

**Asymmetric conjugate additions of 2-substituted benzofuran-3(2*H*)-ones with α ,
 β -unsaturated ketone catalyzed by chiral copper complexes**

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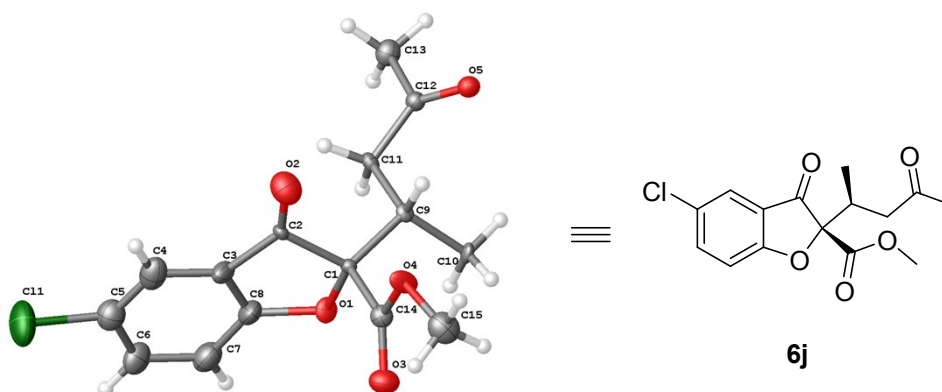
Supporting Information

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1、 Crystal data and structure refinement for enantiopure **6j**

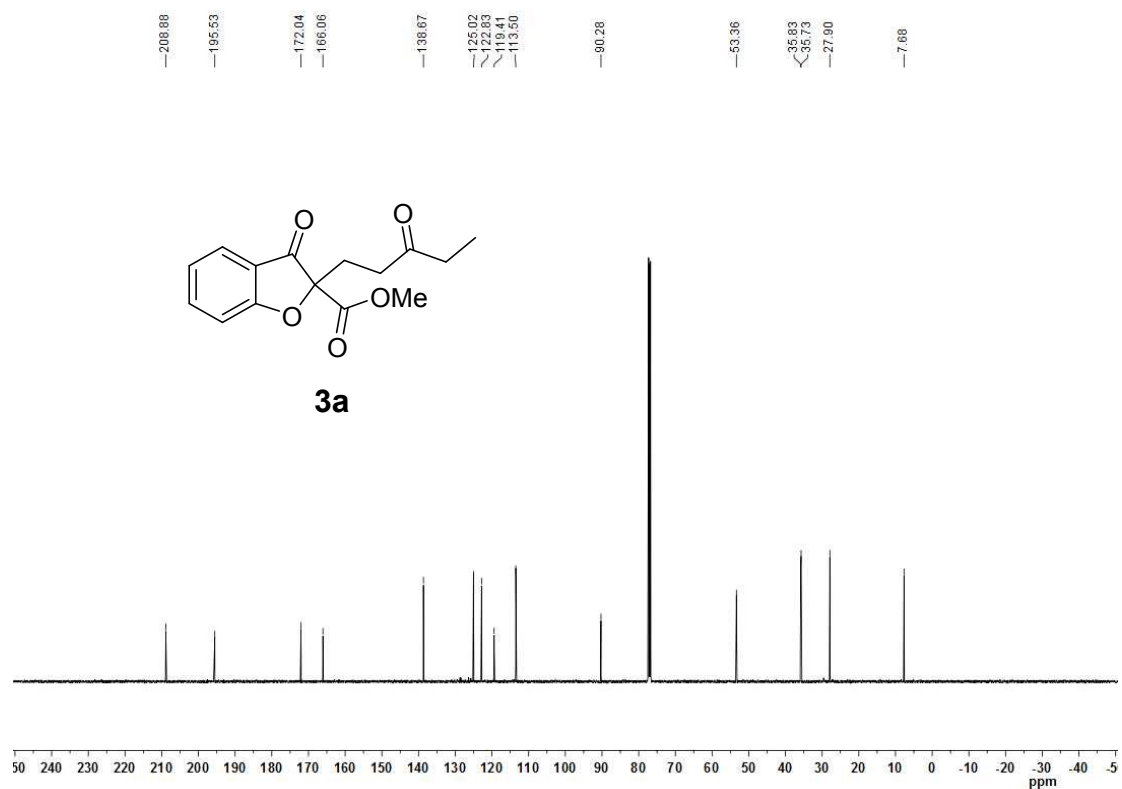
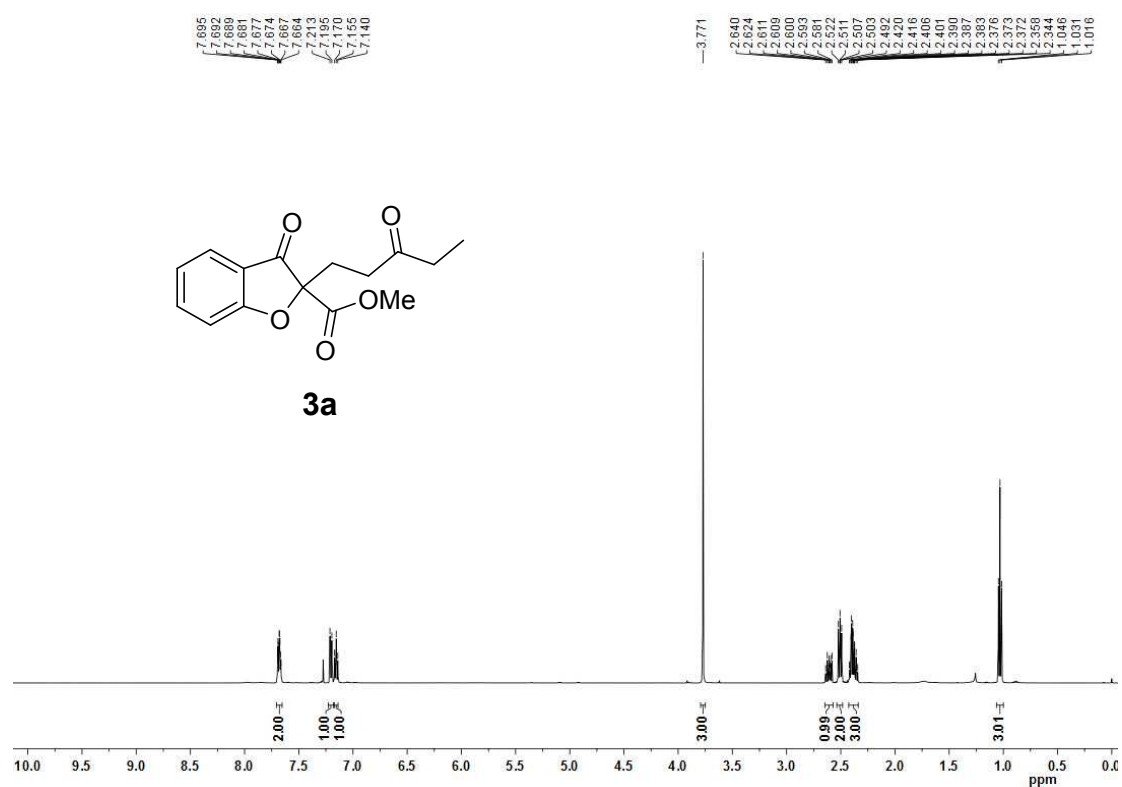
Crystals of enantiopure **6j** suitable for X-ray analysis were obtained from crystallization in a solution of dichloromethane/*n*-hexane.

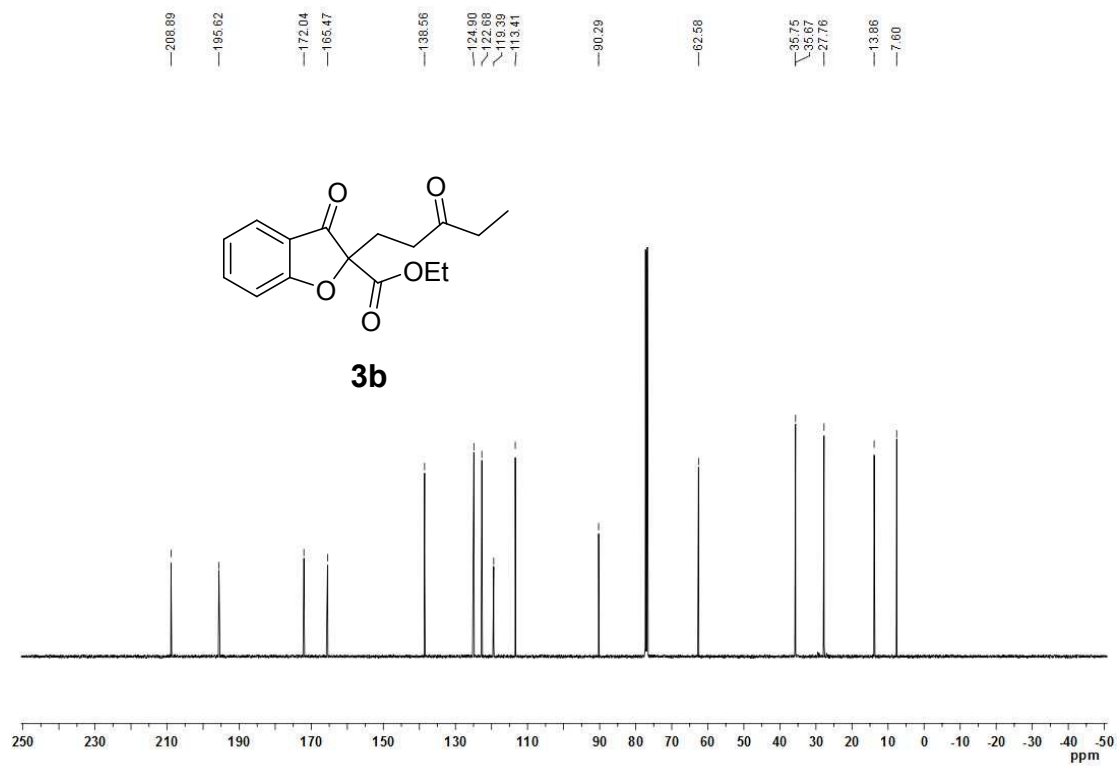
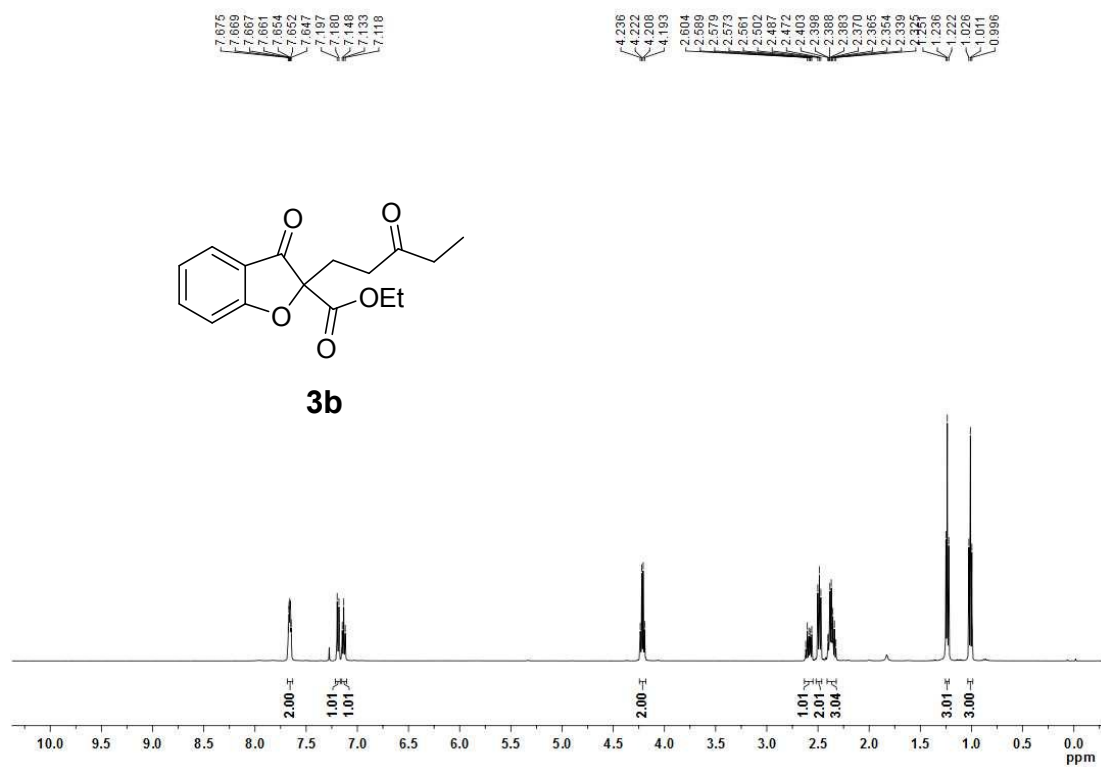


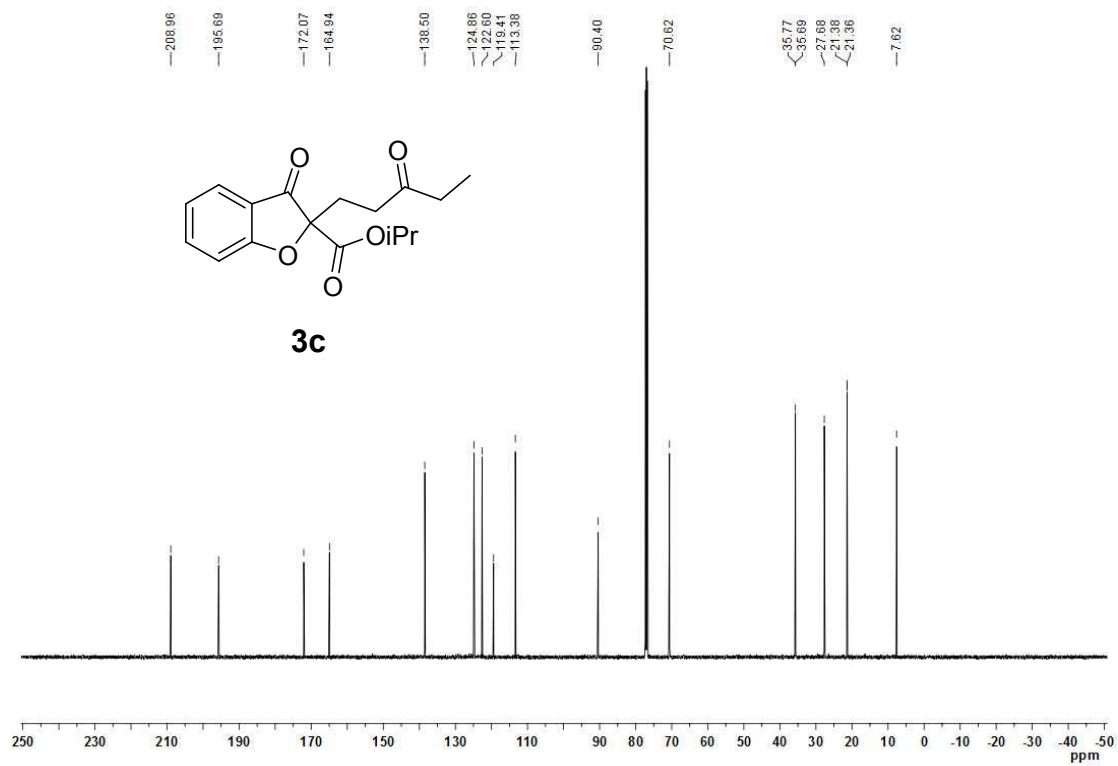
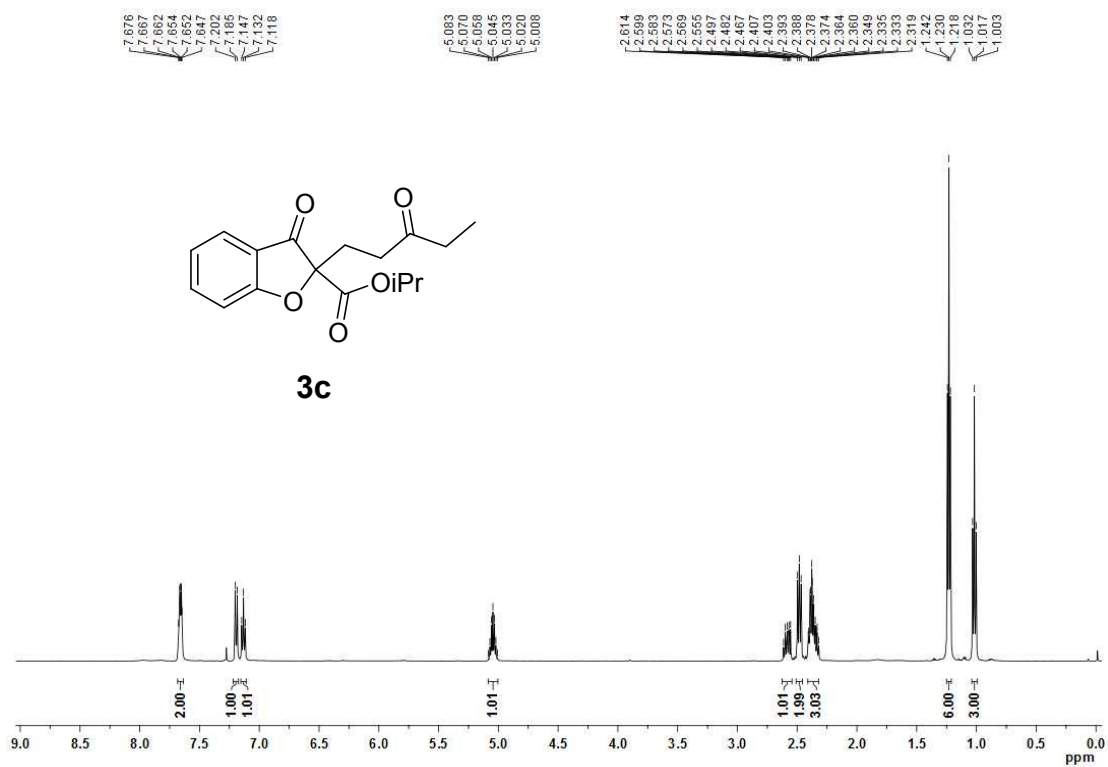
Identification code	mo_dm16873_0m	
Empirical formula	C ₁₅ H ₁₅ ClO ₅	
Formula weight	310.72	
Temperature	130 K	
Wavelength	0.71073 Å	
Crystal system	Hexagonal	
Space group	P 65	
Unit cell dimensions	a = 15.8422(17) Å	∠ = 90°.
	b = 15.8422(17) Å	∠ = 90°.
	c = 11.1926(12) Å	∠ = 120°.
Volume	2432.7(6) Å ³	
Z	6	
Density (calculated)	1.273 Mg/m ³	
Absorption coefficient	0.252 mm ⁻¹	
F(000)	972	
Crystal size	0.05 x 0.05 x 0.05 mm ³	
Theta range for data collection	1.48 to 27.50°.	

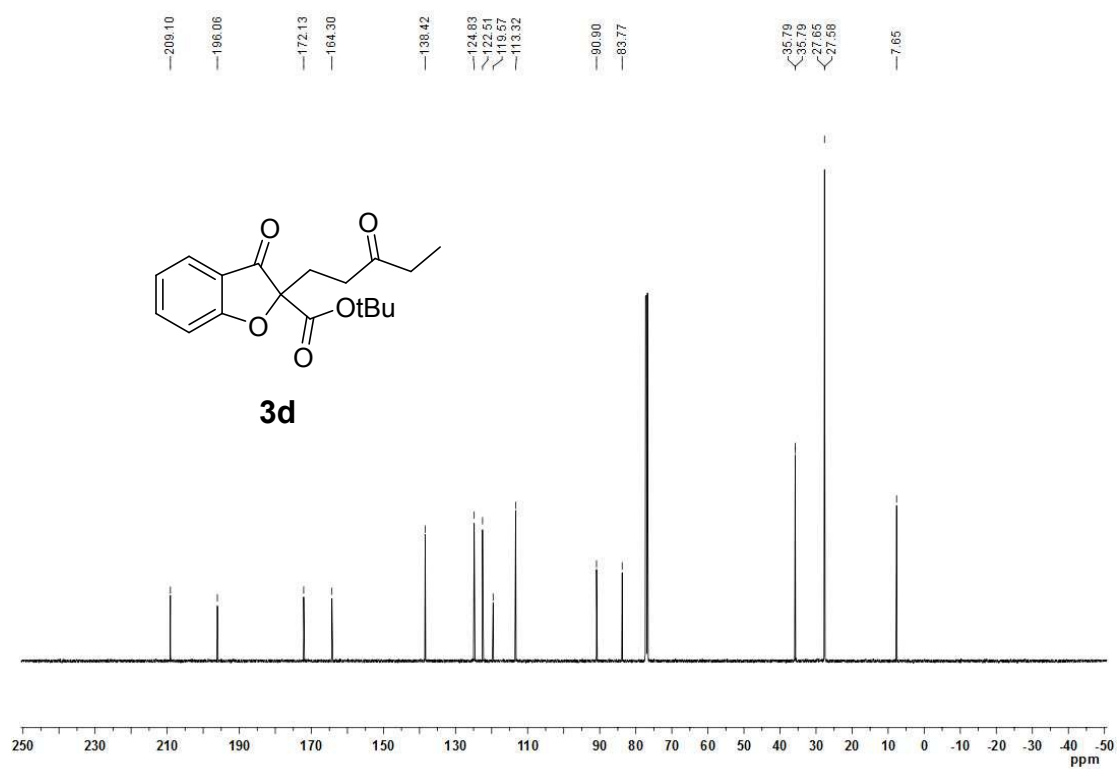
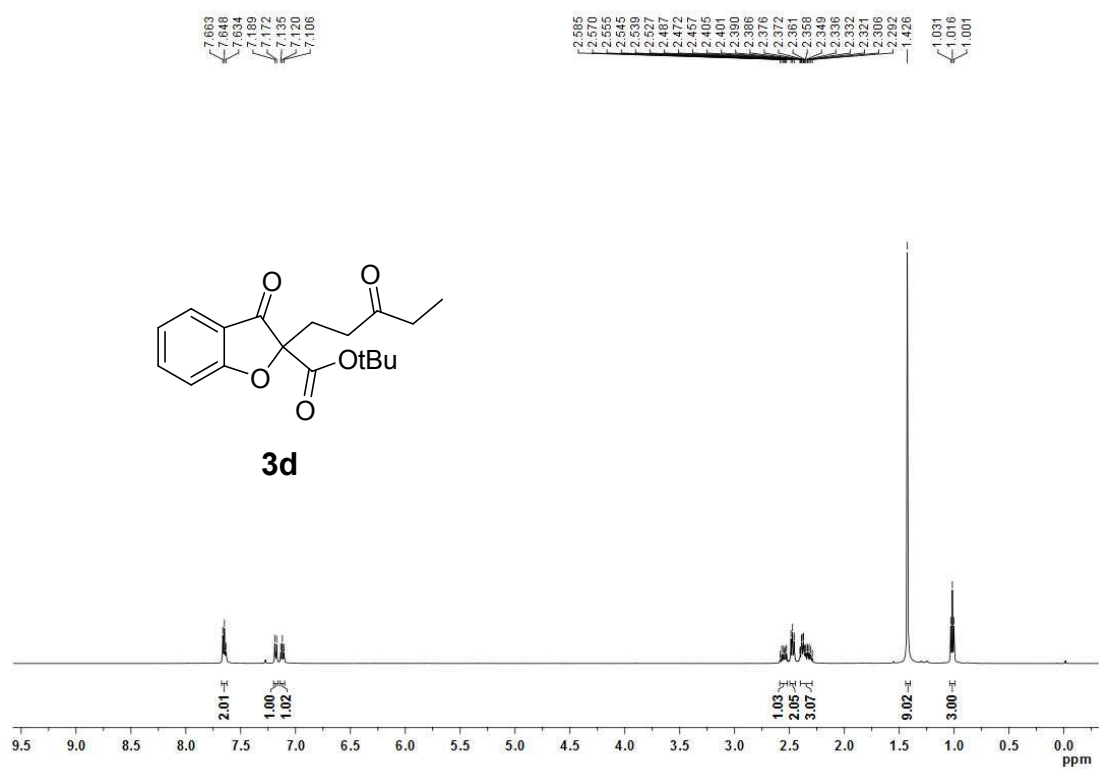
Index ranges	-20<=h<=20, -20<=k<=20, -14<=l<=14
Reflections collected	20466
Independent reflections	3011 [R(int) = 0.0768]
Completeness to theta = 25.242°	100.0 %
Absorption correction	Semi-empirical from equivalents
Max. and min. transmission	0.7456 and 0.6540
Refinement method	Full-matrix least-squares on F ²
Data / restraints / parameters	3011 / 265 / 322
Goodness-of-fit on F ²	0.924
Final R indices [I>2sigma(I)]	R1 = 0.0535, wR2 = 0.1541
R indices (all data)	R1 = 0.0719, wR2 = 0.1733
Absolute structure parameter	0.03(4)
Extinction coefficient	n/a
Largest diff. peak and hole	0.232 and -0.291 e.Å ⁻³

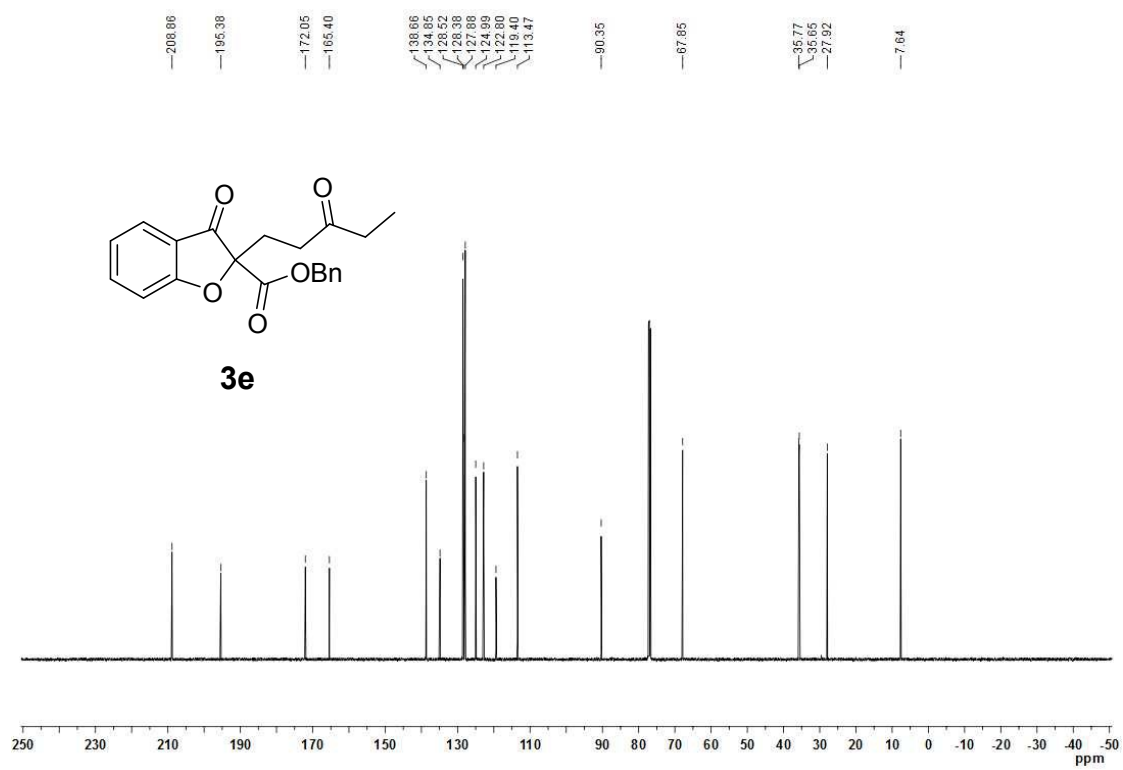
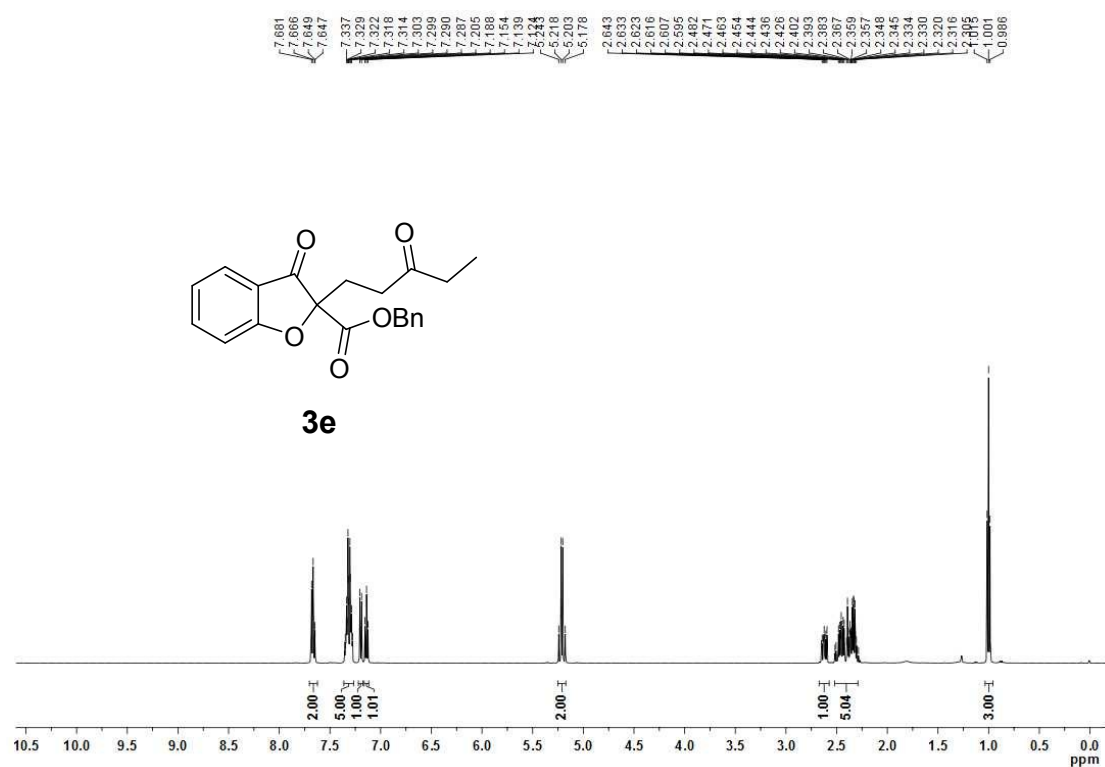
2、 ^1H and ^{13}C NMR spectra of all products

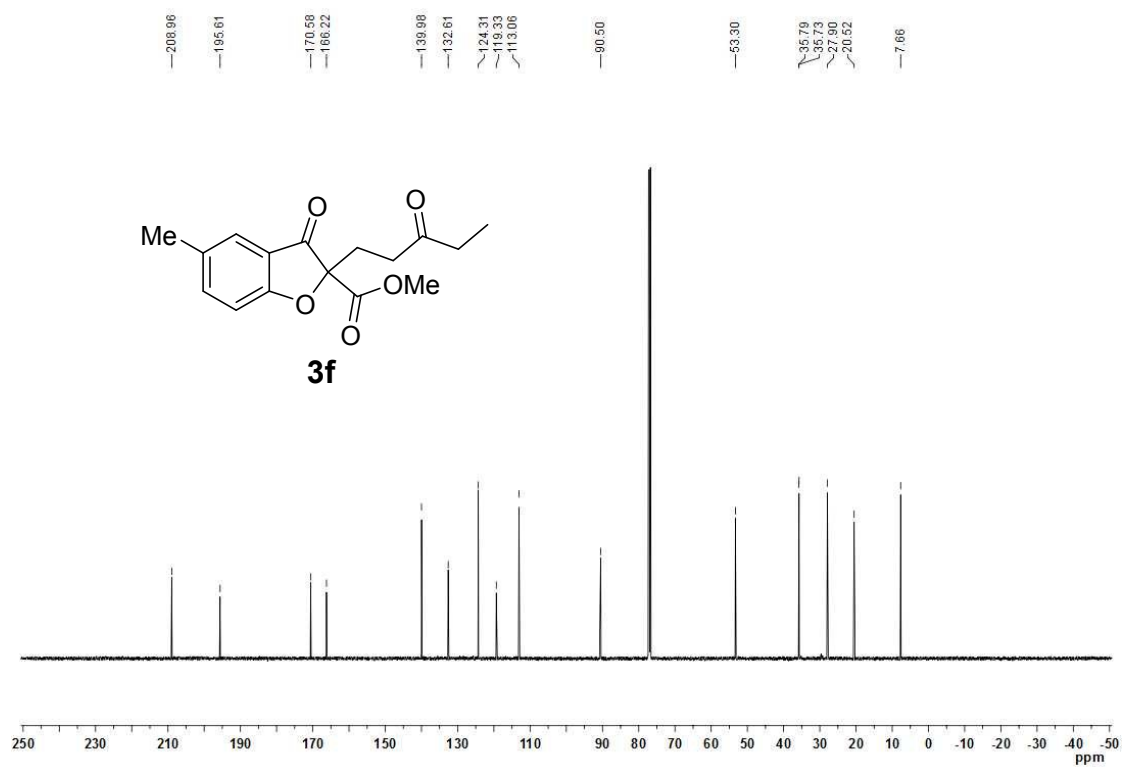
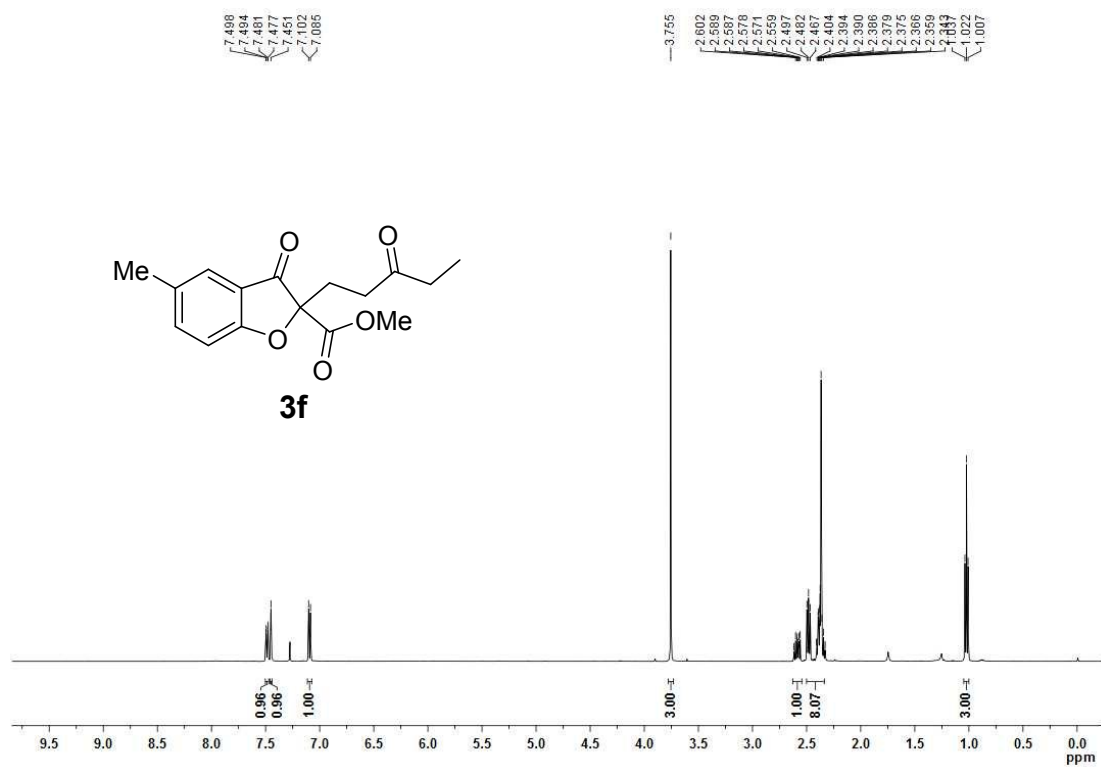


