

Figure S1. ESI MS for $[\text{Co}(\text{dmgBF}_2)_2(\text{H}_2\text{O})_2]$.
Low resolution

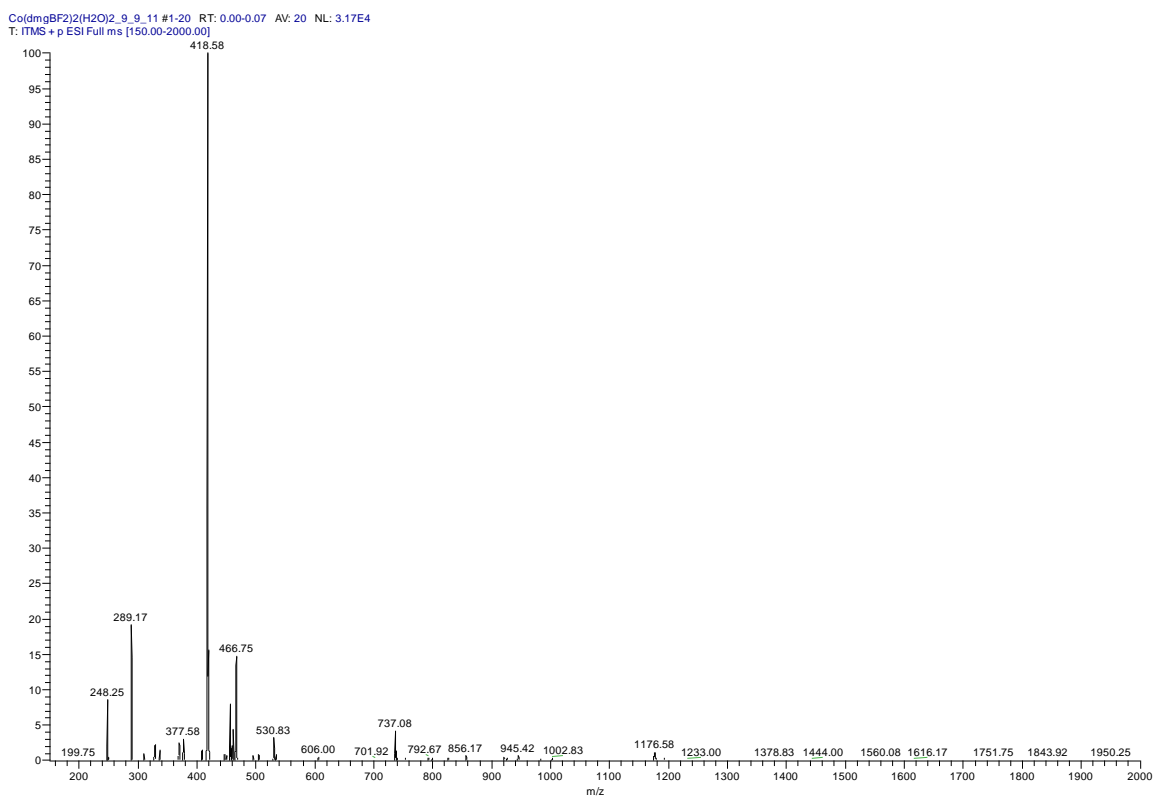
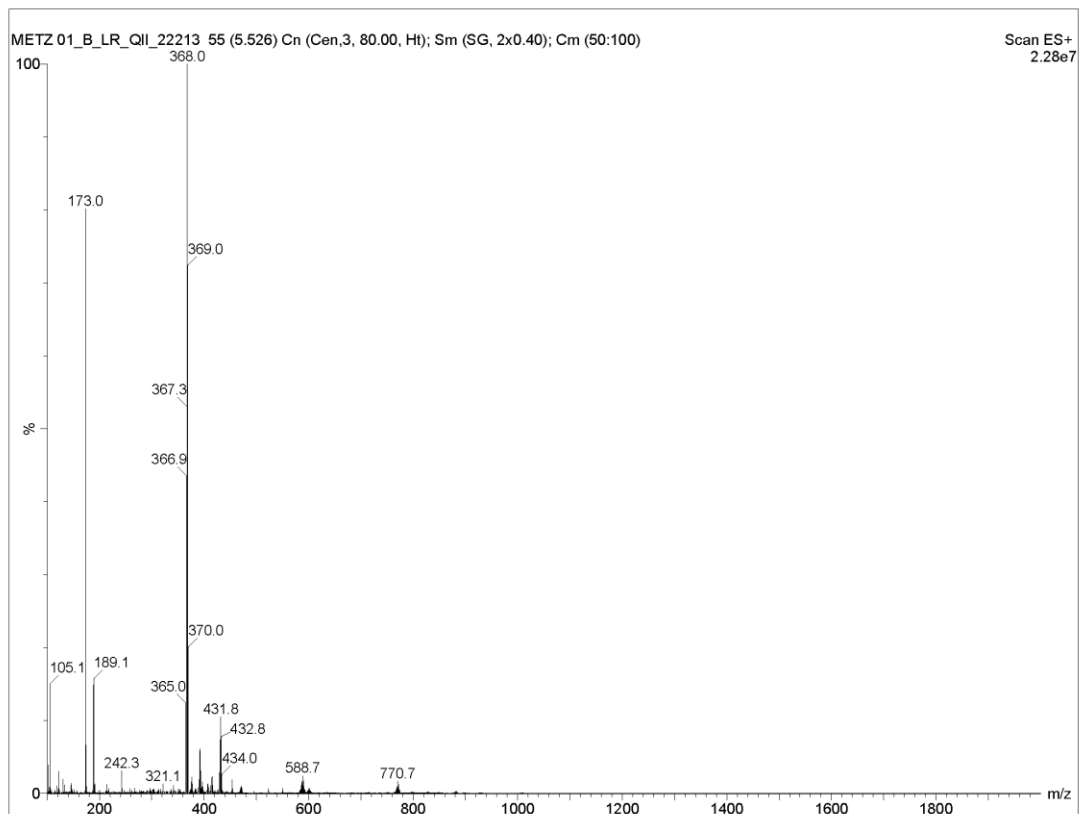
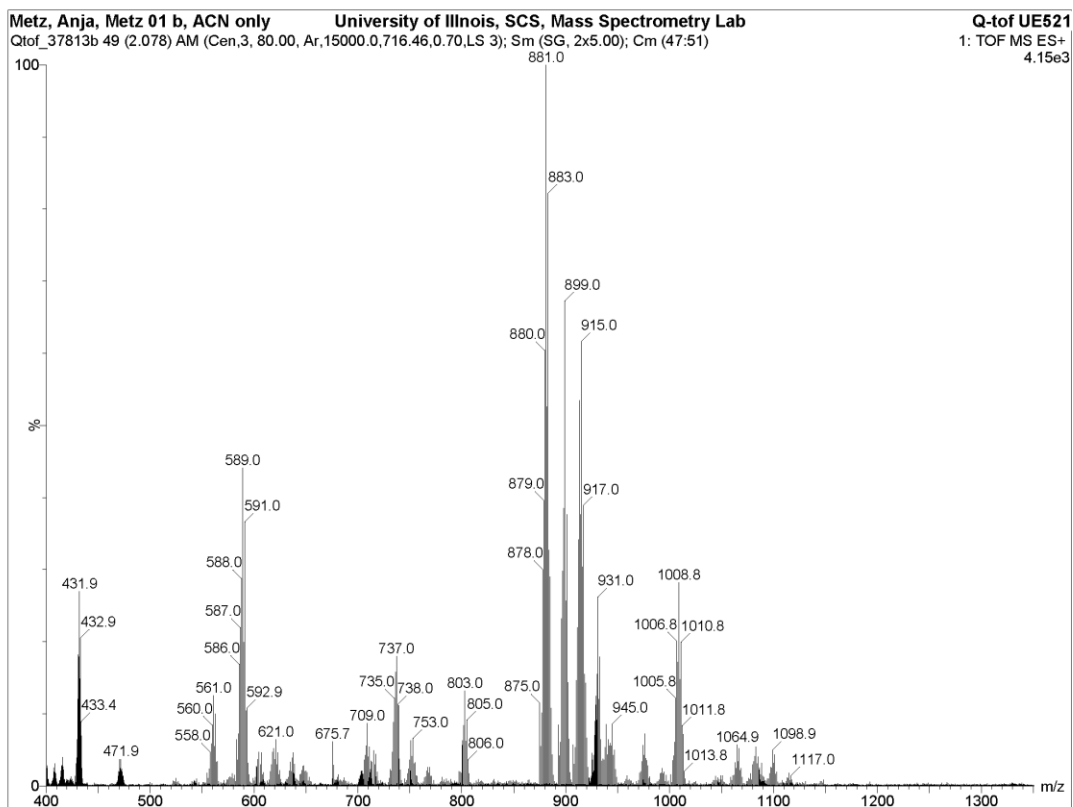


