# [2.2]Paracyclophane-based carbene-copper catalyst tuned by transannular

## electronic effects for asymmetric boration

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### **Electronic Supplementary Information**

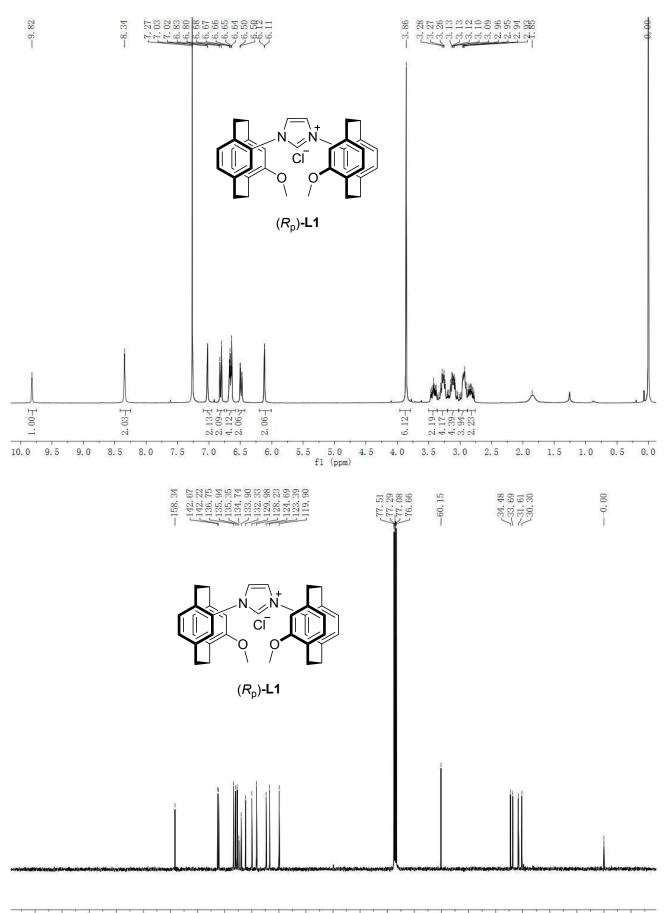
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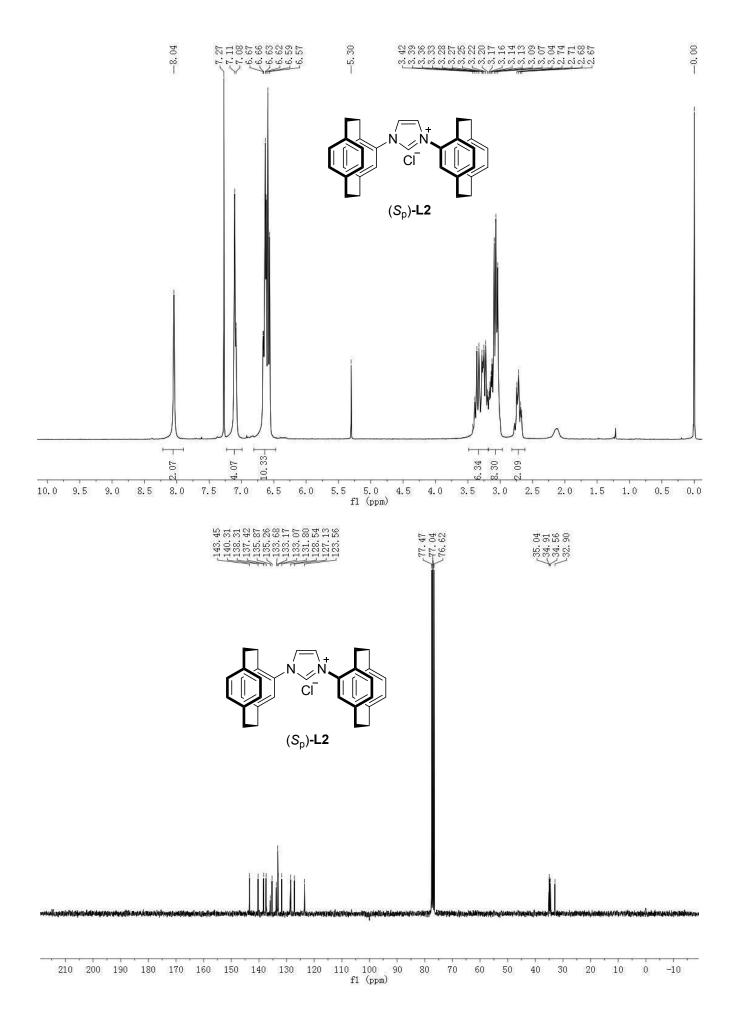
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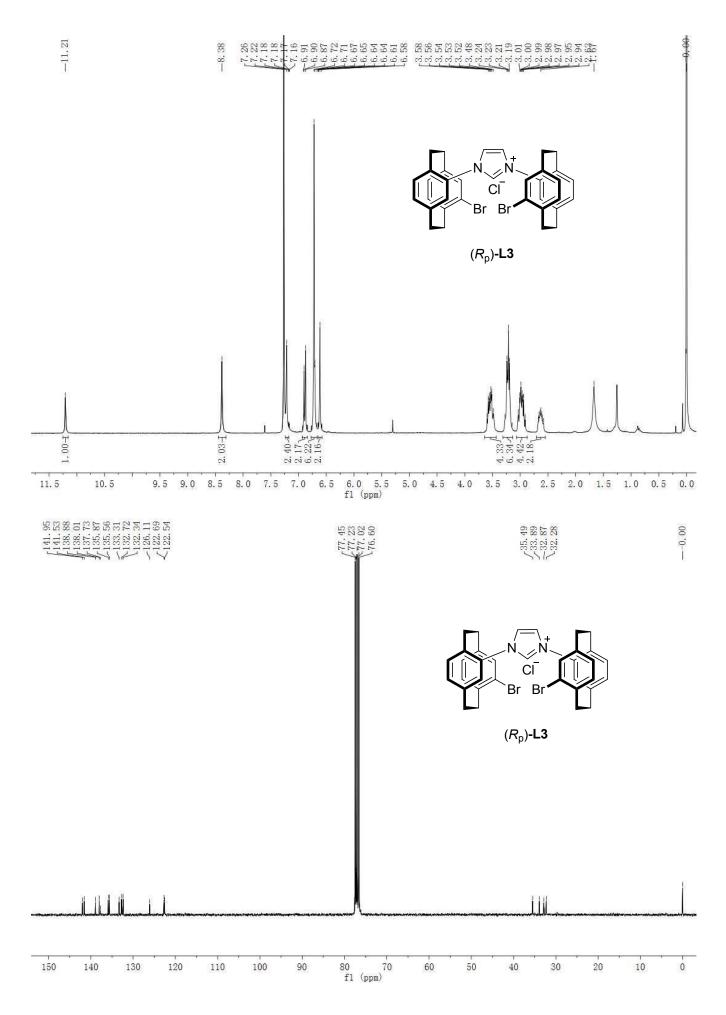
#### **1. General Information**

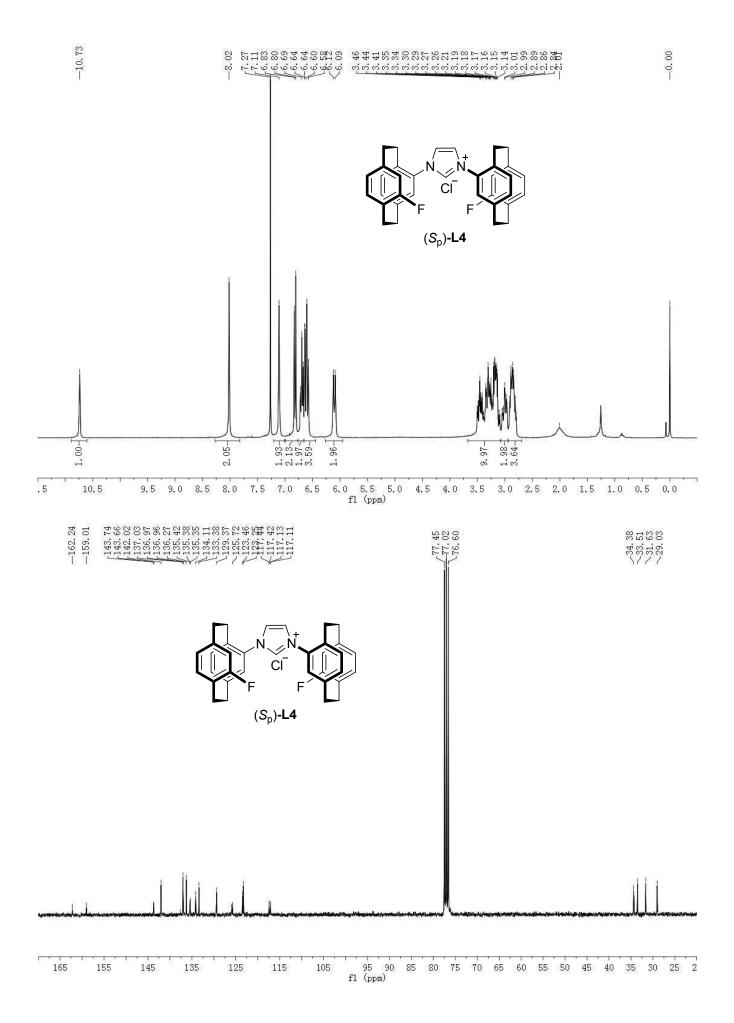
All manipulations of oxygen- and moisture-sensitive materials were conducted with a standard Schlenk technique under a purified argon atmosphere. Solvents were reagent grade and purified by standard techniques. <sup>1</sup>H and <sup>13</sup>C NMR spectra were recorded on a 300 MHz spectrometer. Chemical shifts were reported as  $\delta$  values in ppm and referenced to using tetramethylsilane as an internal. Data are reported as (s = singlet, d = doublet, dd = doublet of doublet, ddd = doublet of doublet of doublet, t = triplet, dt = doublet of triplet, td = triplet of doublet, q = quartet, qd = quartet of doublet, m = multiplet, brs = broad singlet). All high resolution mass spectra were obtained on an Accurate-Mass LC-ESI-Q-TOF-MS. All reactions were monitored by TLC with silica gel-coated plates and visualized with a UV light at 254 nm. Optical rotations were reported as follows:  $[\alpha]_D^T$  (c = g/100 mL, solvent).