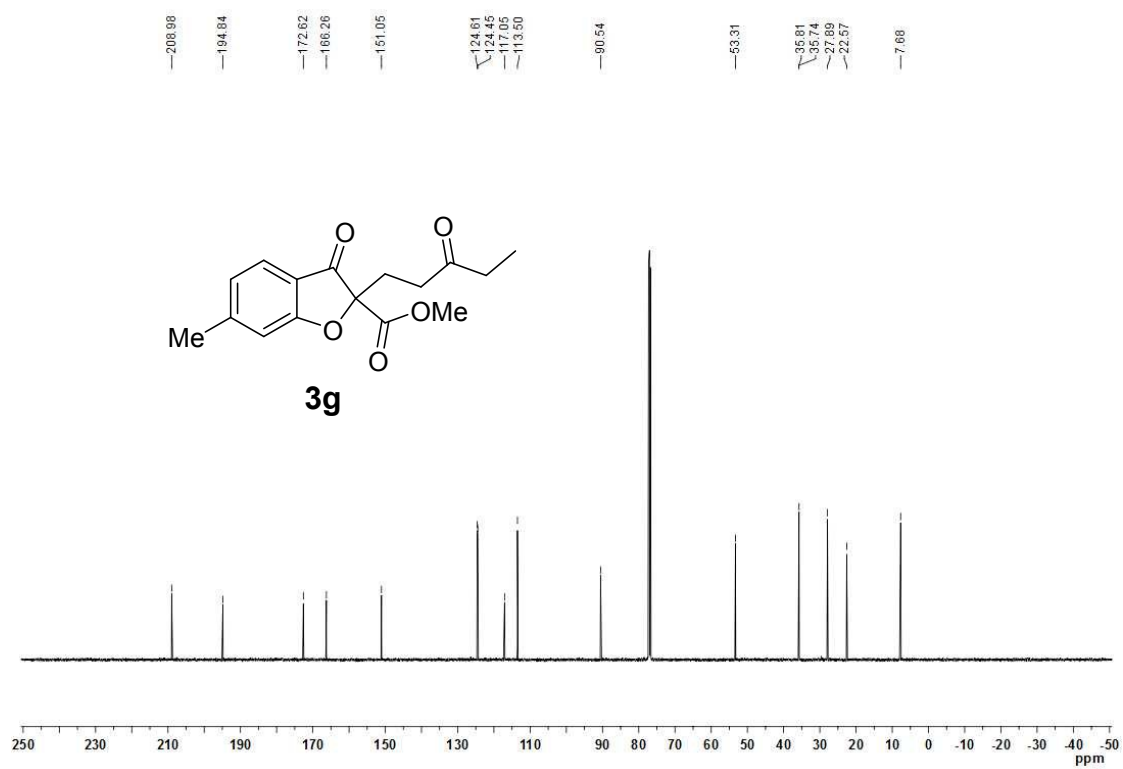
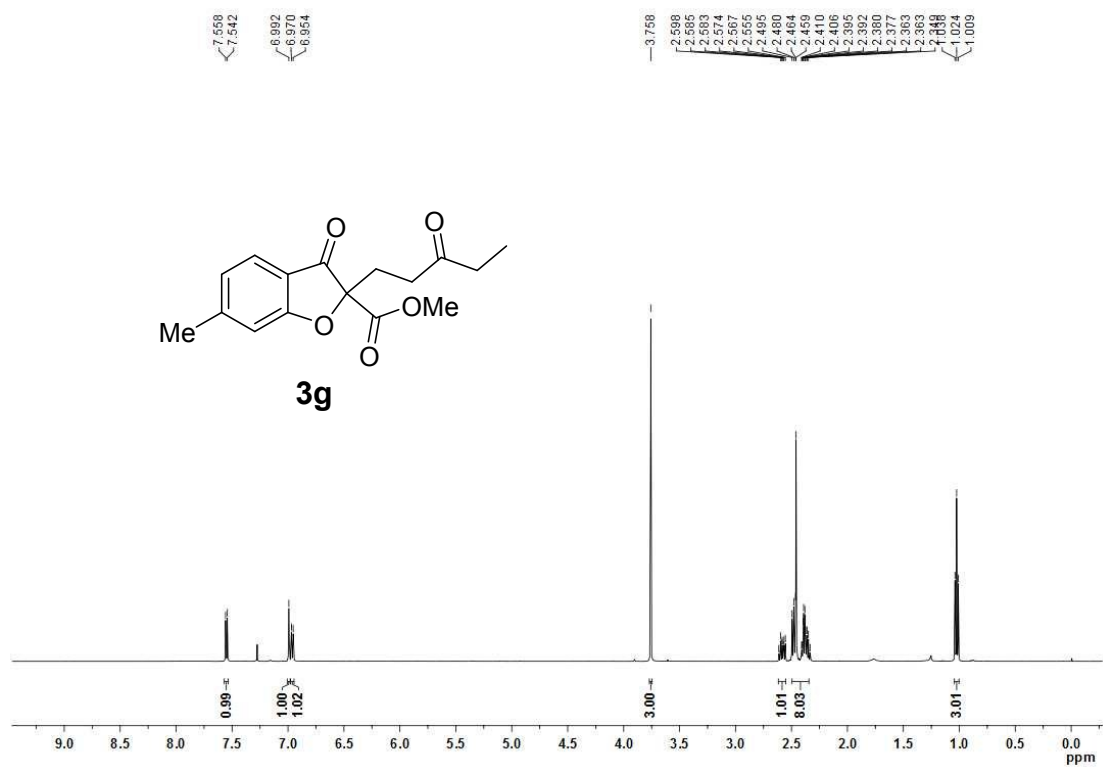


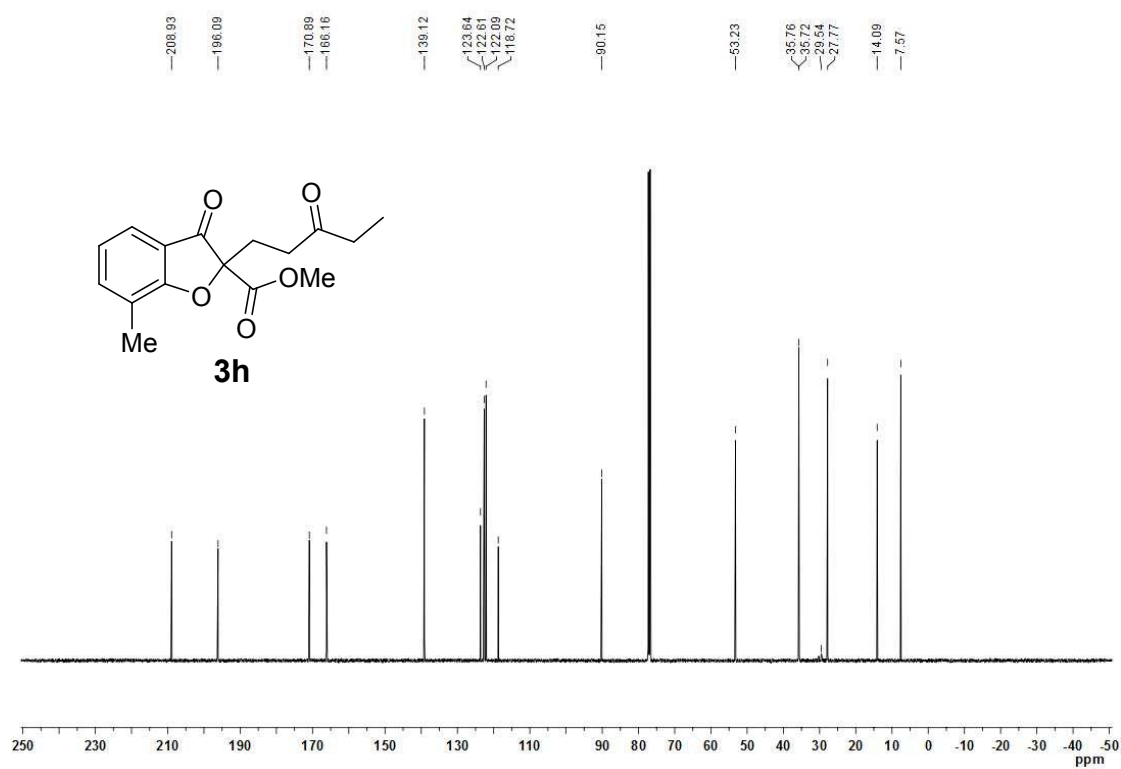
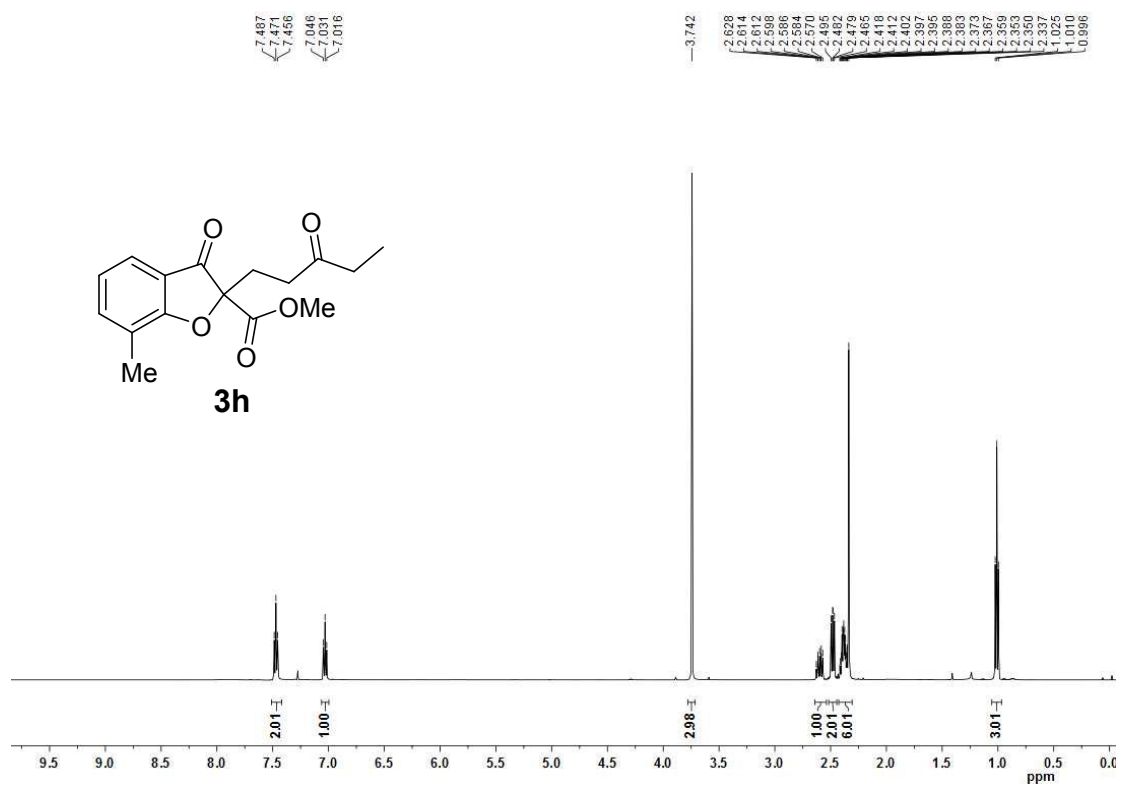


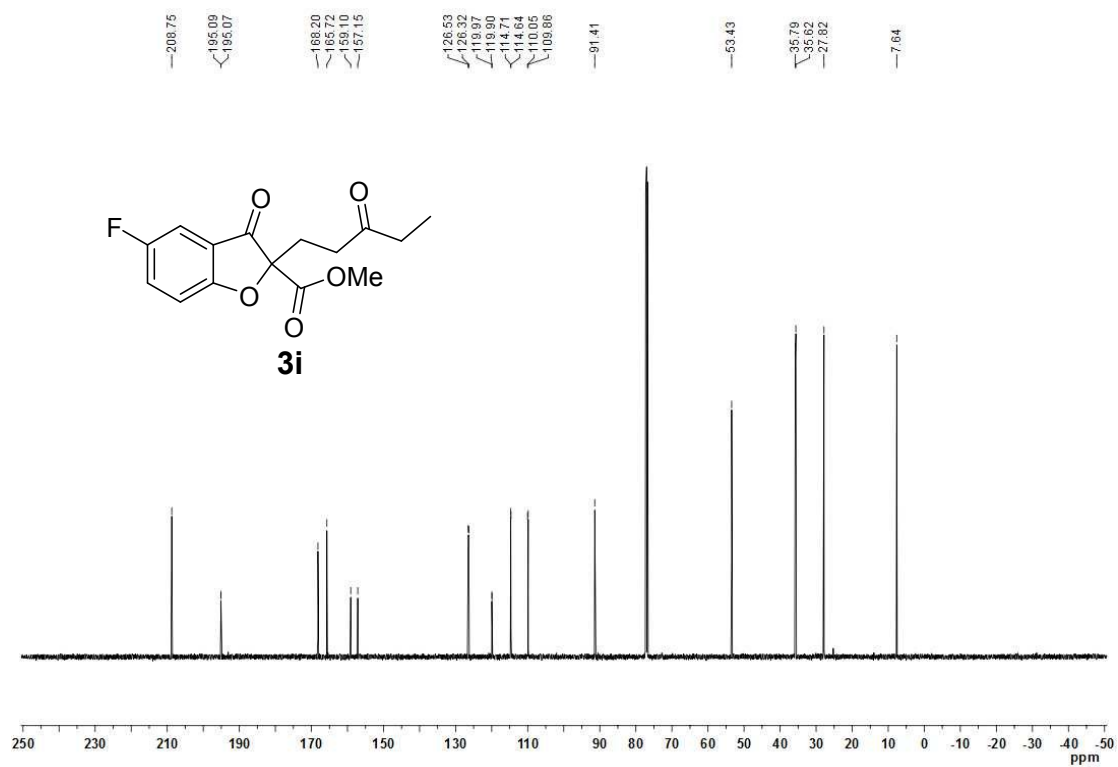
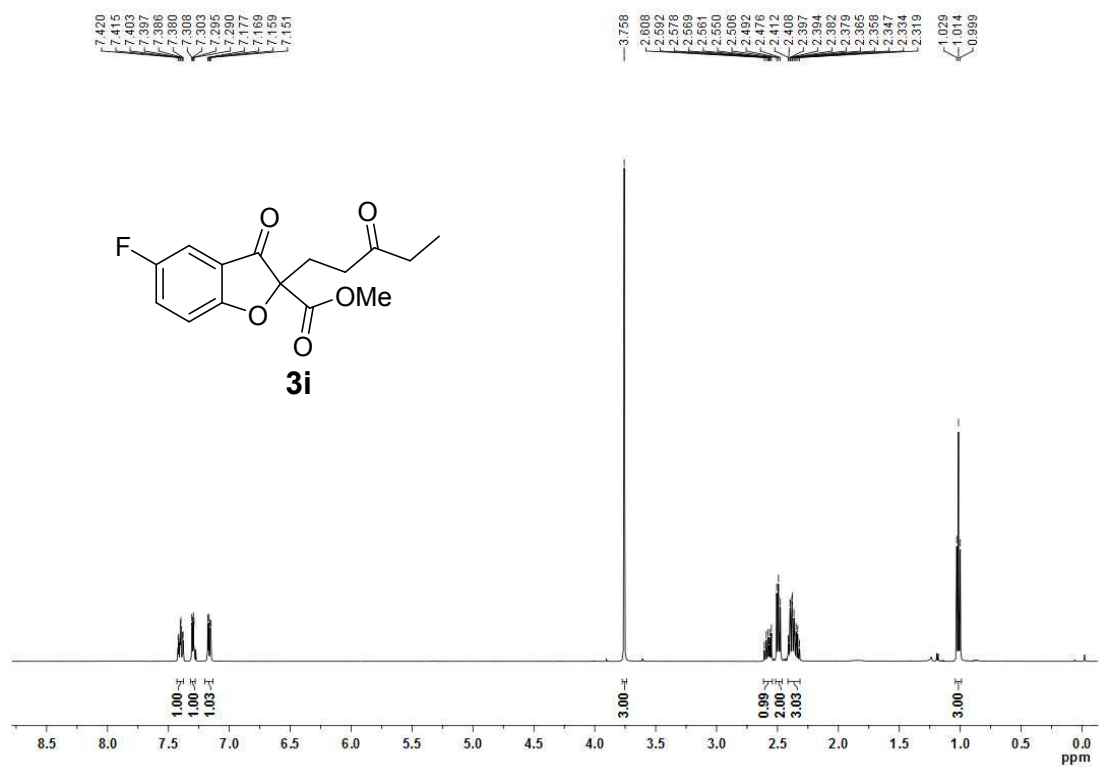


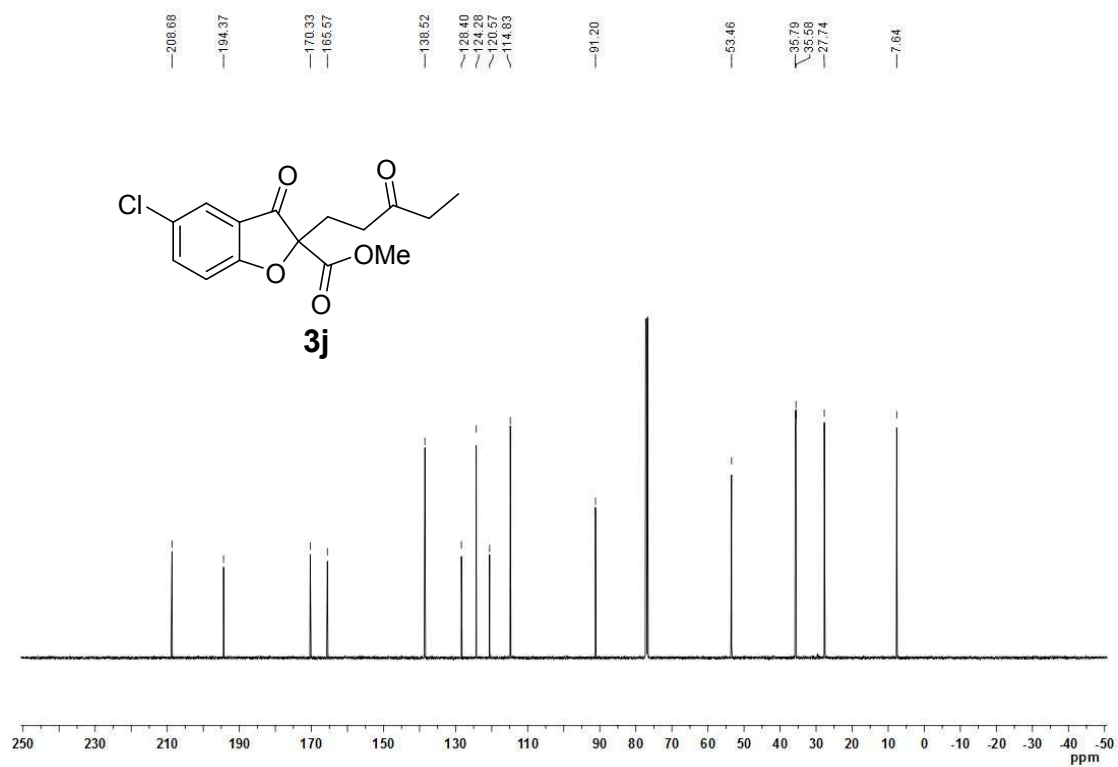
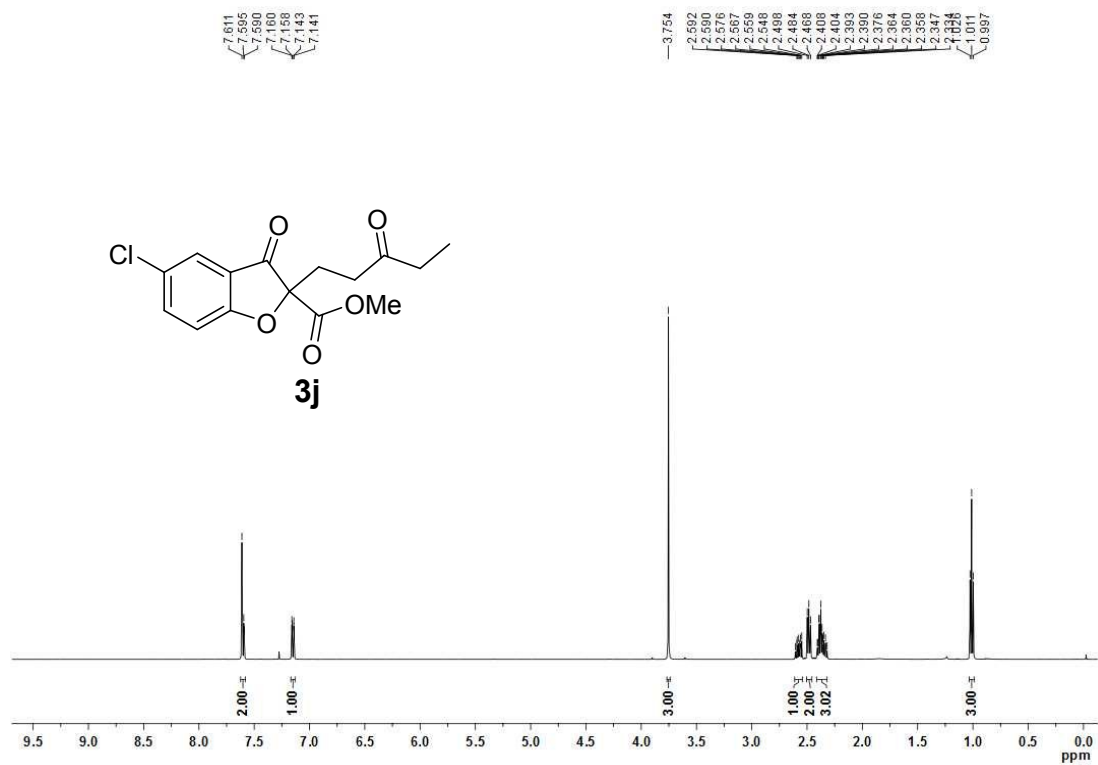


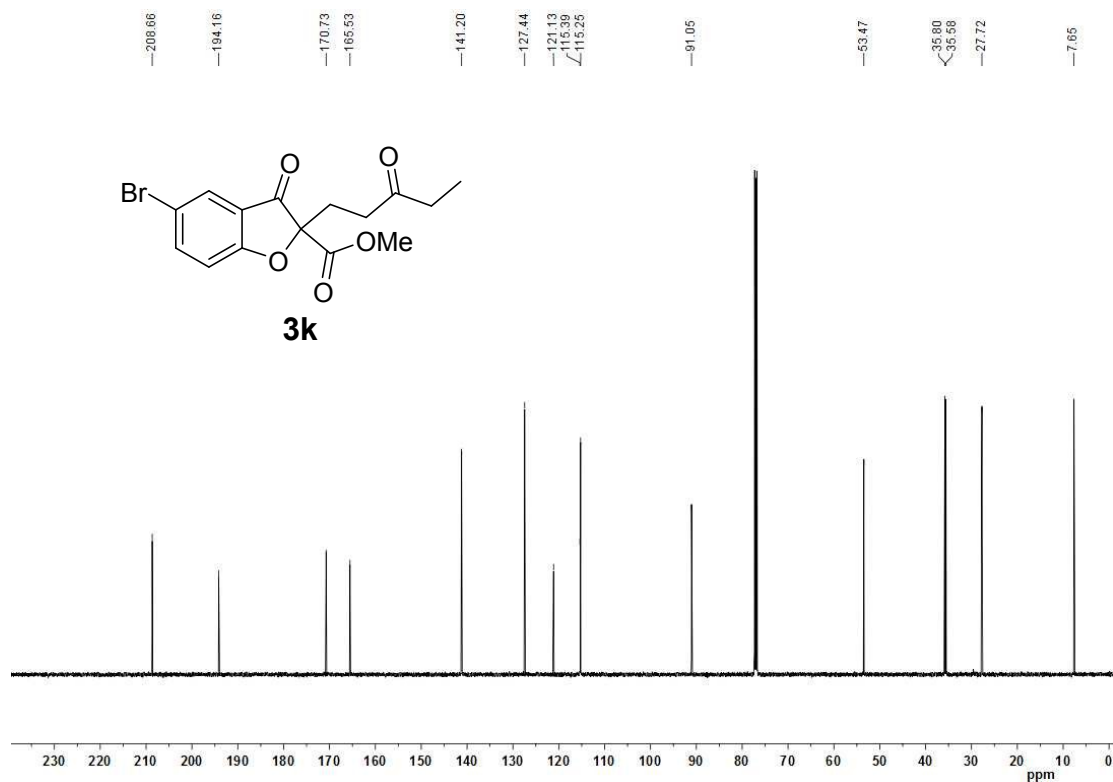
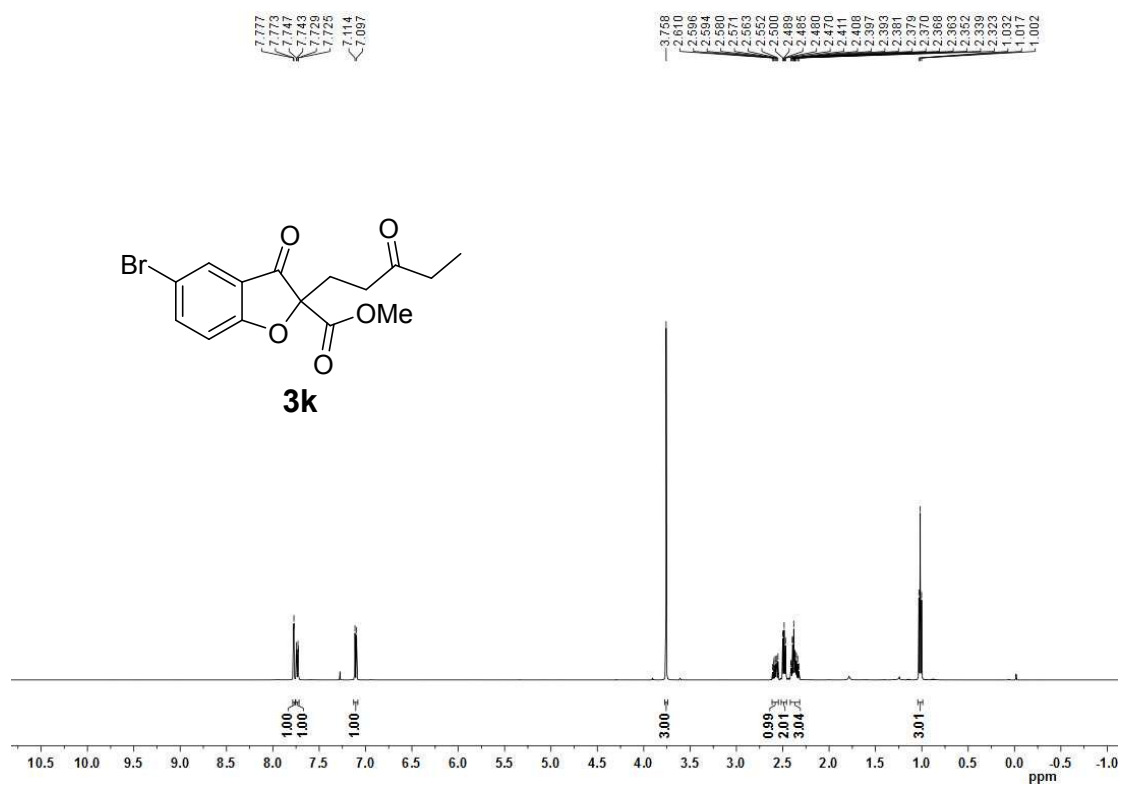


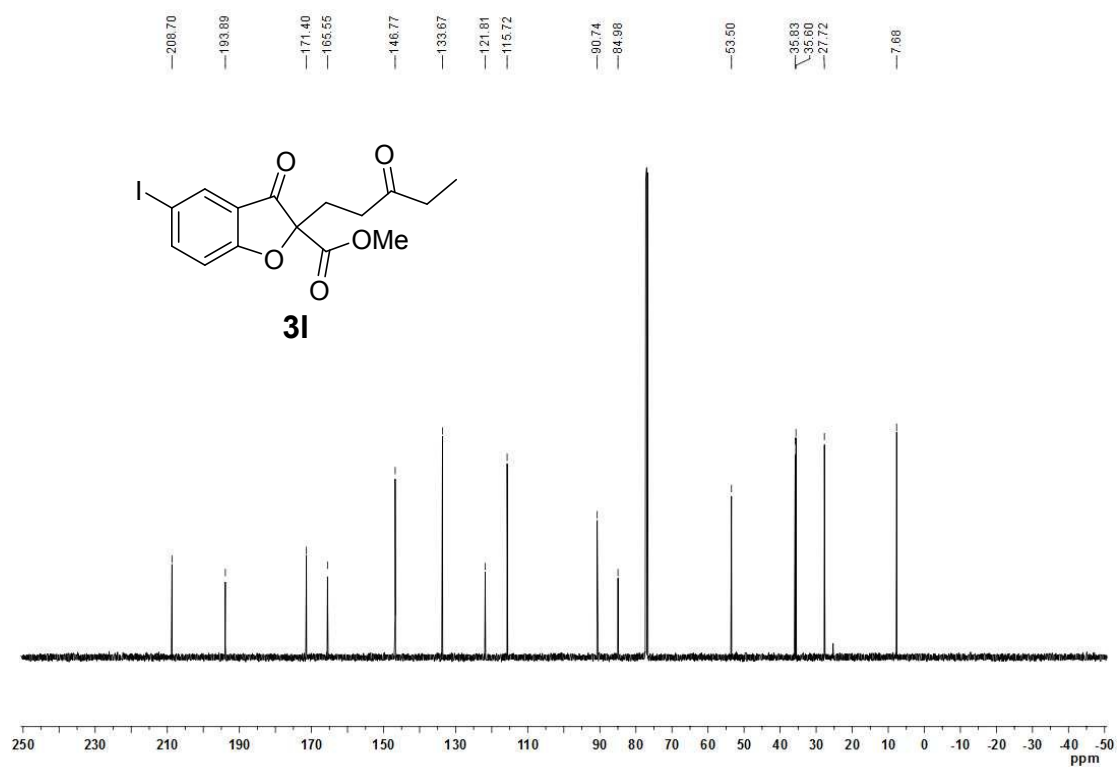
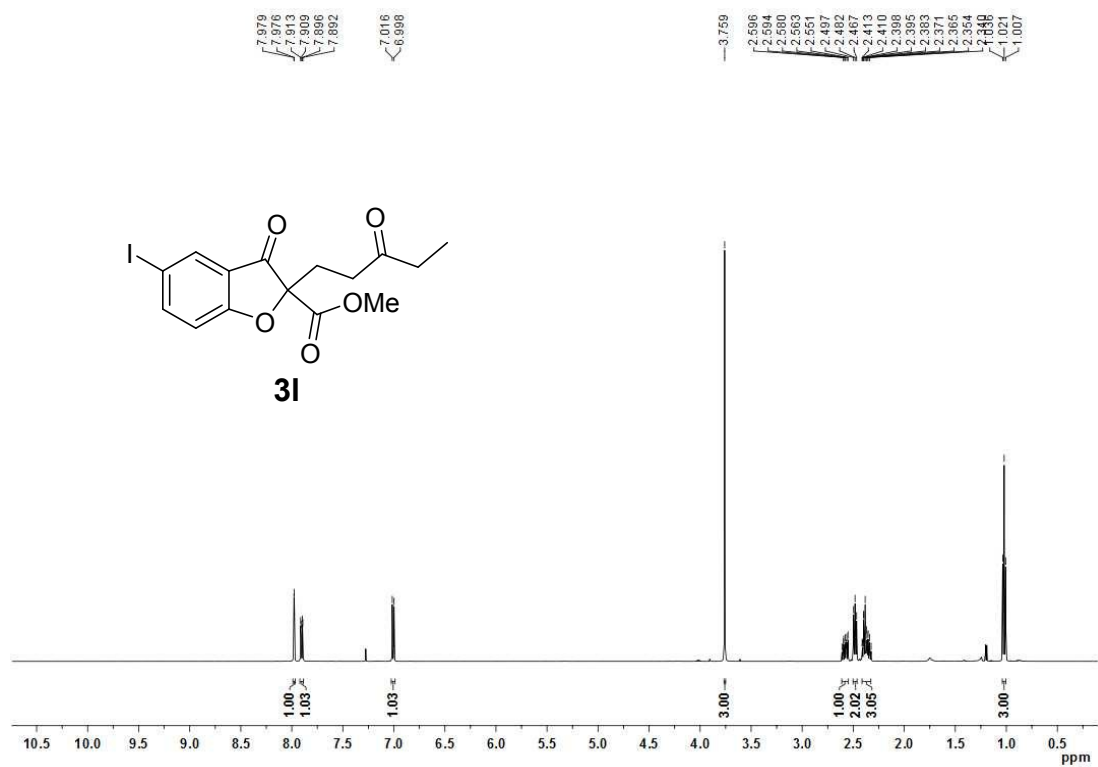