Figure S2. ESI MS for compound **4**.
Low resolution





High Resolution

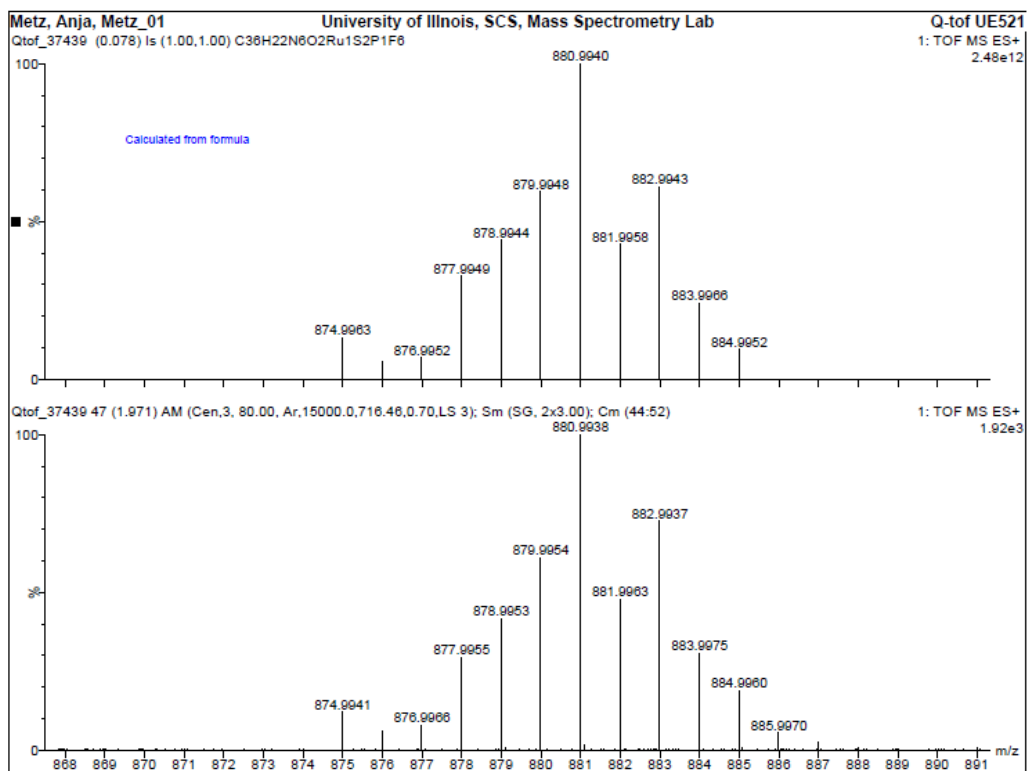
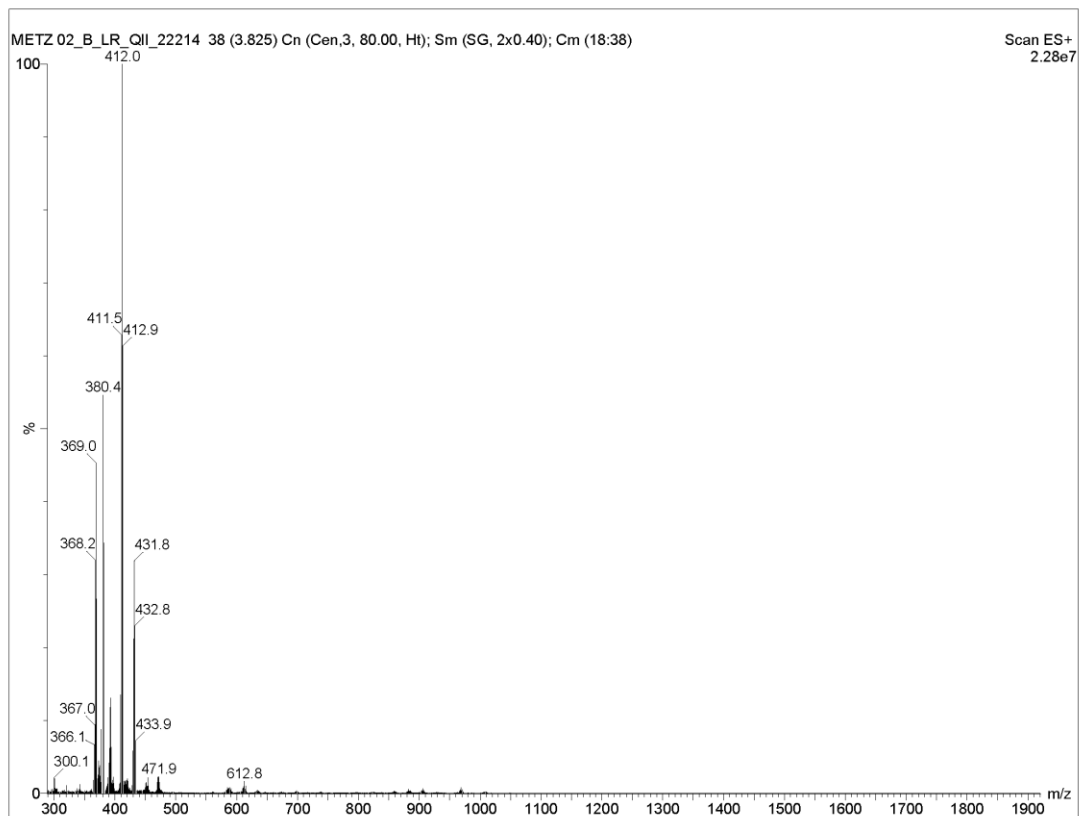
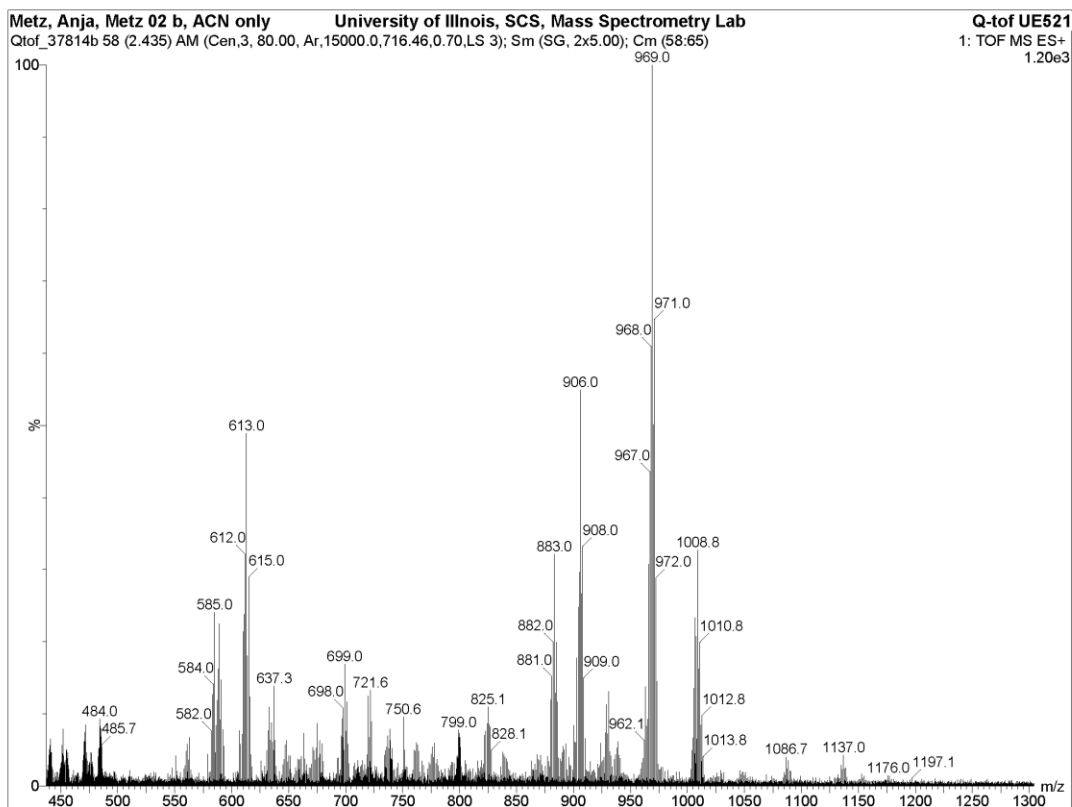