#### 2. NMR spectras and HPLC chromatograms

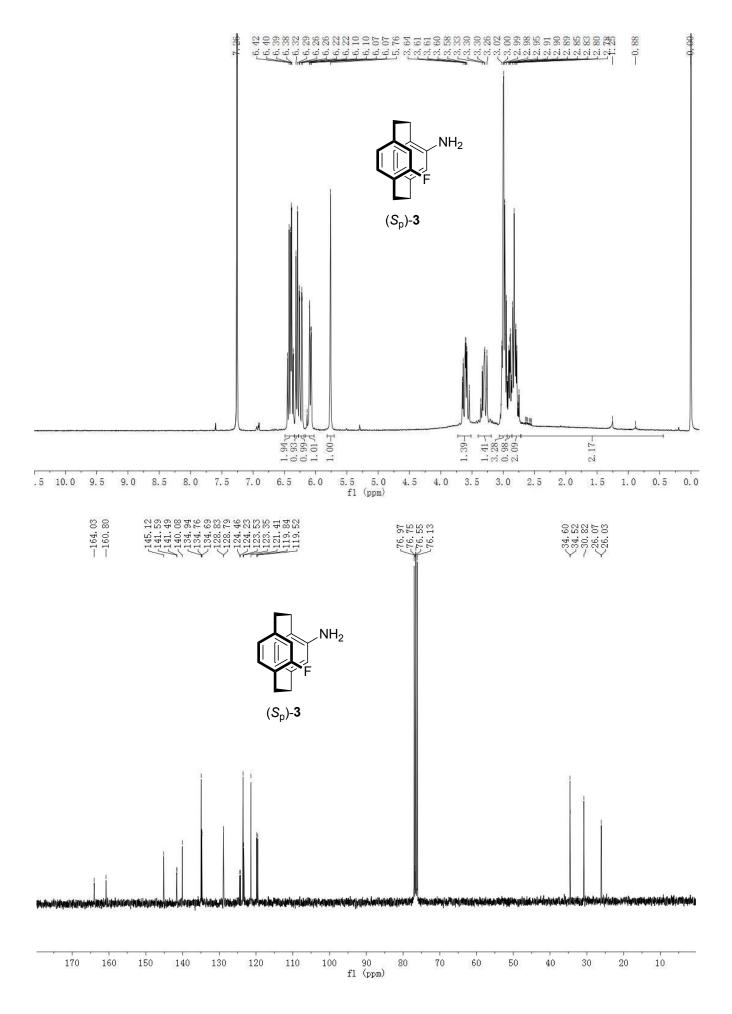


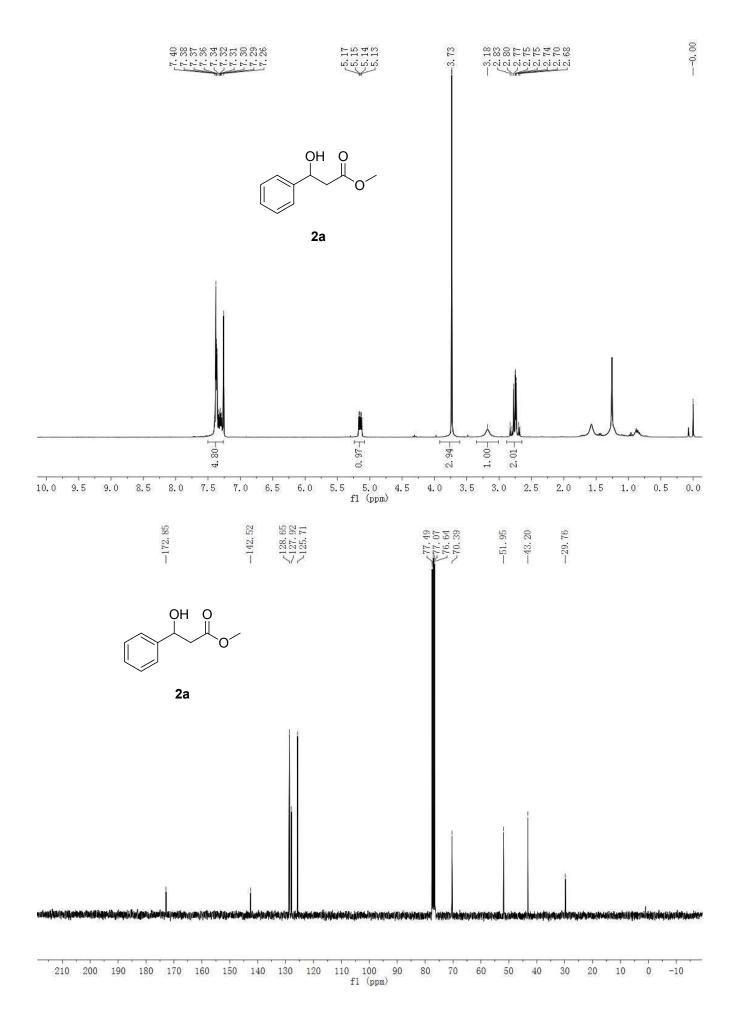


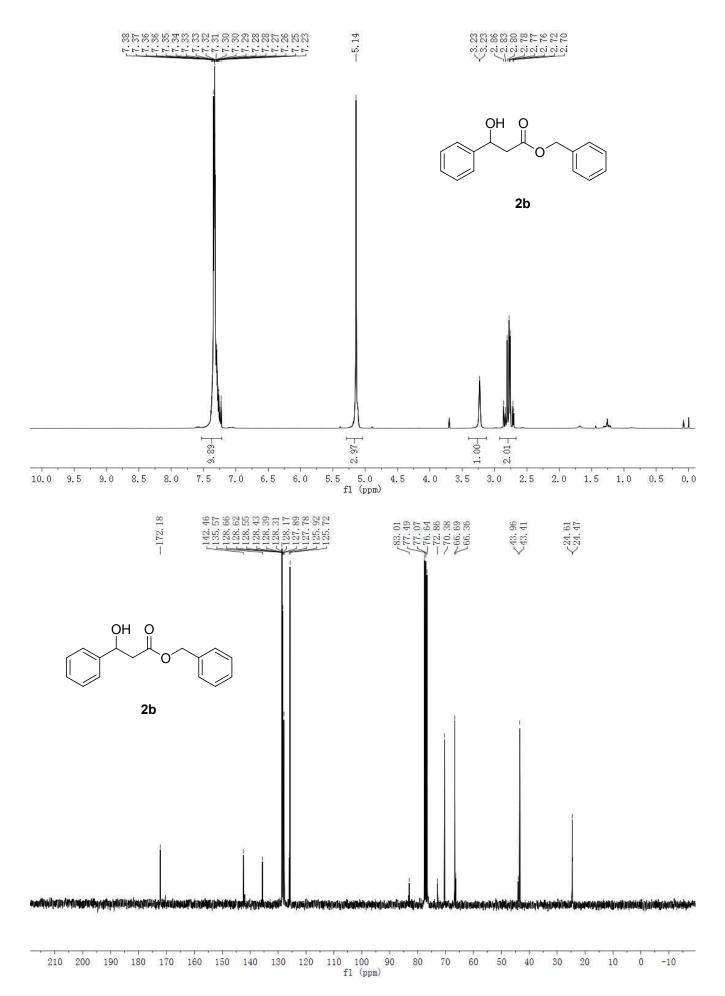


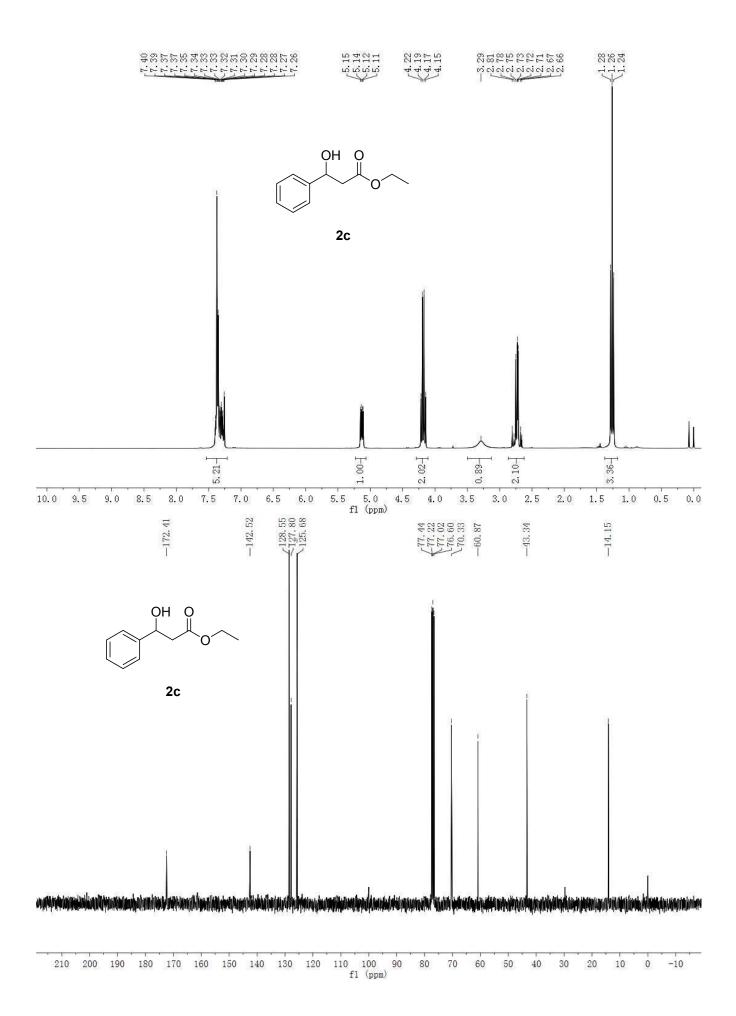


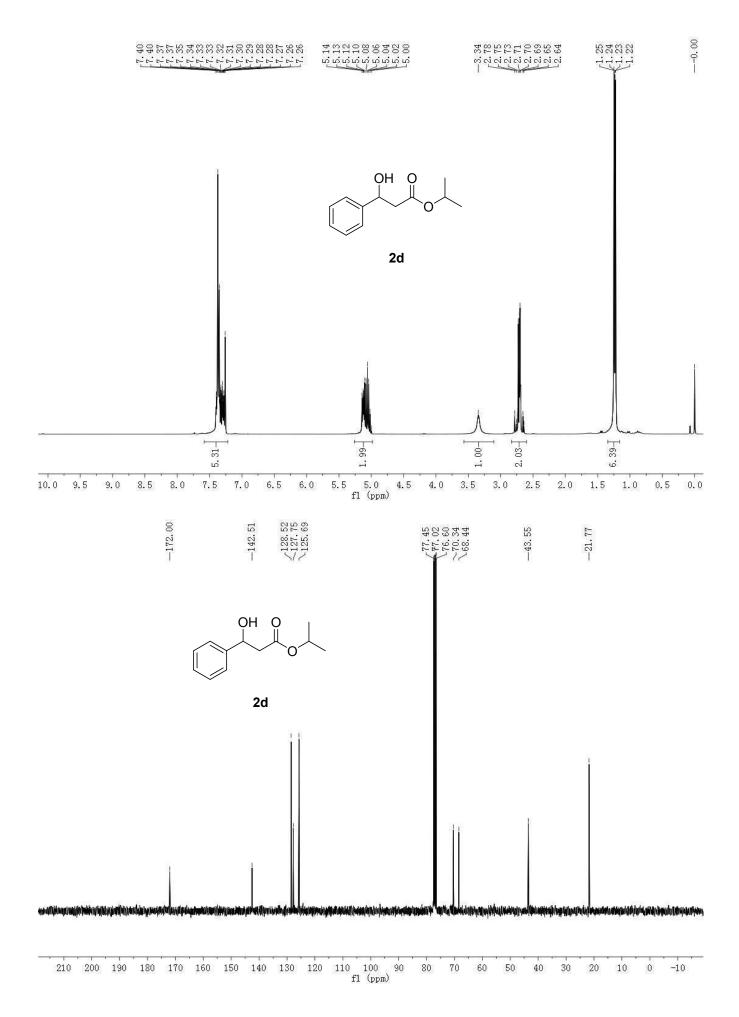
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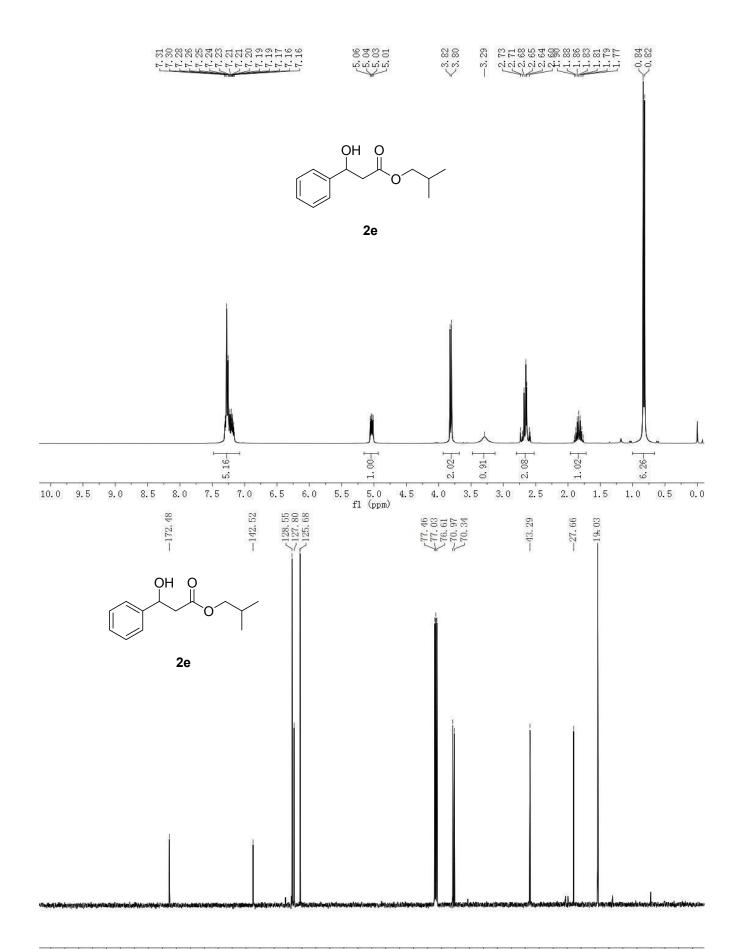