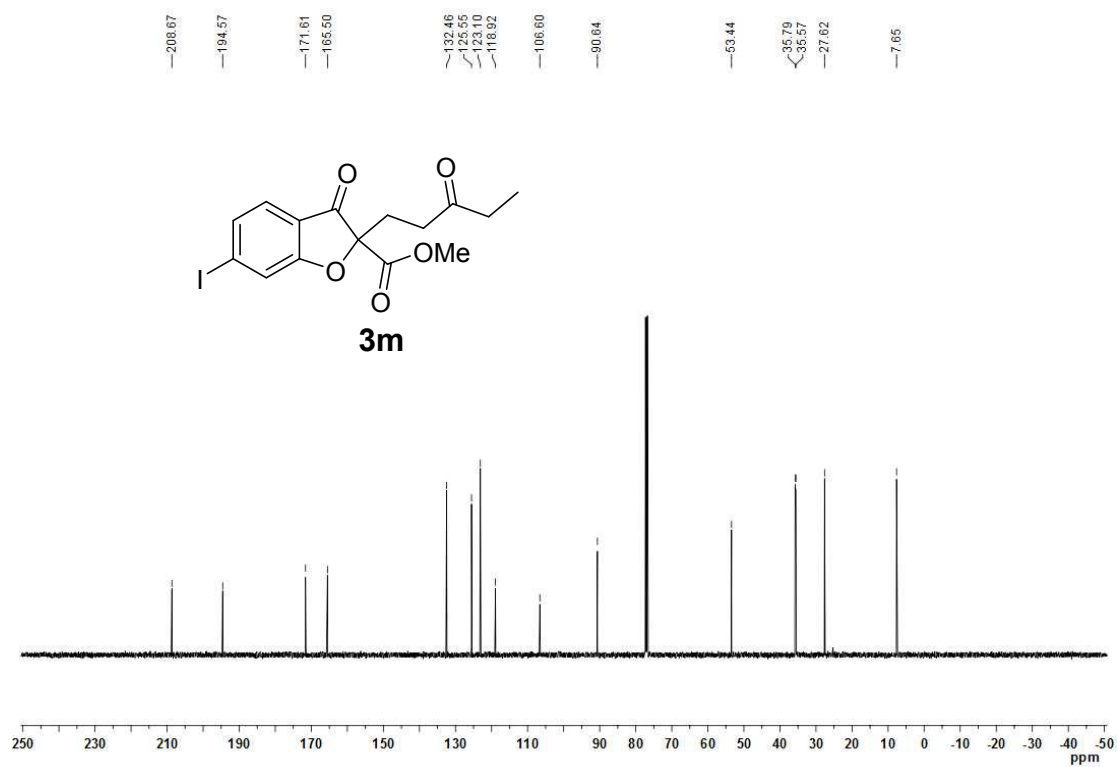
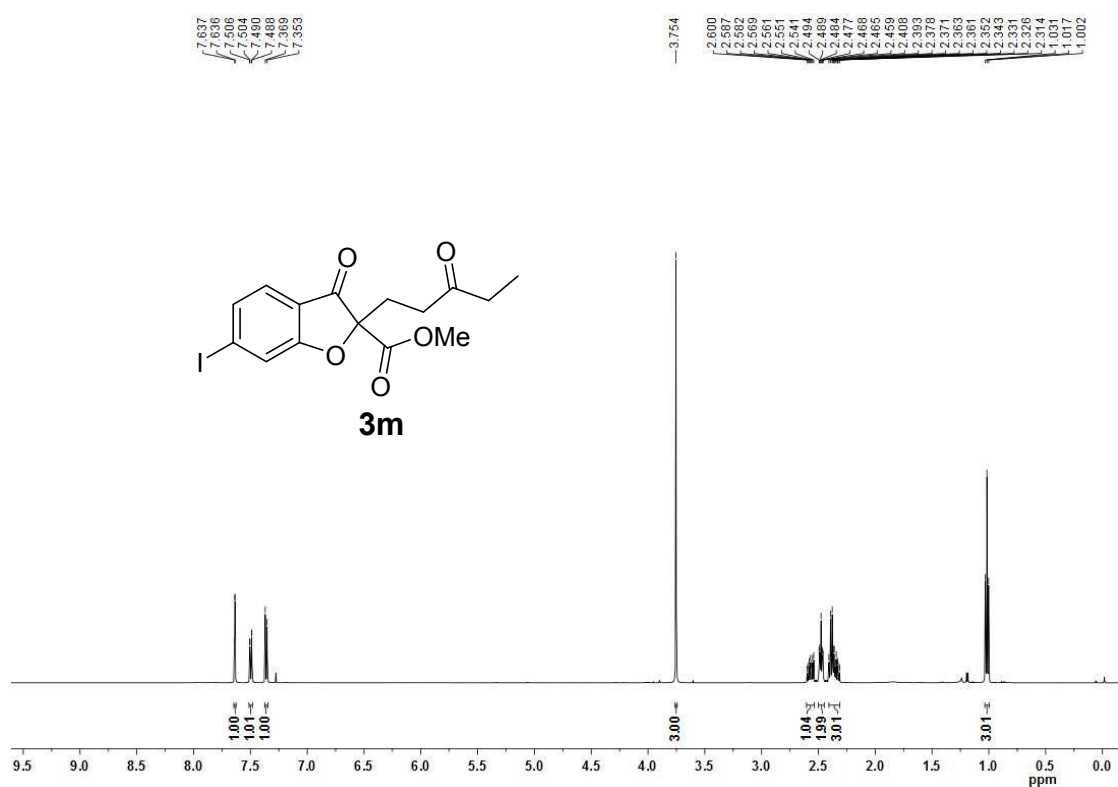


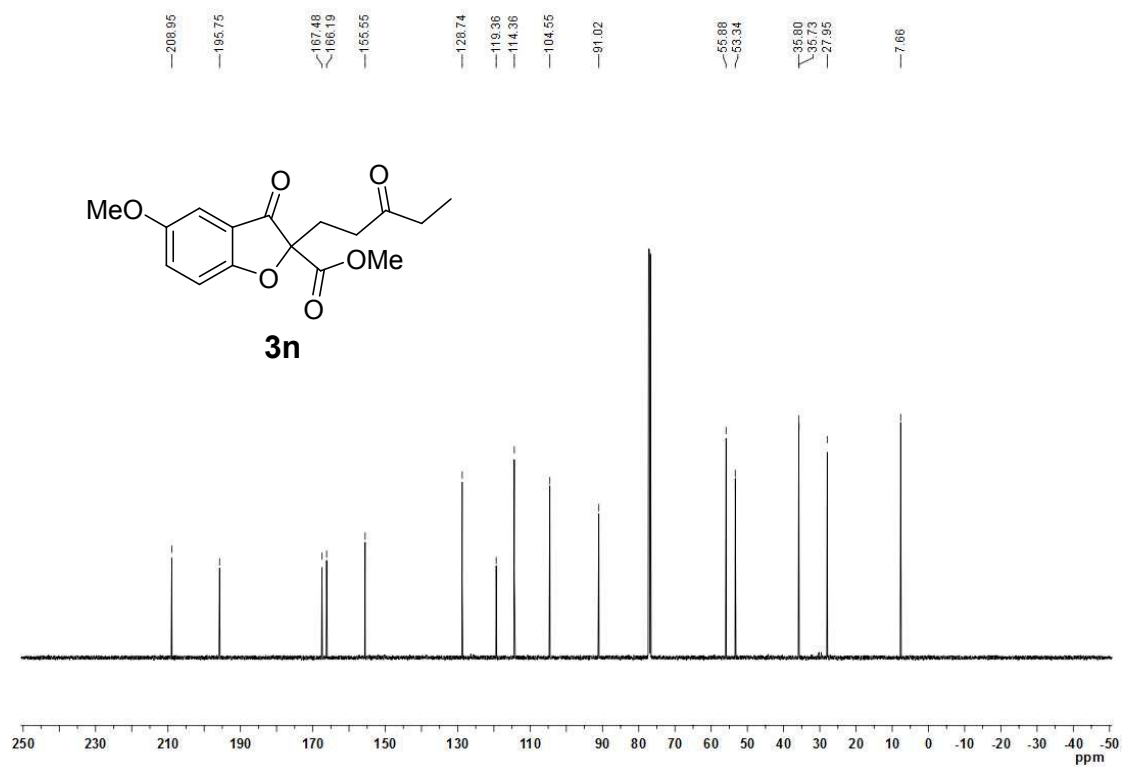
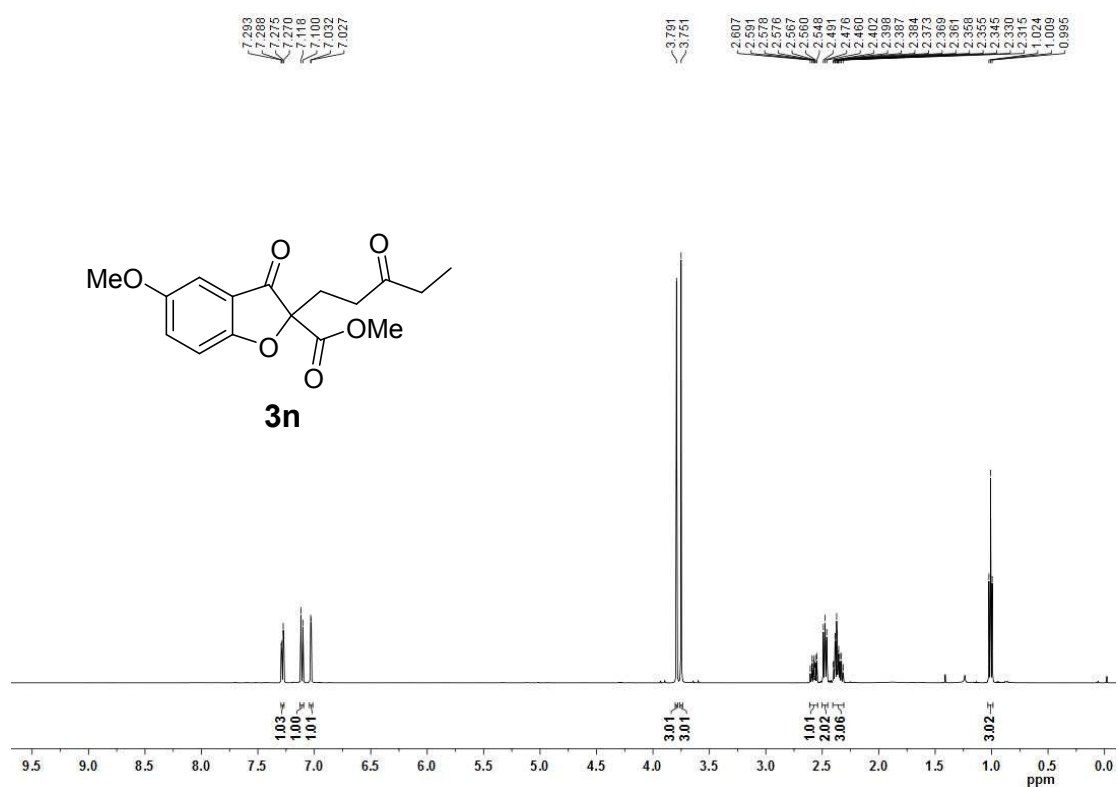


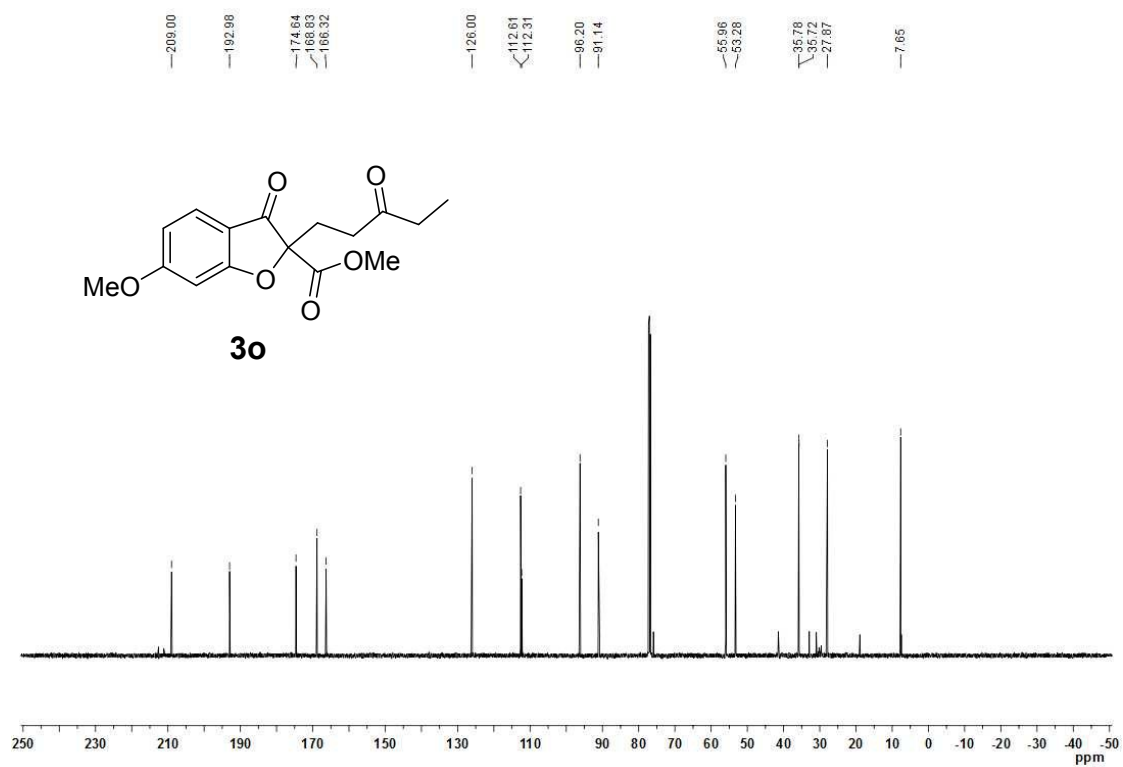
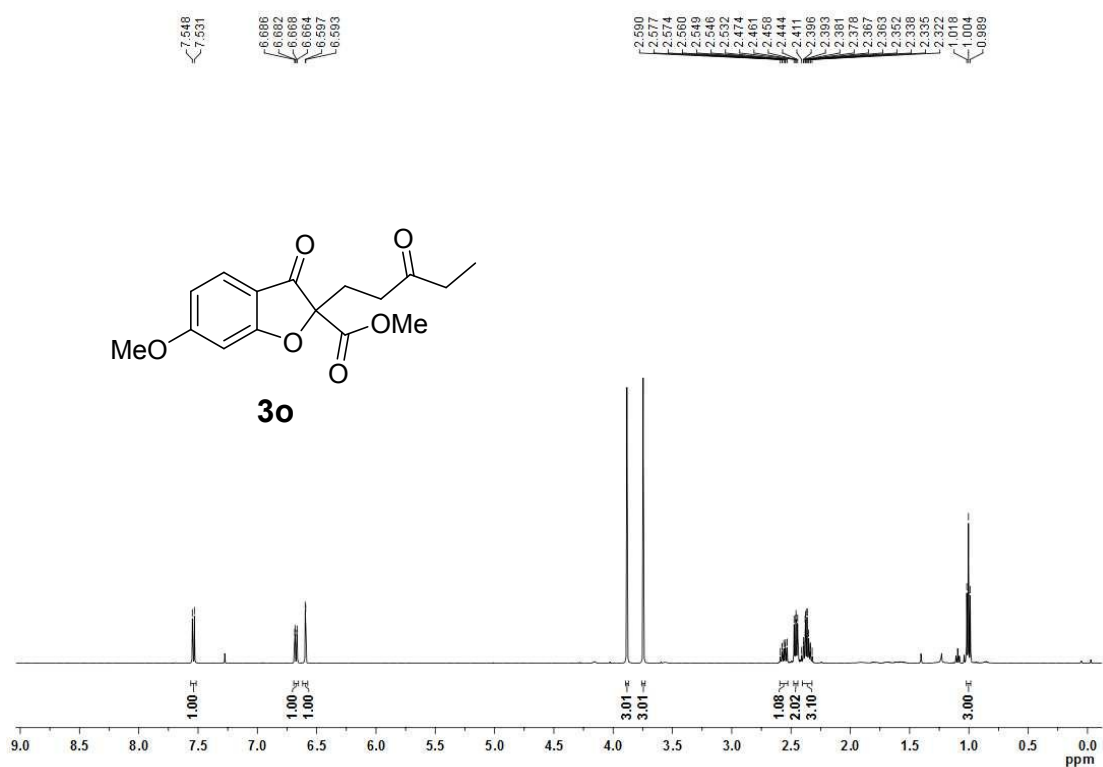


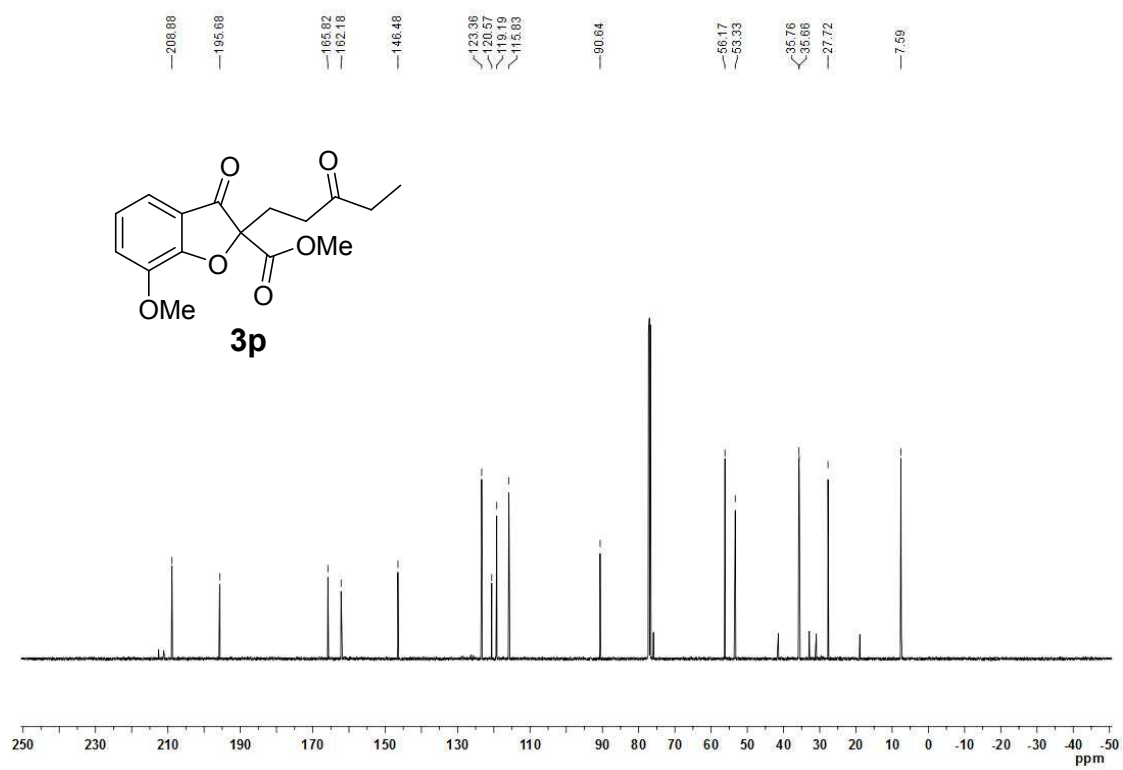
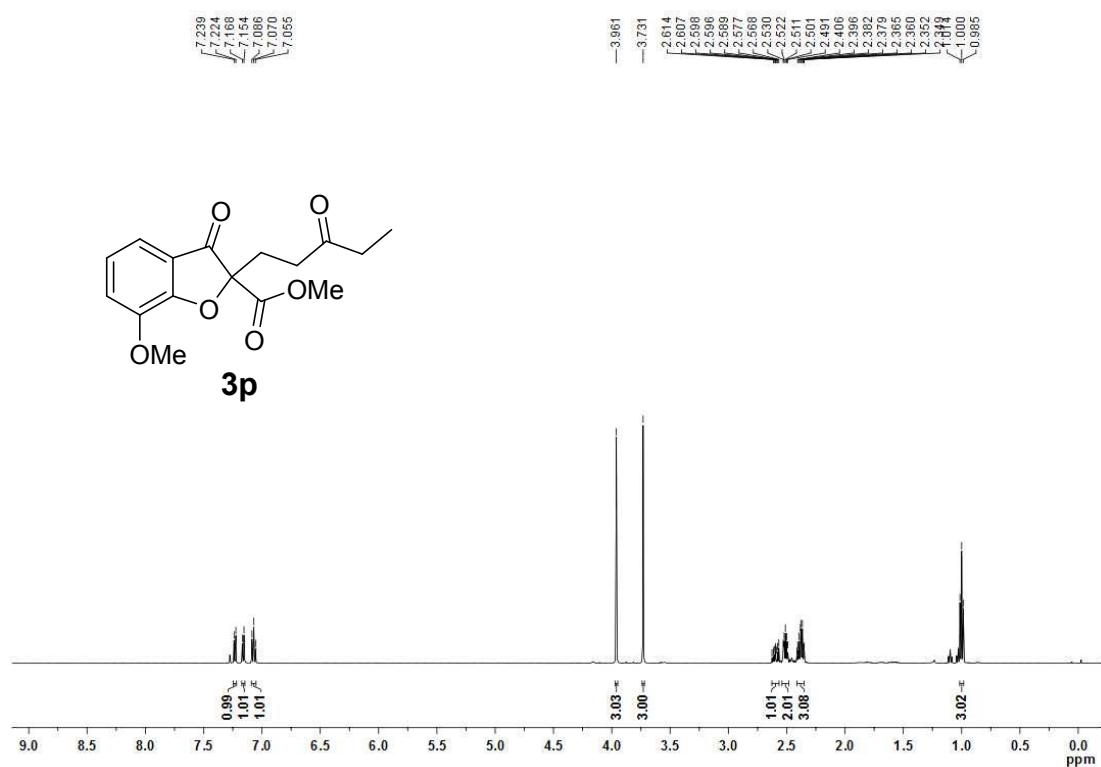


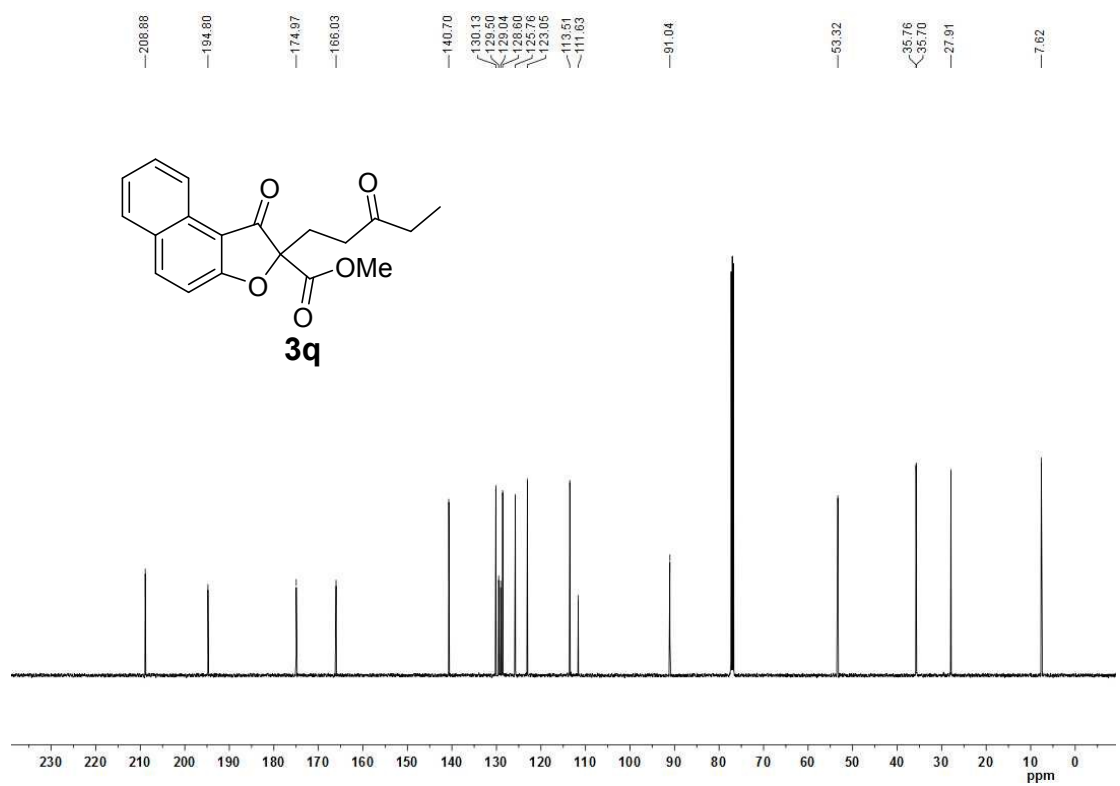
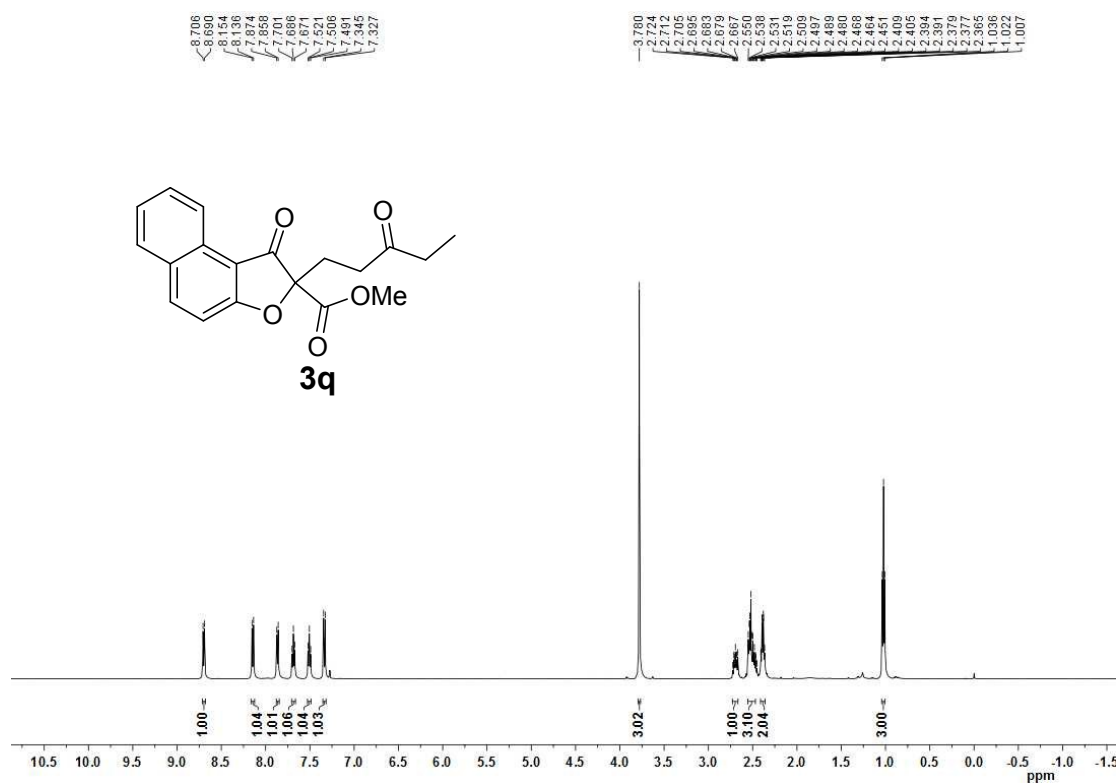


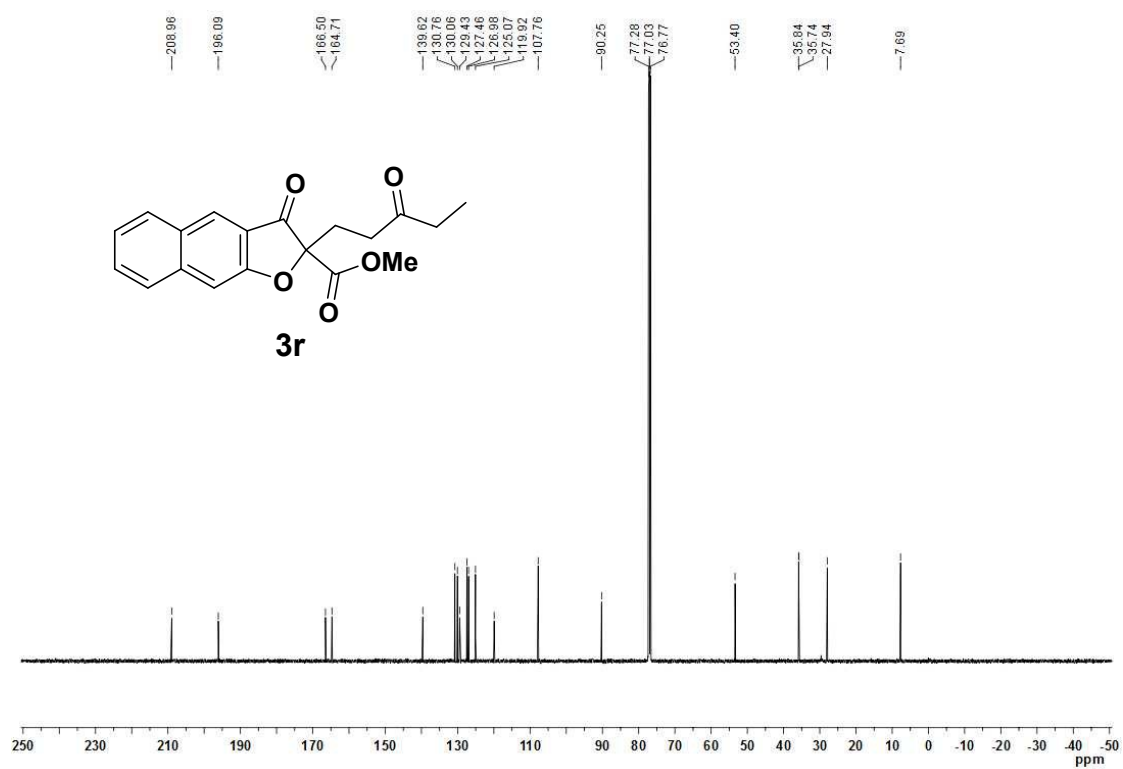
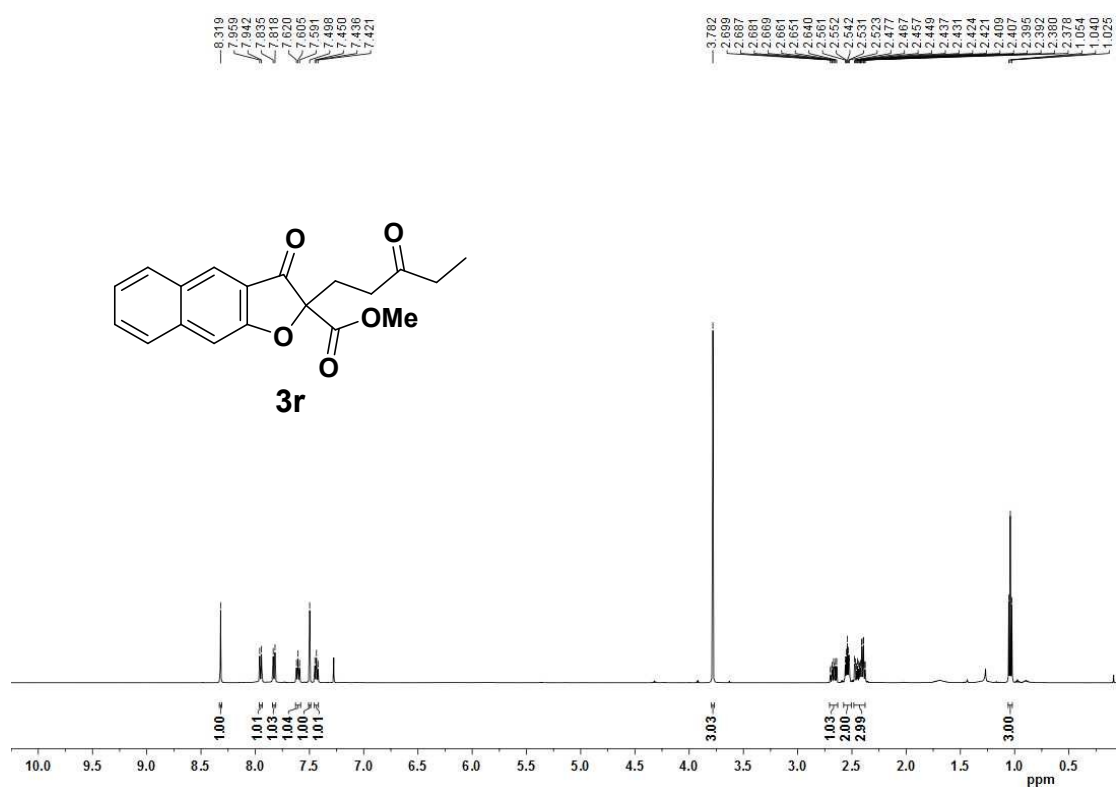