Figure S3. ESI MS for compound **5**.
Low resolution





High resolution

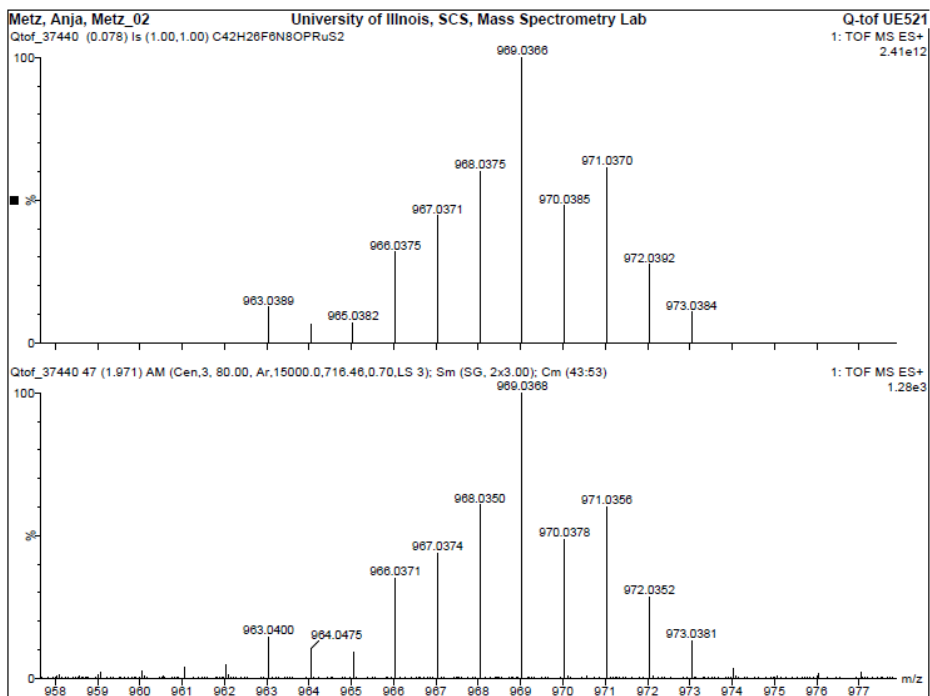
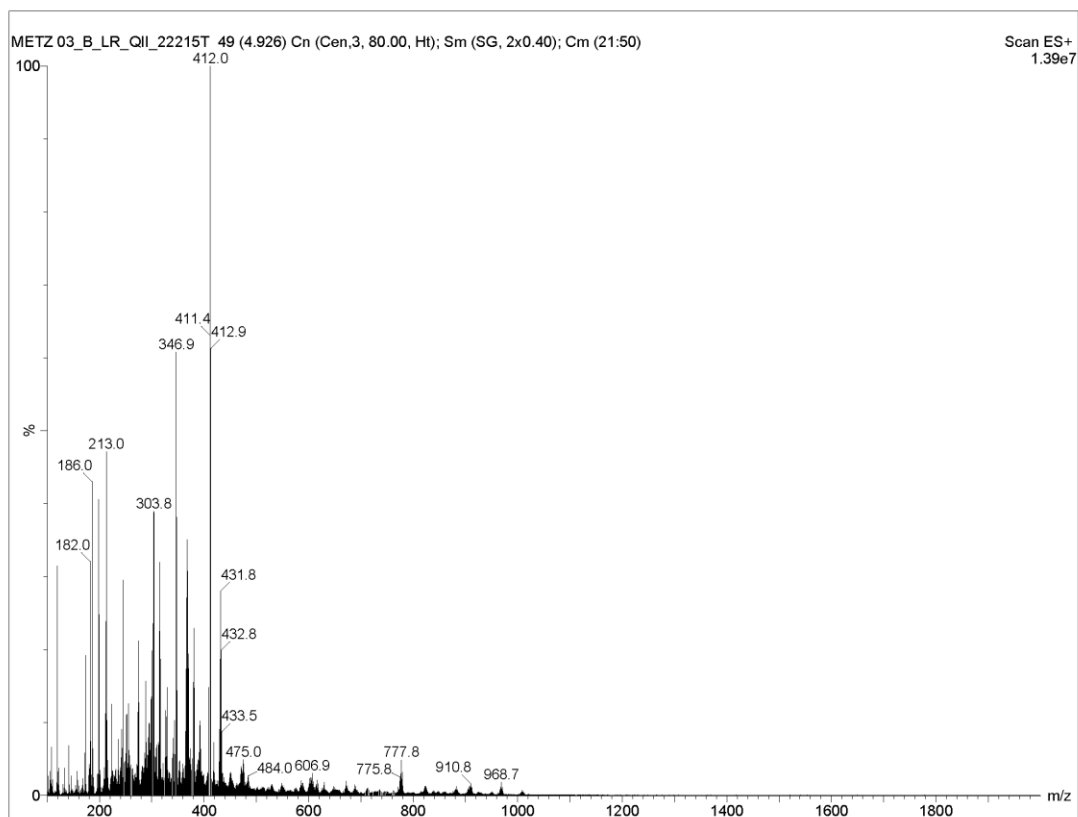
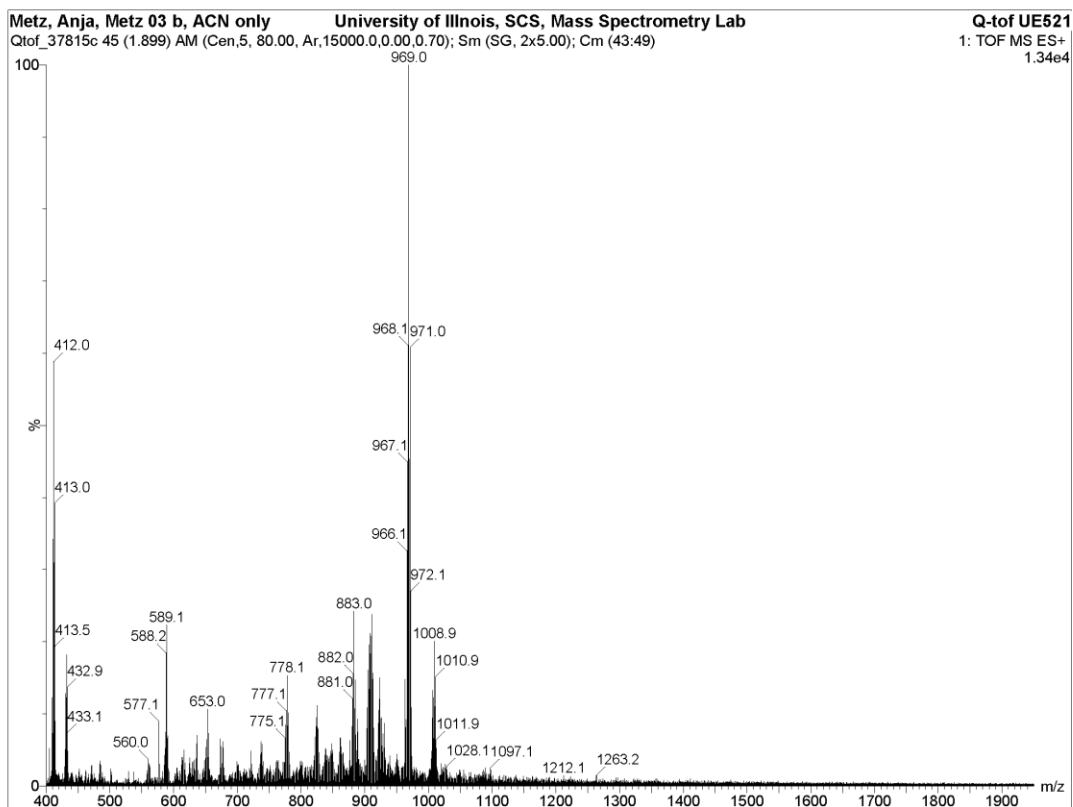


