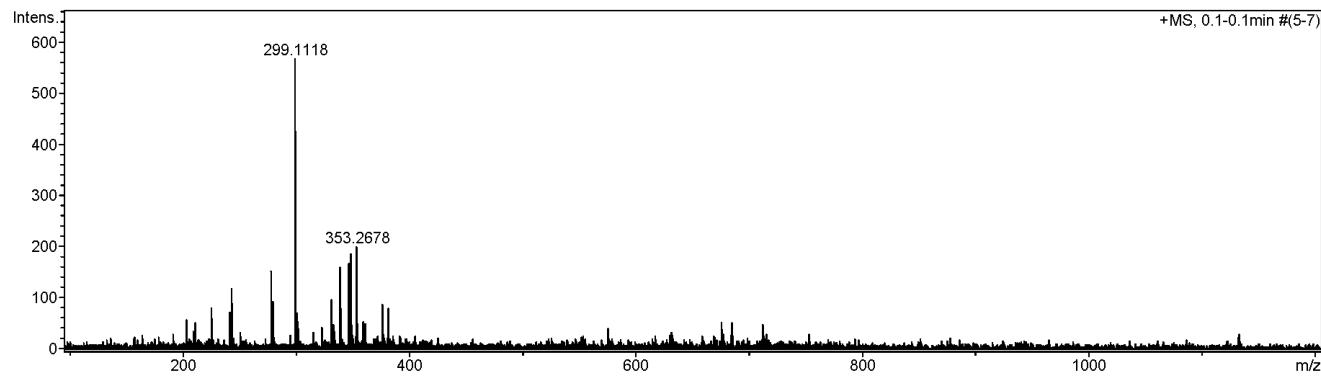
(E)-5-(3-(4-fluorophenyl)acryloyl)-2-hydroxybenzamide (a1)



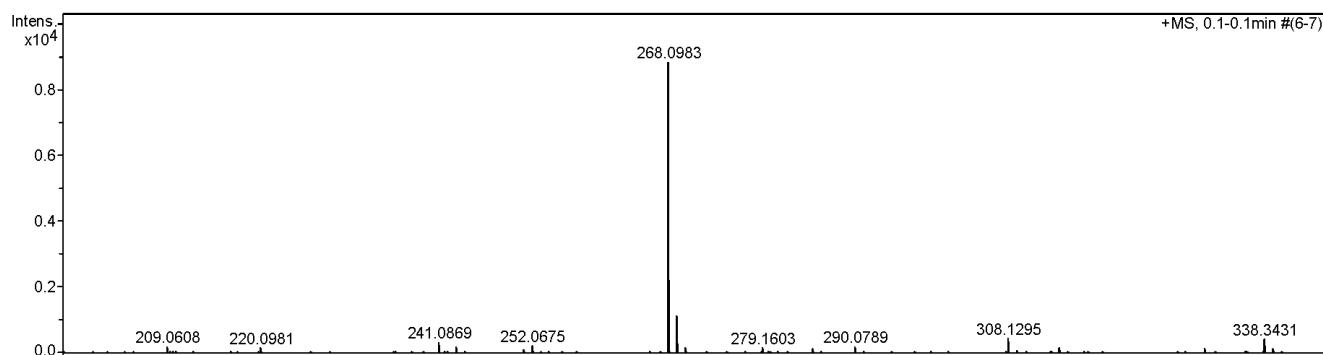
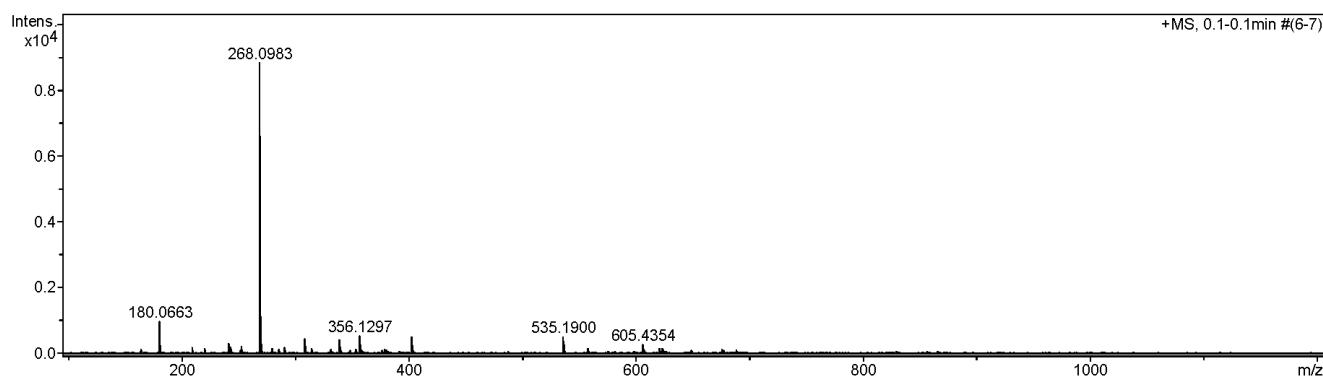
(E)-5-(3-(3-chlorophenyl)acryloyl)-2-hydroxybenzamide (a2)



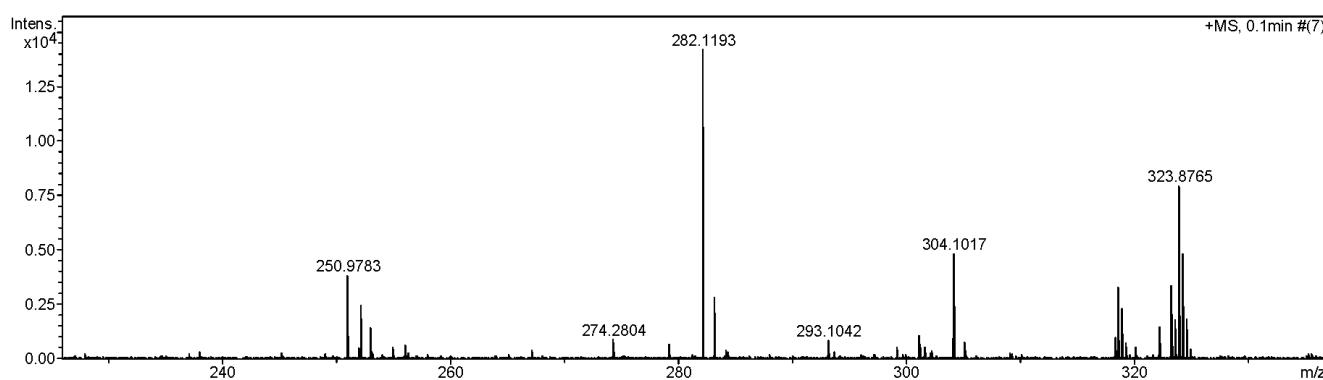
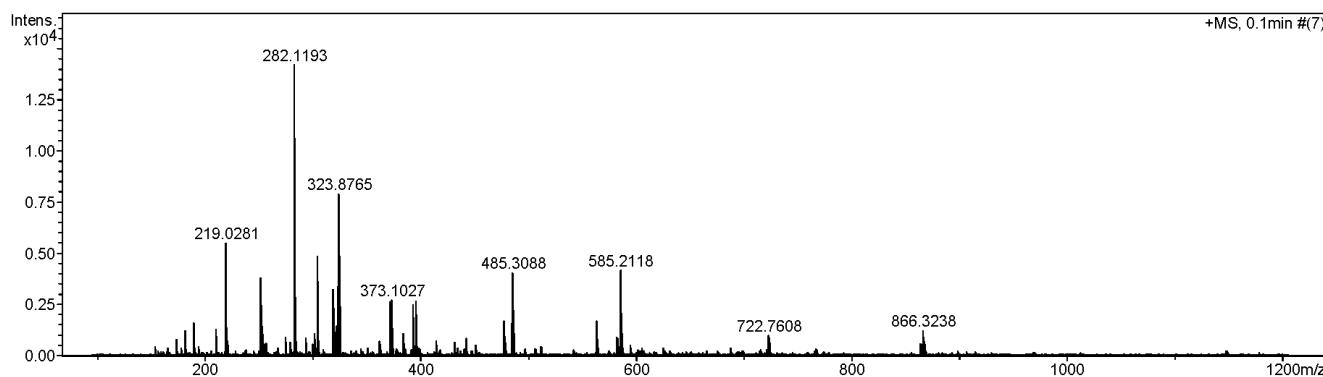
(E)-5-(3-(3-bromophenyl)acryloyl)-2-hydroxybenzamide (a3)