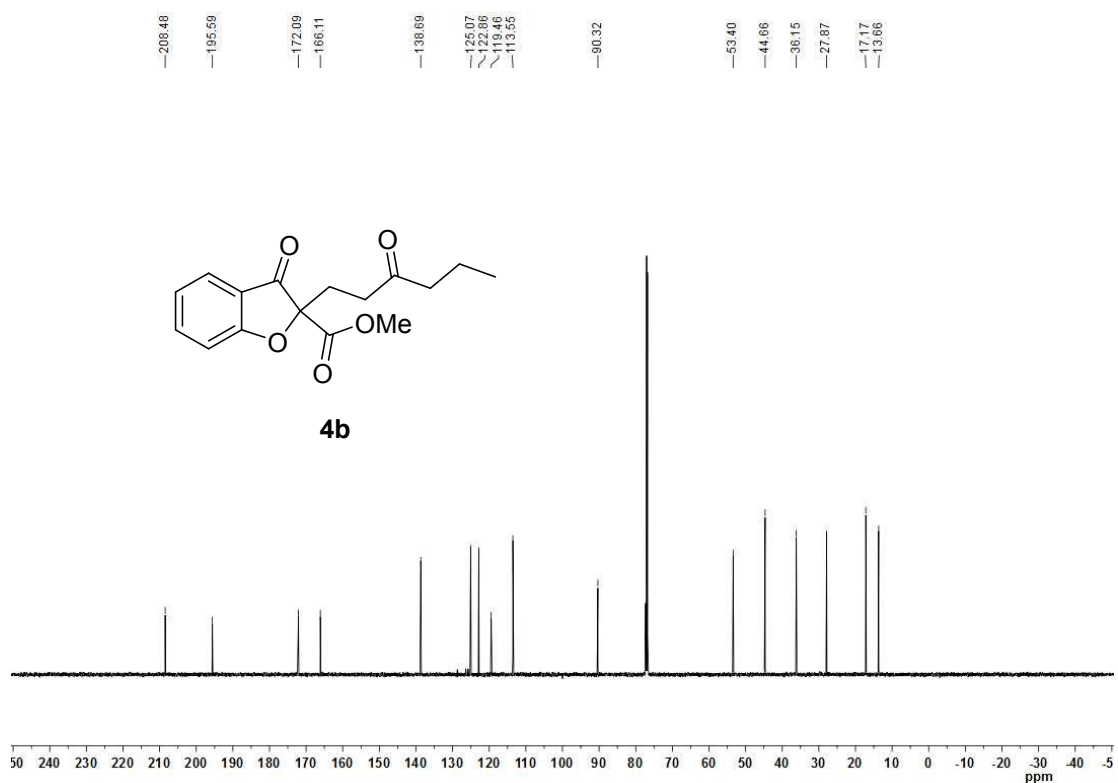
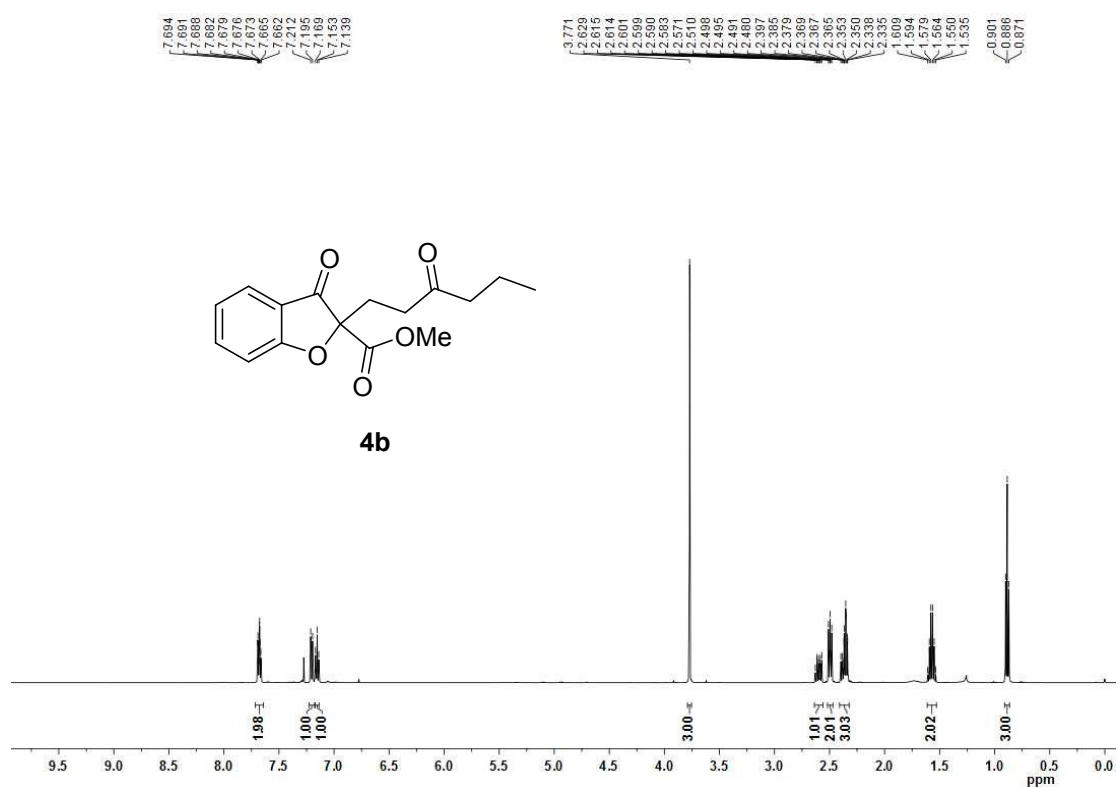


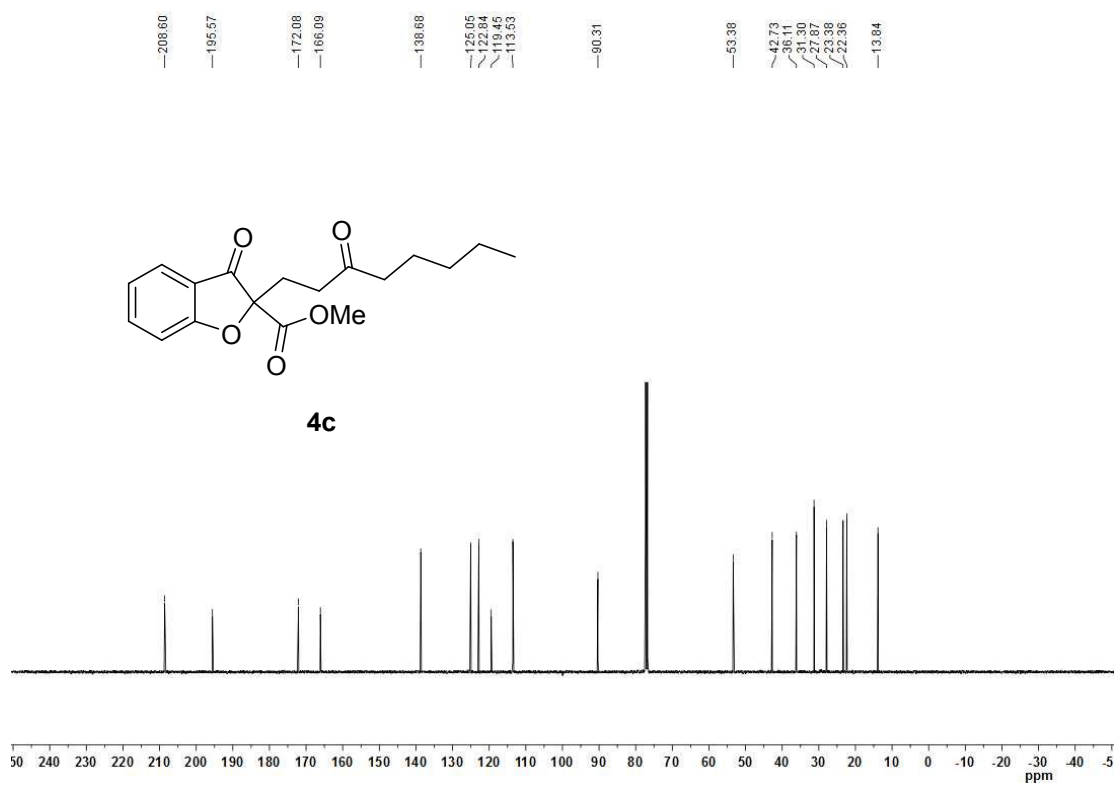
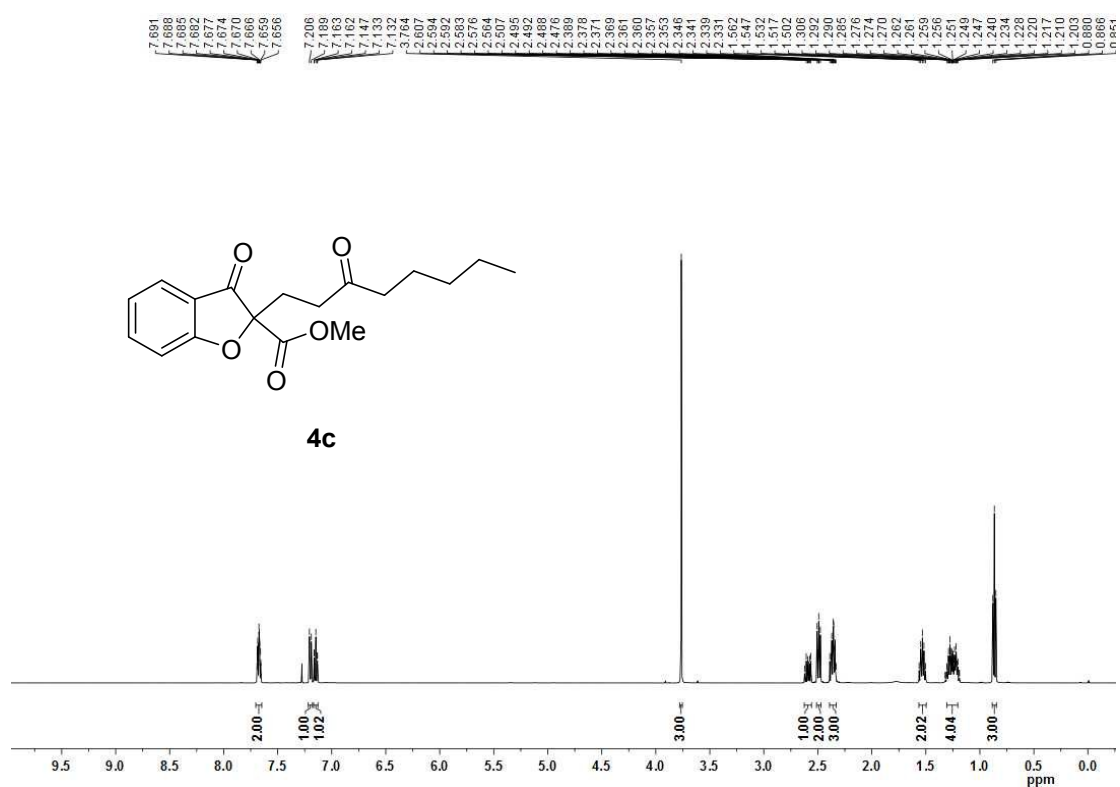


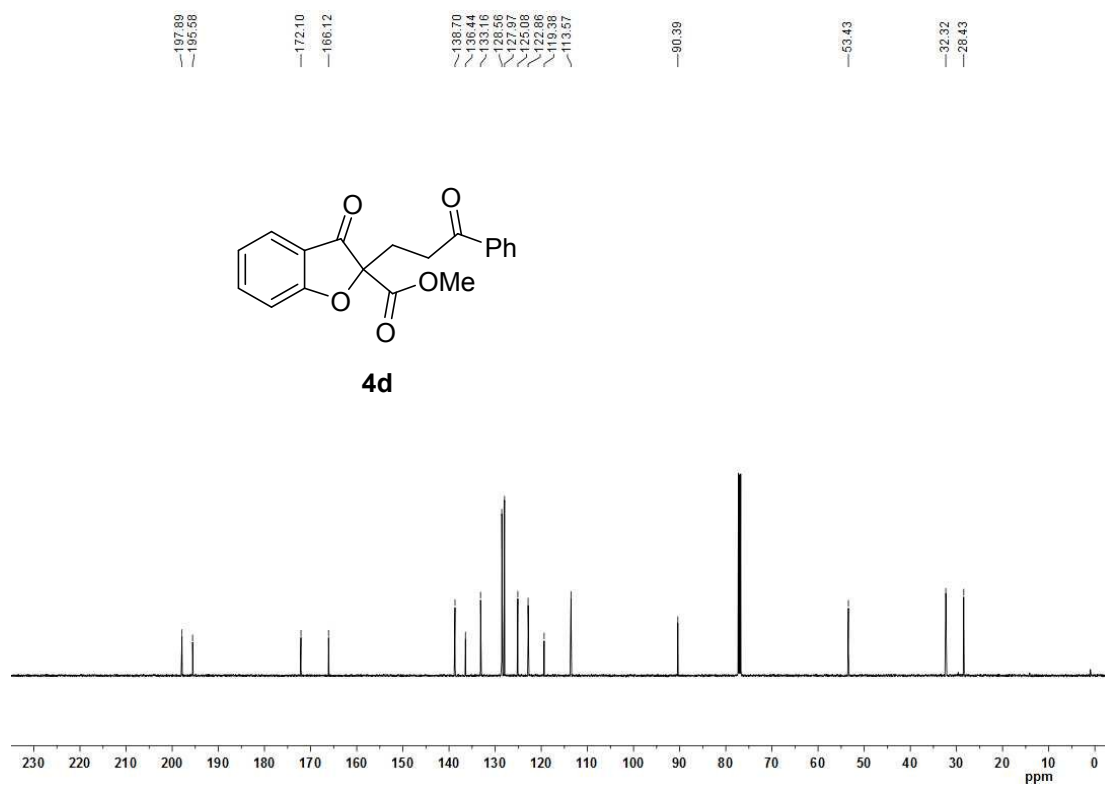
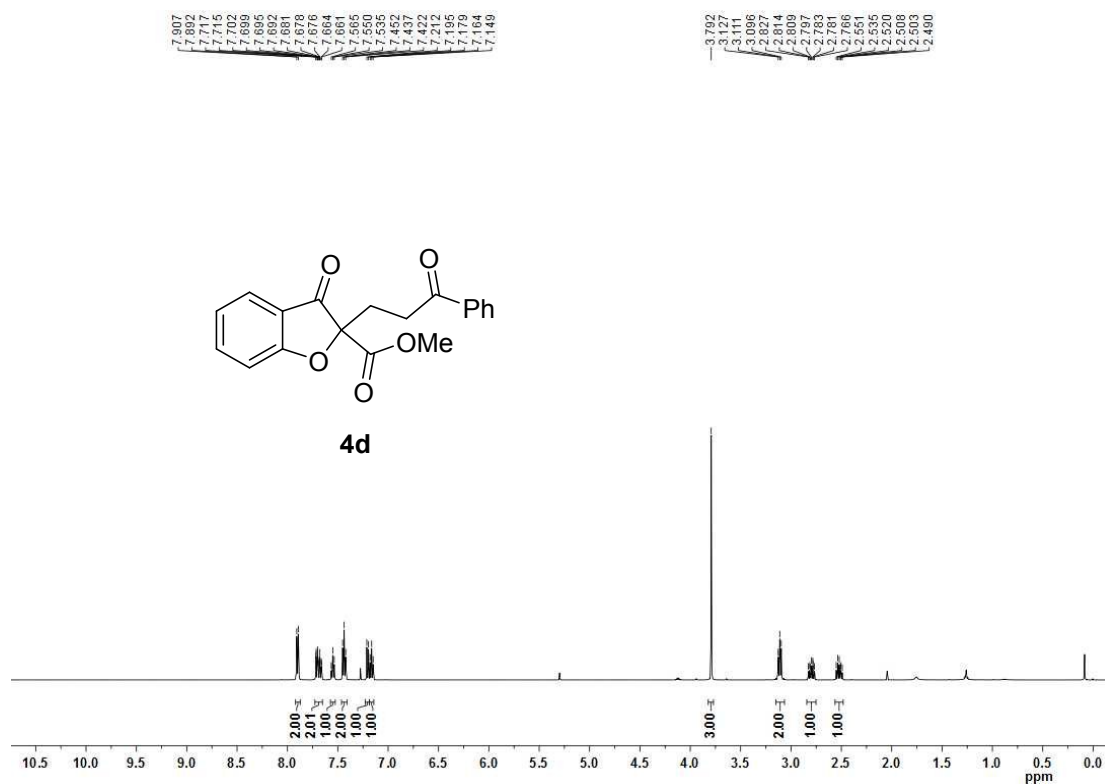


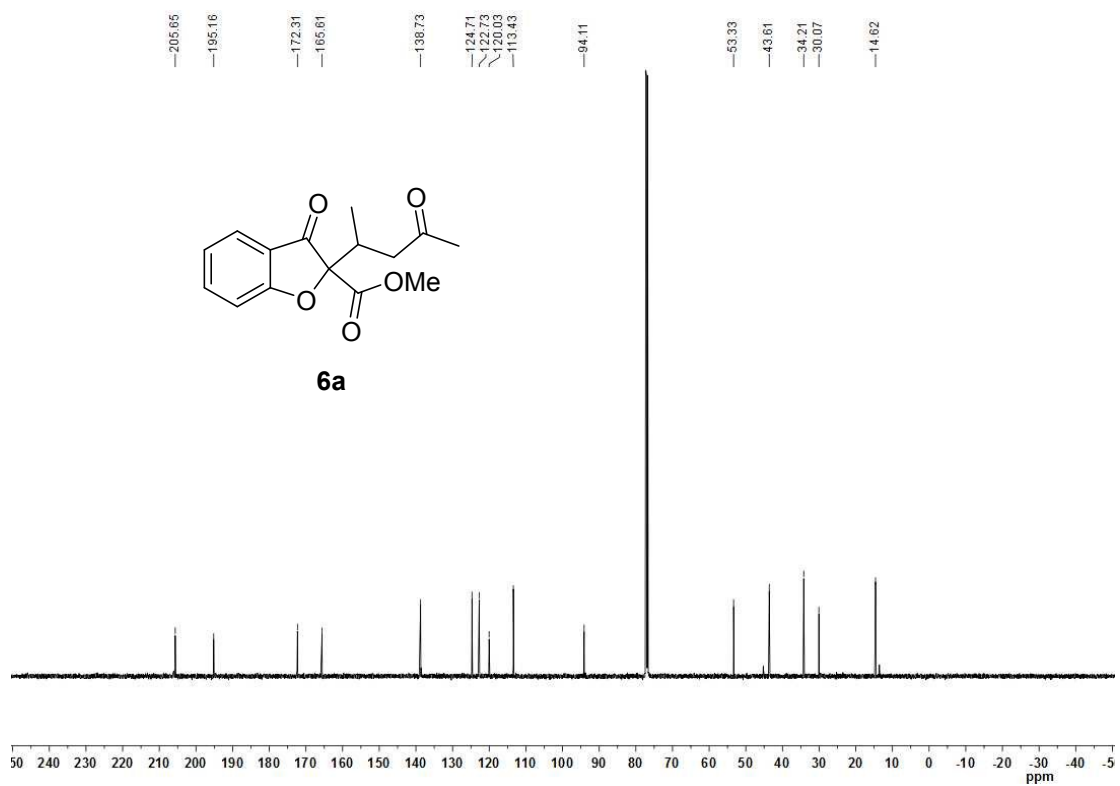
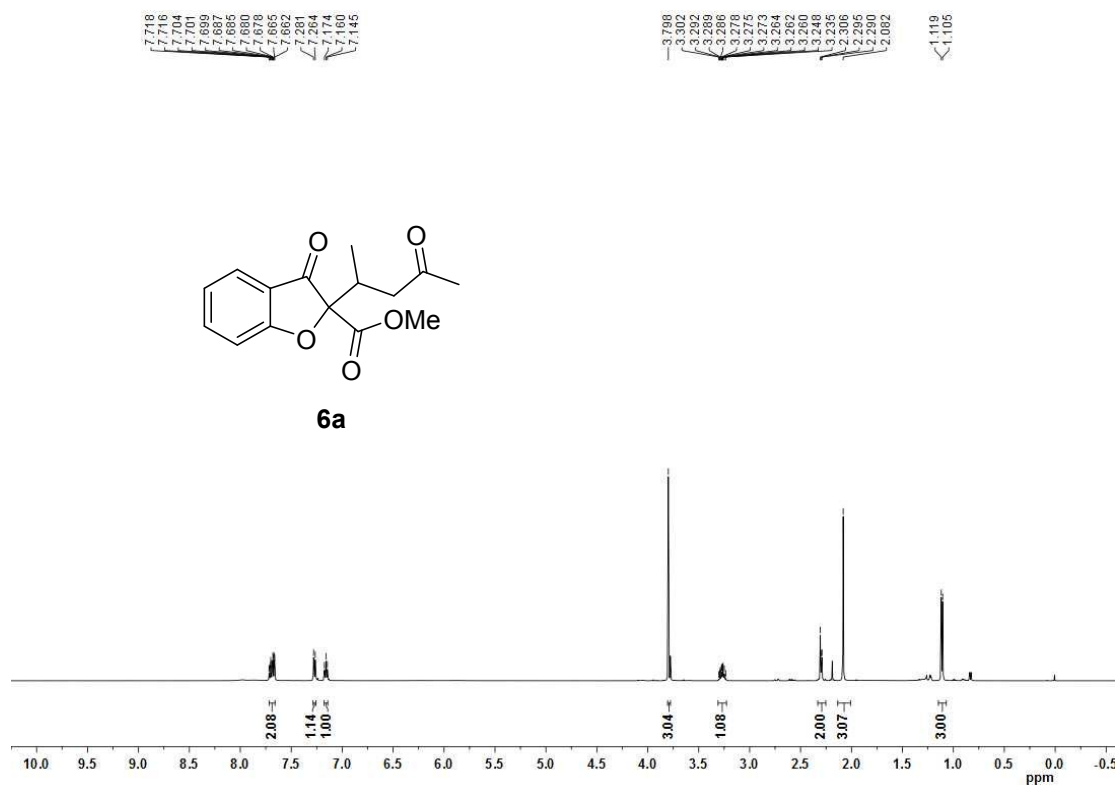


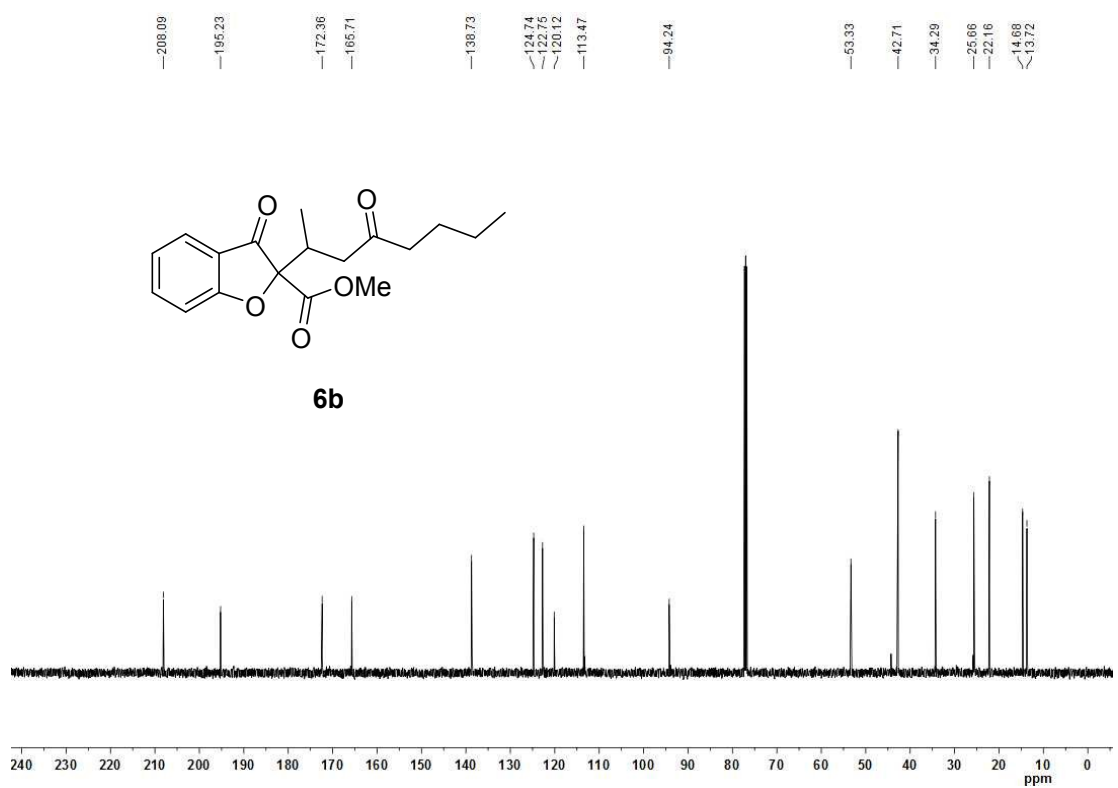
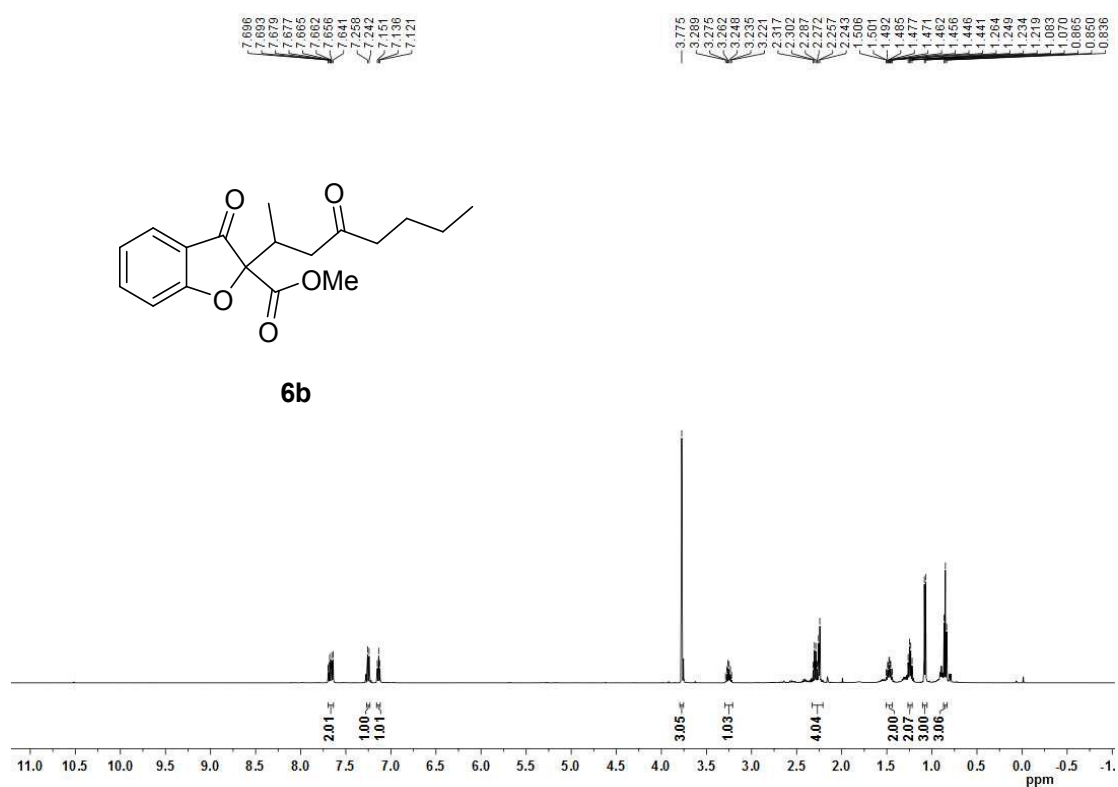


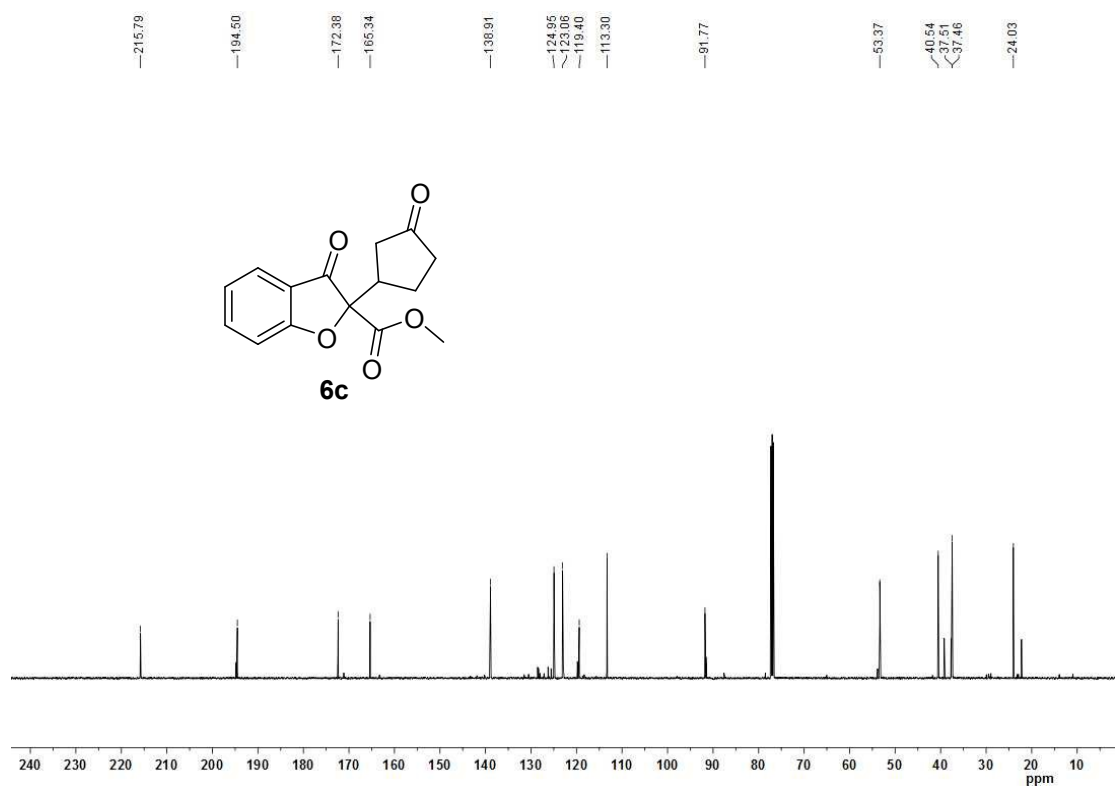
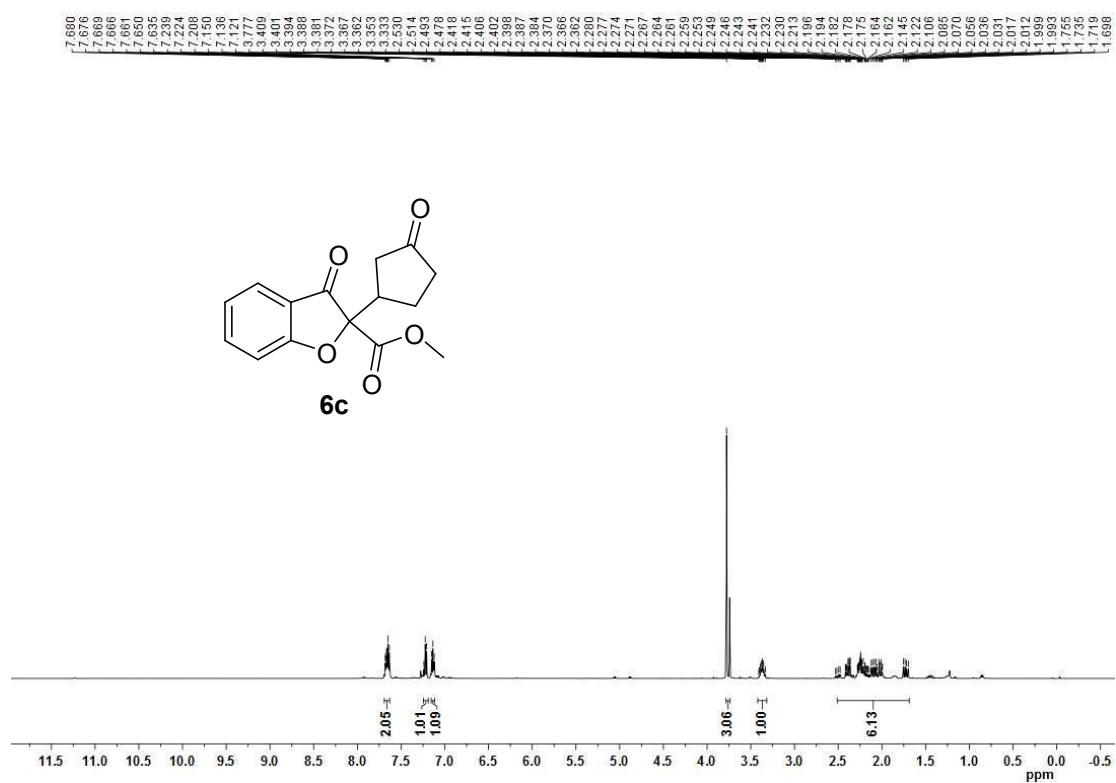


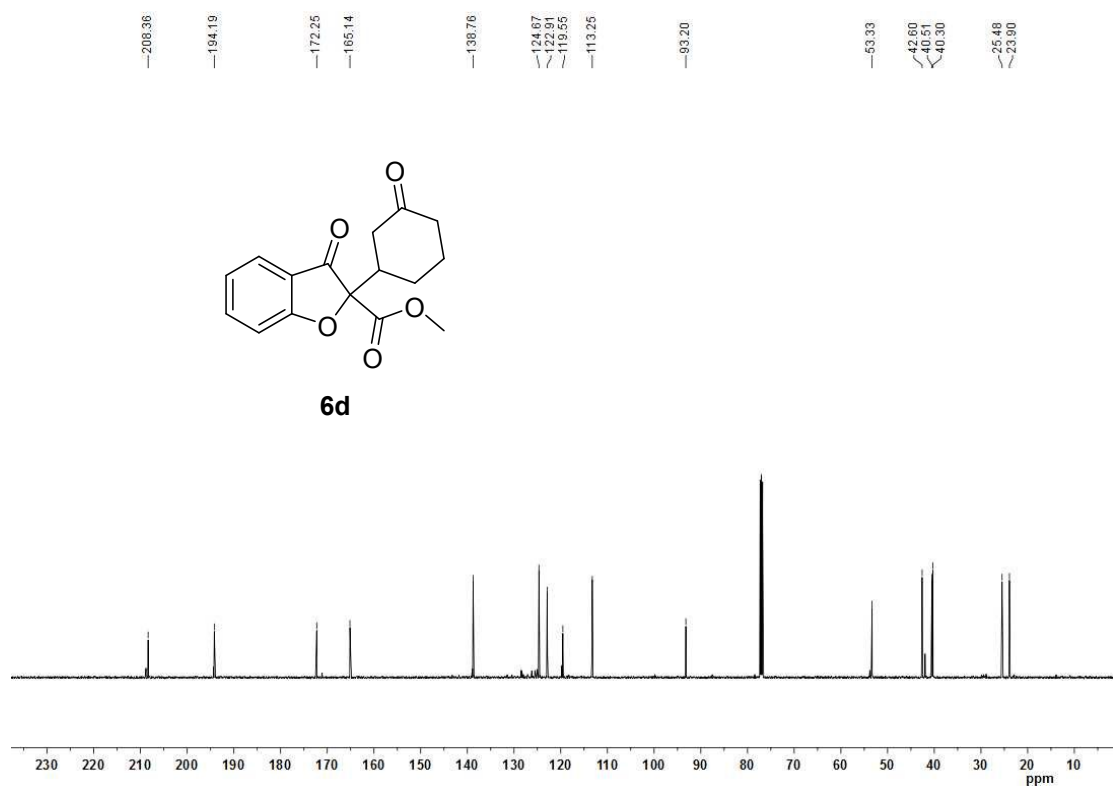
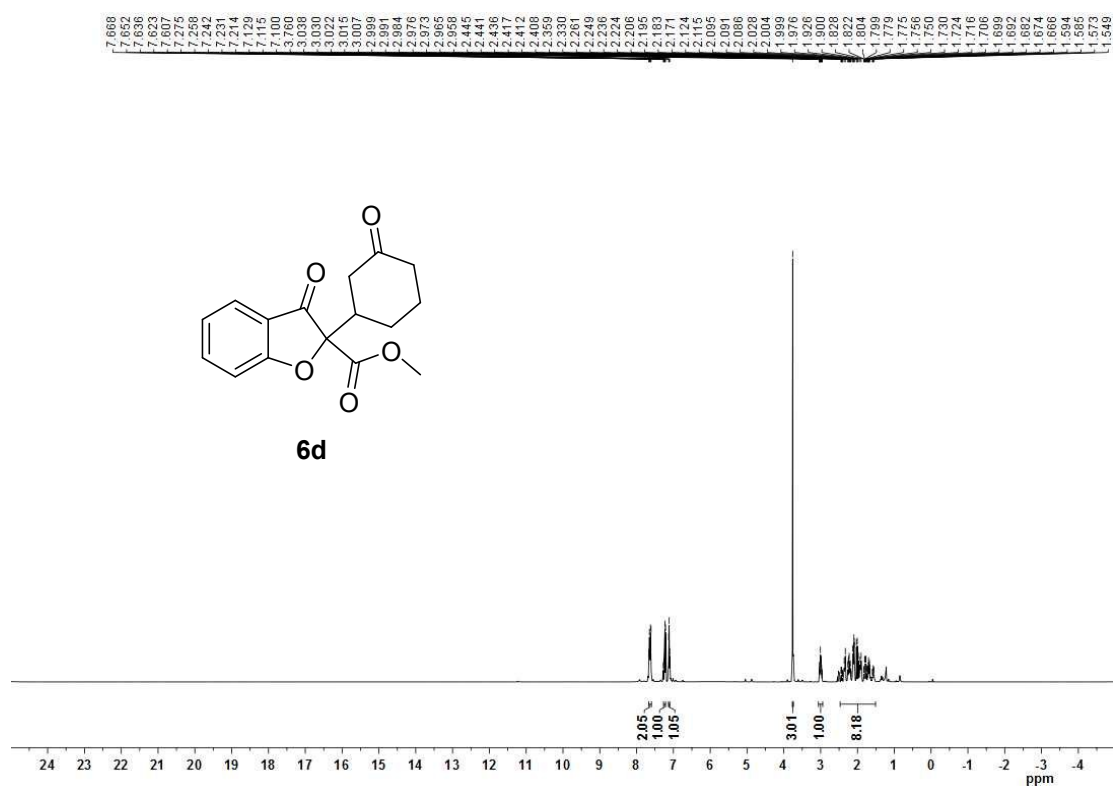