Figure S4. ESI MS for compound **6**.

Low resolution





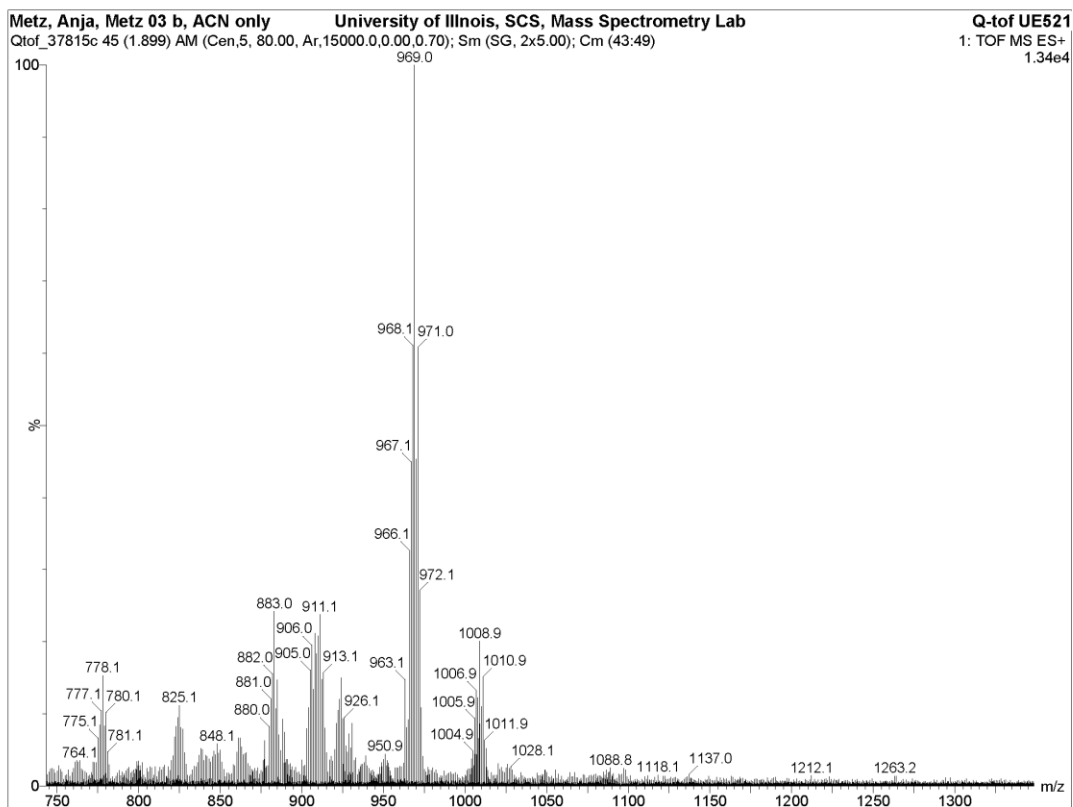


Figure S5. FT IR spectrum of $[\text{Co}(\text{dmgBF}_2)_2(\text{H}_2\text{O})_2]$.

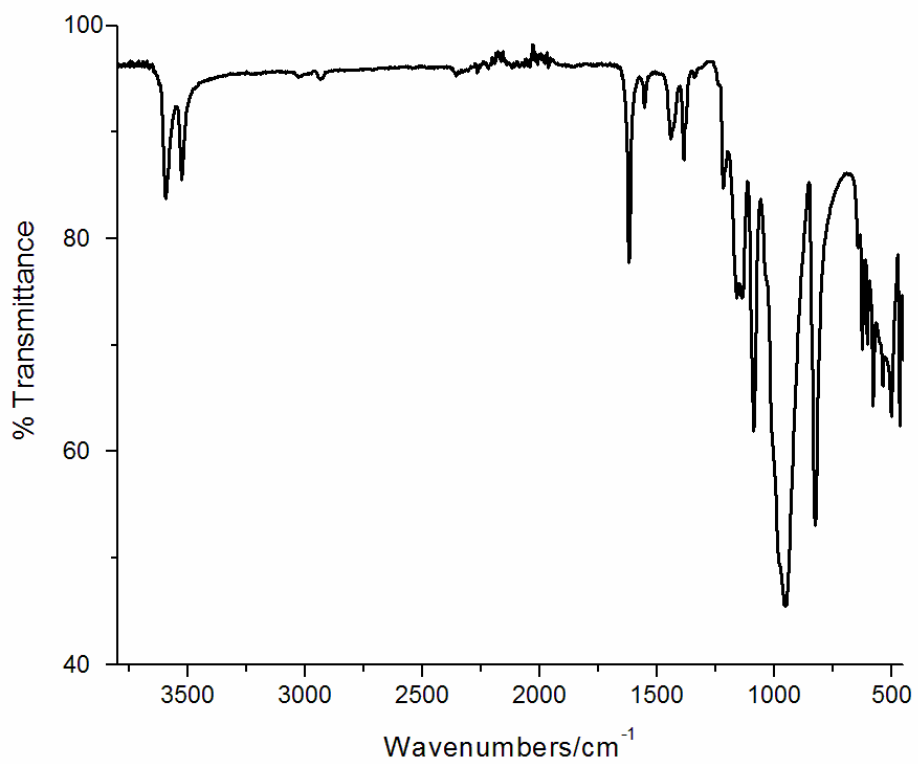


Figure S6. FT IR spectrum of compound 4.