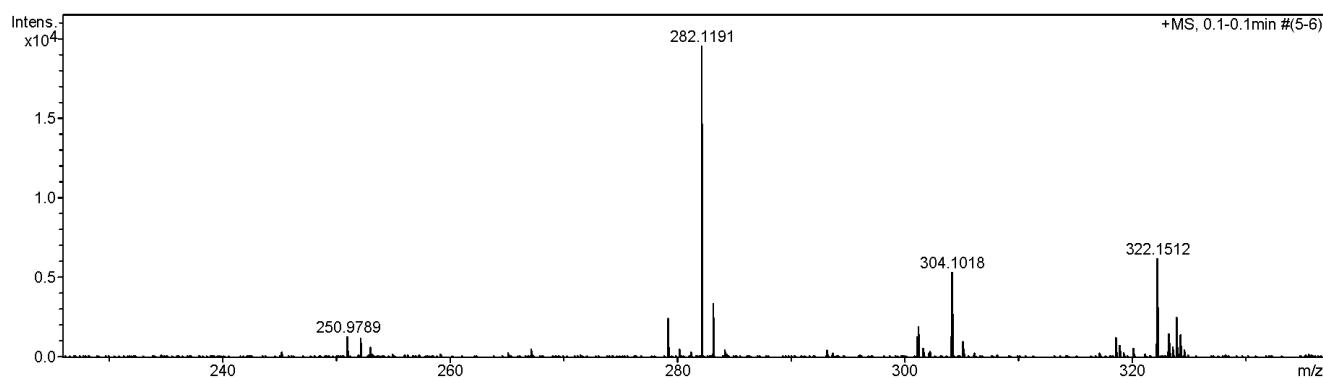
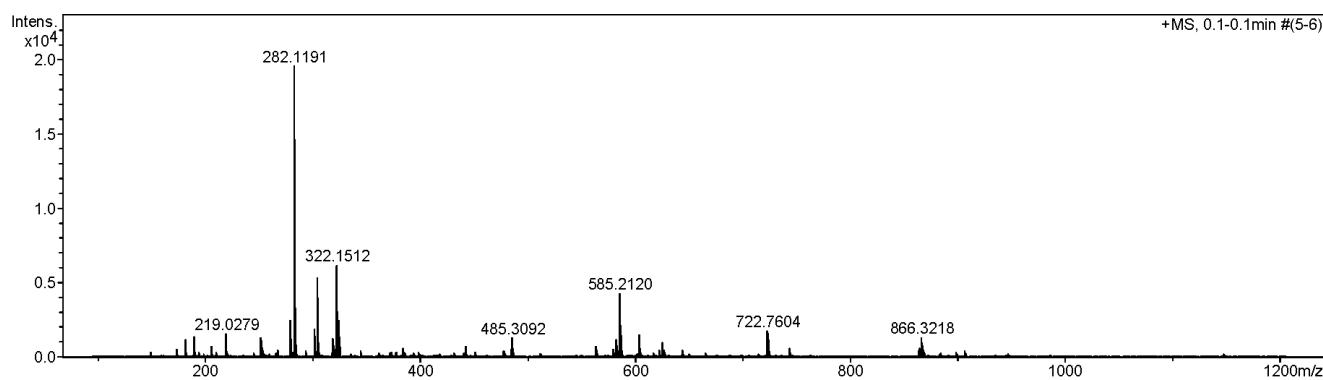
(E)-5-(3-(4-bromophenyl)acryloyl)-2-hydroxybenzamide (a4)



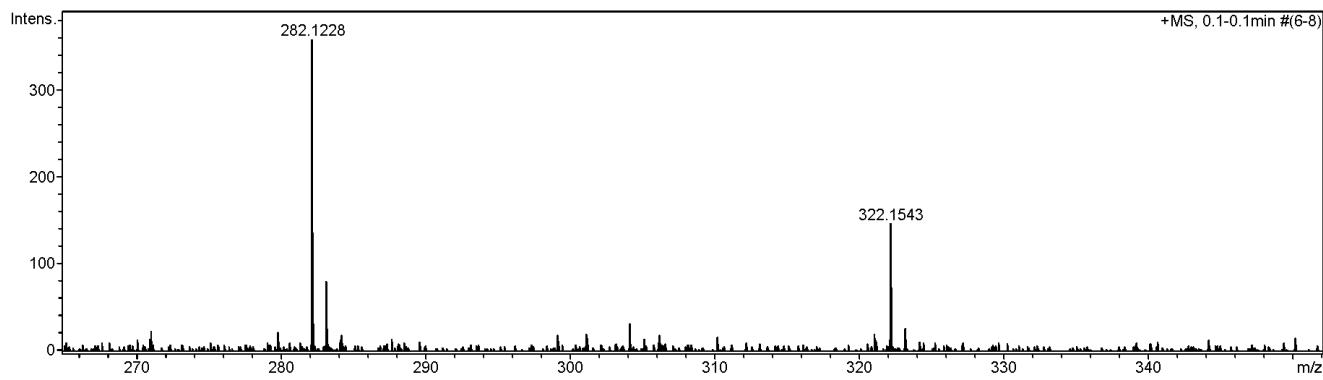
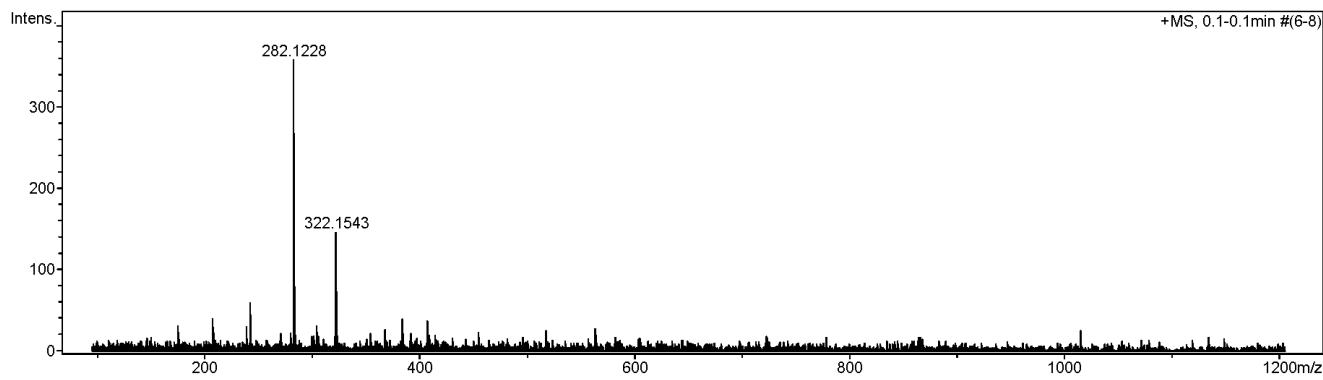
(E)-5-cinnamoyl-2-hydroxybenzamide (**a5**)