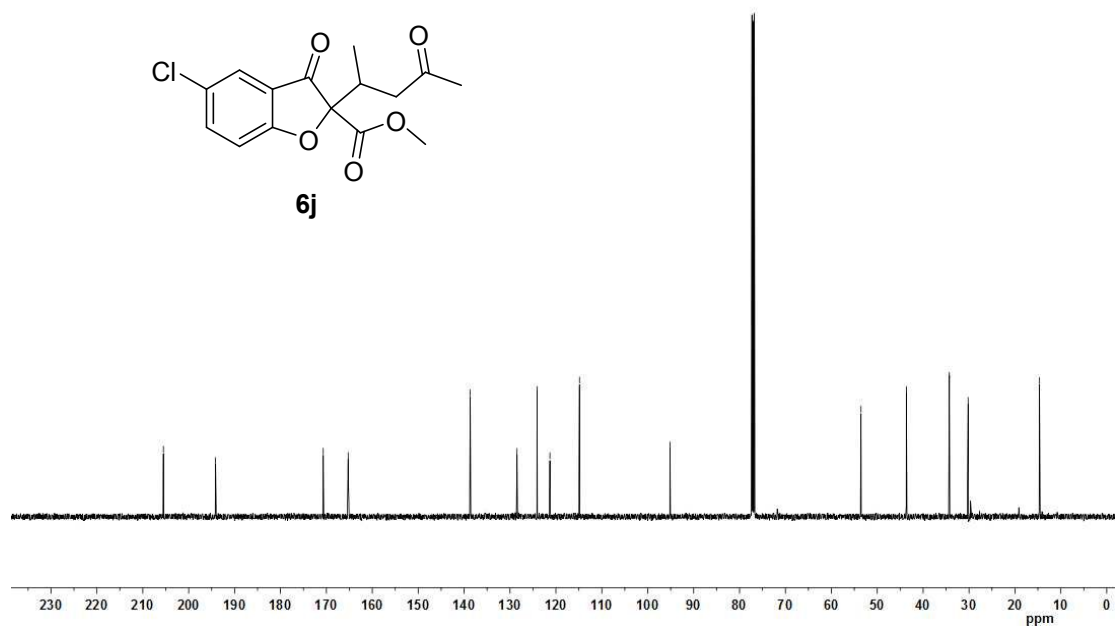
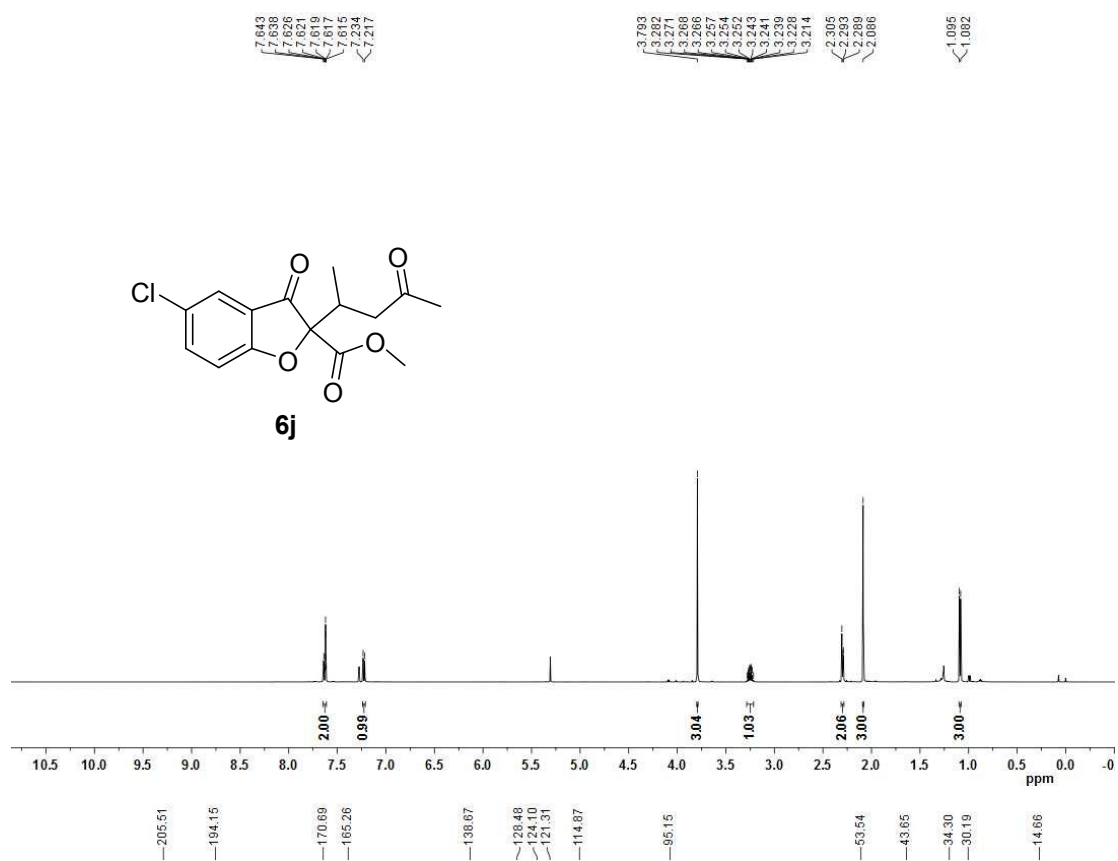


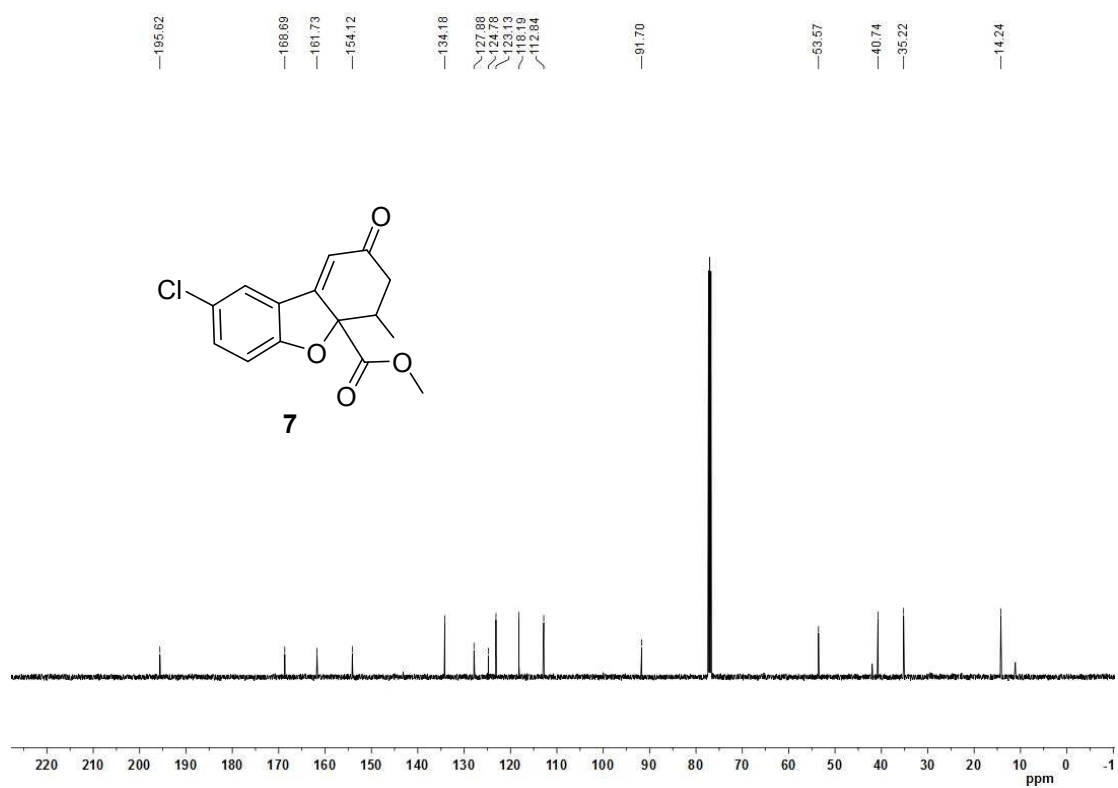
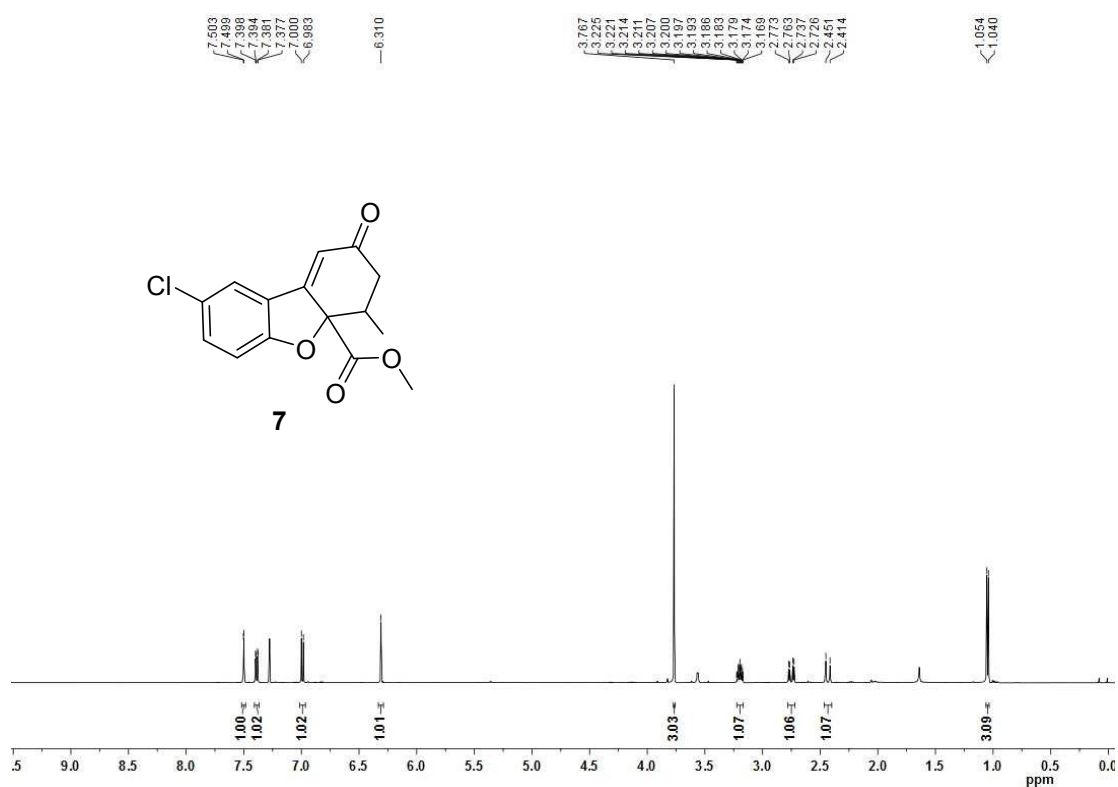




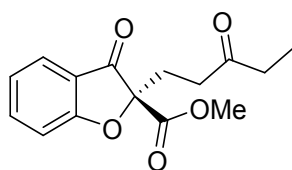




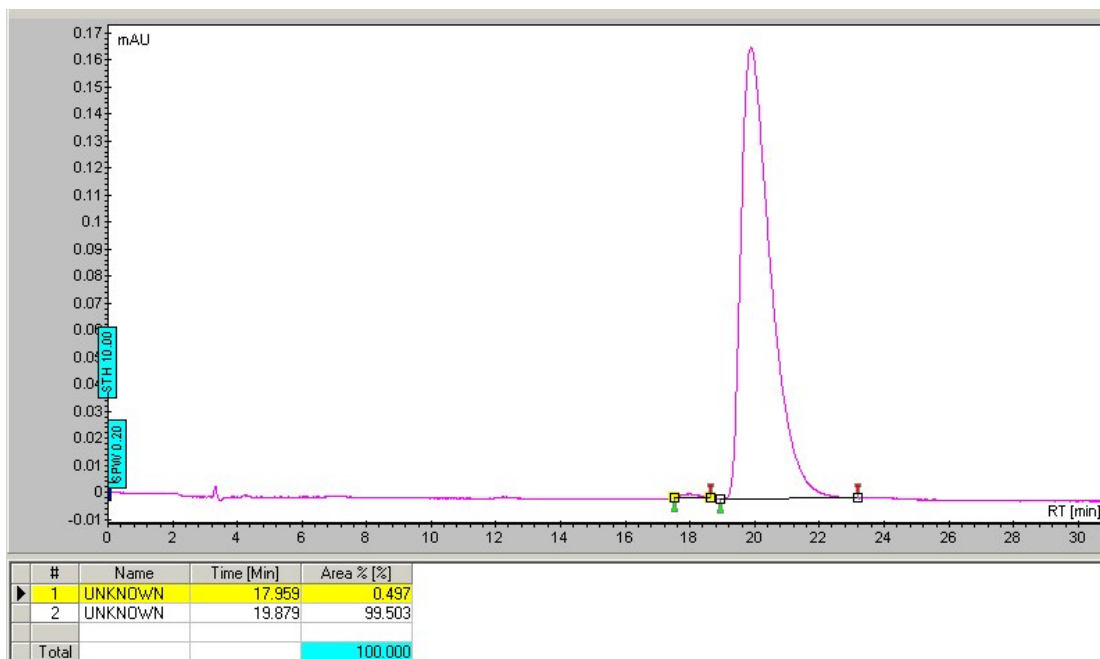
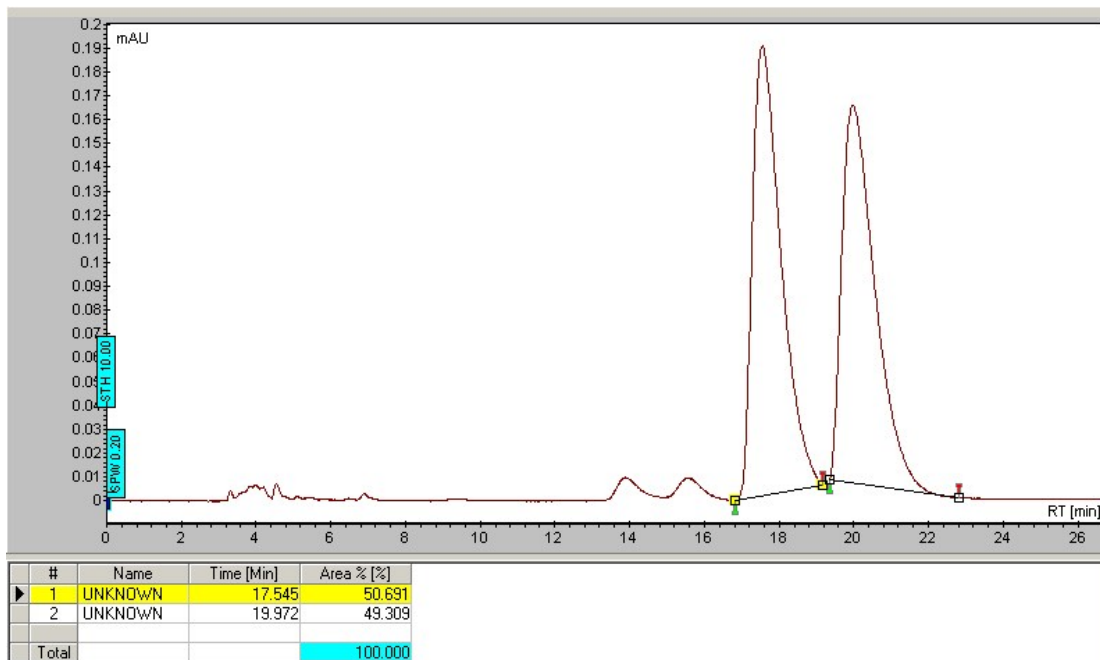


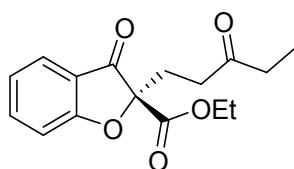


3、Chiral HPLC data of all products

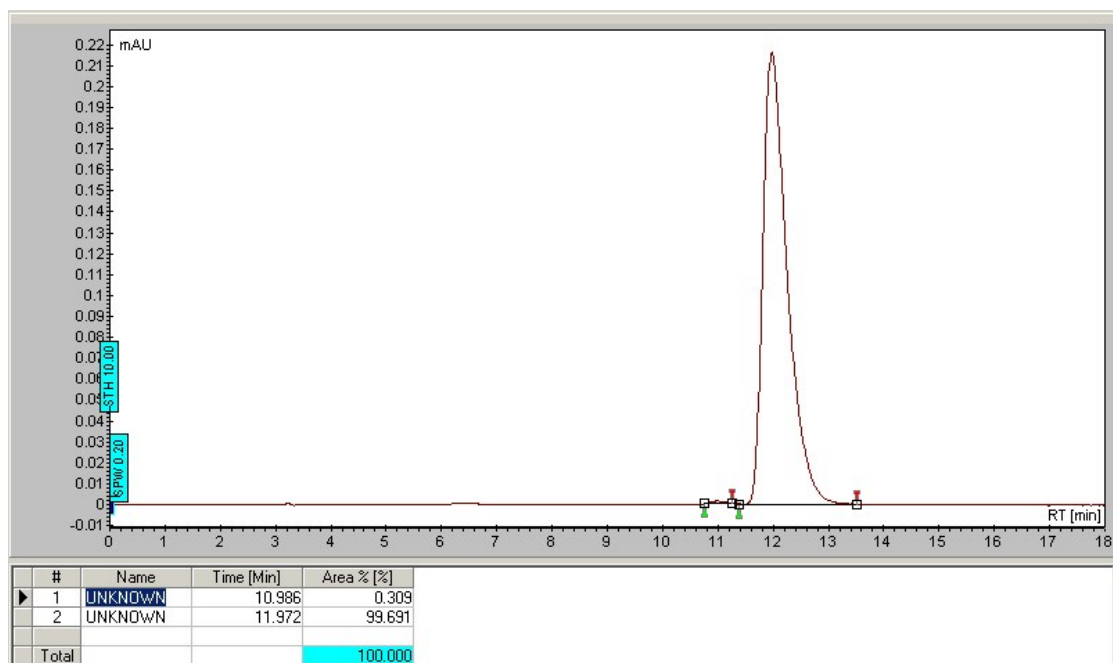
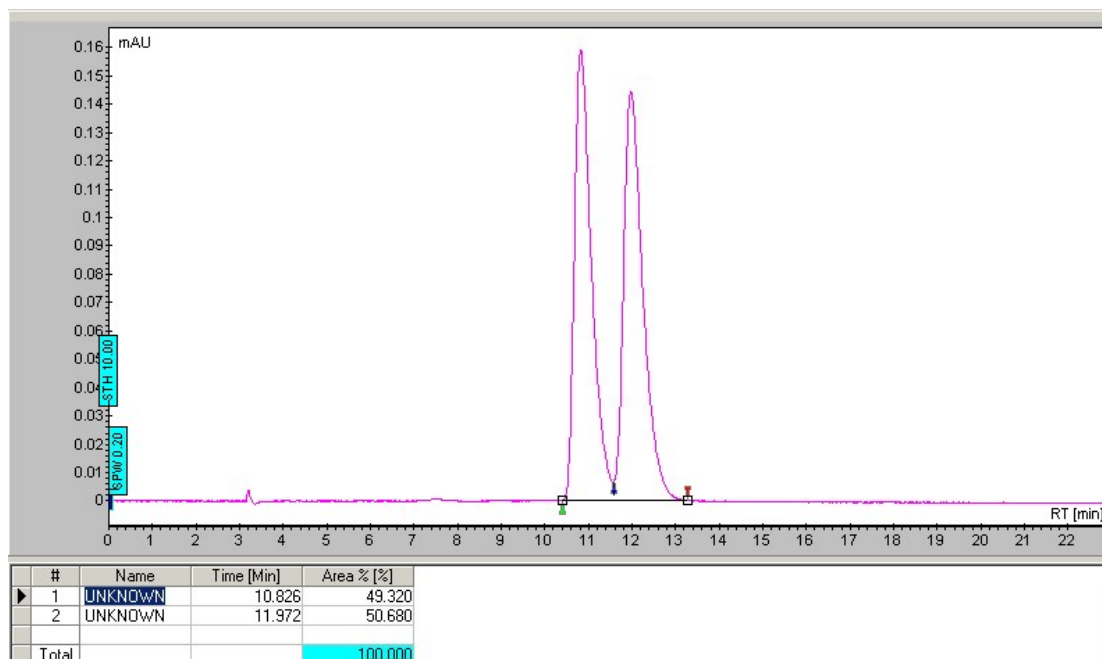


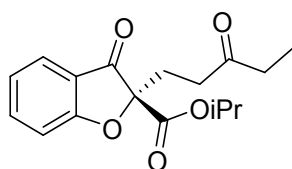
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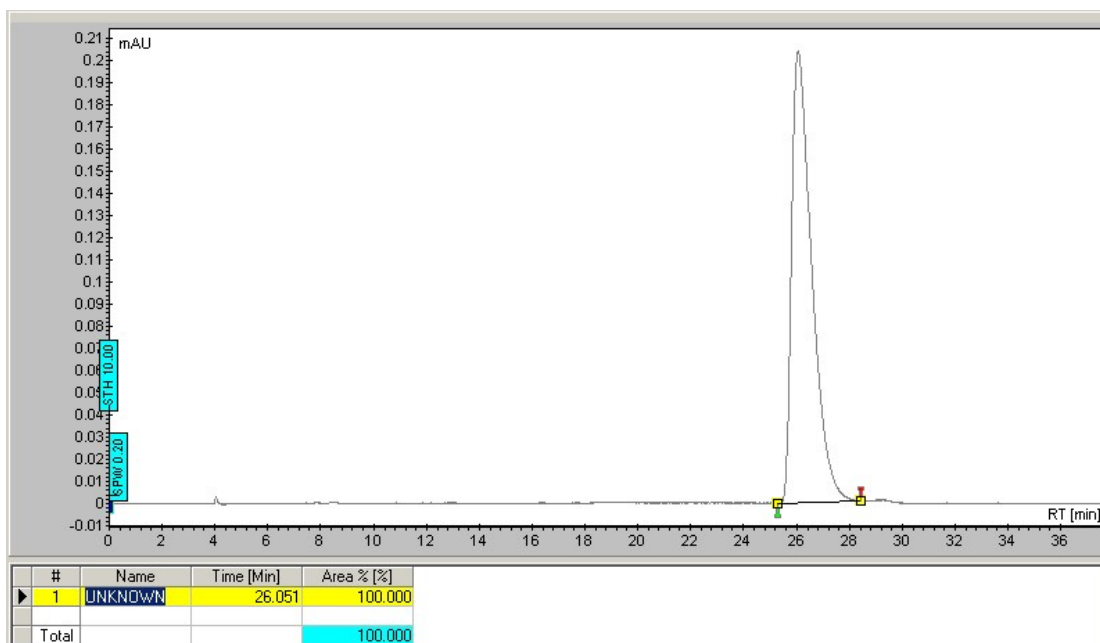
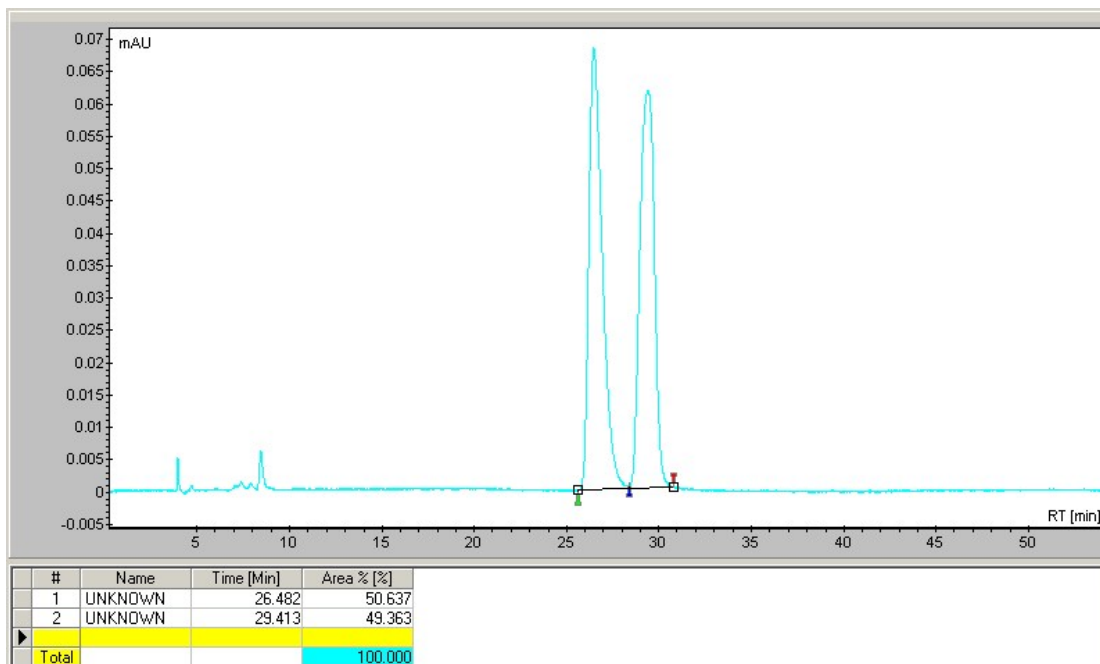


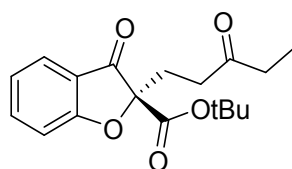
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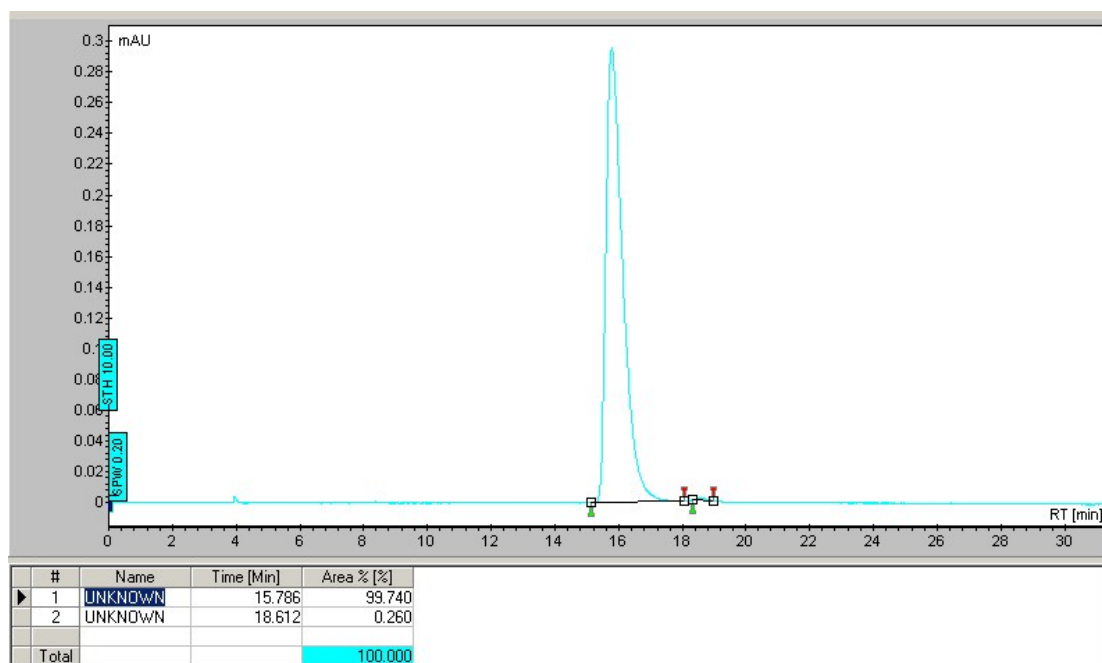
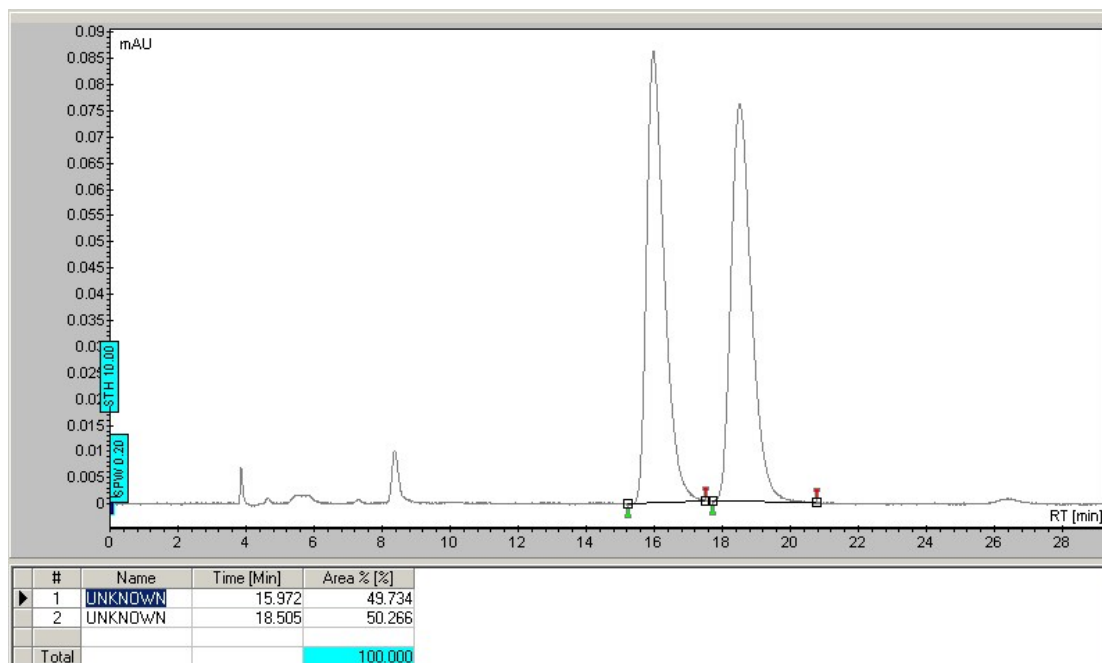


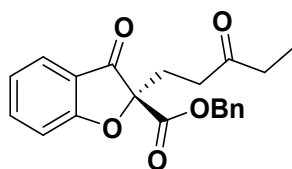
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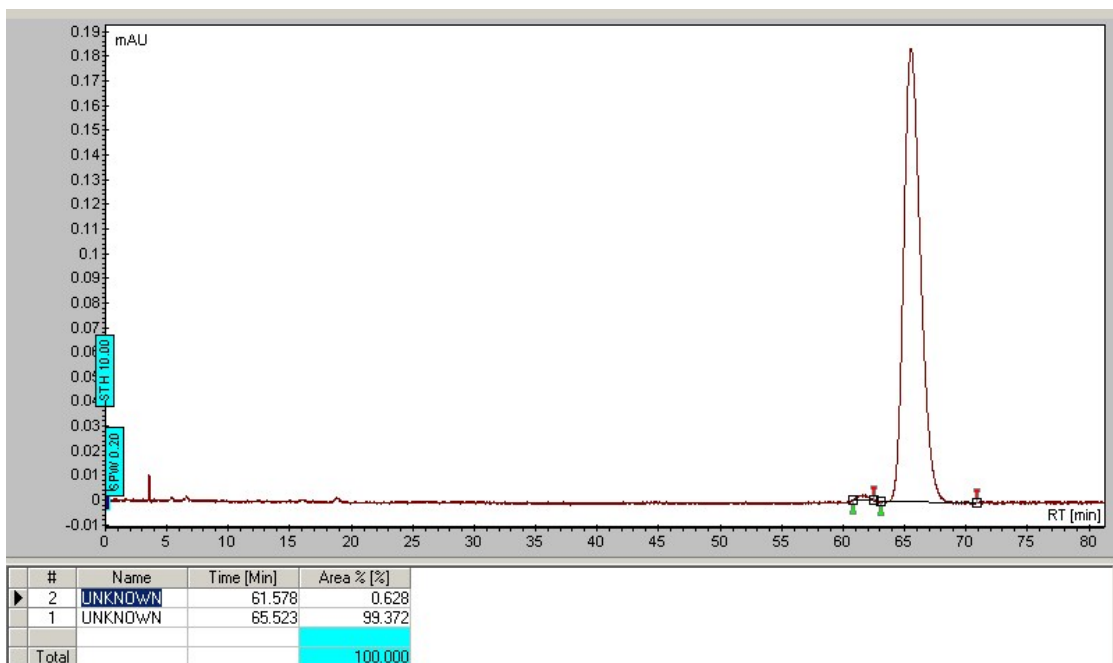
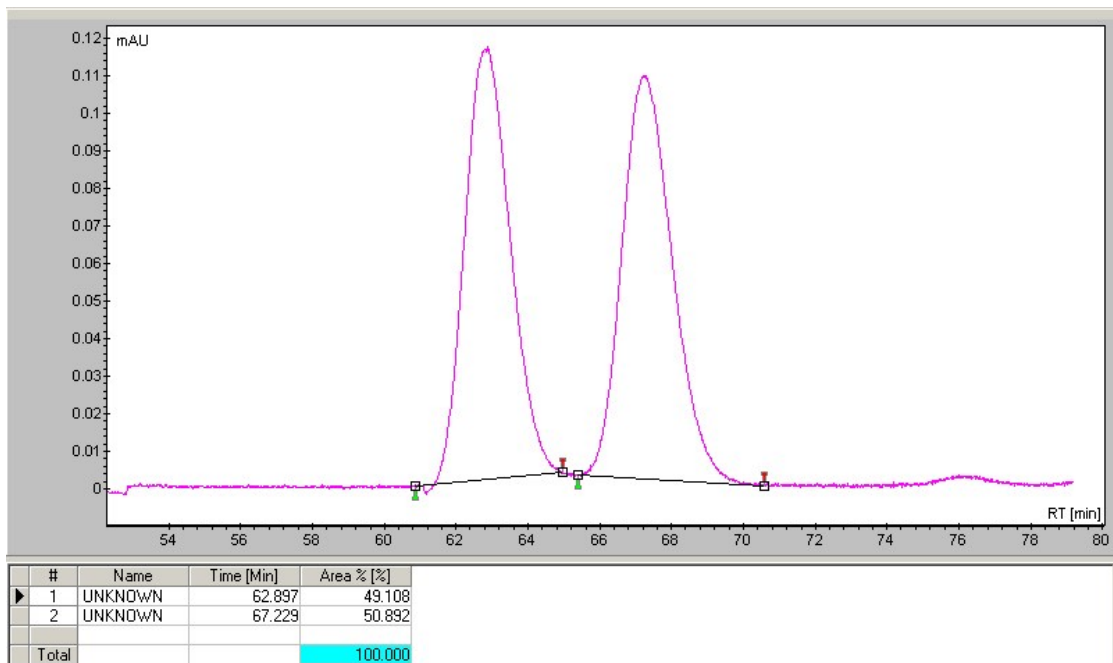