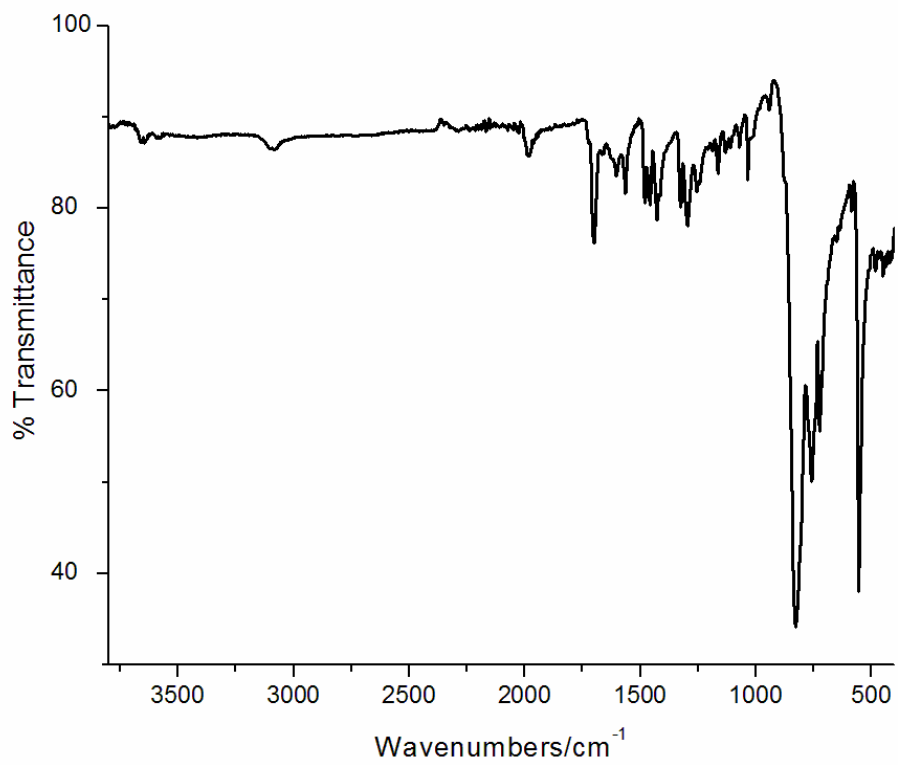


Figure S7. FT IR spectrum of compound **5**.

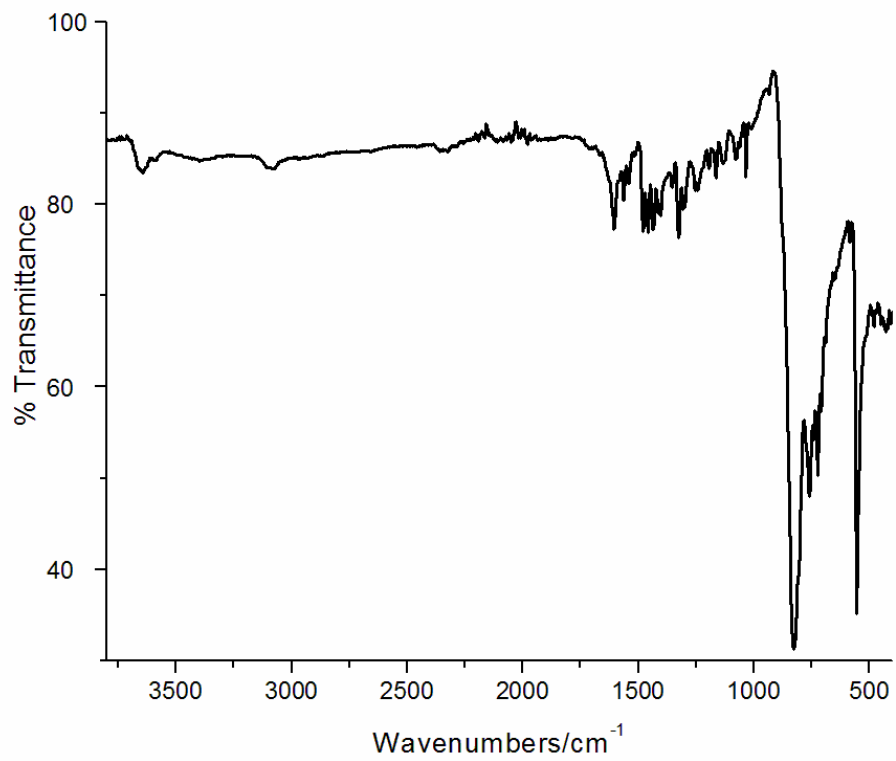


Figure S8. FT IR spectrum of compound **6**.