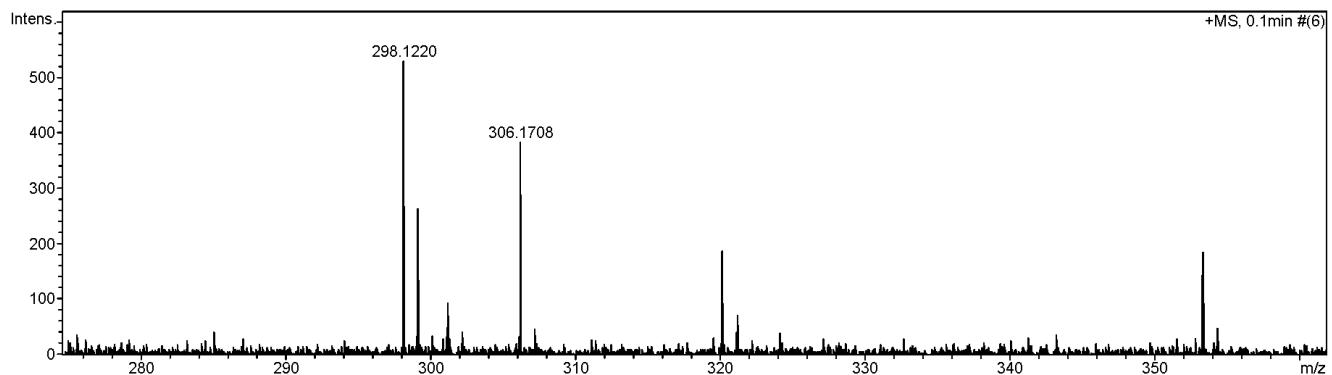
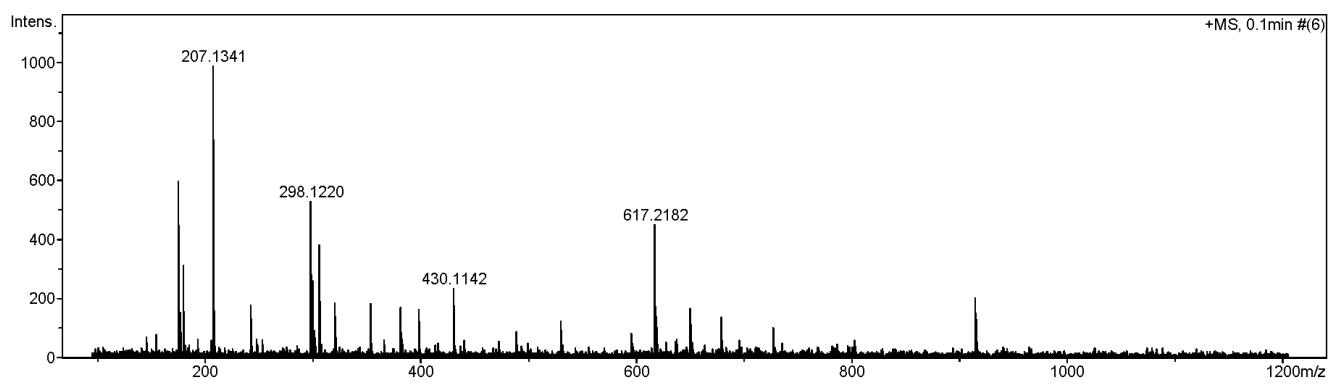
(E)-2-hydroxy-5-(3-(o-tolyl)acryloyl)benzamide (**a6**)



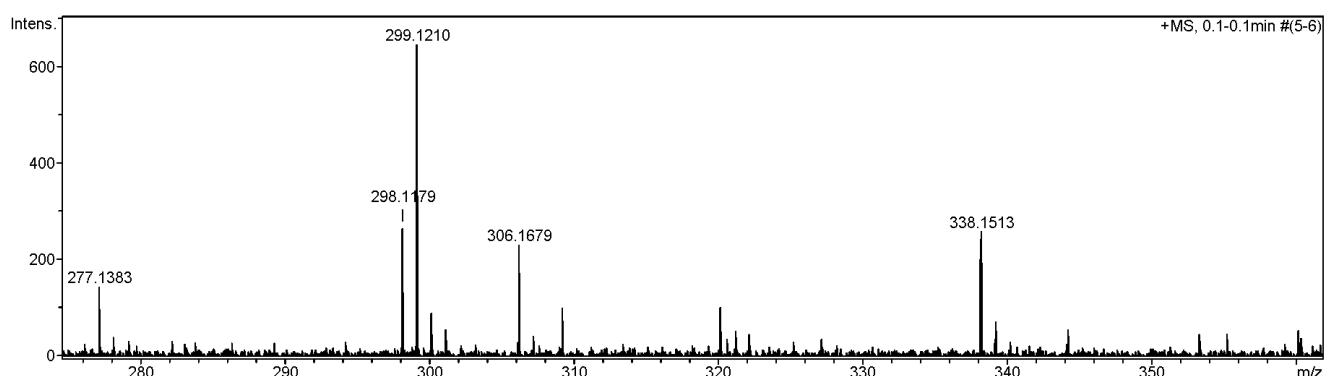
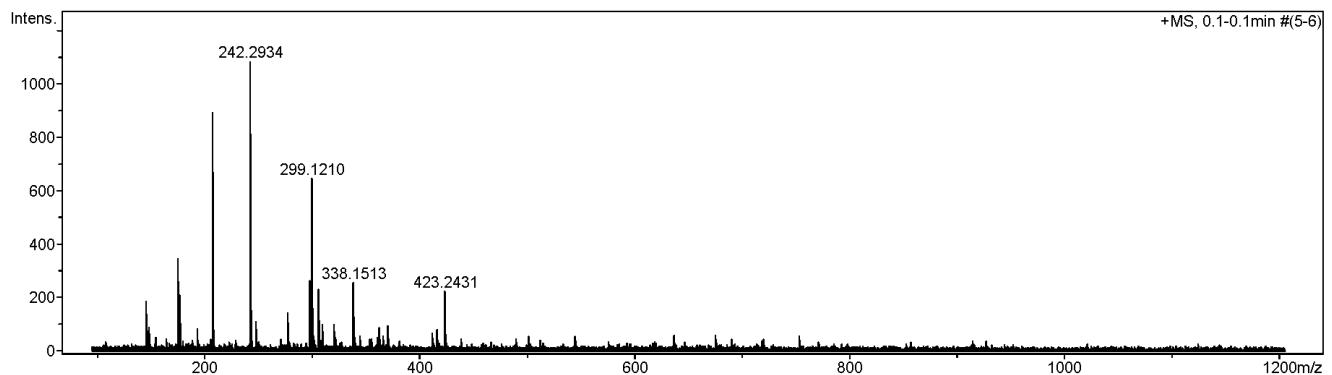
(E)-2-hydroxy-5-(3-(m-tolyl)acryloyl)benzamide (a7)



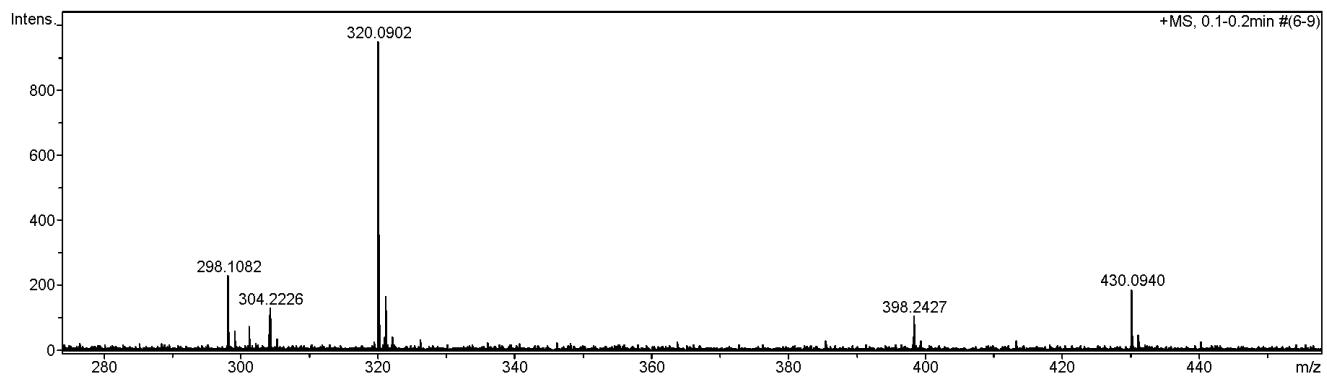
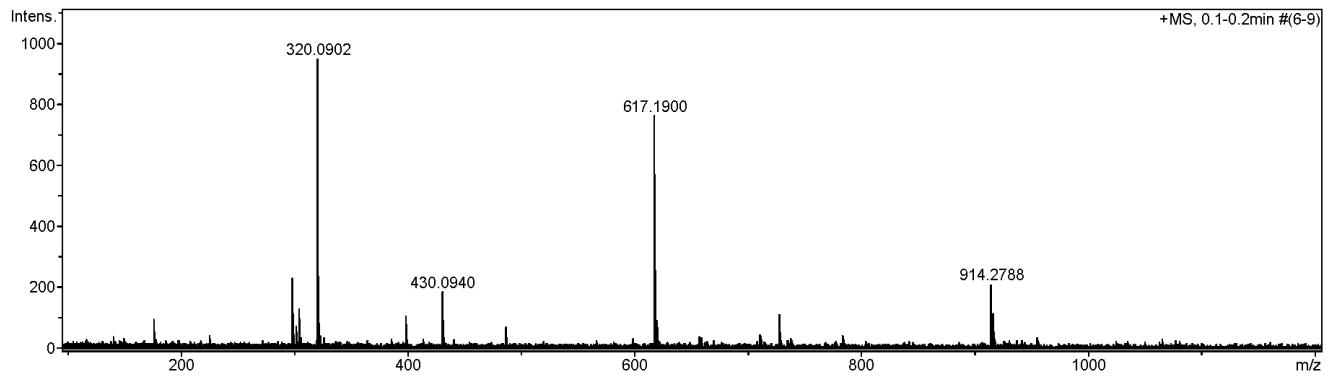
(E)-2-hydroxy-5-(3-(p-tolyl)acryloyl)benzamide (a8)