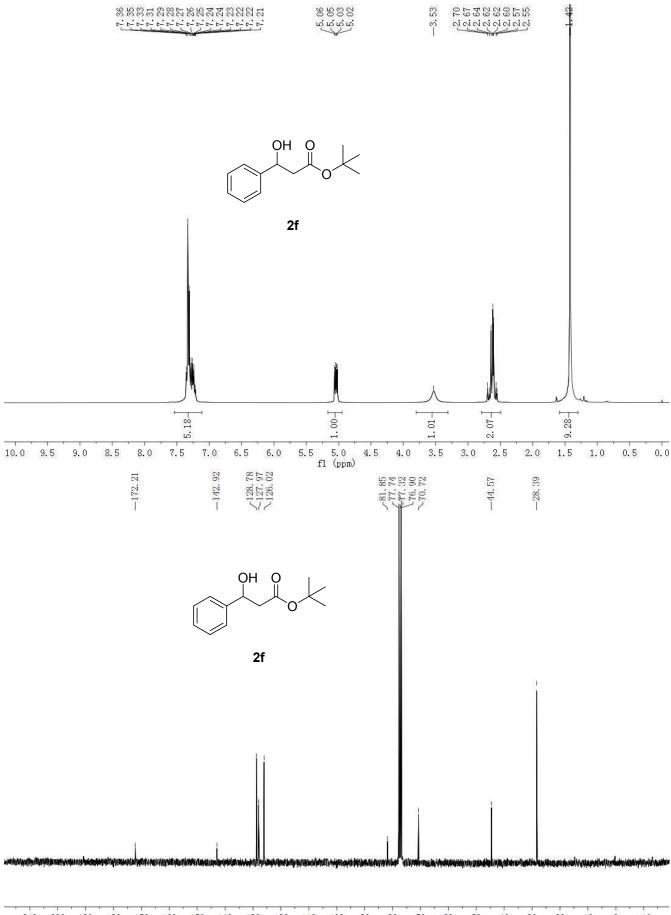


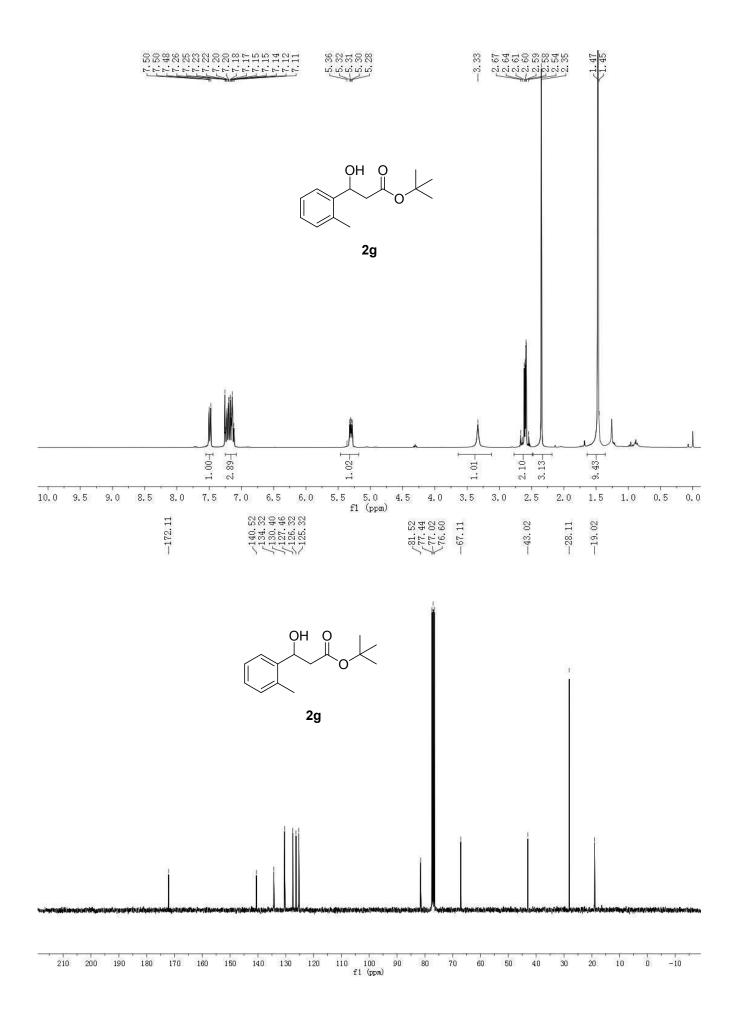




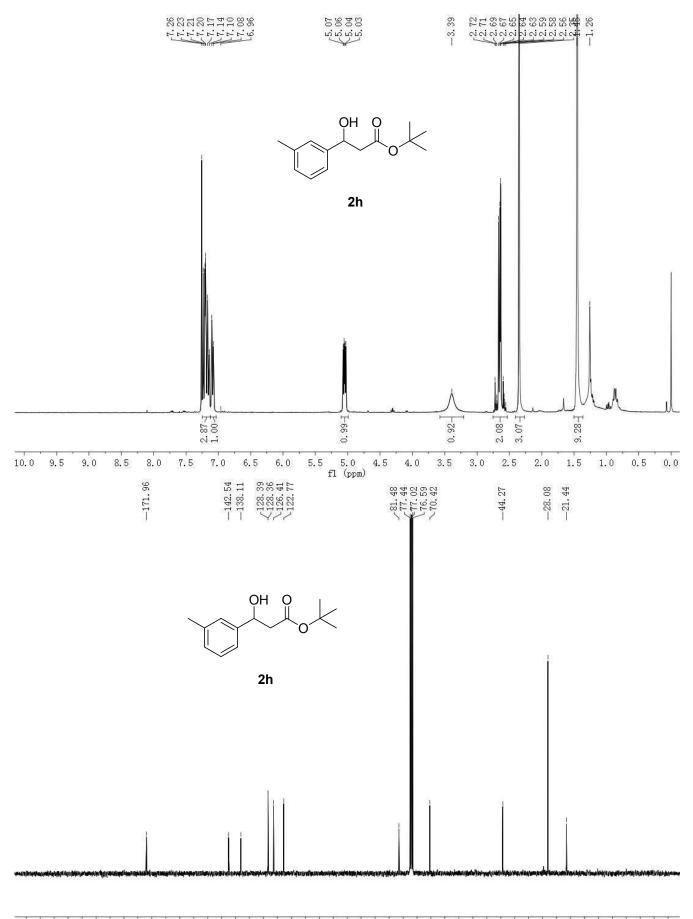




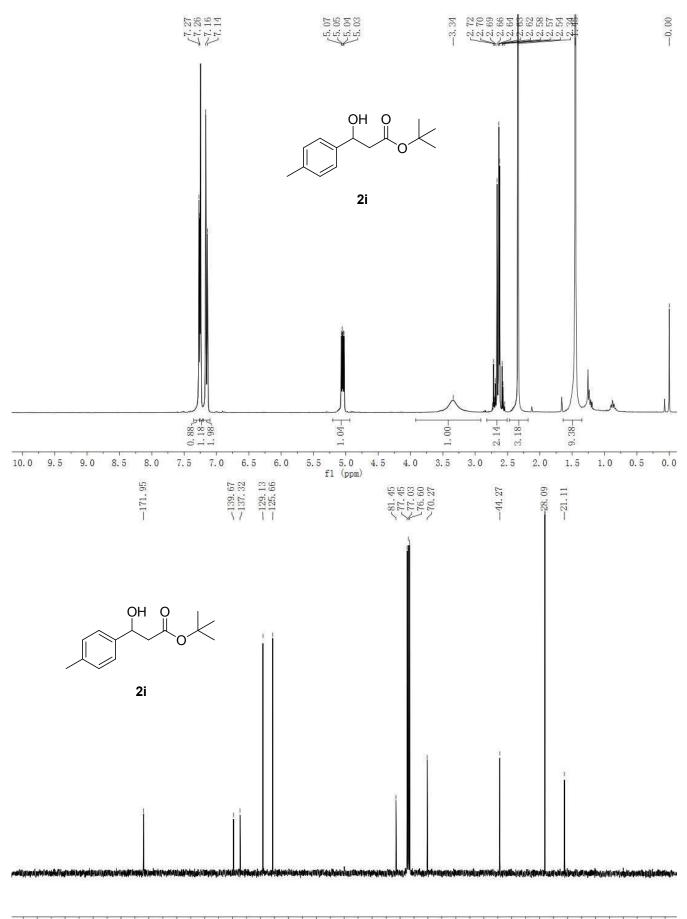