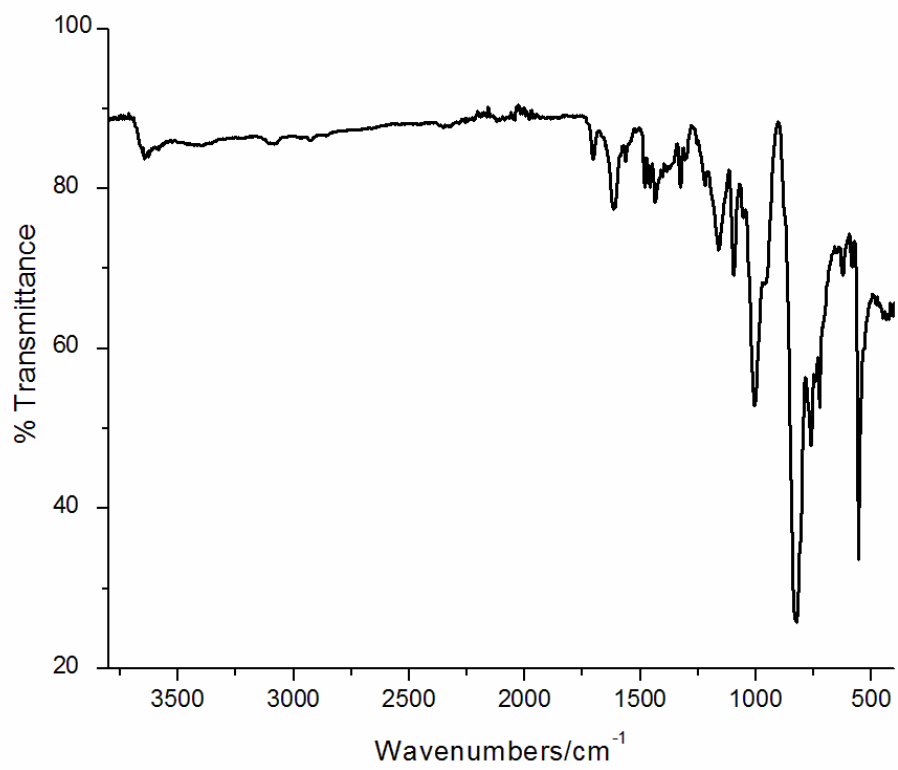
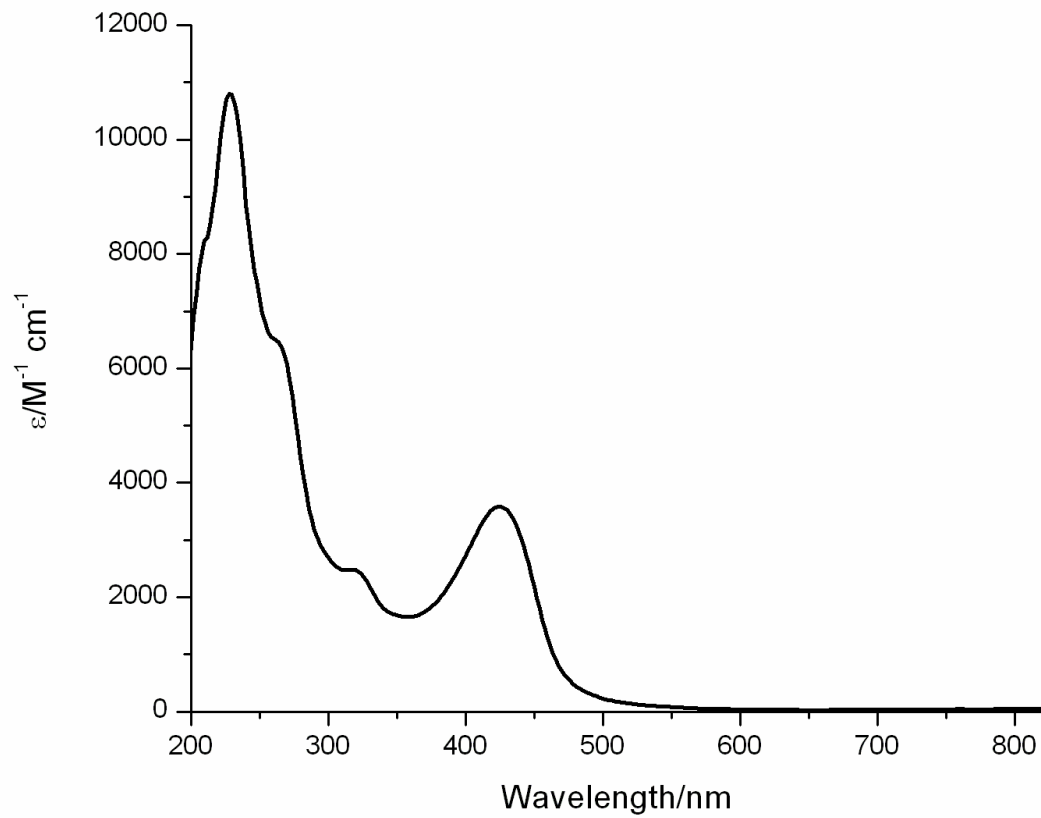


Figure S9. UV-visible spectra of dilute solutions of $[\text{Co}(\text{dmgBF}_2)_2(\text{H}_2\text{O})_2]$, compounds **4**, **5**, and **6** in CH_3CN .

(A) $[\text{Co}(\text{dmgBF}_2)_2(\text{H}_2\text{O})_2]$



(B) Compounds 4, 5, and 6.