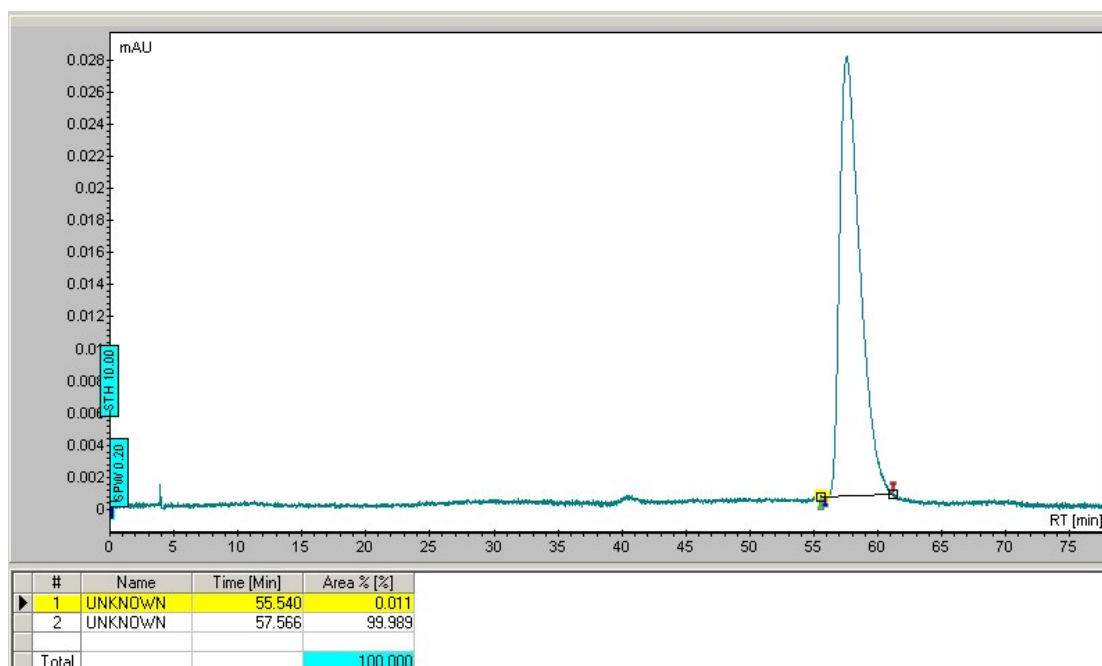
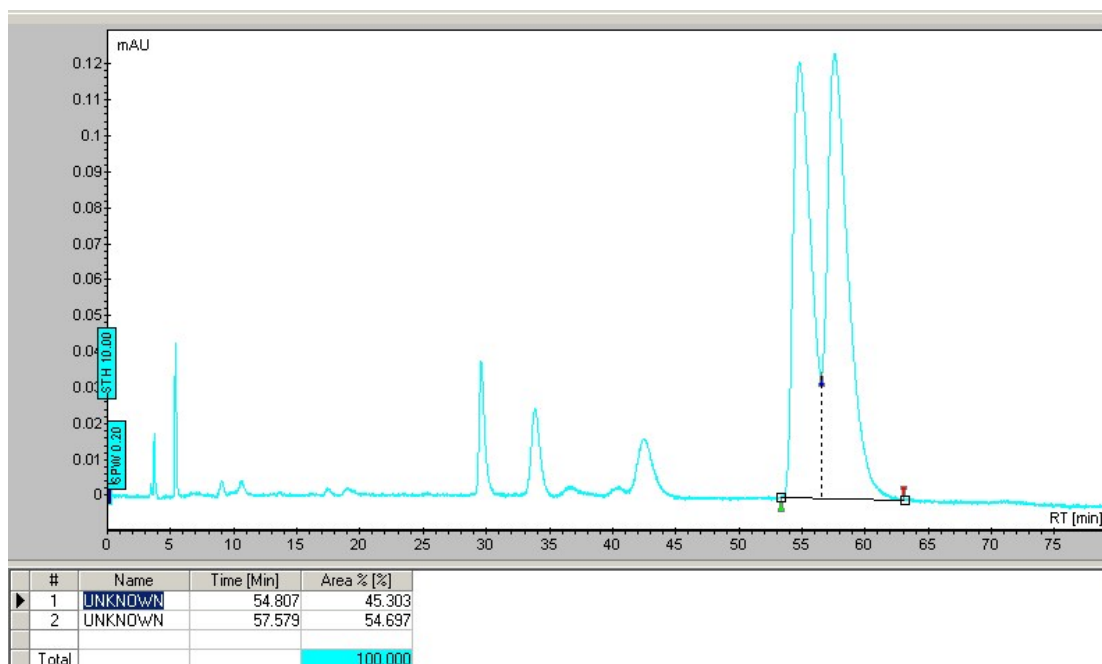
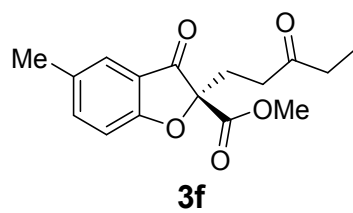
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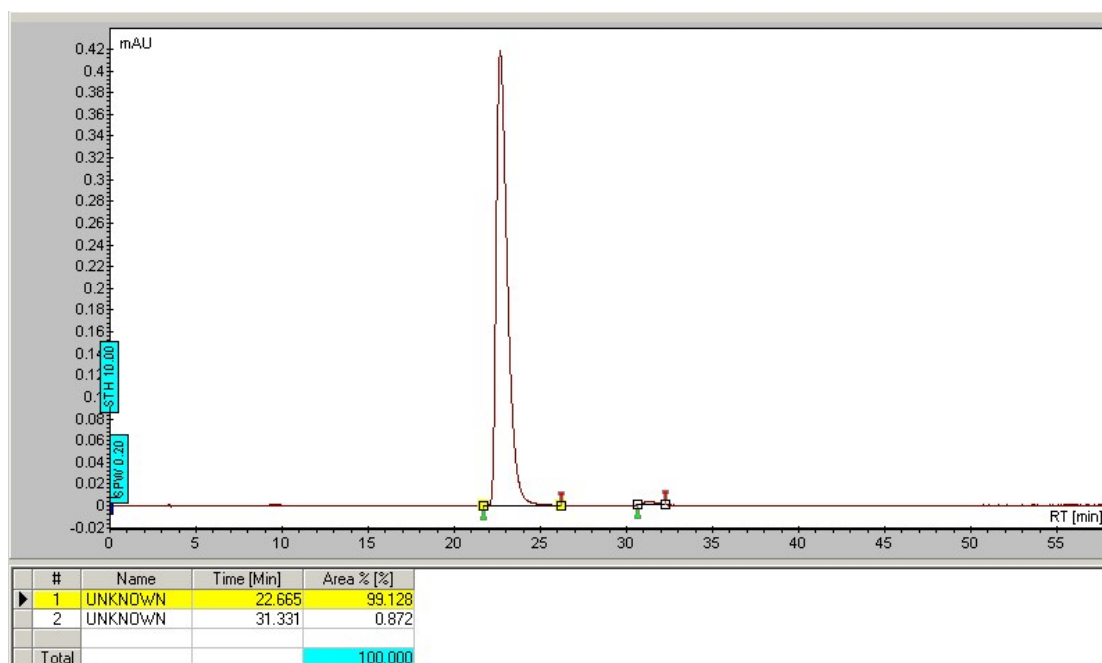
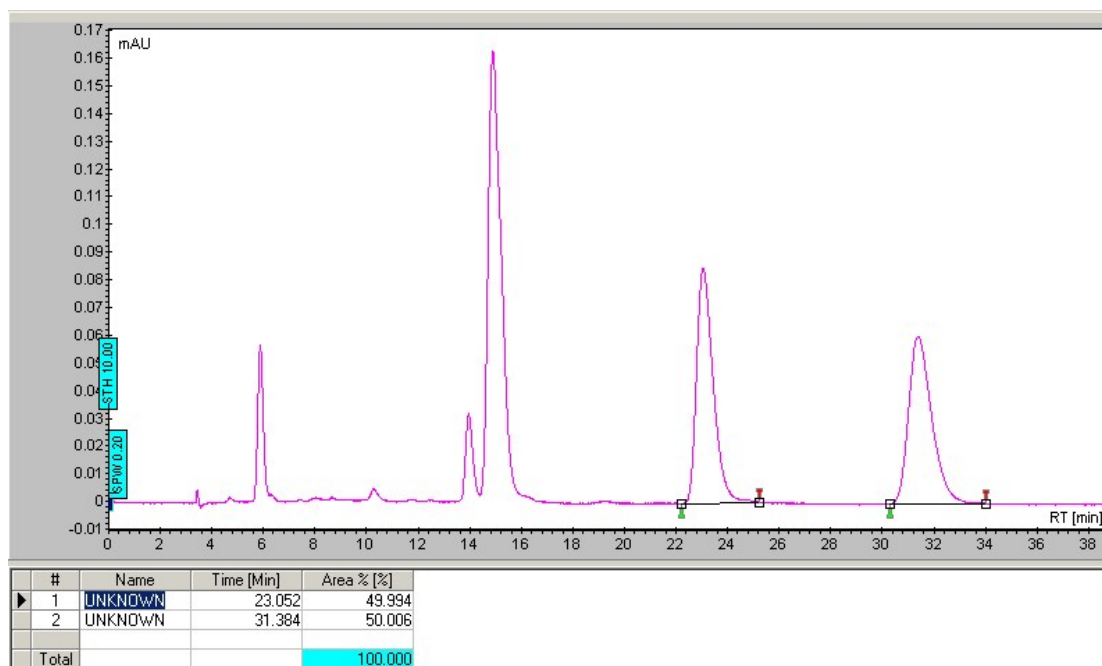
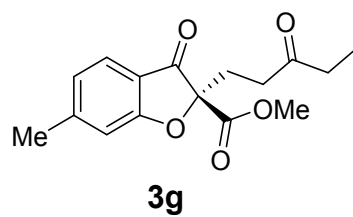


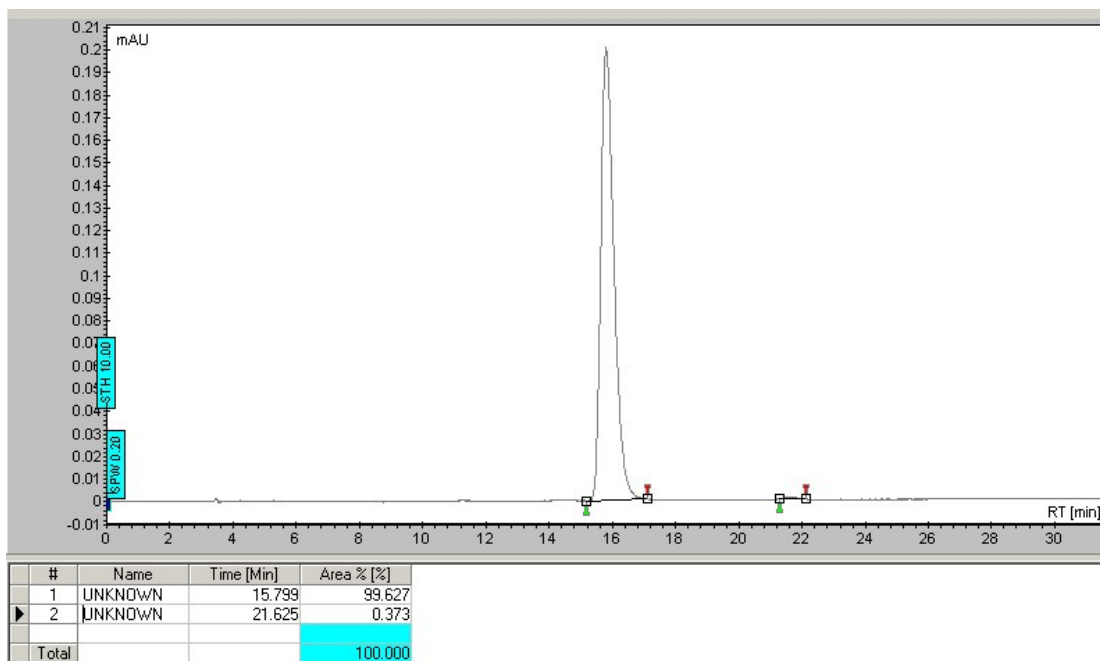
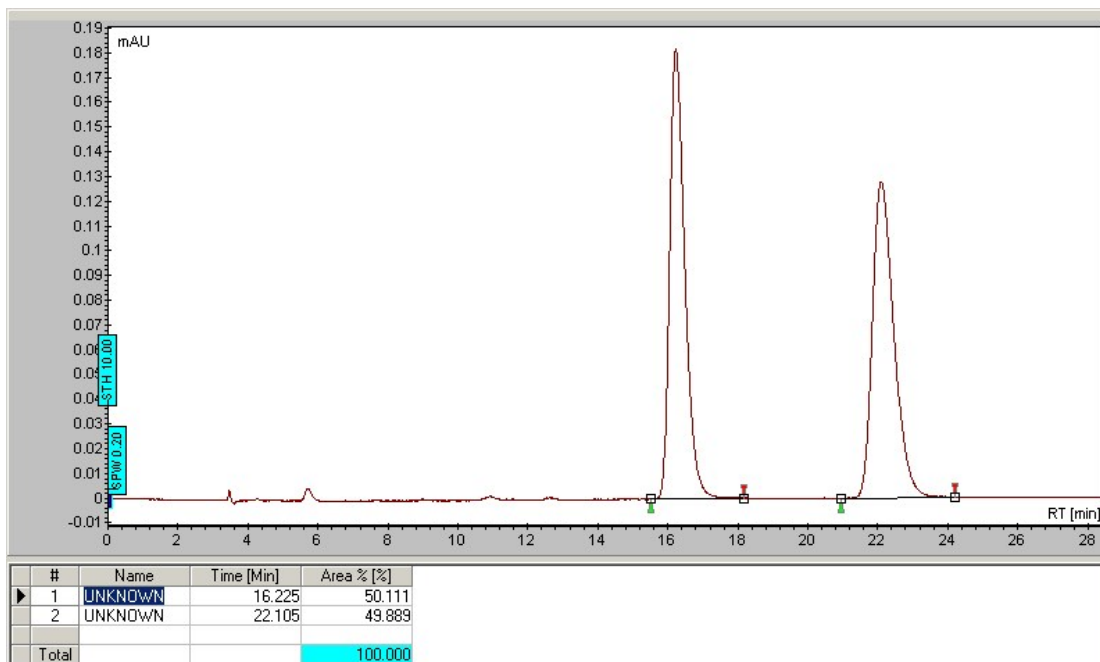
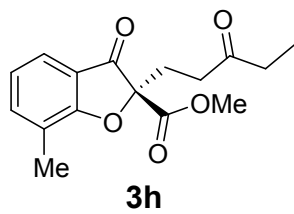


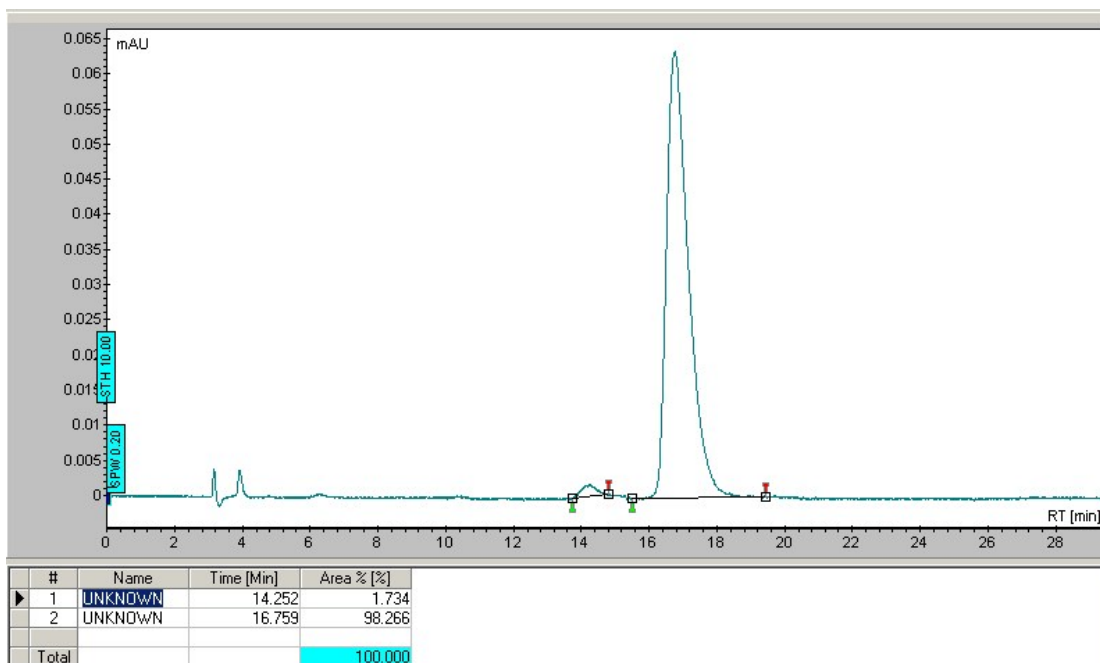
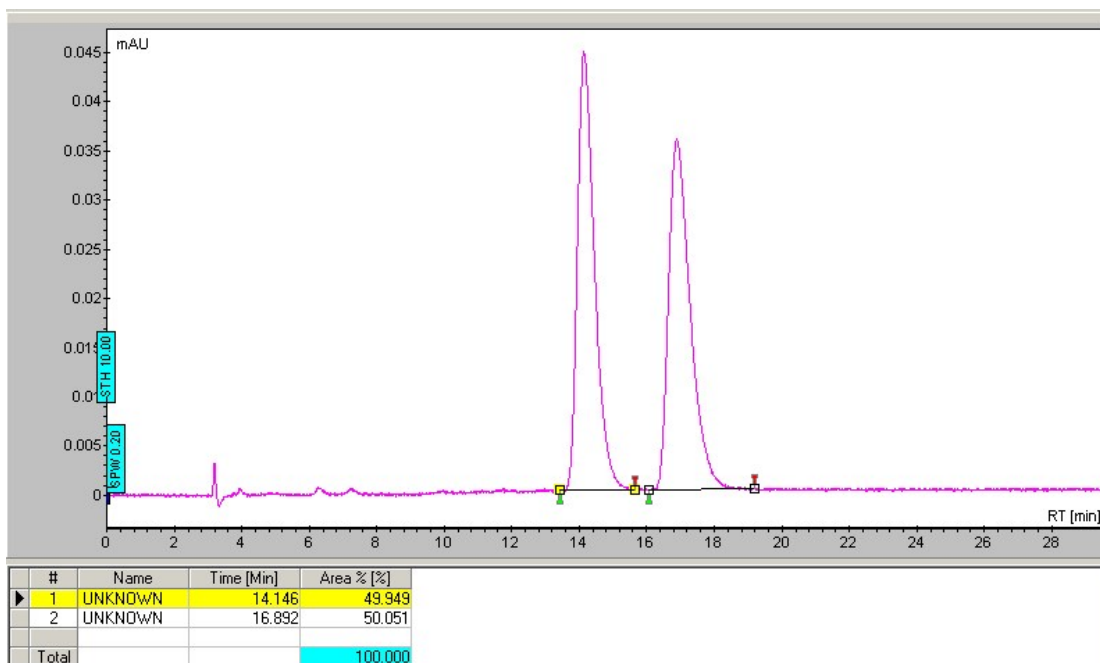
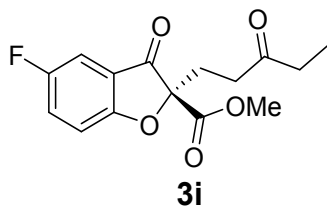
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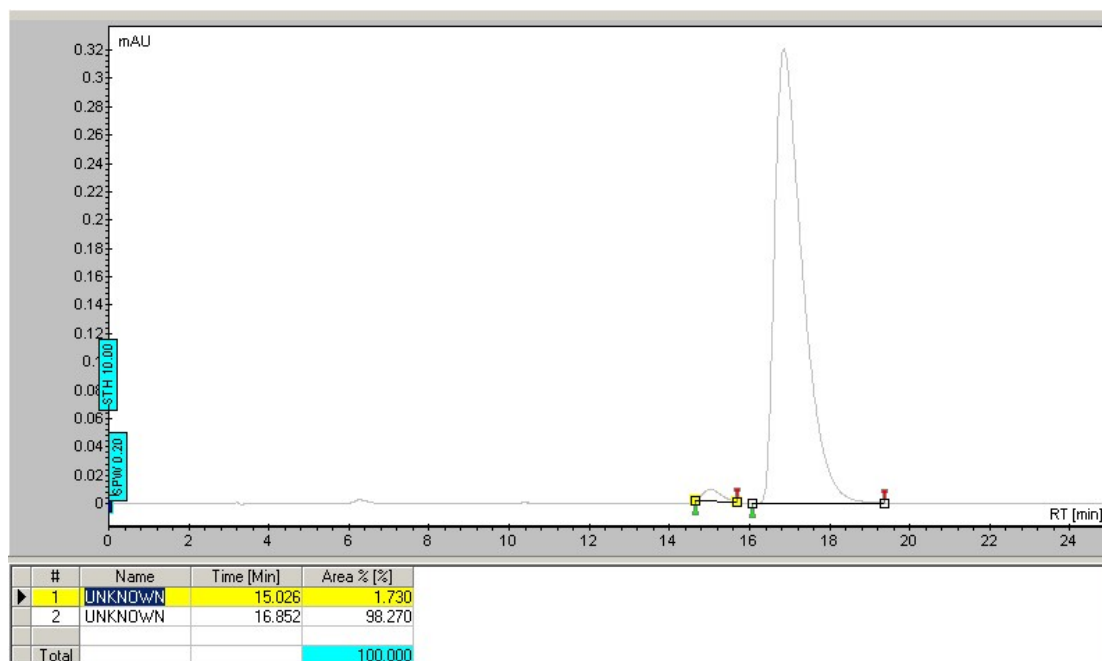
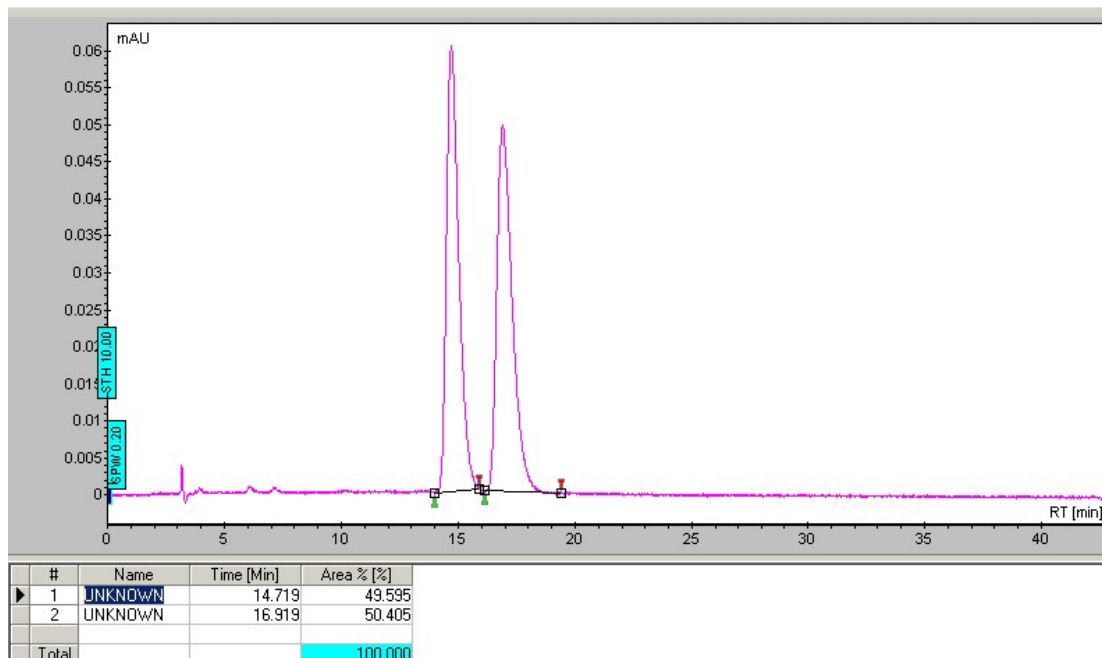
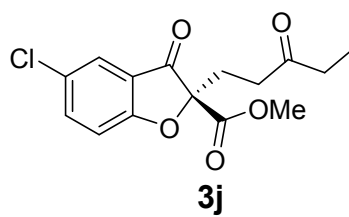


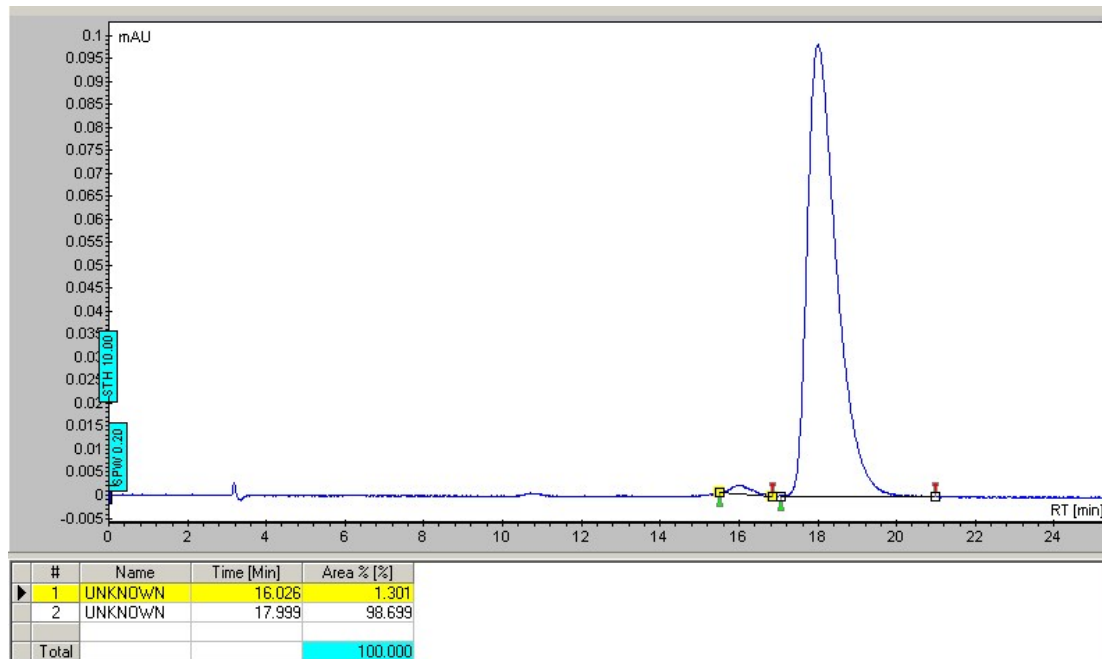
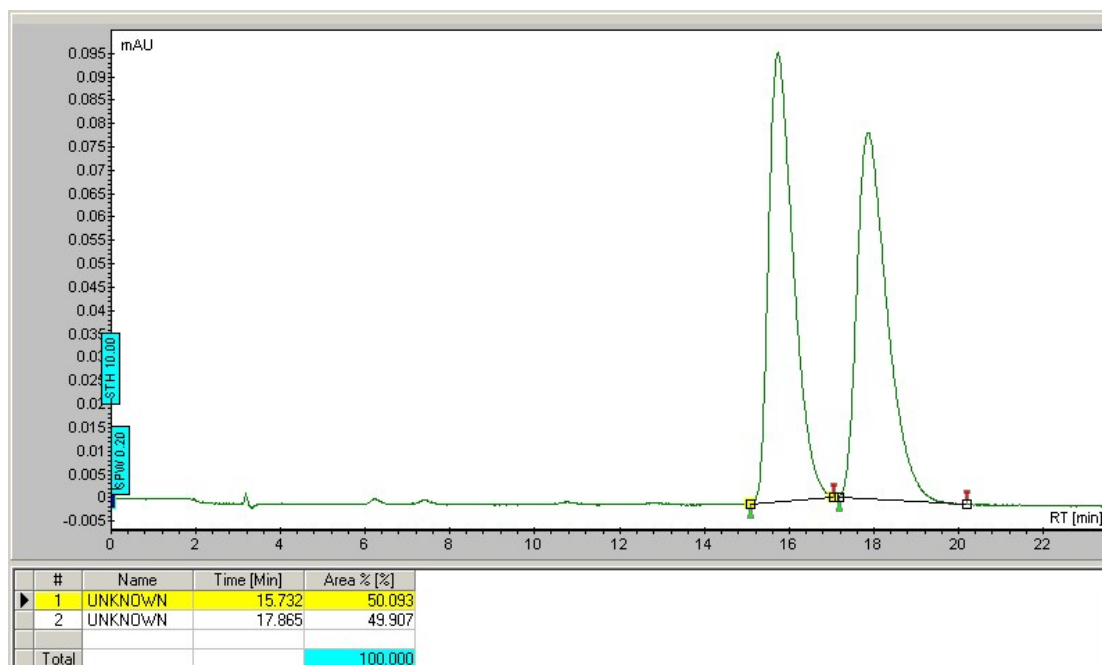
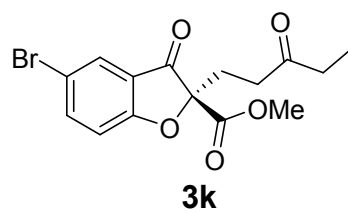


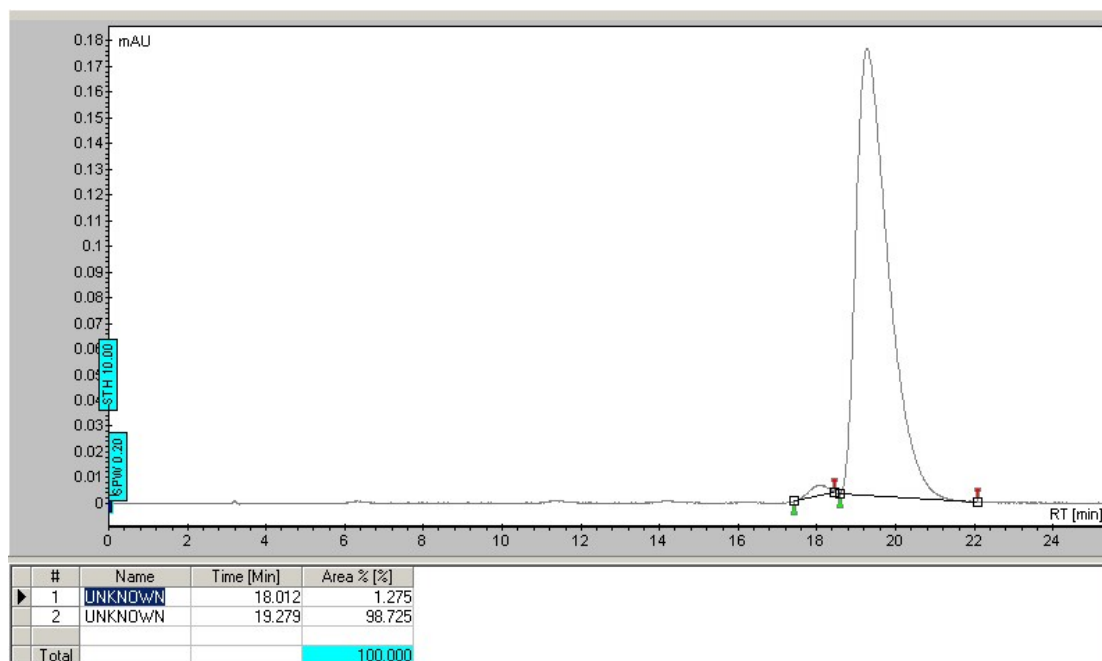
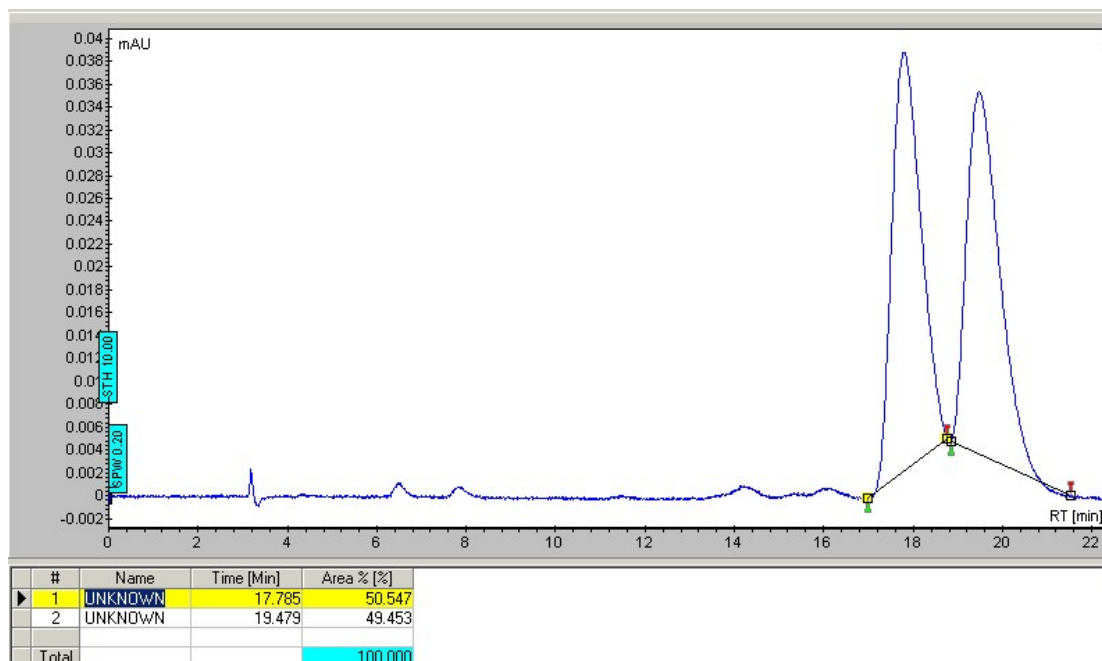
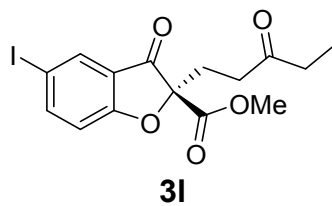