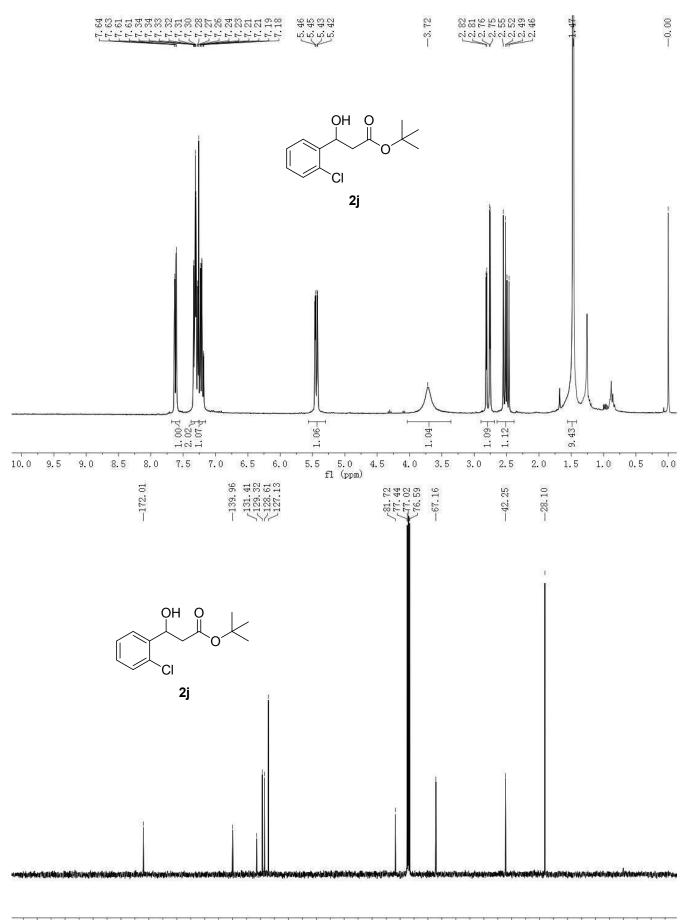


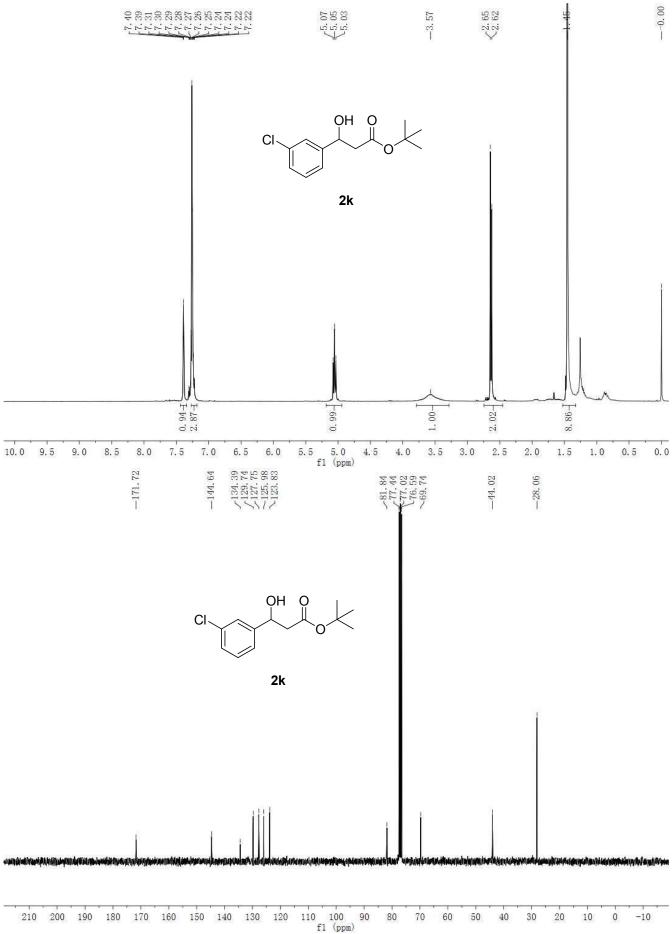
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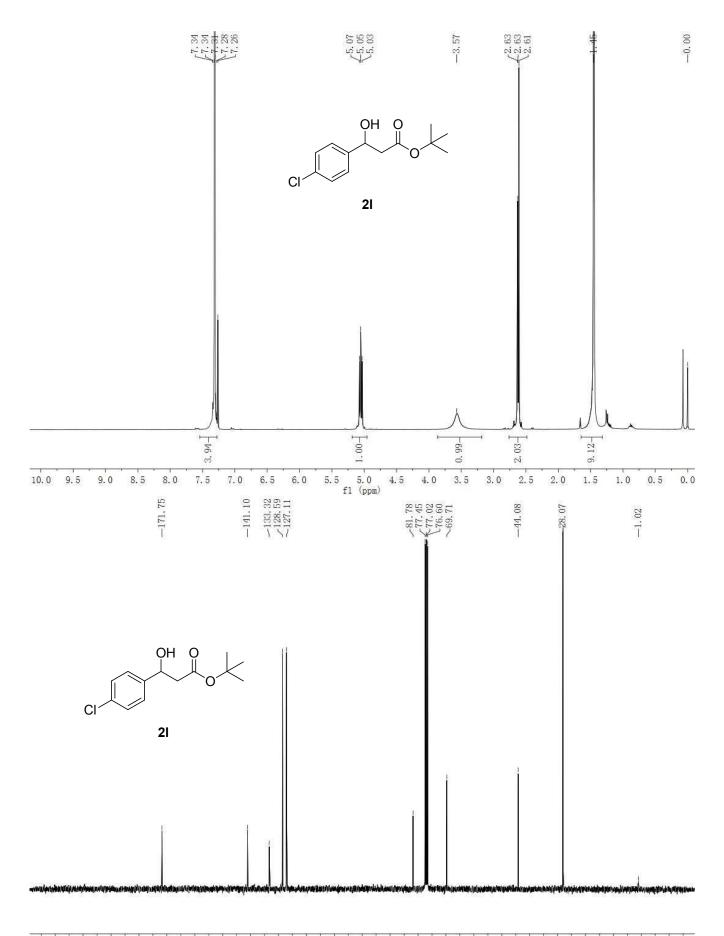