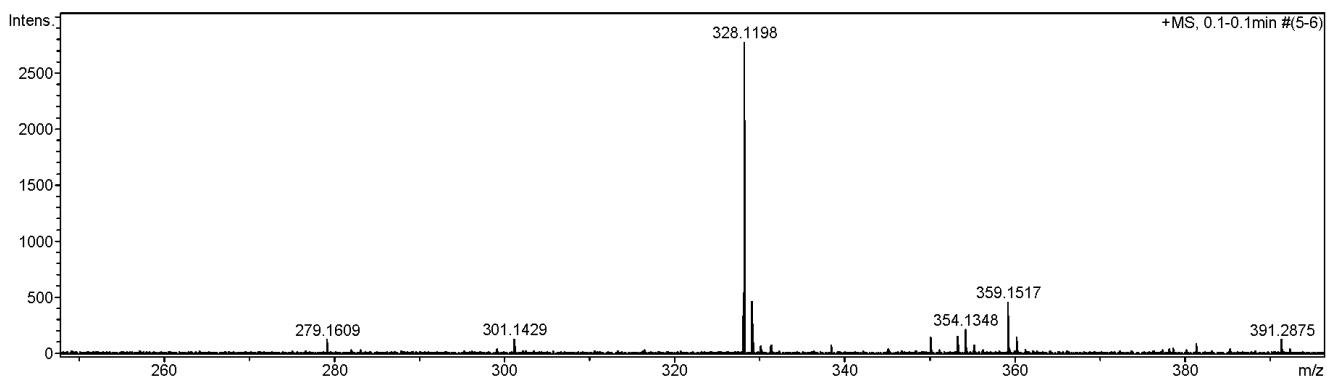
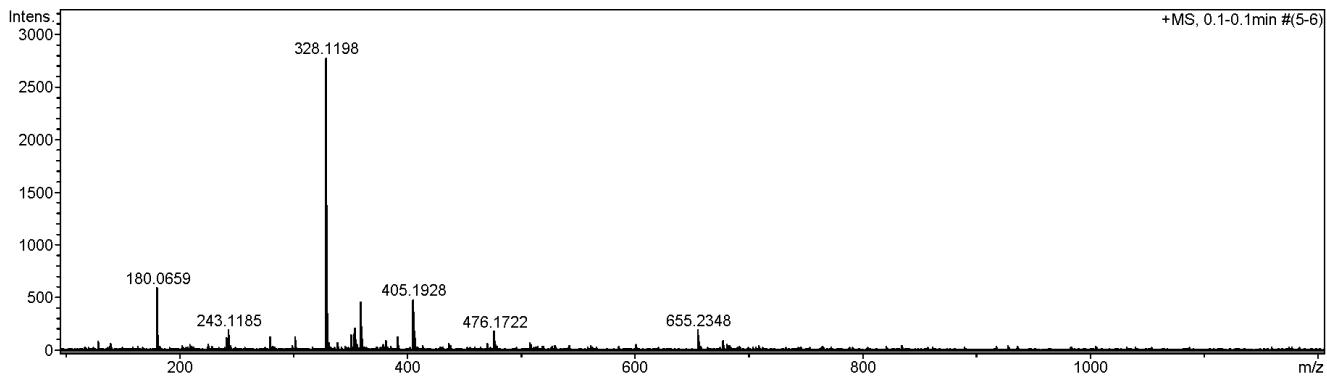
(E)-2-hydroxy-5-(3-(2-methoxyphenyl)acryloyl)benzamide (**a9**)



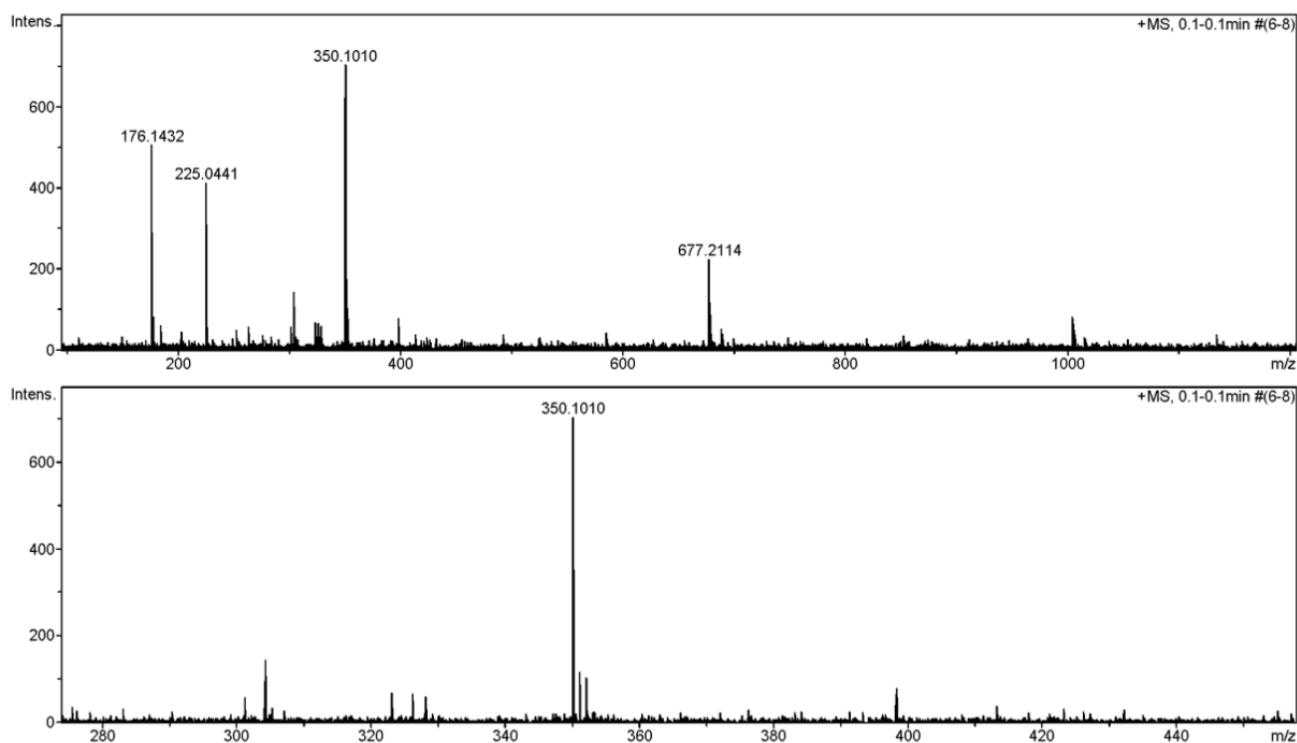
(E)-2-hydroxy-5-(3-(3-methoxyphenyl)acryloyl)benzamide(**a10**)