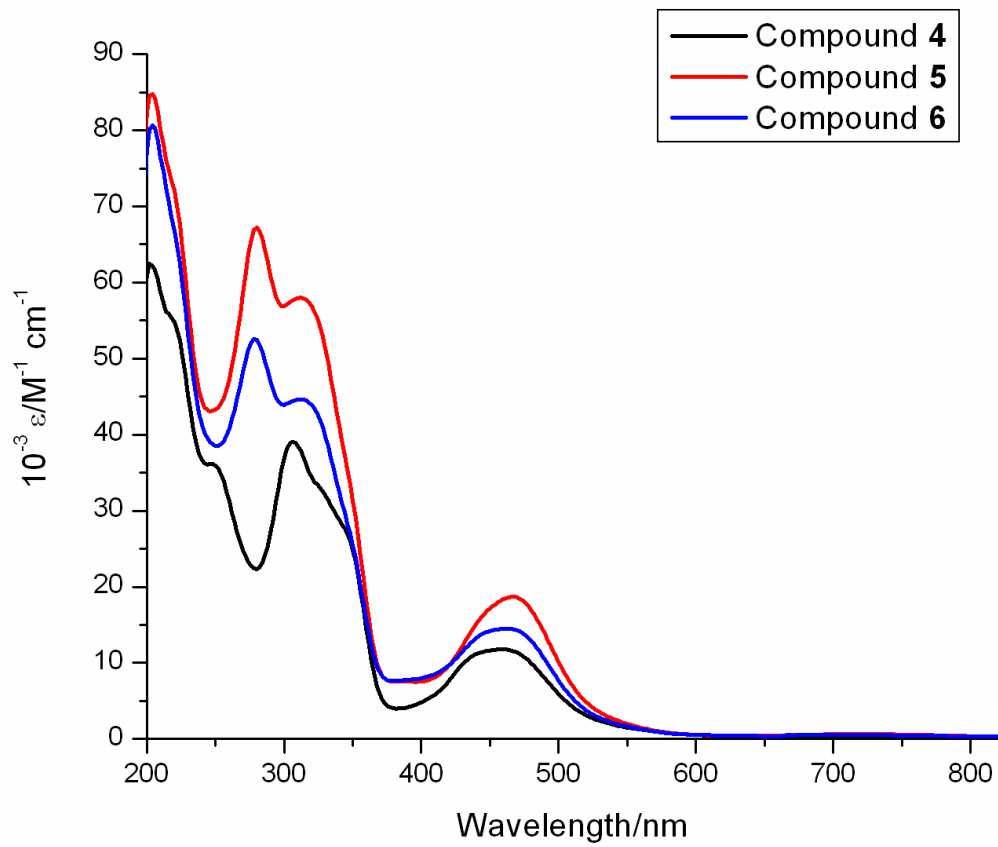


Figure S10. Q-band ESR spectra simulations of $[\text{Co}(\text{dmgBF}_2)_2(\text{OH}_2)_2]$ and compound **6** in DMSO at 100 K. Instrument settings were as follows: centre field = 11.5 kG, sweep width = 3500.00 G, microwave frequency = 34.06 GHz, microwave power = 2.00 mW, modulation frequency = 100 kHz, and modulation amplitude = 10.0 G.

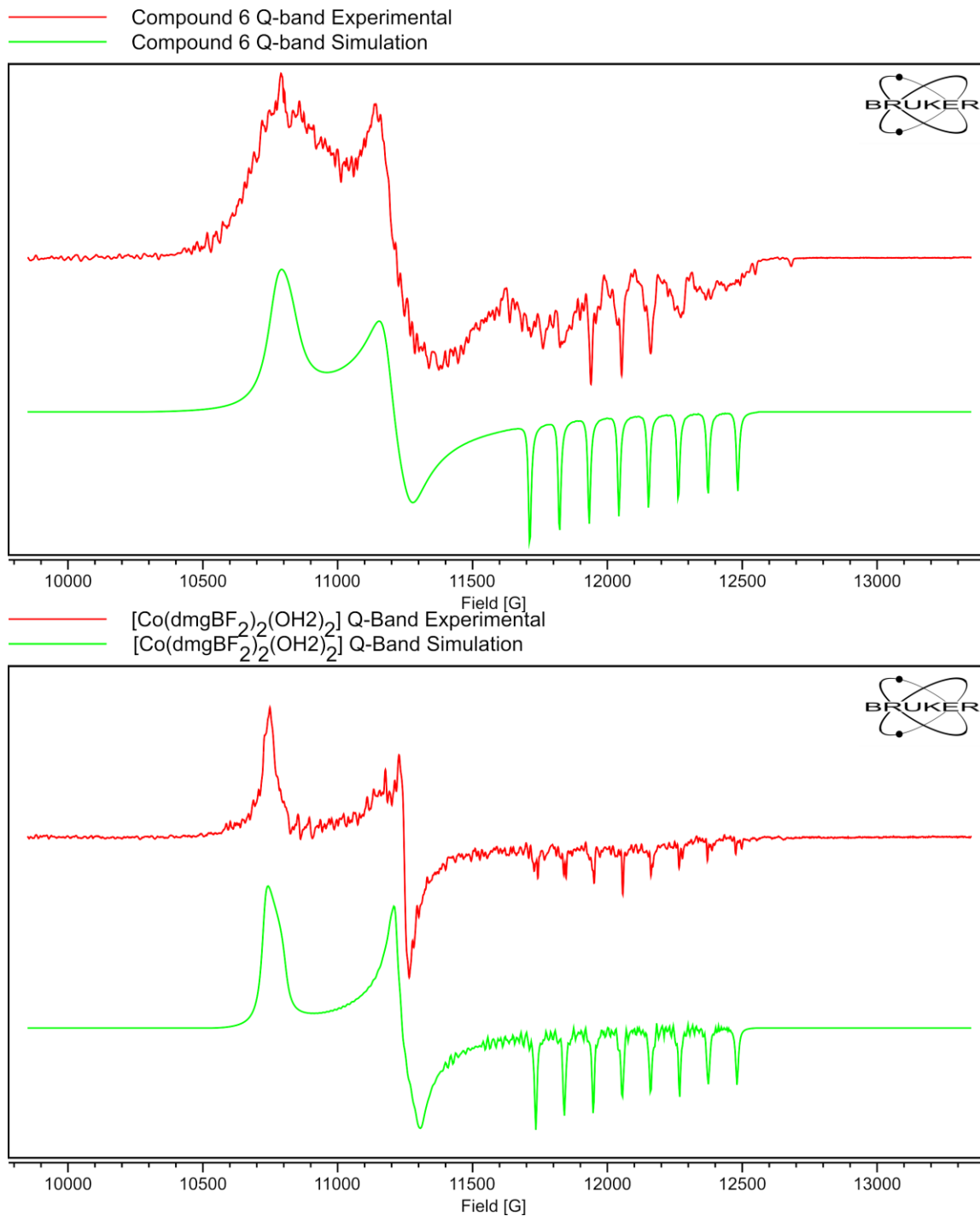


Figure S11. An X-band ESR spectrum and simulation of compound **8** in DMSO at 101 K. Instrument settings were as follows: centre field = 3200.00 G, sweep width = 1500.00 G, static field = 2450 G, microwave frequency = 9.46 GHz, microwave power = 0.22 mW, receiver gain = 5.02×10^4 , modulation frequency = 100 kHz, modulation amplitude = 2.00 G, modulation phase = 0° , time constant = 40.96 ms, conversion time = 40.00 ms, and sweep time = 96.00 s.