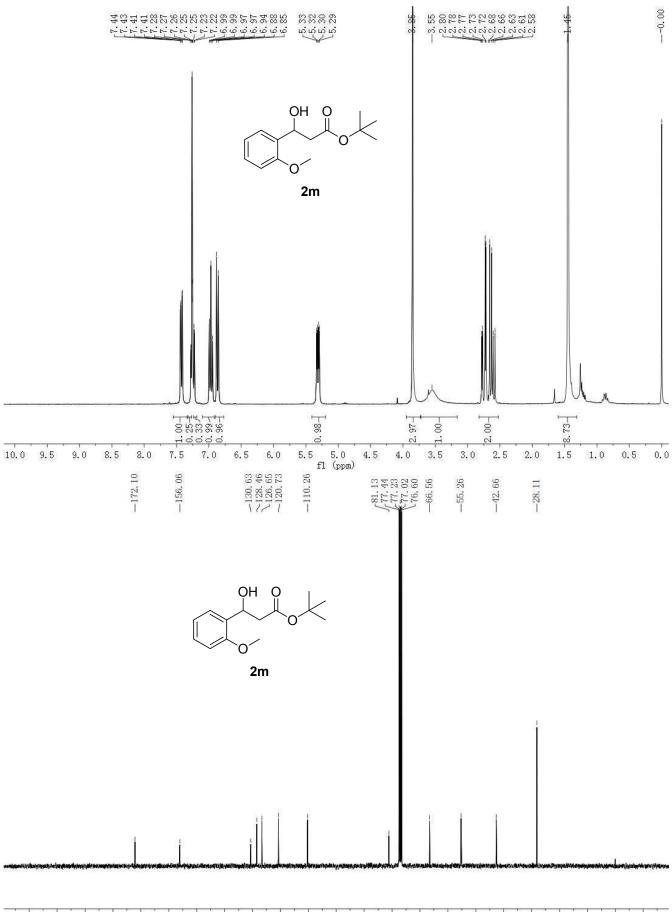
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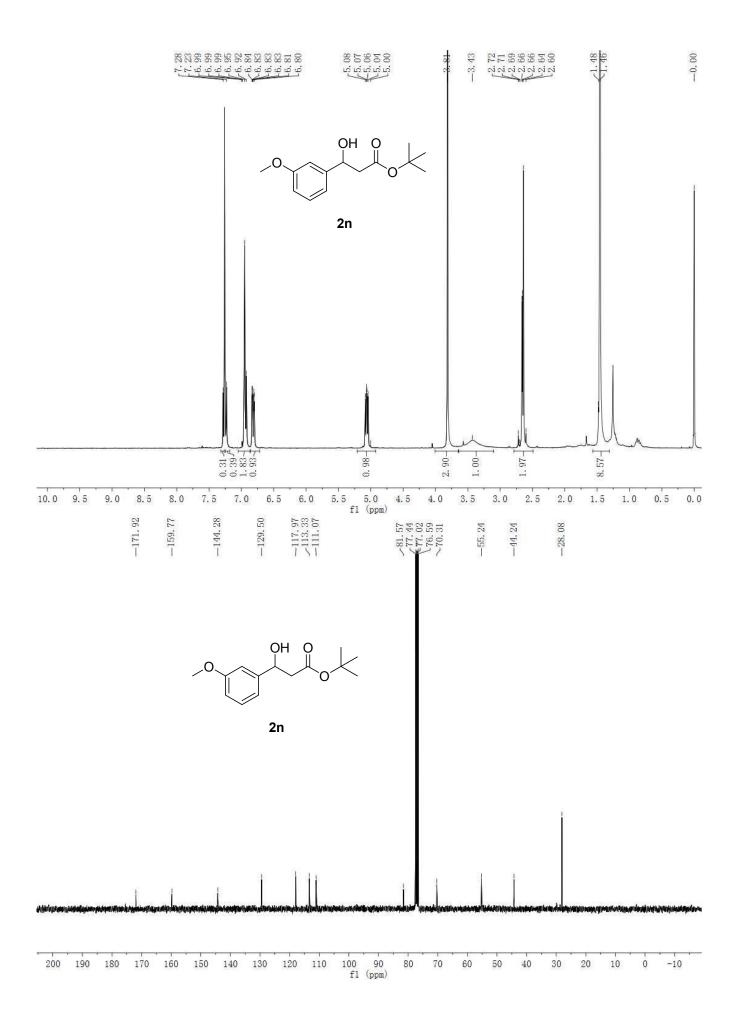