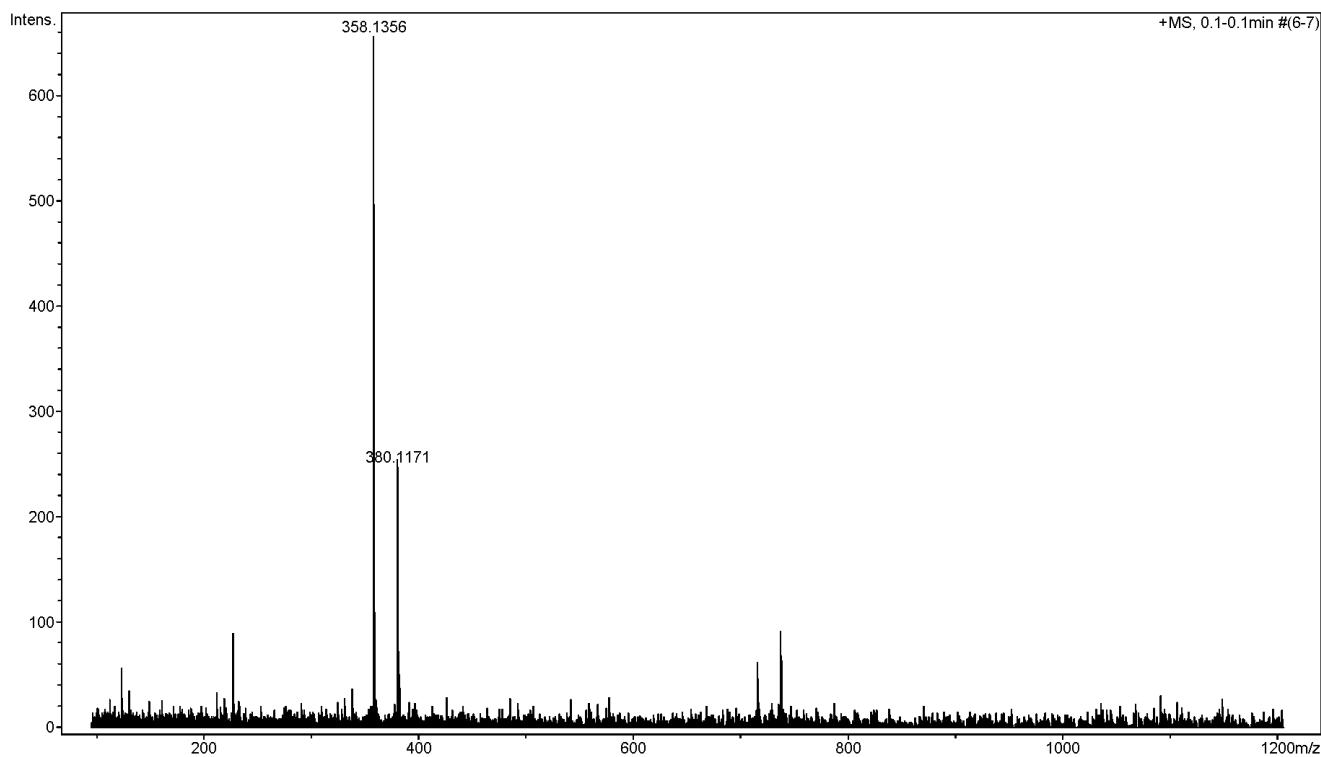
(E)-2-hydroxy-5-(3-(4-methoxyphenyl)acryloyl)benzamide (**a11**)



(E)-5-(3-(2,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (**a12**)

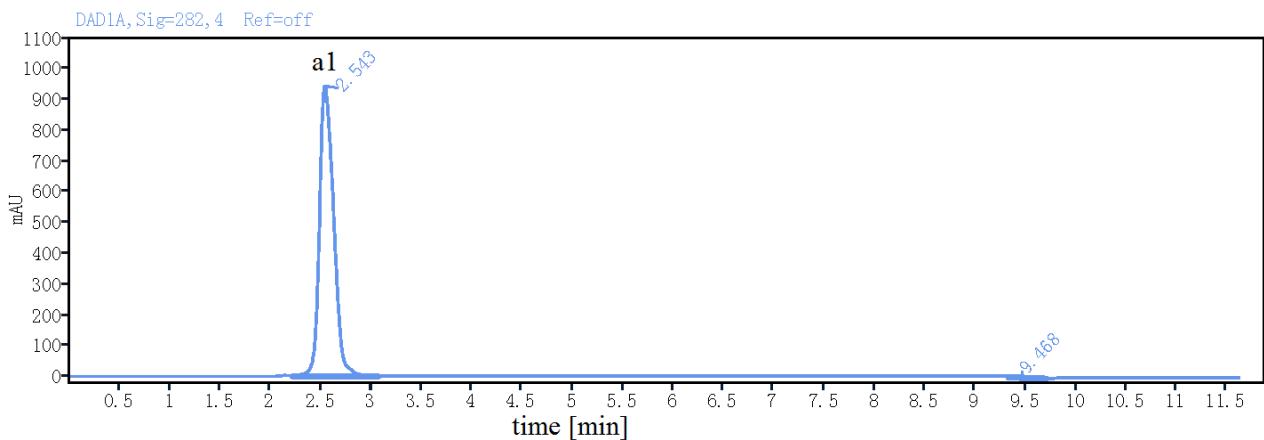


(E)-5-(3-(3,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (a13)



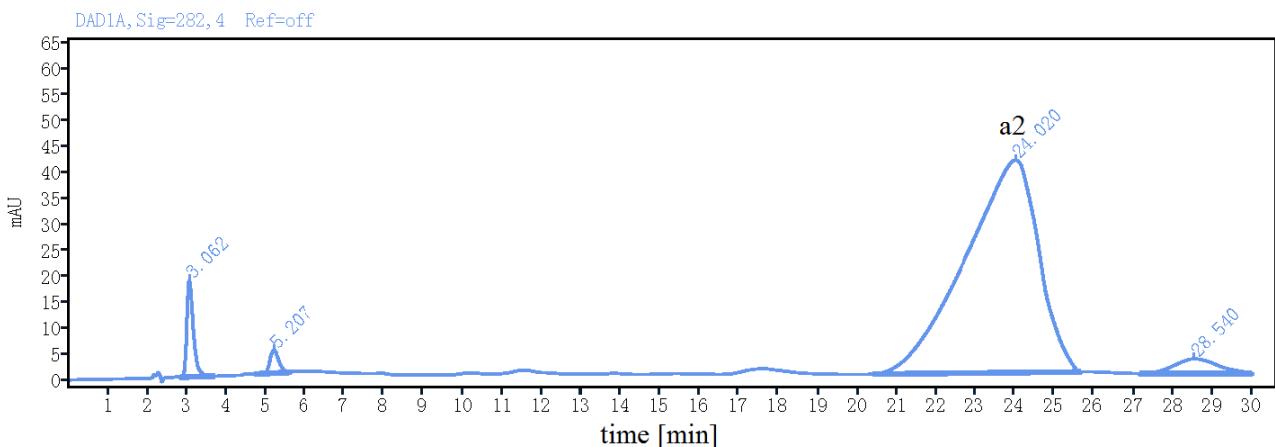
(F)-2-hydroxy-5-(3-(3,4,5-trimethoxyphenyl)acryloyl)benzamide (a14)

5. The purity information information of the synthesized compounds.



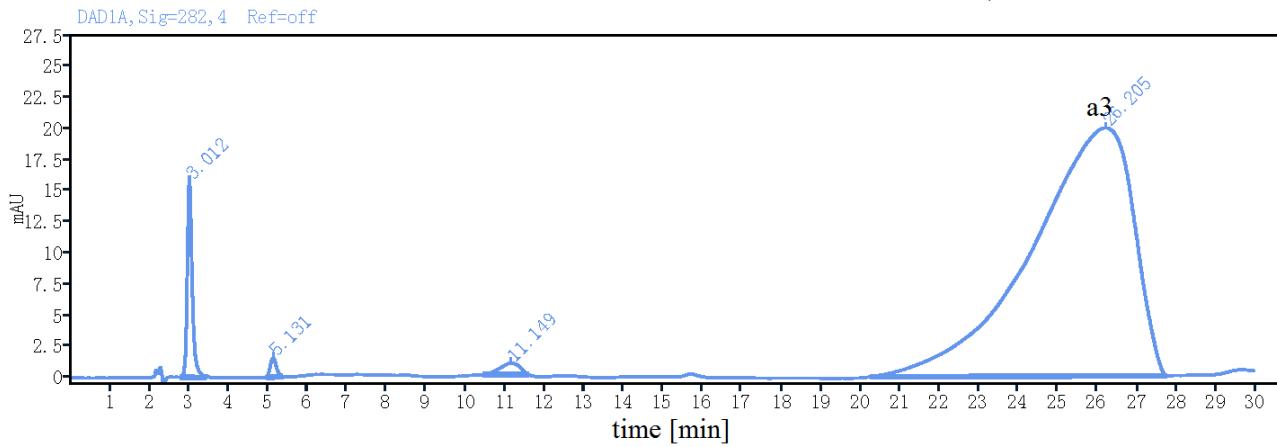
Retention time[min]	Area	Height	Area(%)
2. 543	9283.22	943.09	98. 97
9. 468	96.94	5.76	1. 03
			9380.17