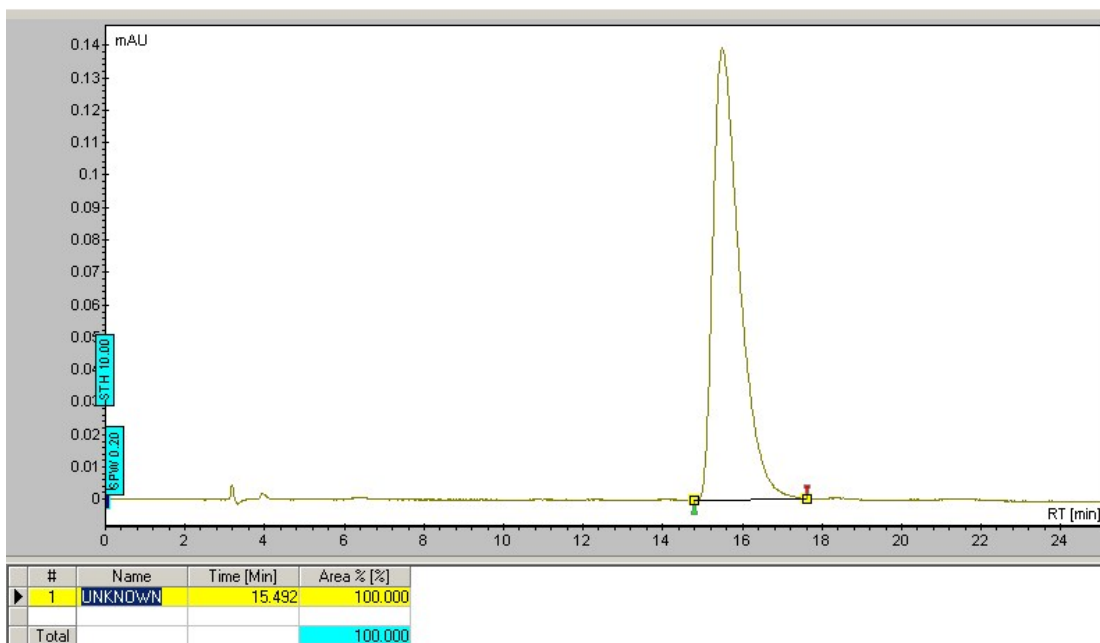
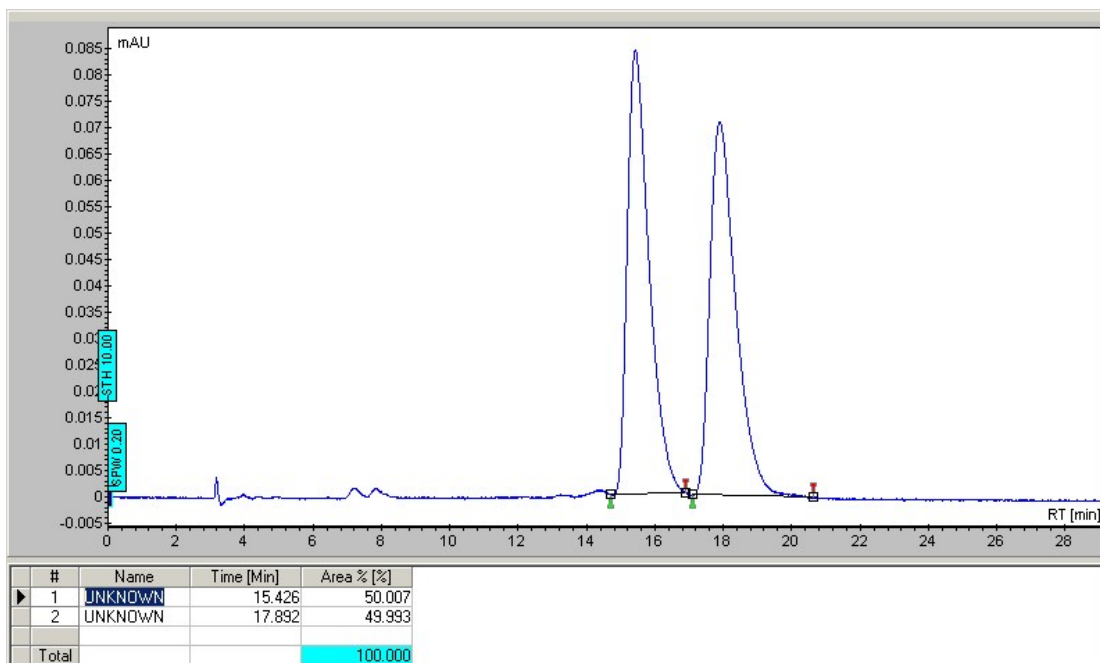
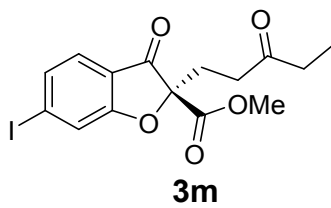


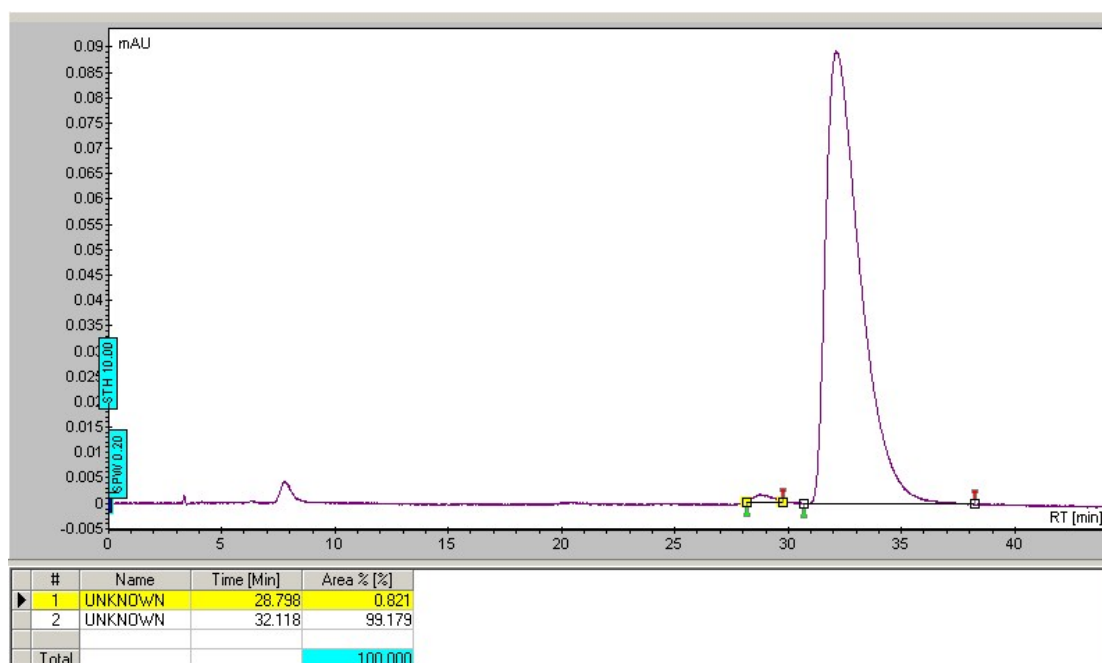
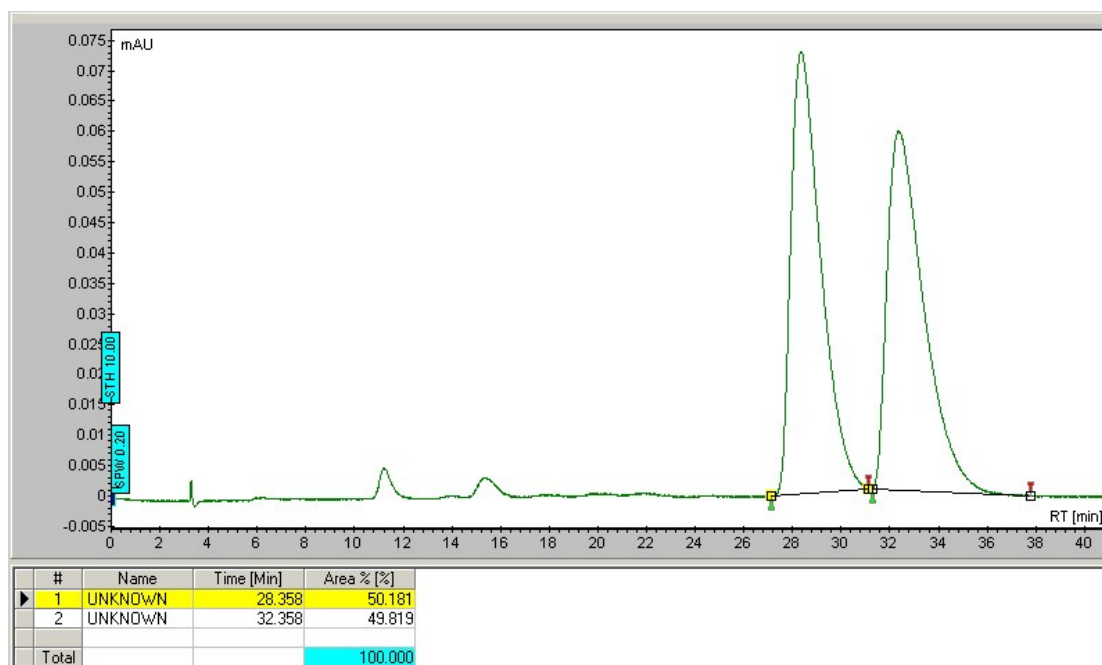
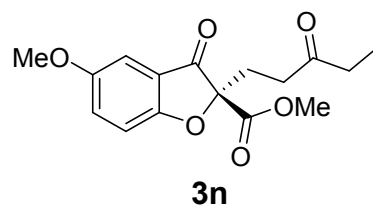


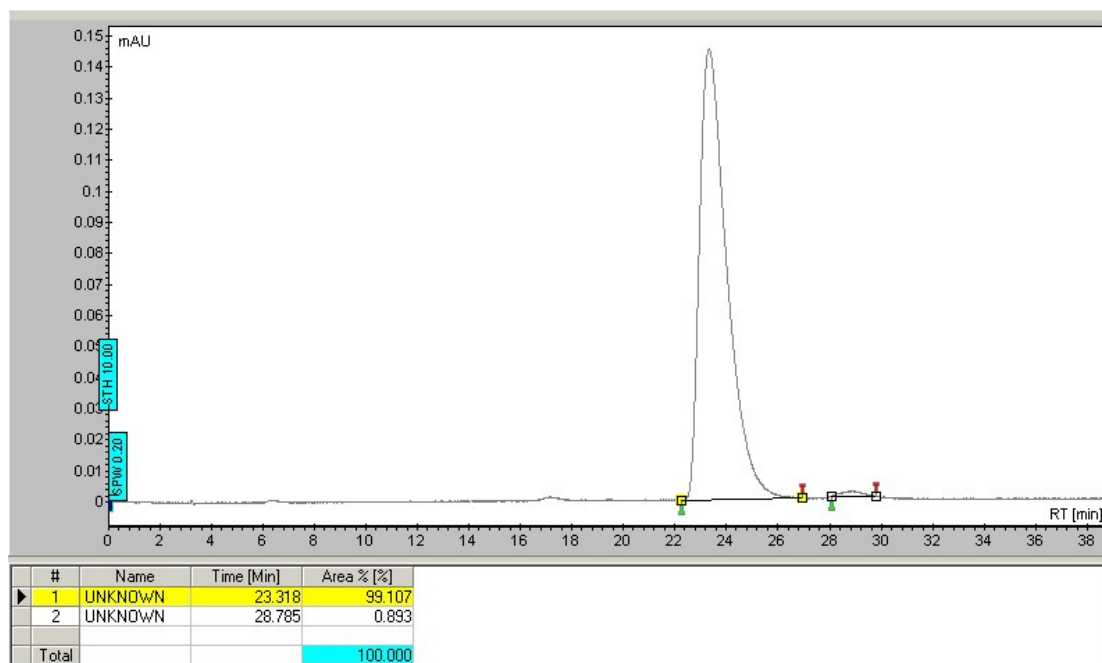
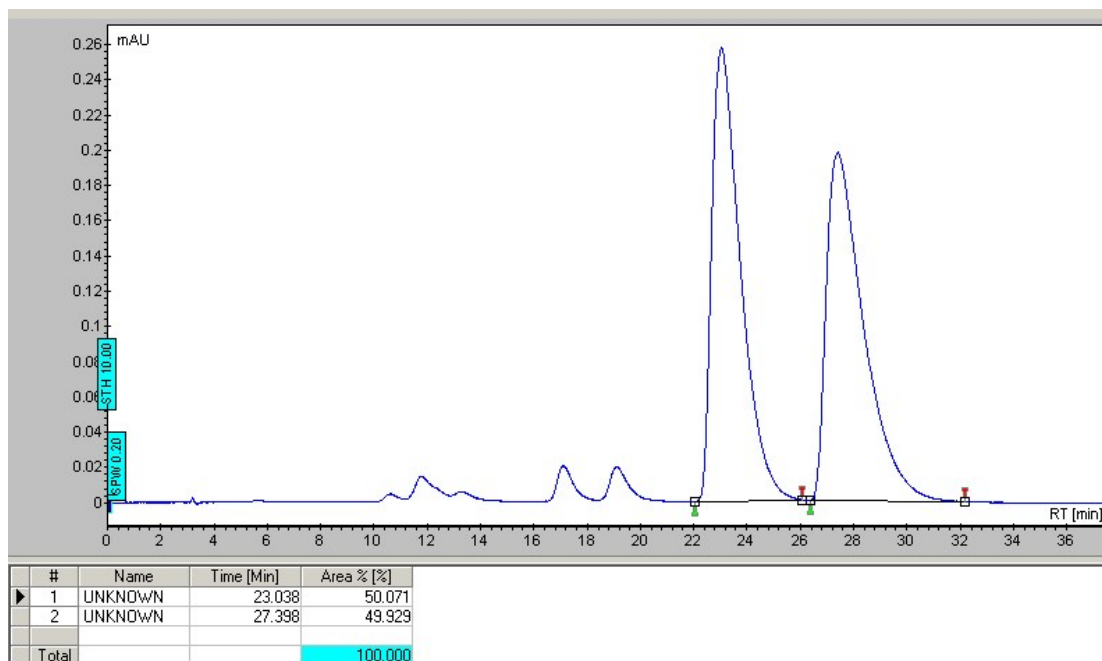
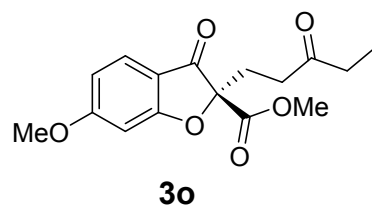


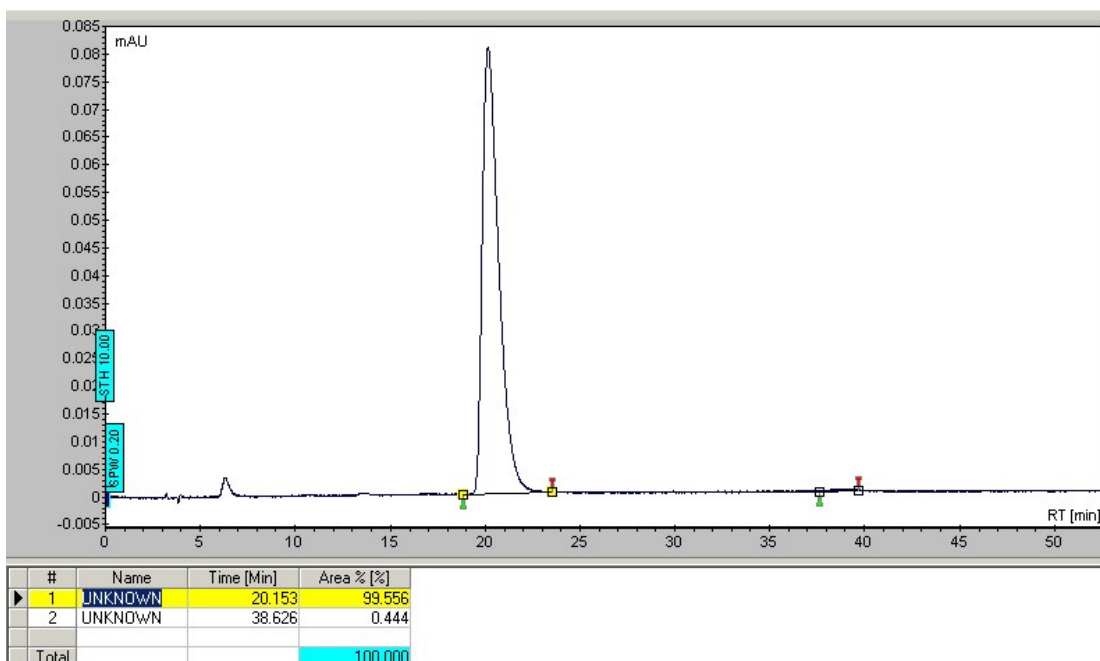
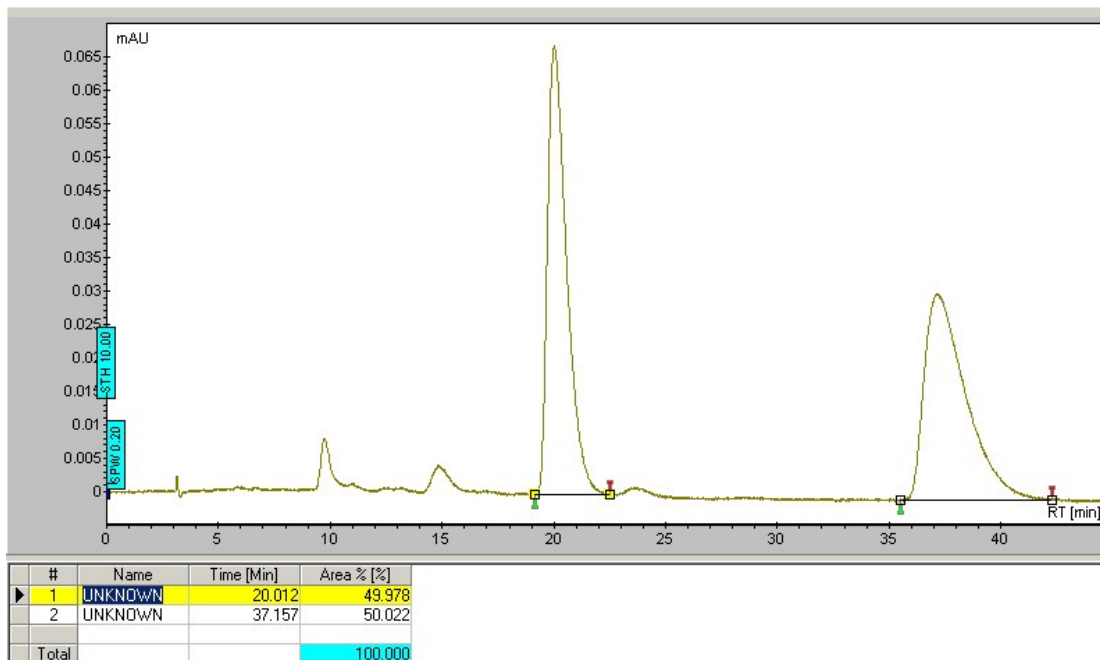
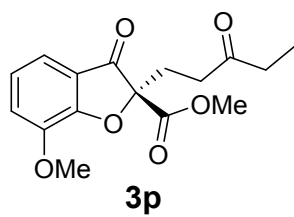


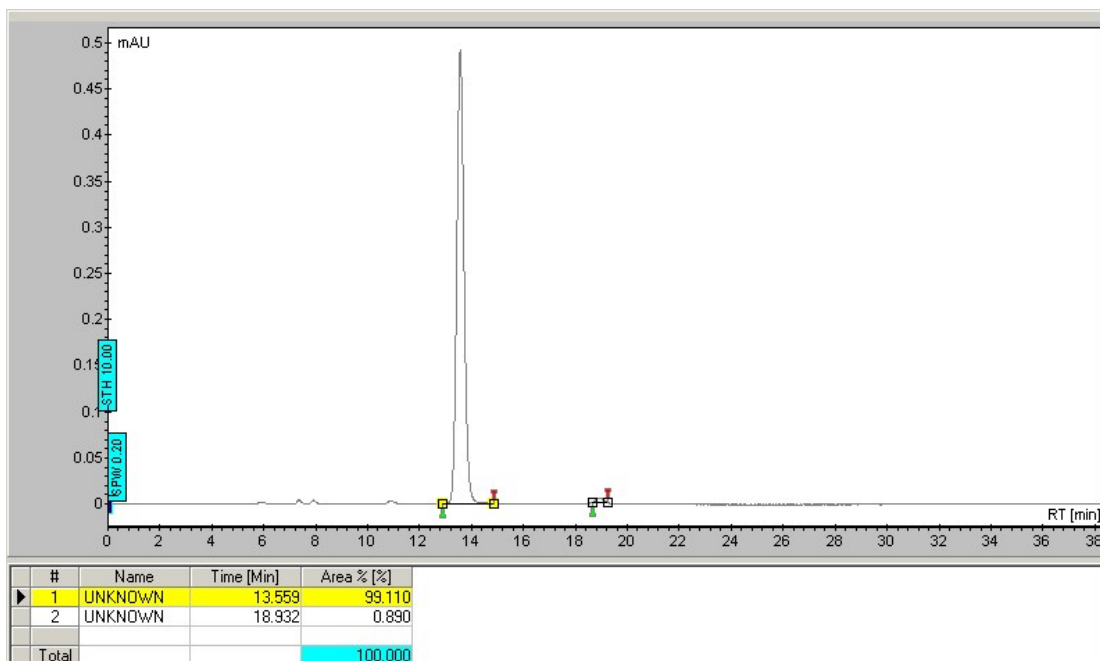
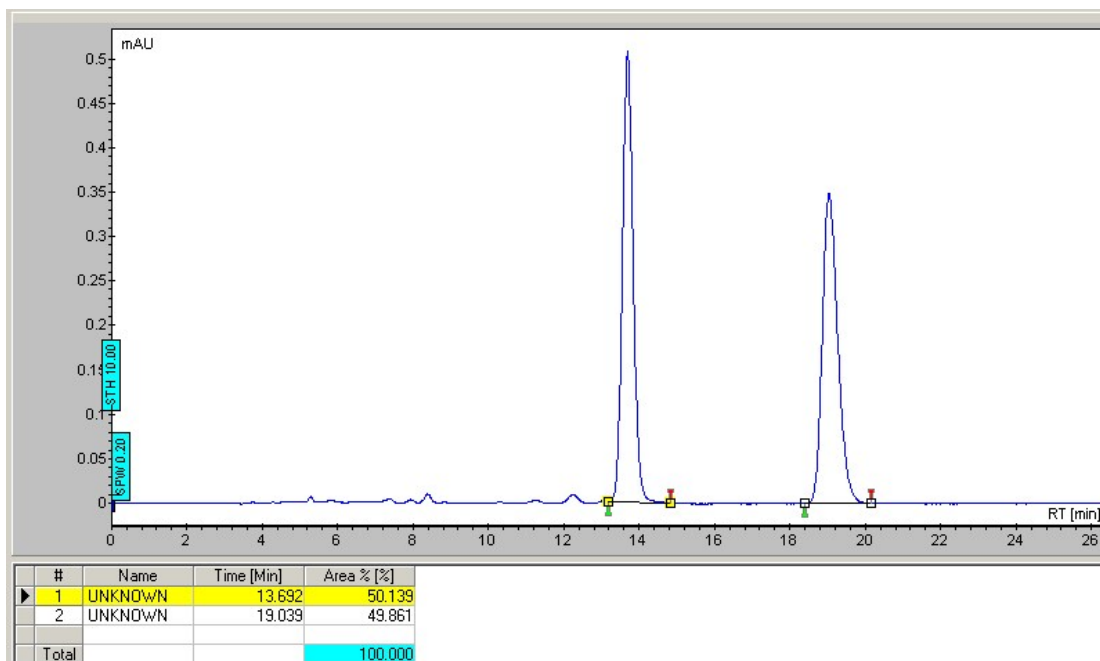
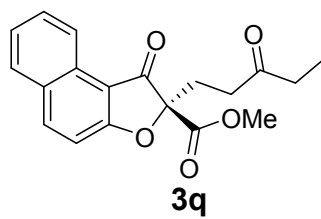


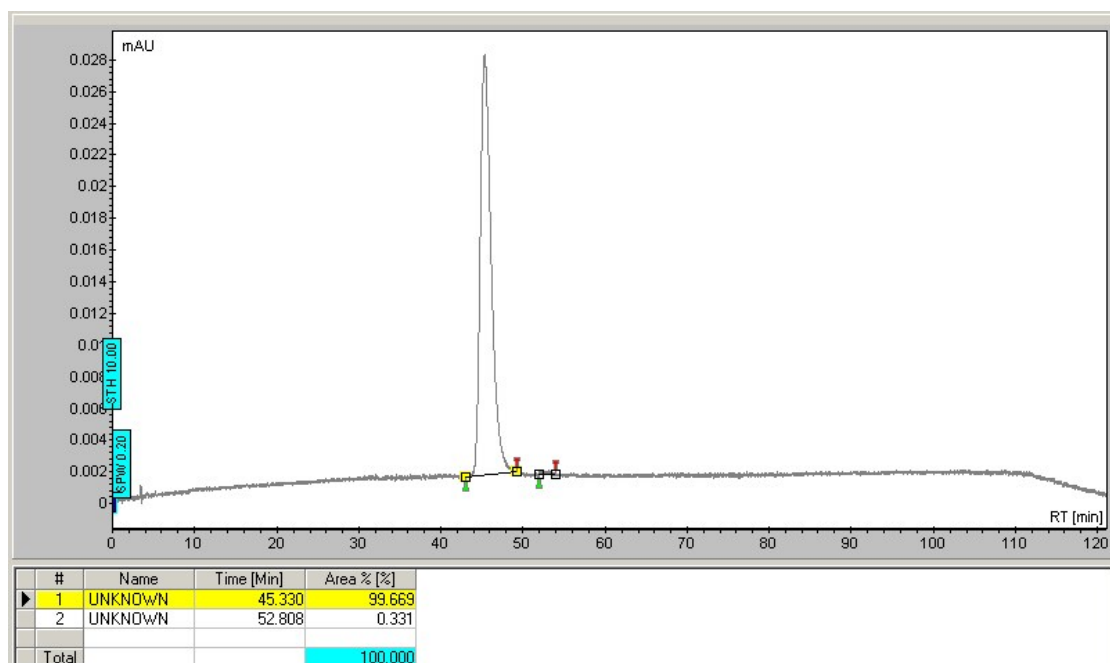
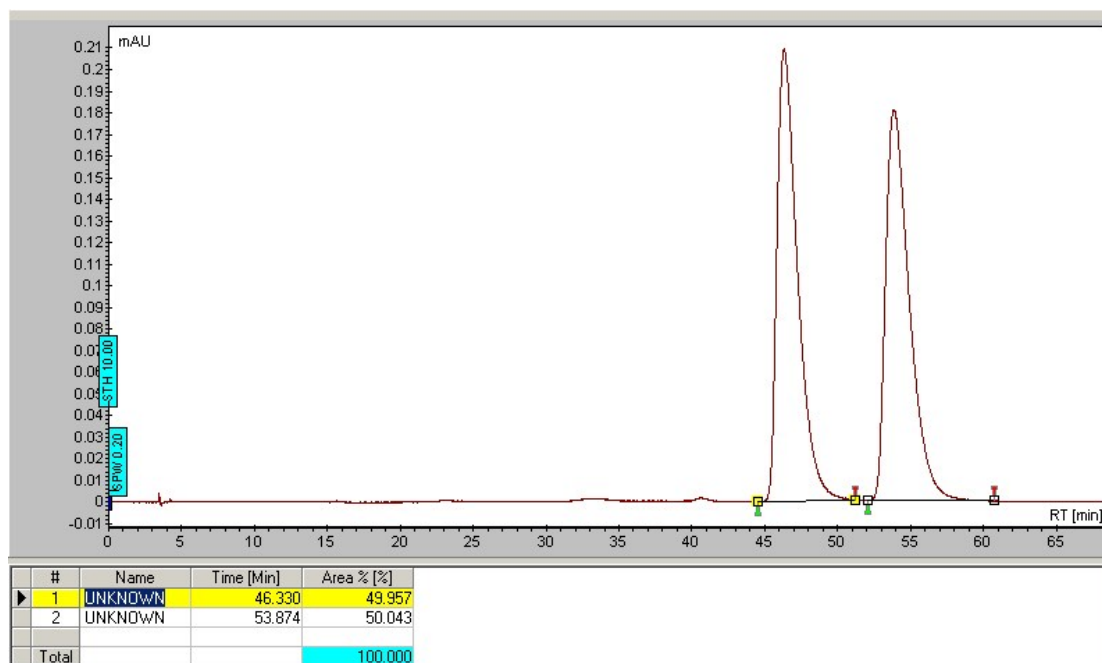
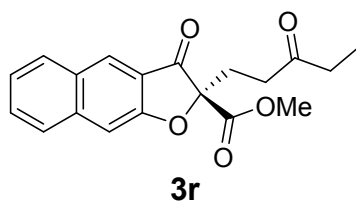


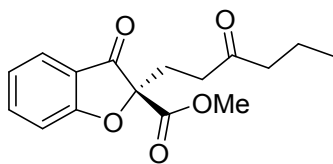




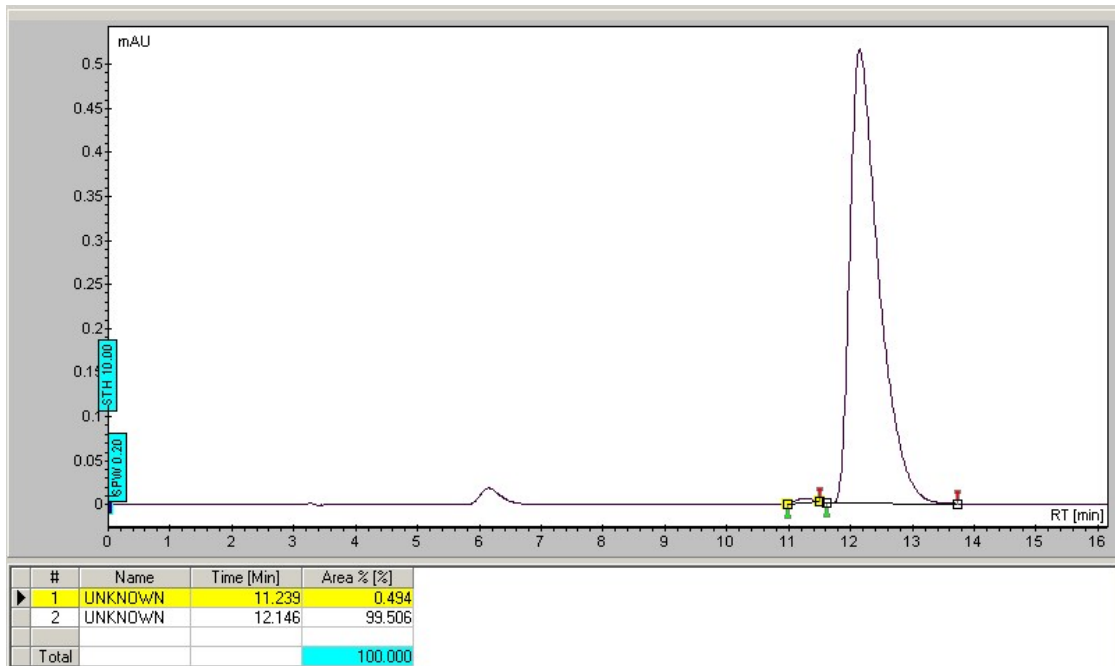
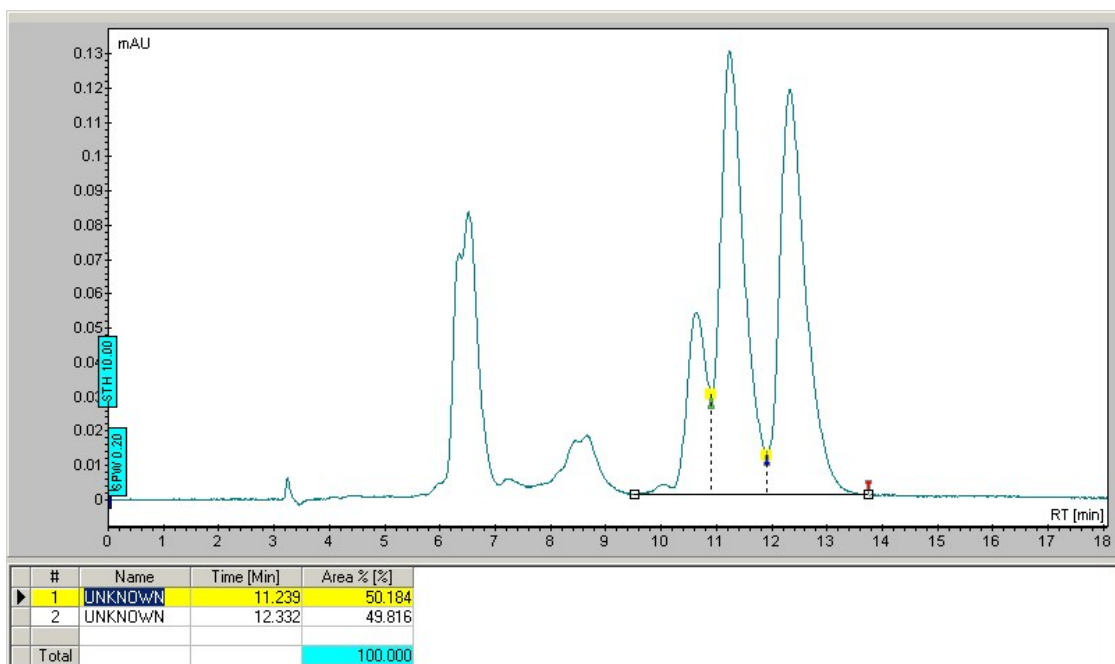