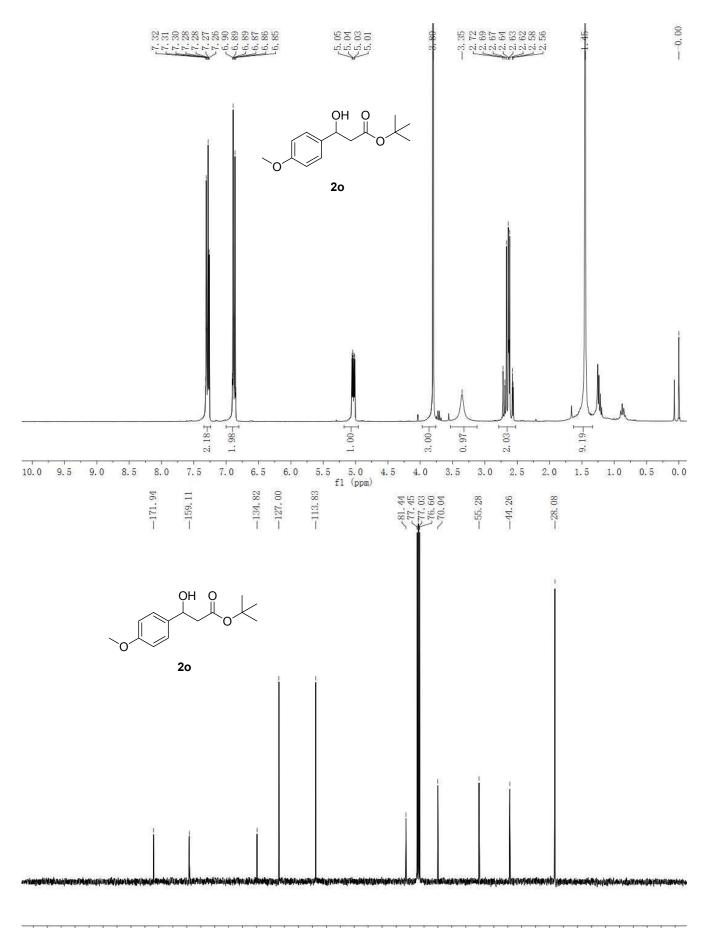


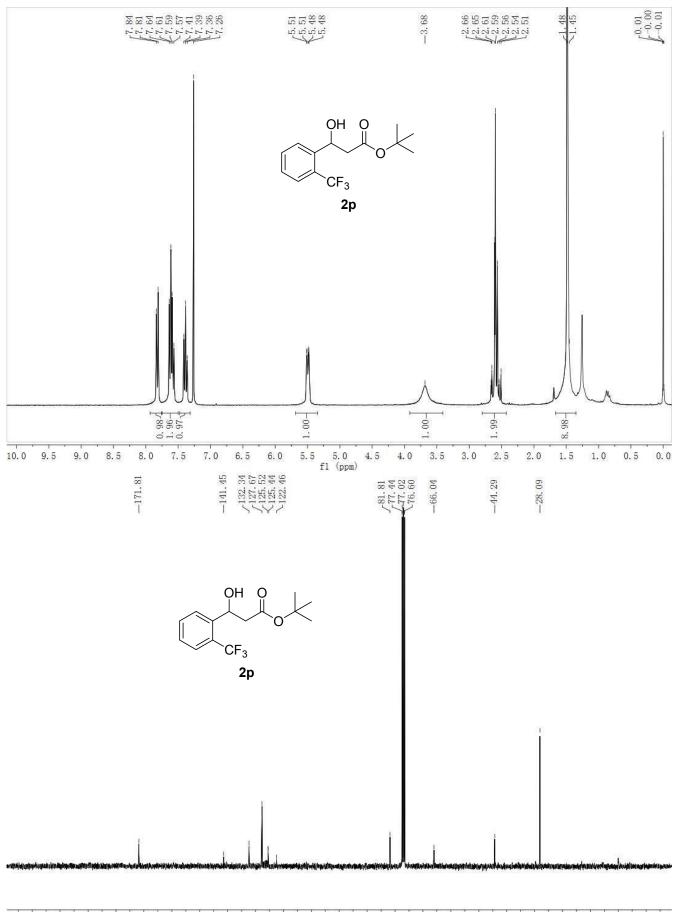




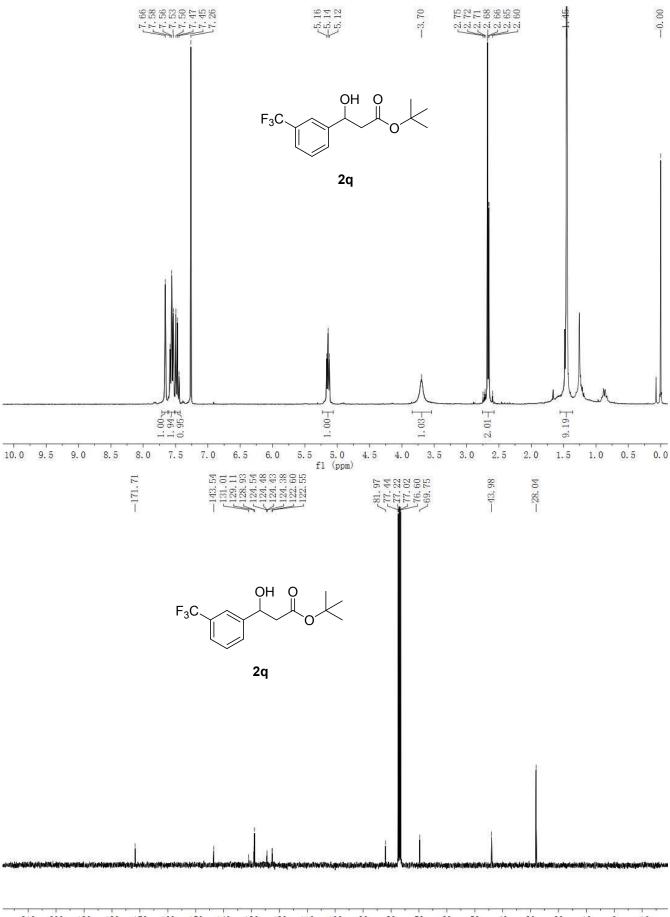


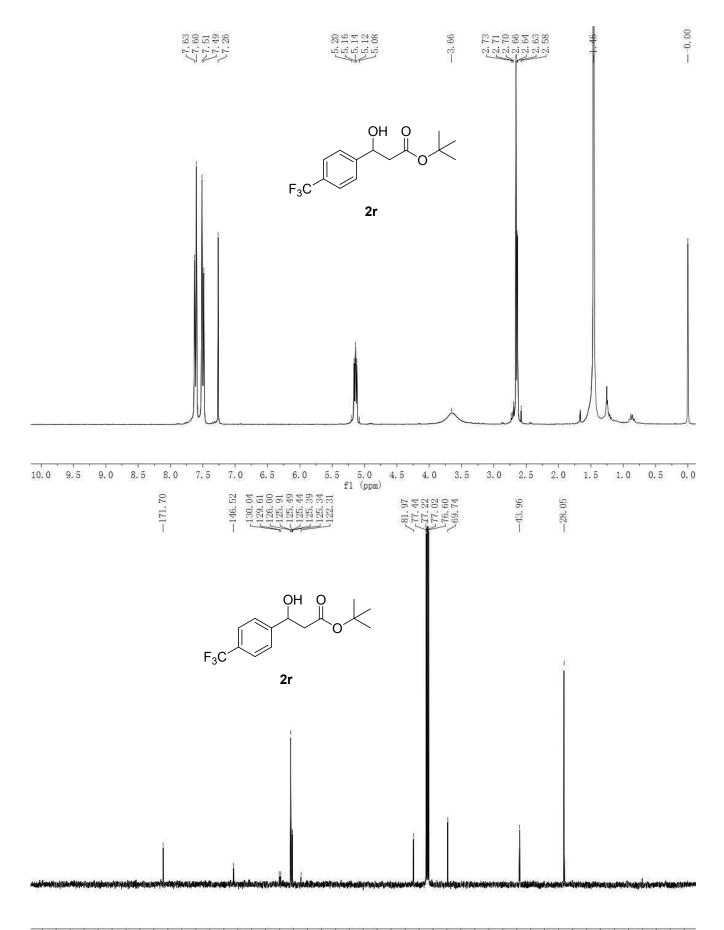


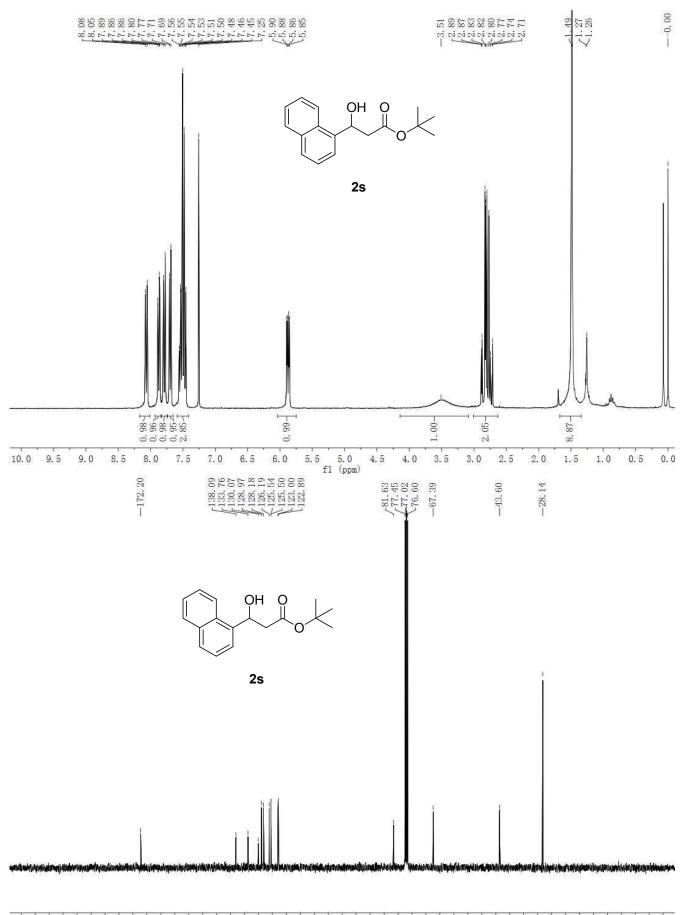


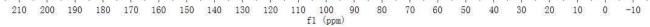


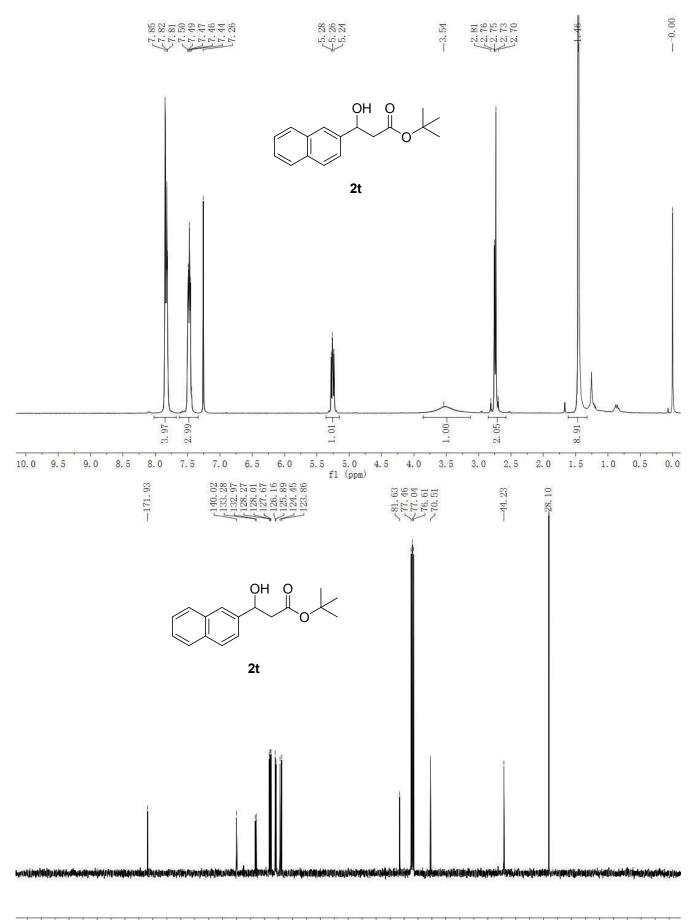
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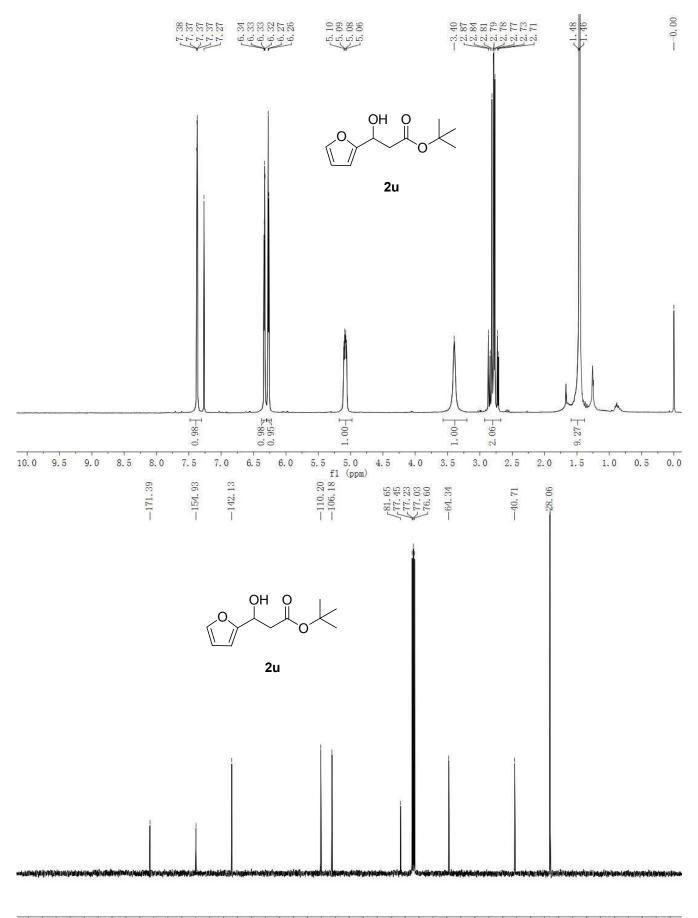