(F)-5-(3-(4-fluorophenyl)acryloyl)-2-hydroxybenzamide (**a1**)



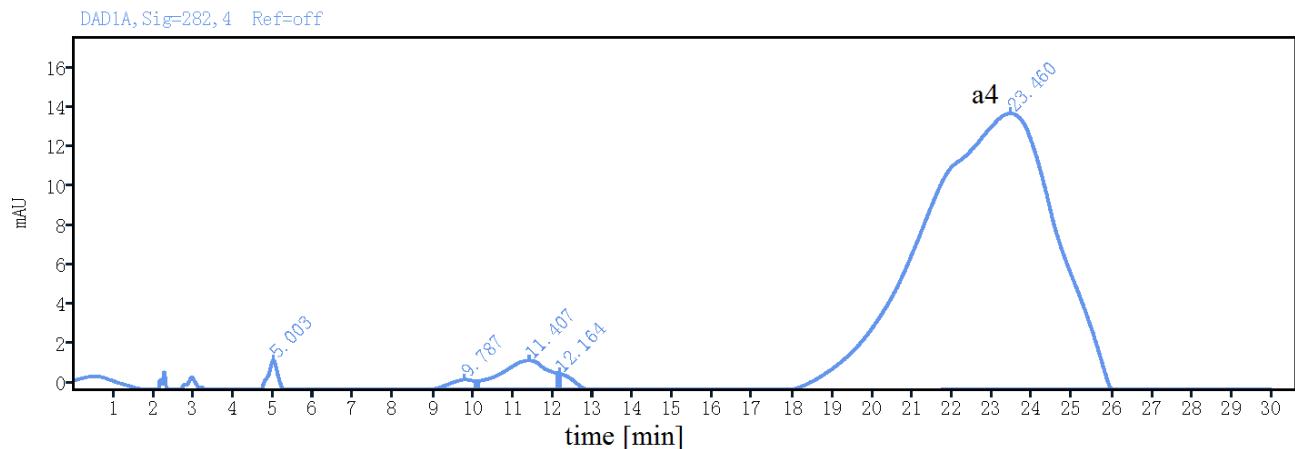
Retention time[min]	Area	Height	Area(%)
3. 062	212.99	18.59	3.66
5. 207	69.97	4.42	1.20
24. 020	5317.79	40.97	91.39
28. 540	218.31	2.84	3.75
			5819.05

(E)-5-(3-(3-chlorophenyl)acryloyl)-2-hydroxybenzamide (**a2**)



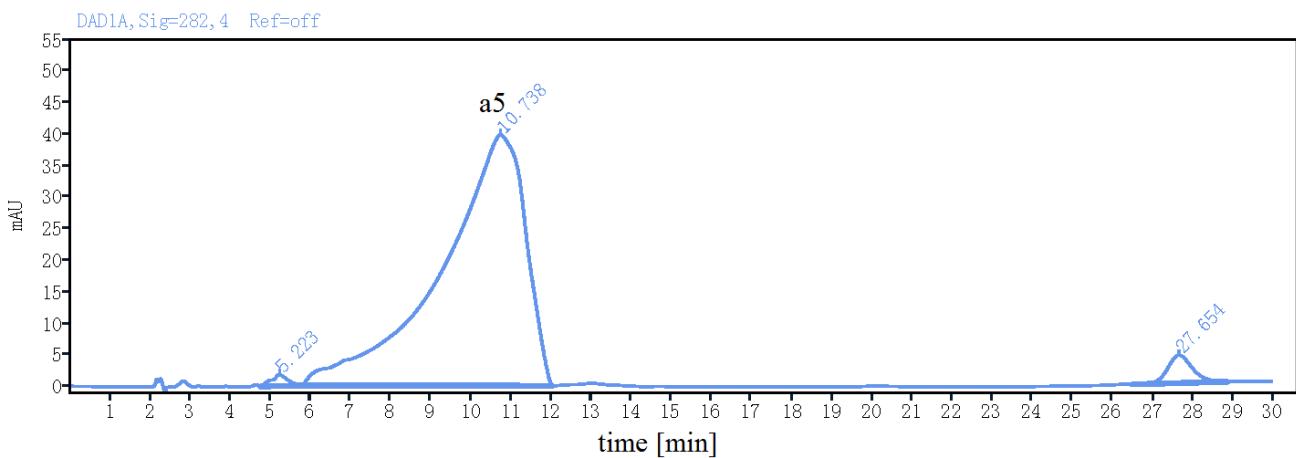
Retention time[min]	Area	Height	Area(%)
3.012	136.93	15.73	3.65
5.131	16.45	1.54	0.44
11.149	32.42	0.93	0.86
26.205	3568.92	19.94	95.05
			3754.72

(F)-5-(3-(3-bromophenyl)acryloyl)-2-hydroxybenzamide (a3)



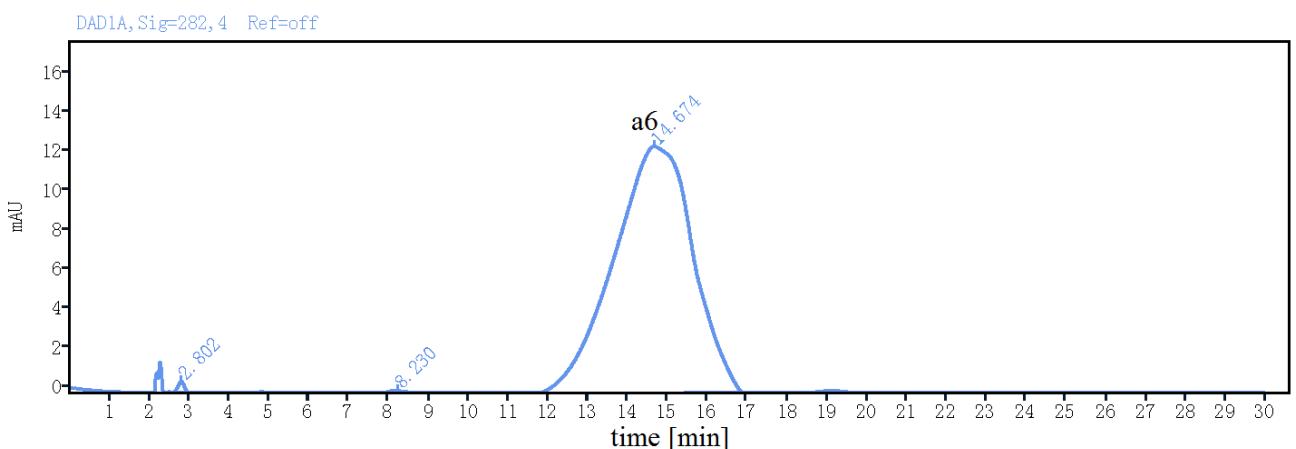
Retention time[min]	Area	Height	Area(%)
5.003	30.77	1.68	0.88
9.787	24.61	0.56	0.70
11.407	126.81	1.51	3.63
12.164	20.02	0.83	0.57
23.460	3295.37	14.09	94.22
			3497.58

(E)-5-(3-(4-bromophenyl)acryloyl)-2-hydroxybenzamide (a4)



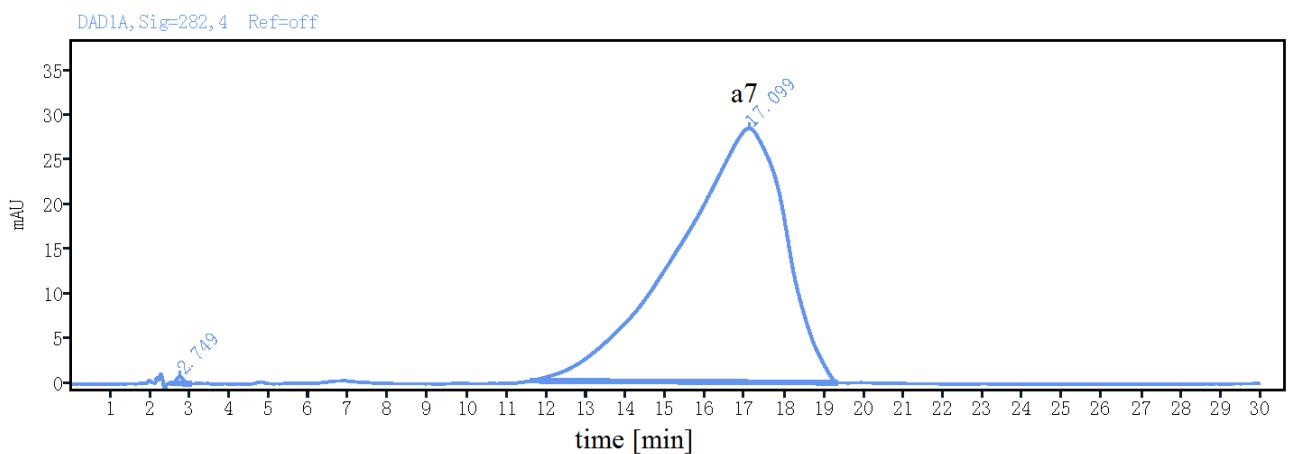
Retention time[min]	Area	Height	Area(%)
5.223	50.87	1.79	0.85
10.738	5744.32	39.75	95.93
27.654	192.87	4.42	3.22
	5988.05		