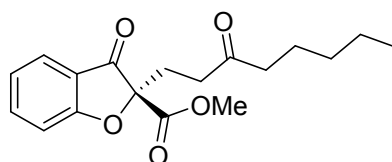




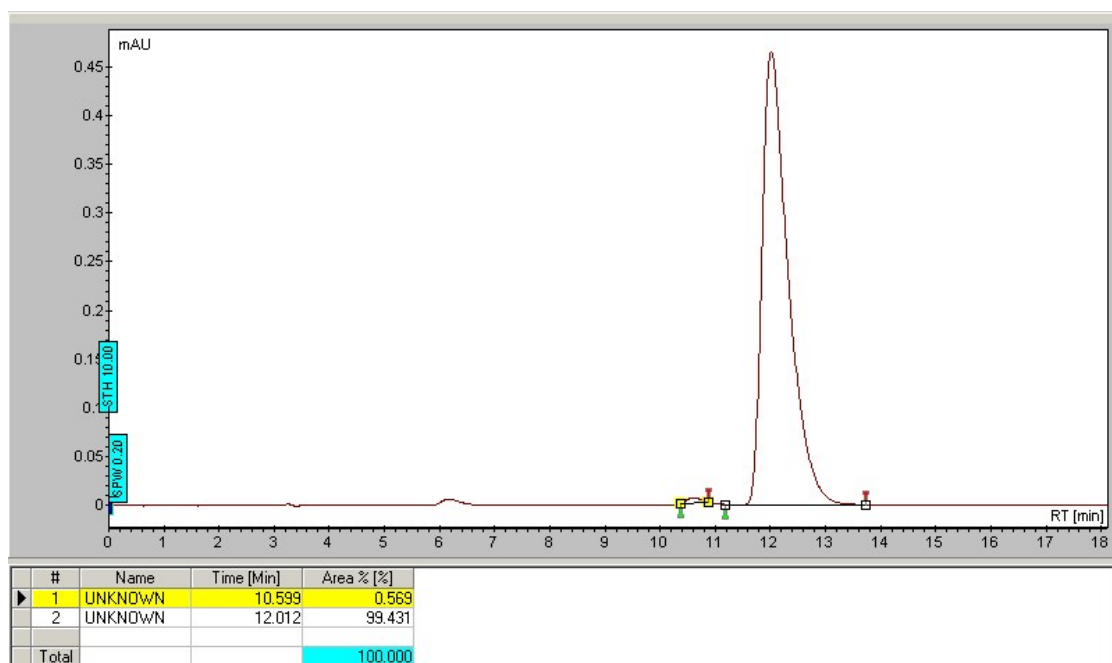
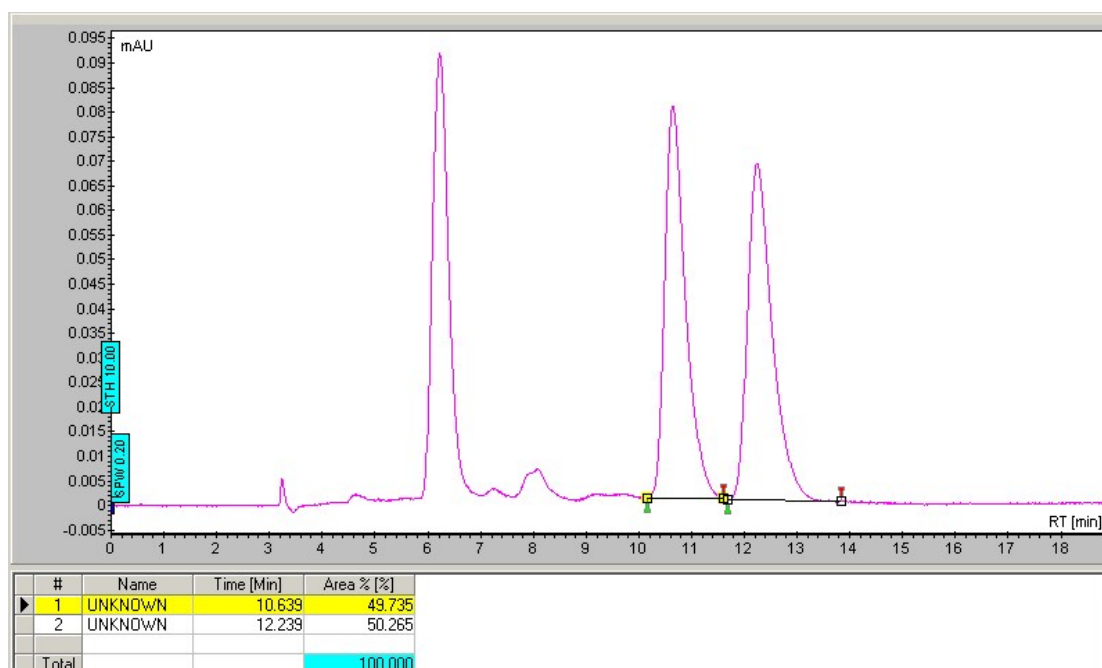


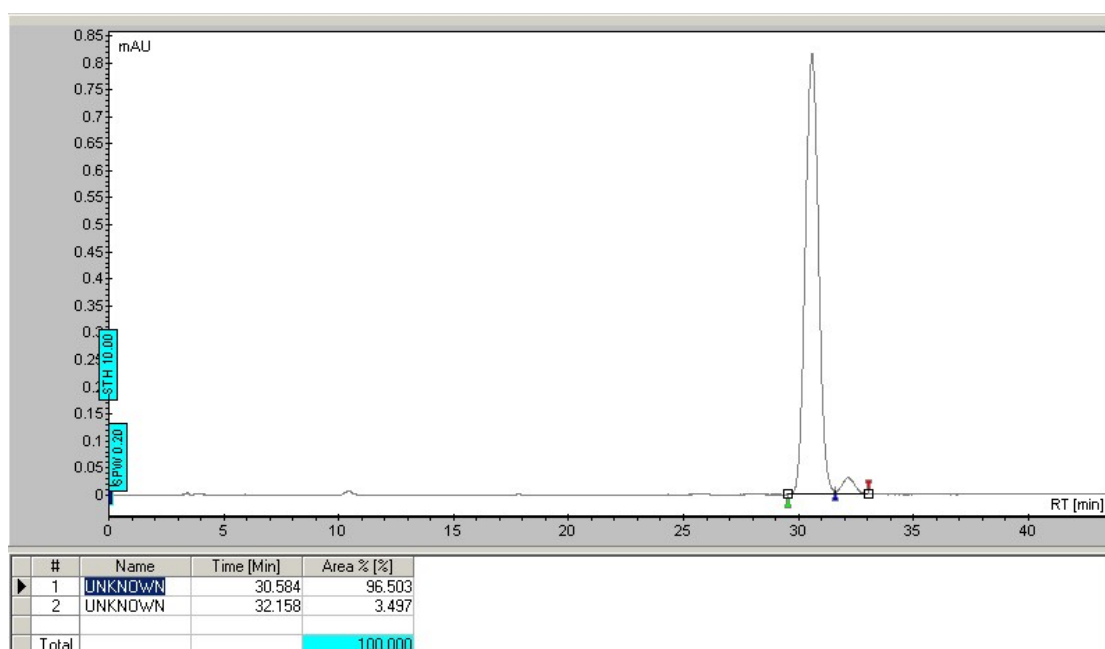
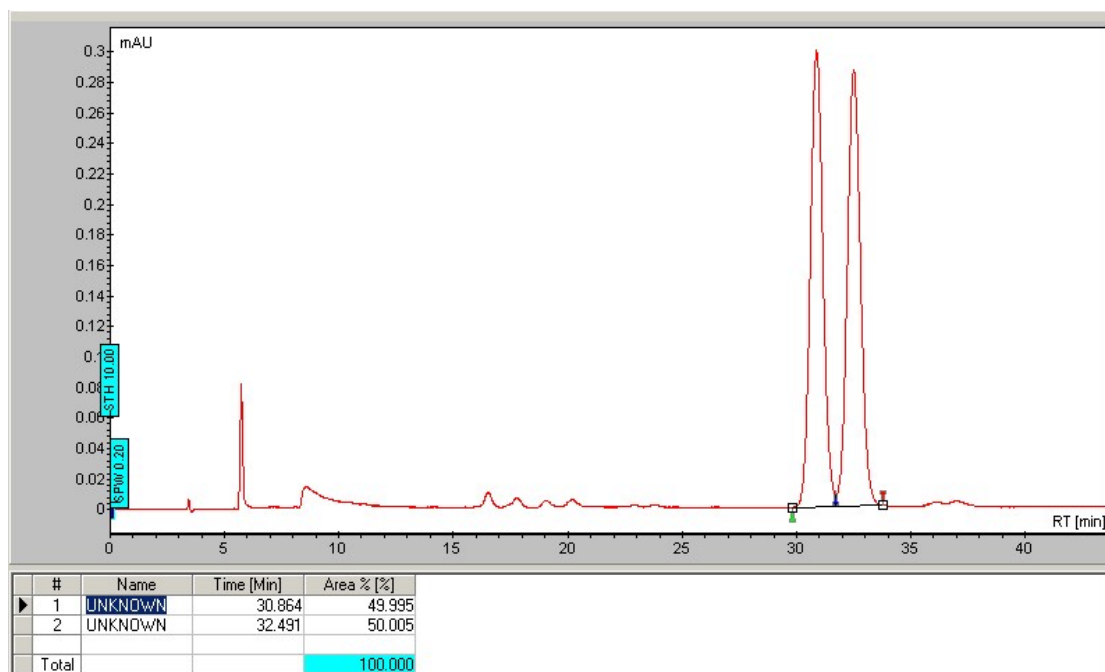
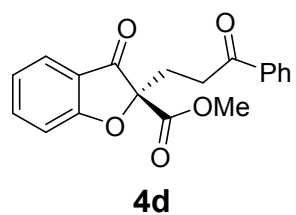
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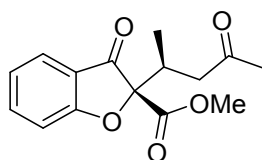




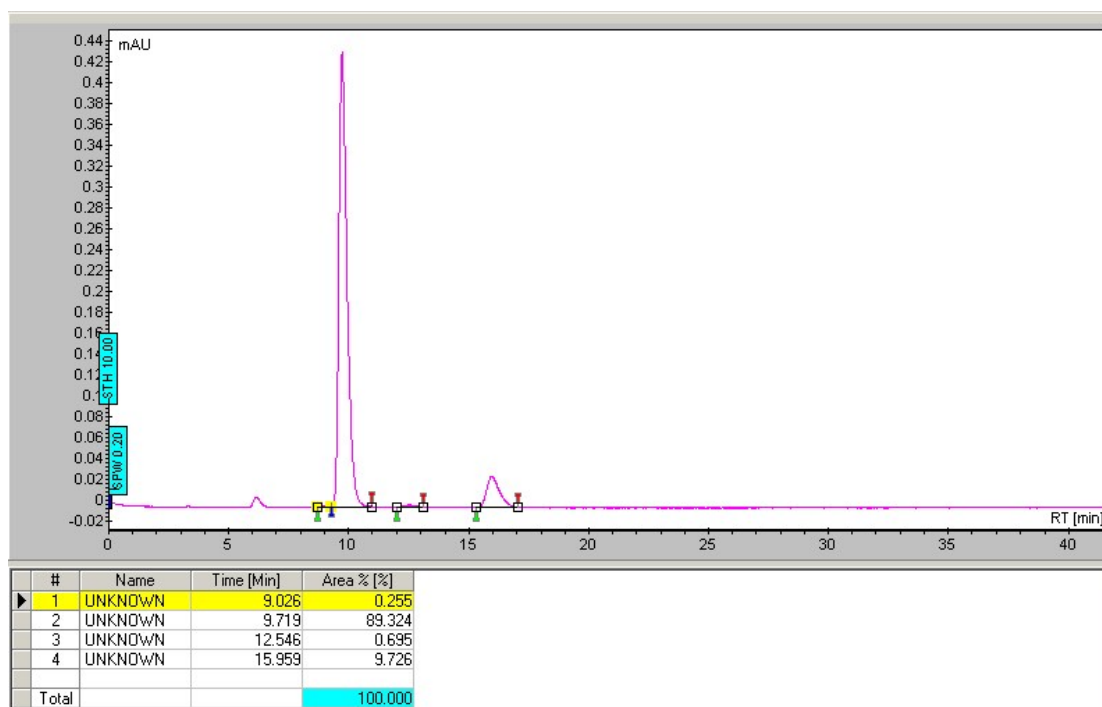
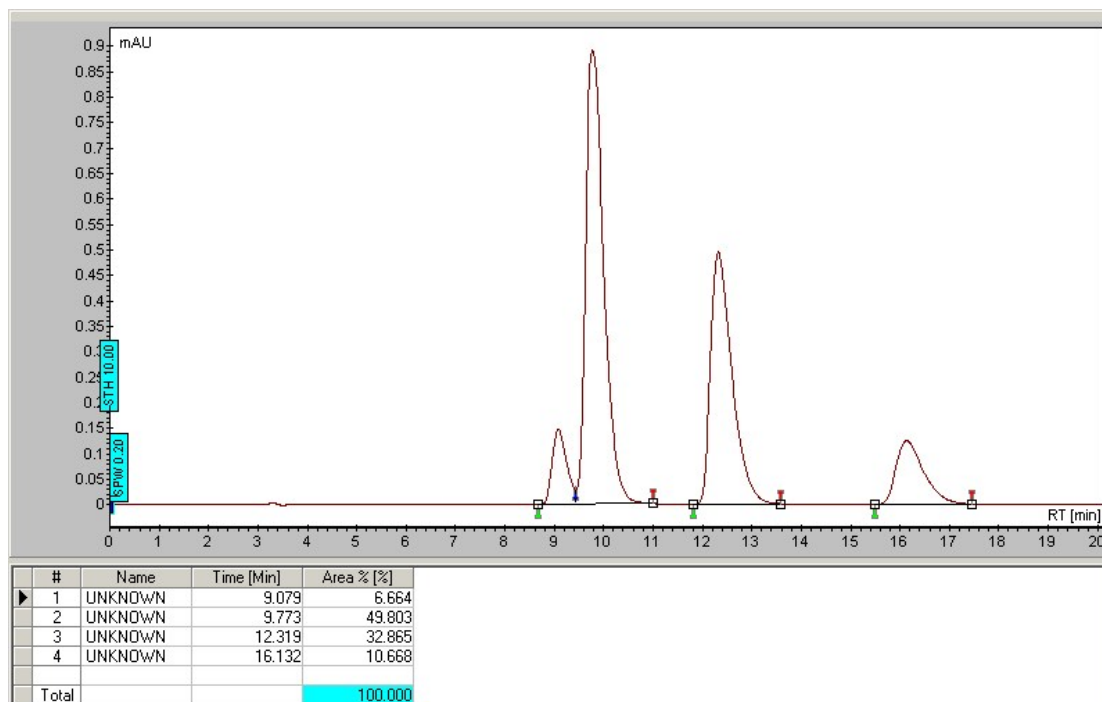
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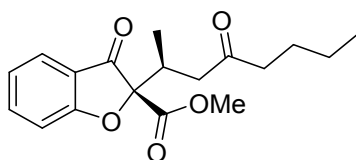




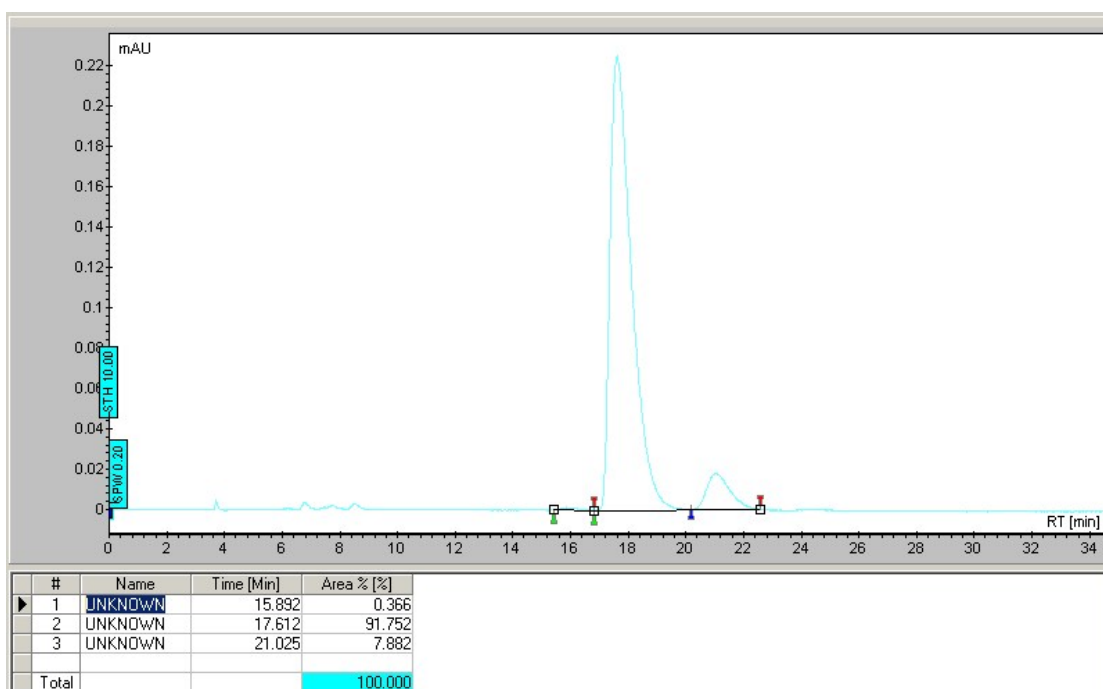
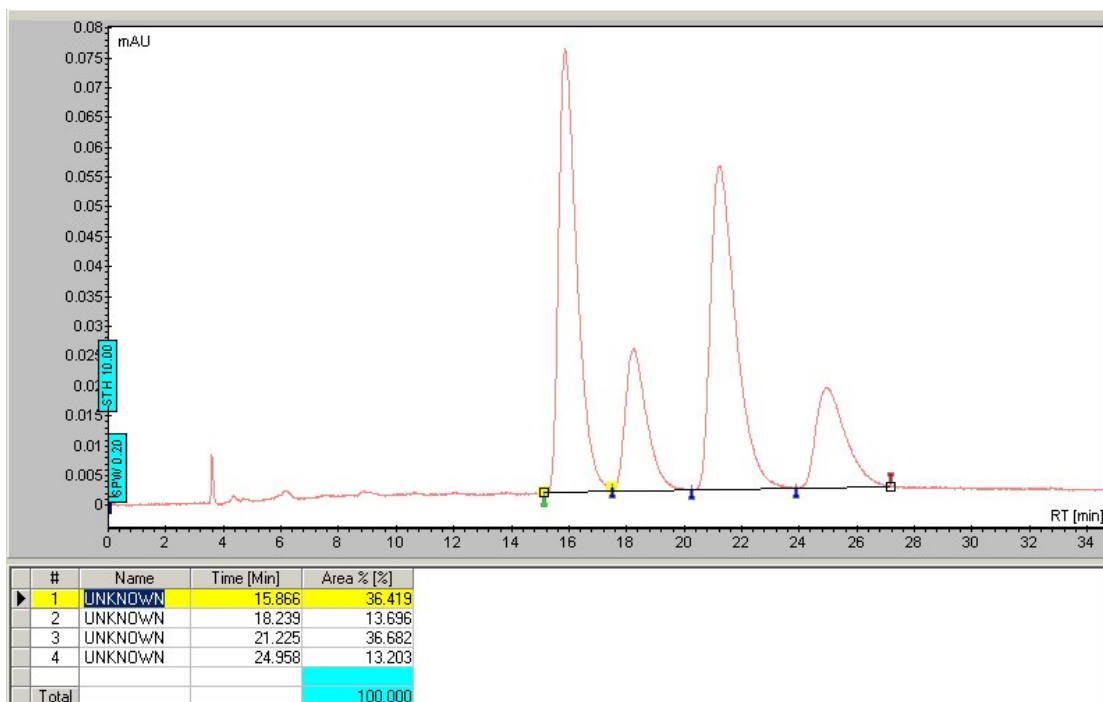


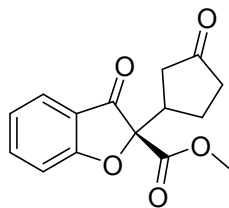
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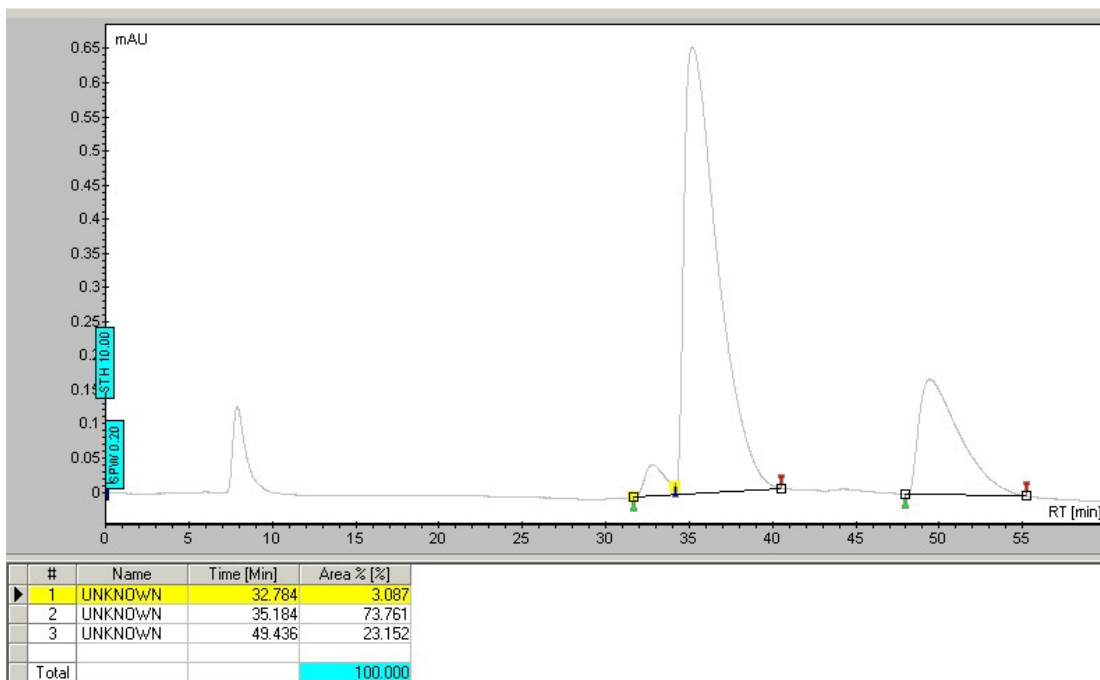
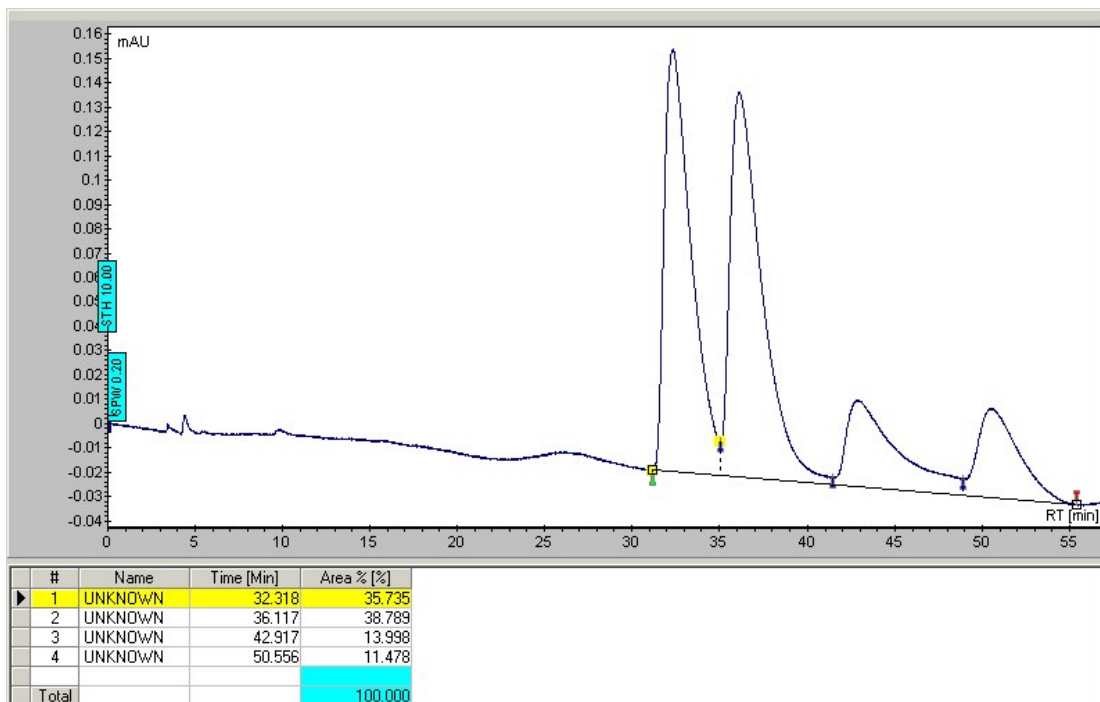


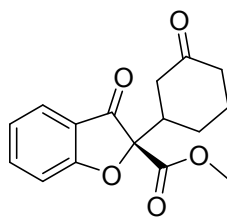
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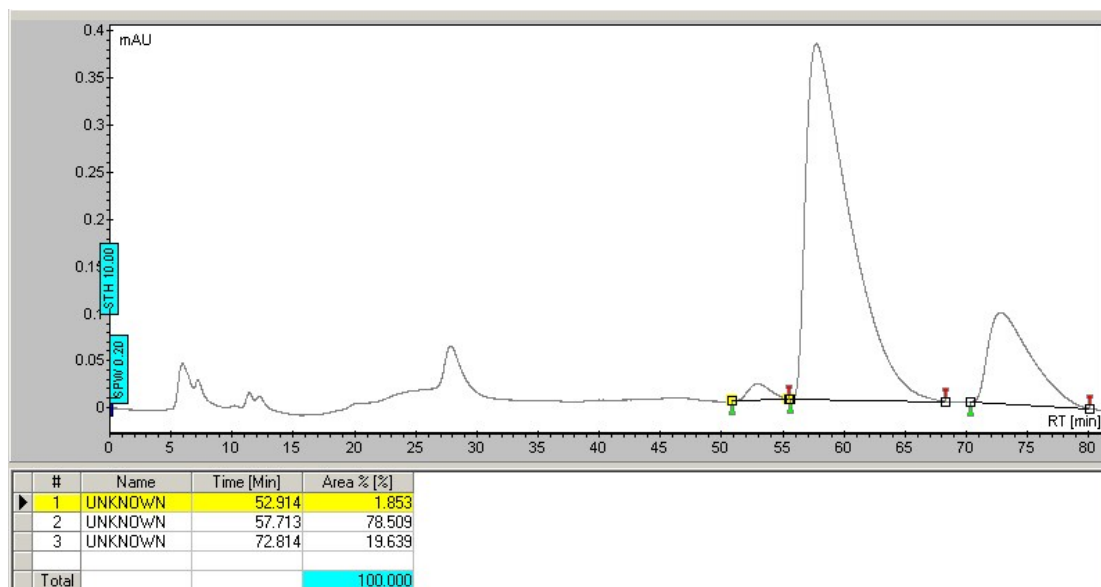
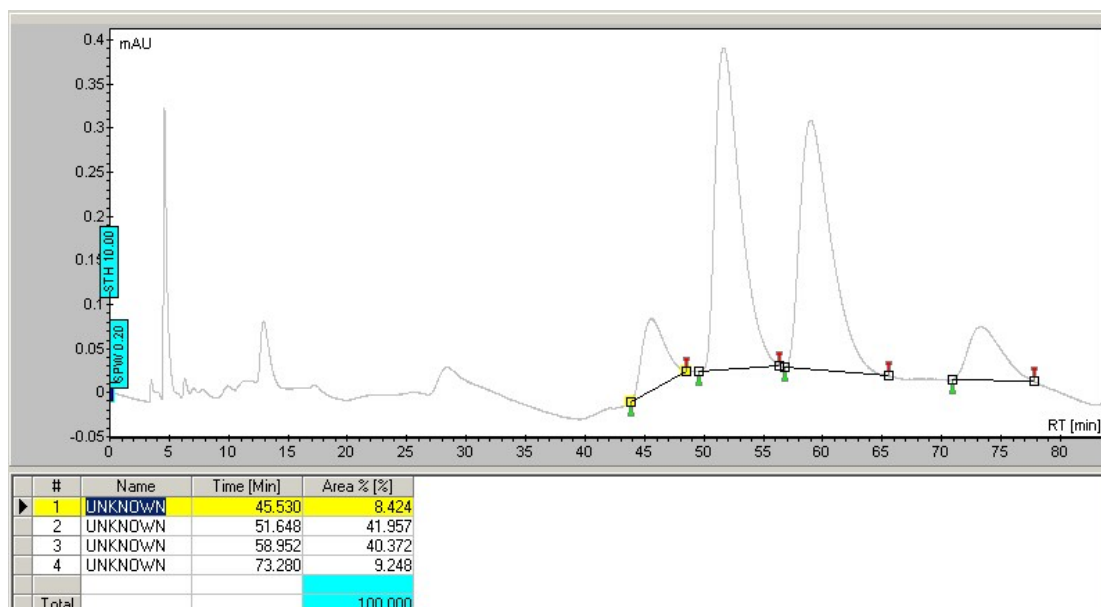


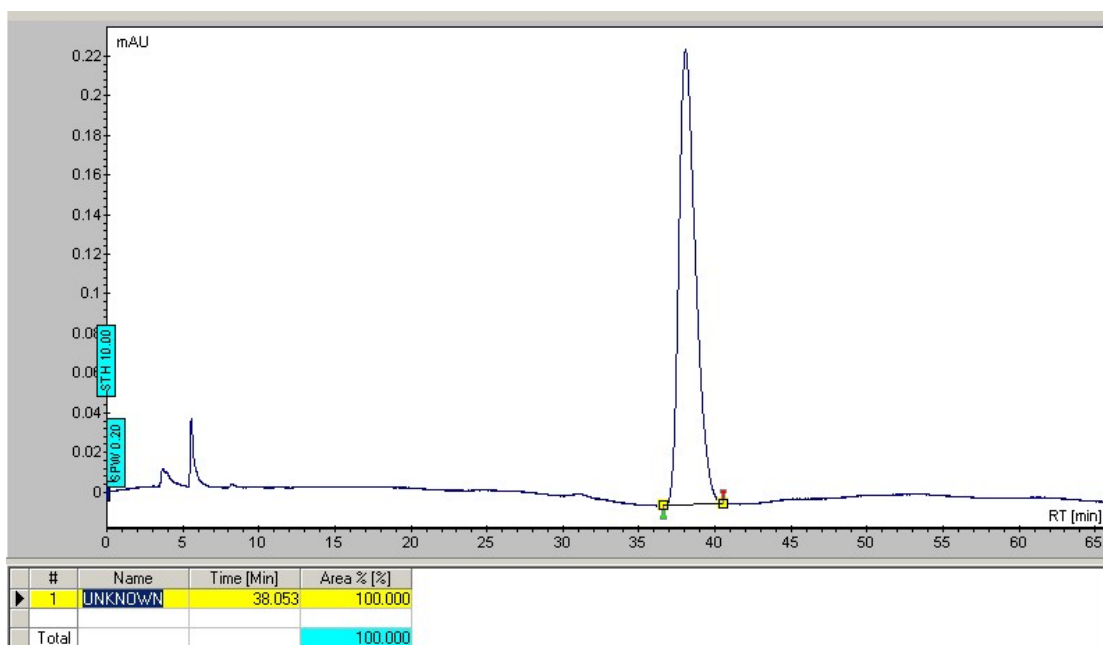
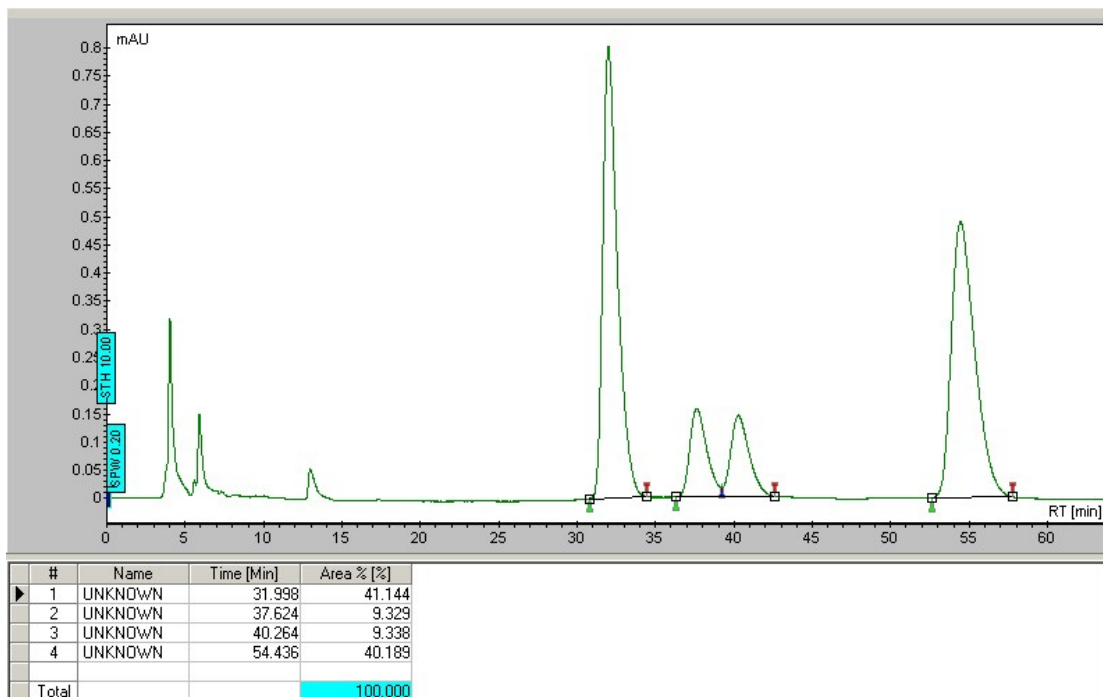
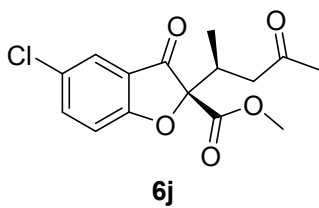
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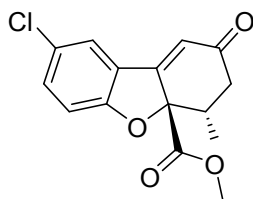




6d







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