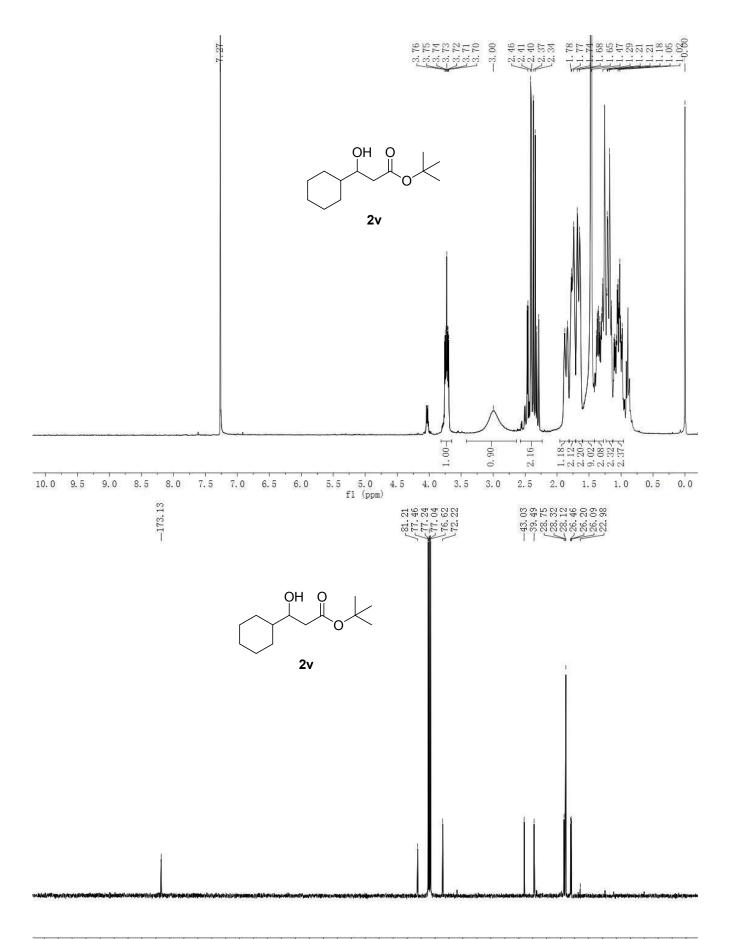


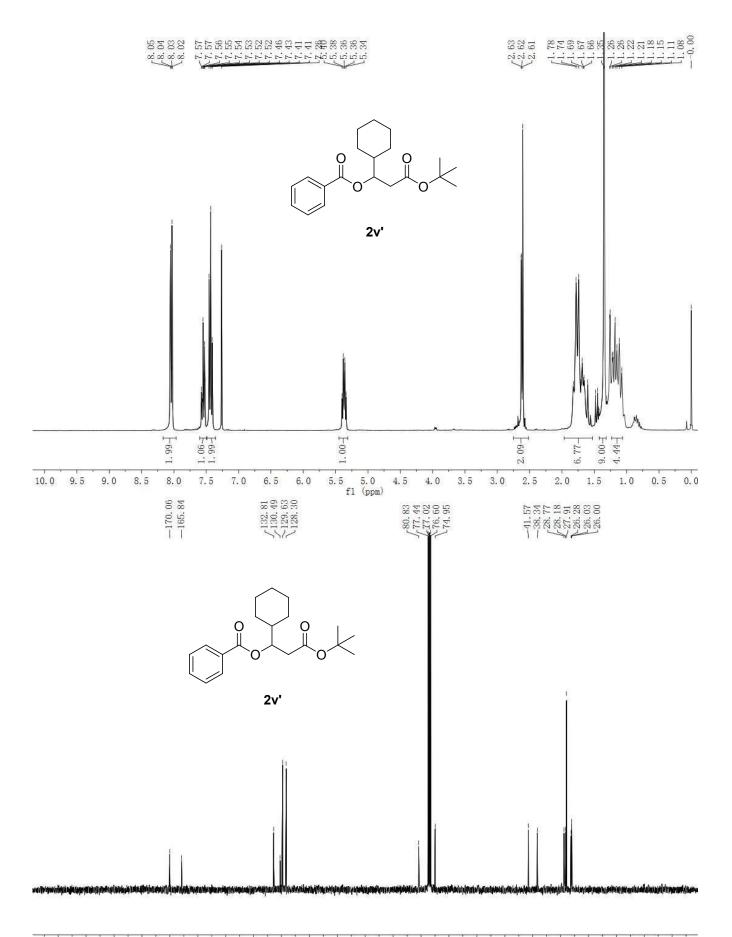


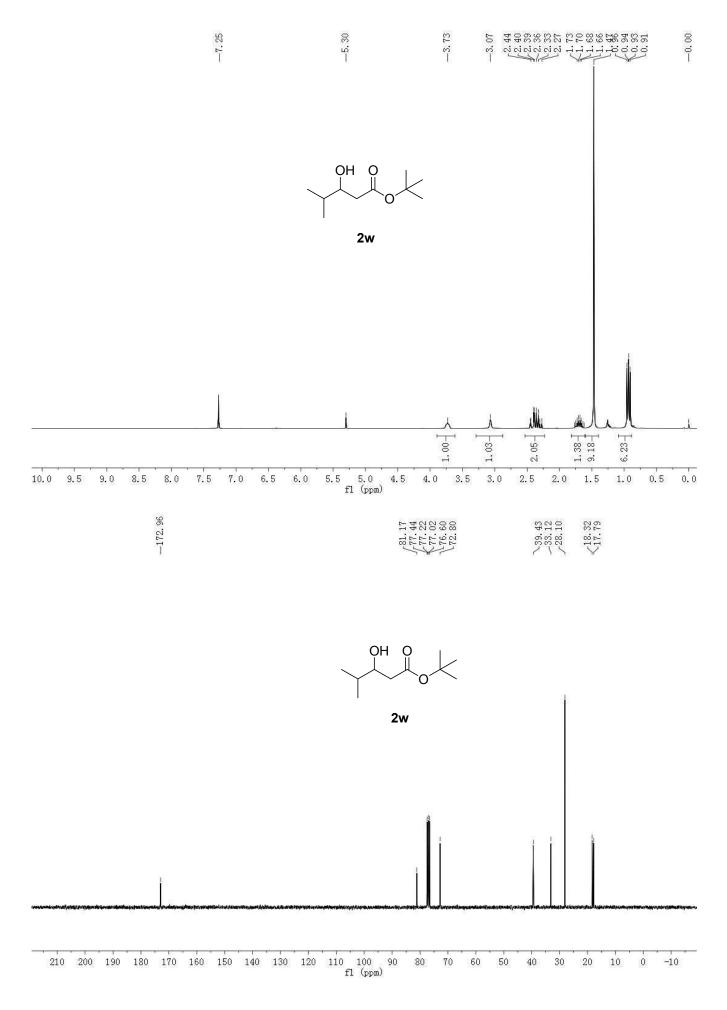


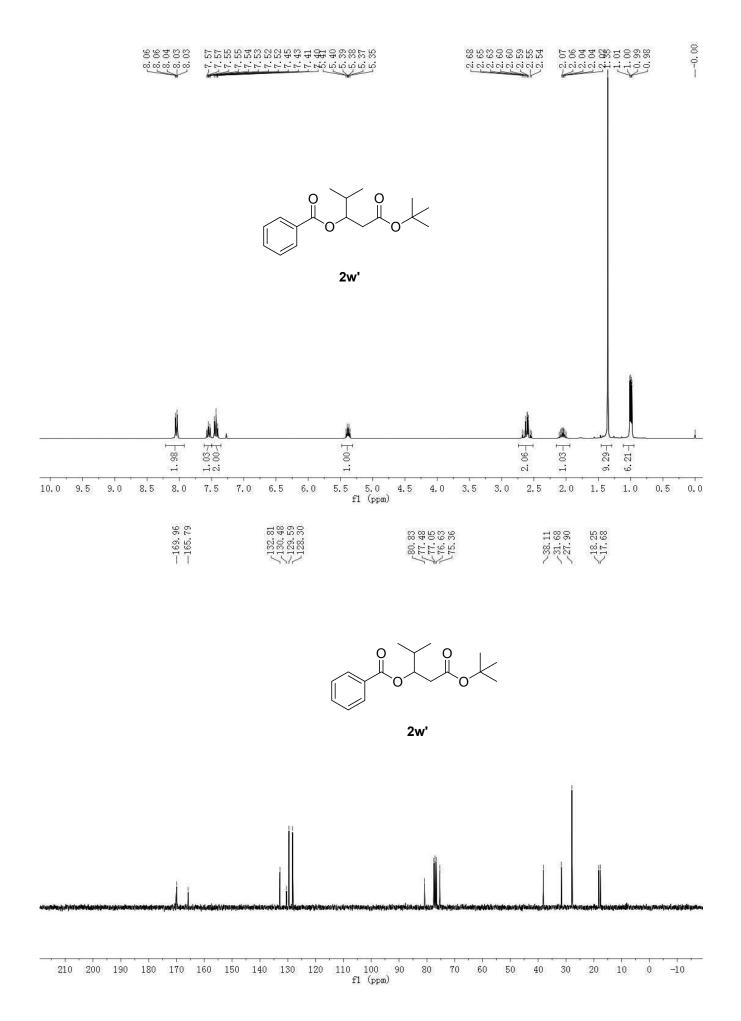


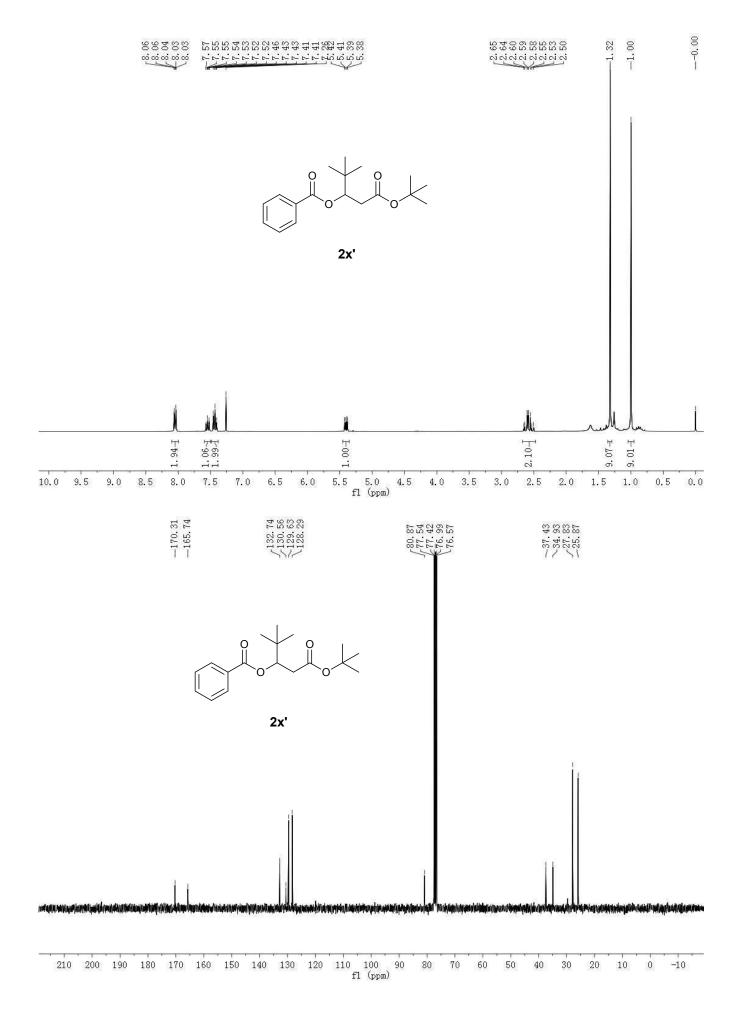


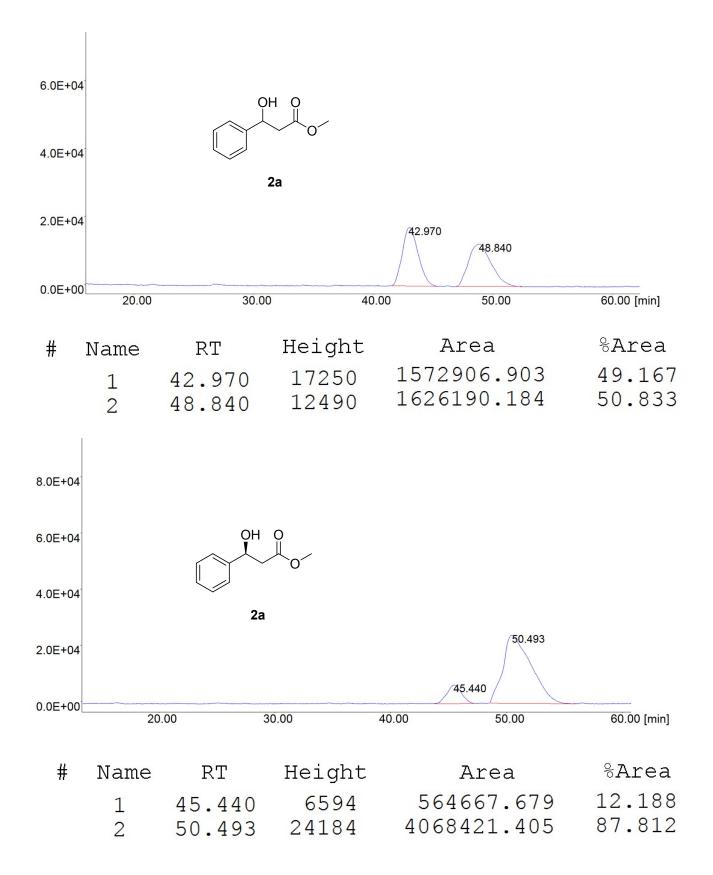