(F)-5-cinnamoyl-2-hydroxybenzamide (**a5**)



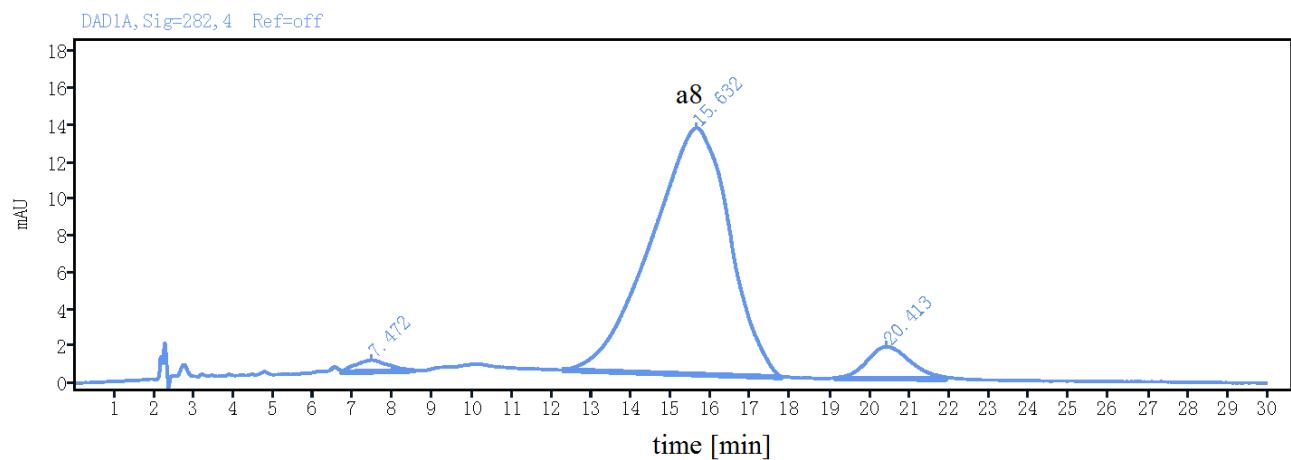
Retention time[min]	Area	Height	Area(%)
2.802	7.06	0.61	0.40
8.230	8.75	0.21	0.49
14.674	1765.65	12.64	99.11
	1781.45		

(E)-2-hydroxy-5-(3-(o-tolyl)acryloyl)benzamide (**a6**)



Retention time[min]	Area	Height	Area(%)
2.749	9.12	0.78	0.17
17.099	5384.61	28.45	99.83
			5393.73

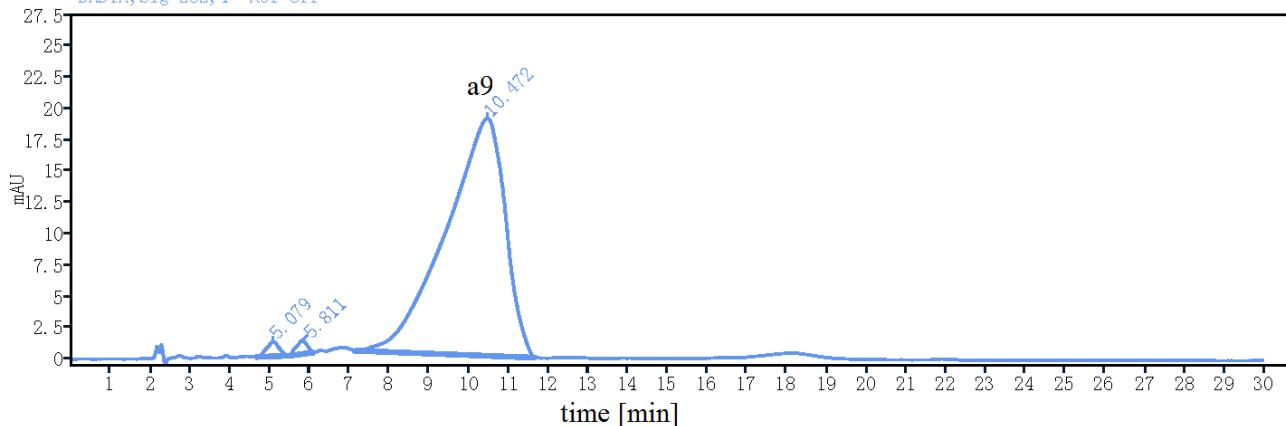
(E)-2-hydroxy-5-(3-(m-tolyl)acryloyl)benzamide (a7)



Retention time[min]	Area	Height	Area(%)
7.472	36.65	0.64	1.83
15.632	1834.75	13.39	91.55
20.413	132.60	1.75	6.62
	2004.00		

(E)-2-hydroxy-5-(3-(p-tolyl)acryloyl)benzamide (a8)

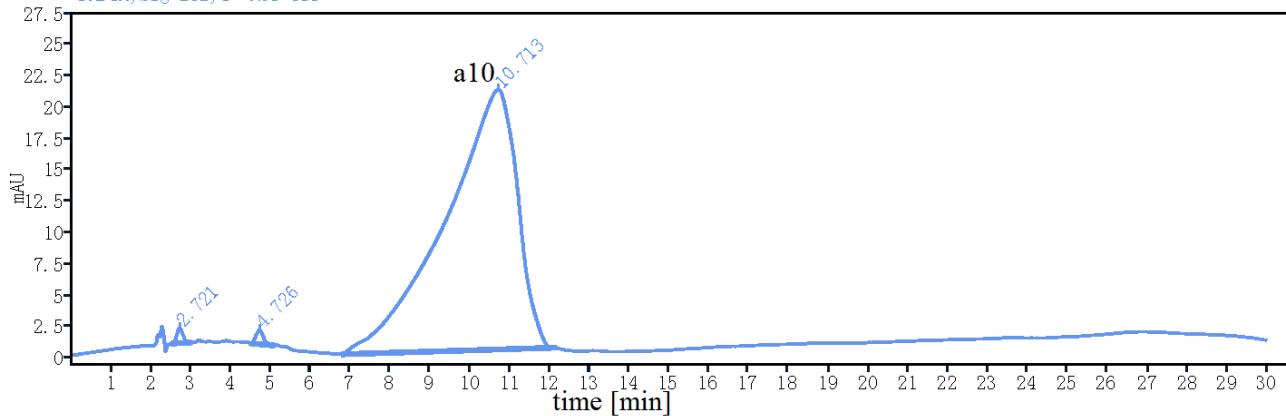
DAD1A, Sig=282, 4 Ref=off



Retention time[min]	Area	Height	Area(%)
5.079	25.23	1.16	1.26
5.811	20.62	1.07	1.03
10.472	1960.23	18.97	97.71
	2006.08		

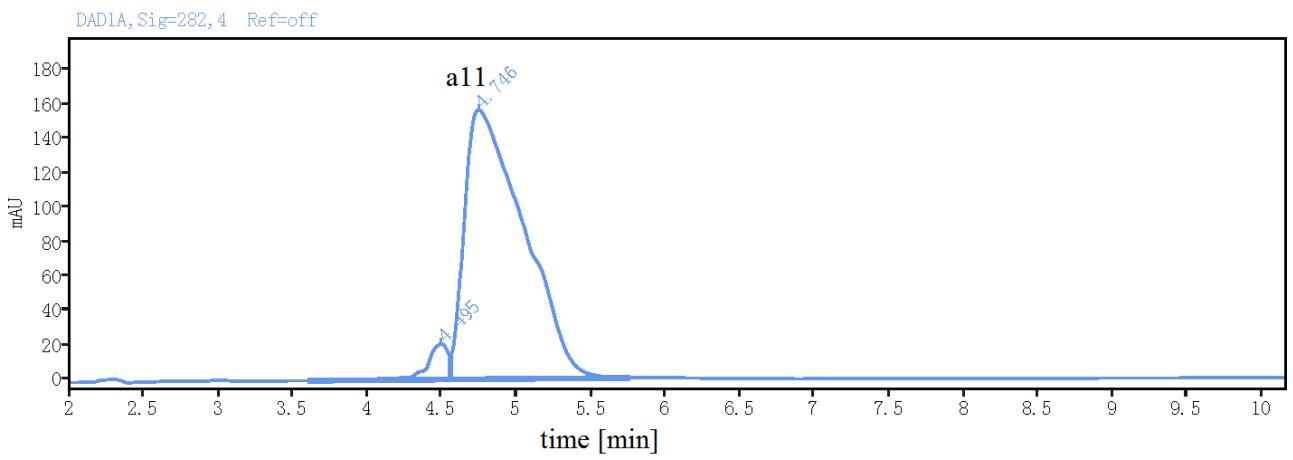
(F)-2-hydroxy-5-(3-(2-methoxyphenyl)acryloyl)benzamide (**a9**)

DAD1A, Sig=282, 4 Ref=off

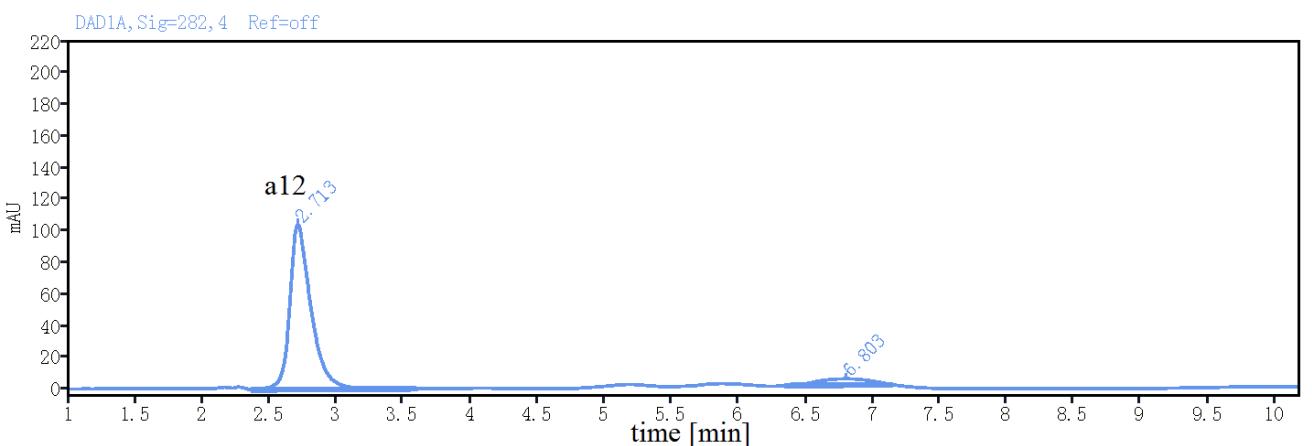


Retention time[min]	Area	Height	Area(%)
2.721	12.74	1.17	0.49
4.726	15.06	1.13	0.58
10.713	2589.88	20.72	98.94
	2617.69		

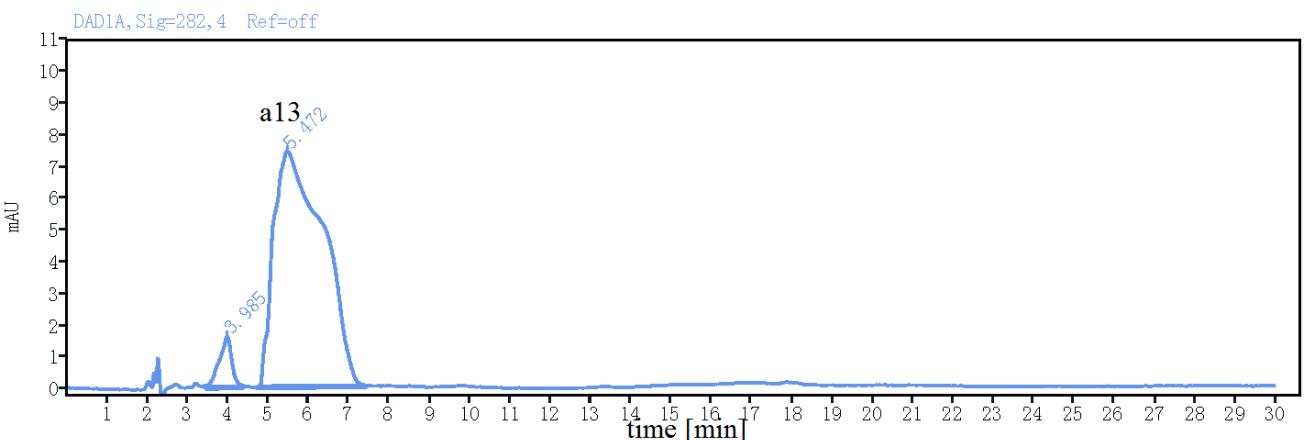
(E)-2-hydroxy-5-(3-(3-methoxyphenyl)acryloyl)benzamide(**a10**)



(F)-2-hydroxy-5-(3-(4-methoxyphenyl)acryloyl)benzamide(**a11**)

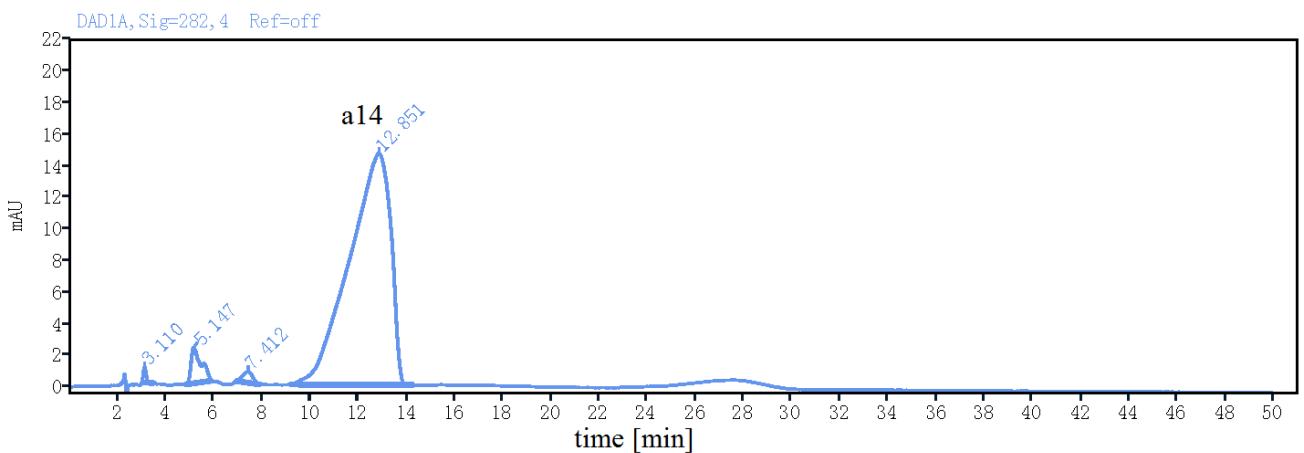


(E)-5-(3-(2,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (**a12**)



Retention time[min]	Area	Height	Area(%)
3. 985	34. 79	1. 61	5. 19
5. 472	635. 41	7. 43	94. 81
		670. 19	

(F) -5-(3,4-dimethoxyphenyl)acryloyl)-2-hydroxybenzamide (a13)



Retention time[min]	Area	Height	Area(%)
3. 110	6. 56	0. 96	0. 36
5. 147	65. 99	2. 26	3. 58
7. 412	20. 02	0. 70	1. 09
12. 851	1751. 58	14. 67	94. 98
		1844. 15	

(E)-2-hydroxy-5-(3-(3,4,5-trimethoxyphenyl)acryloyl)benzamide (a14)