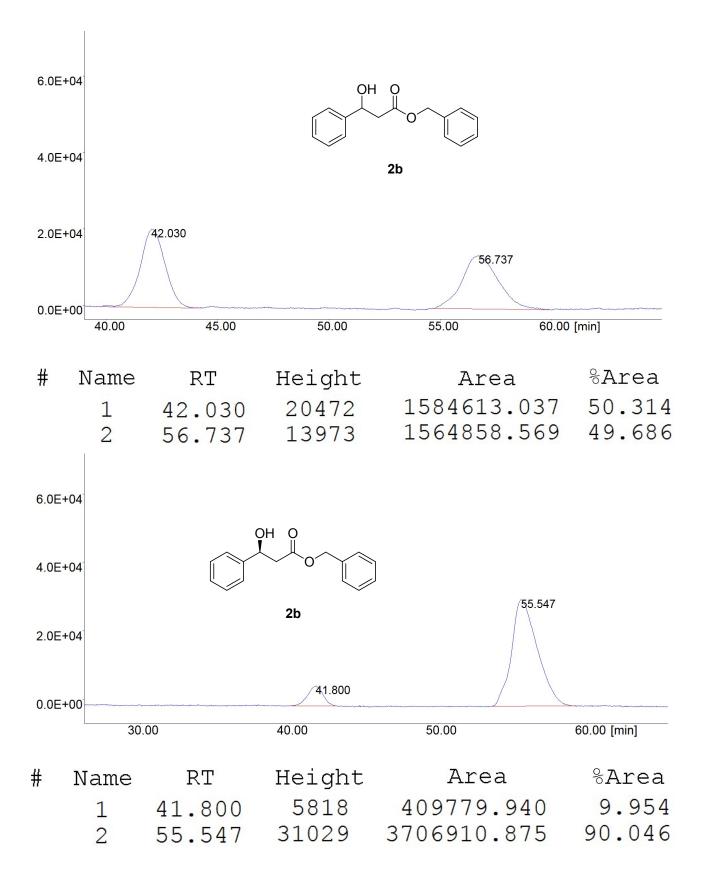


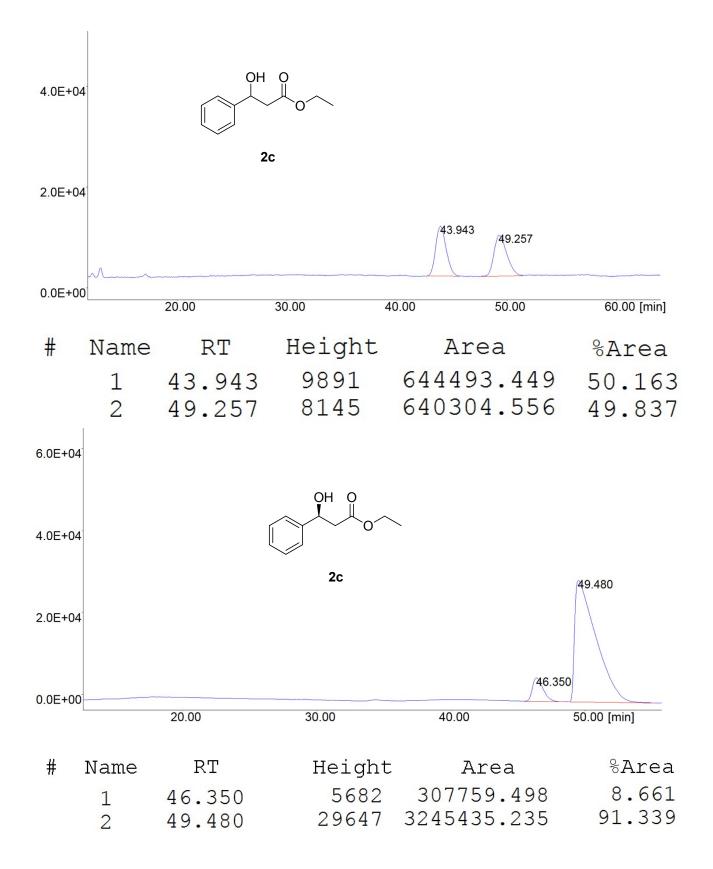


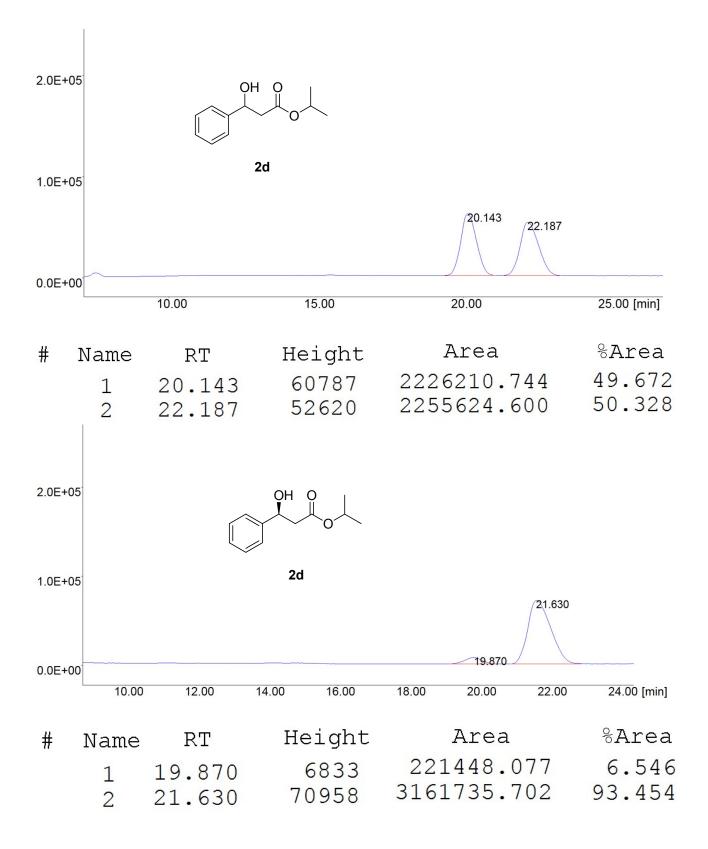


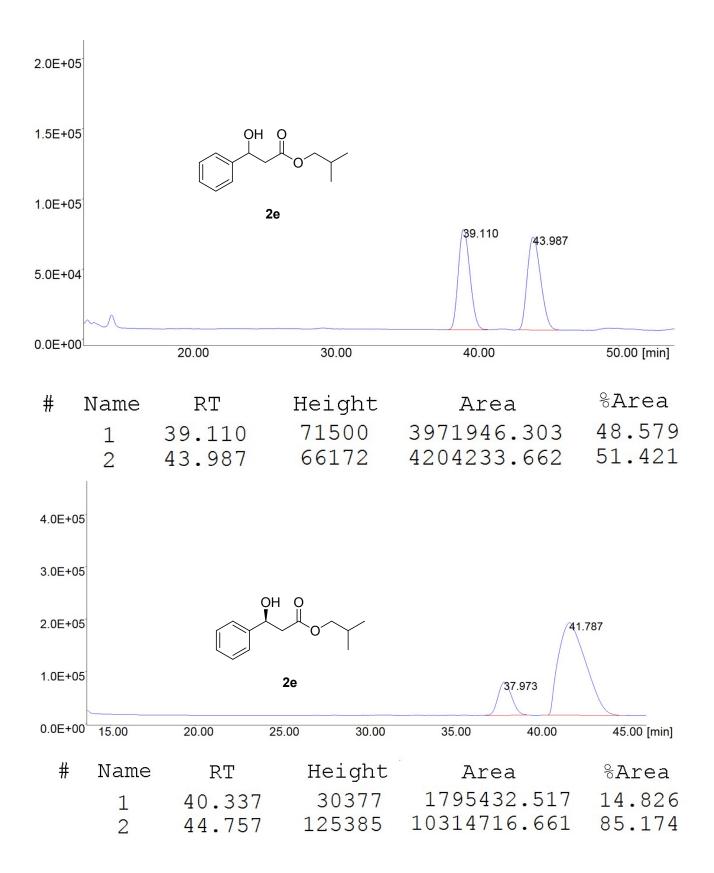


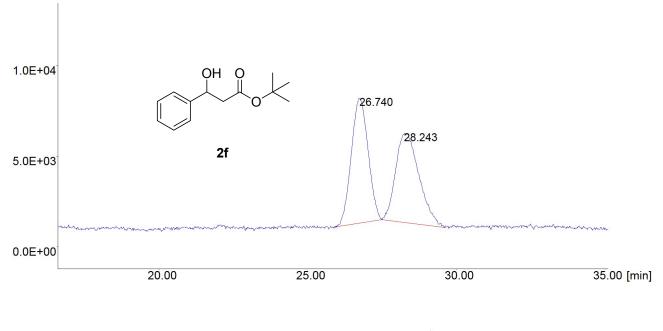




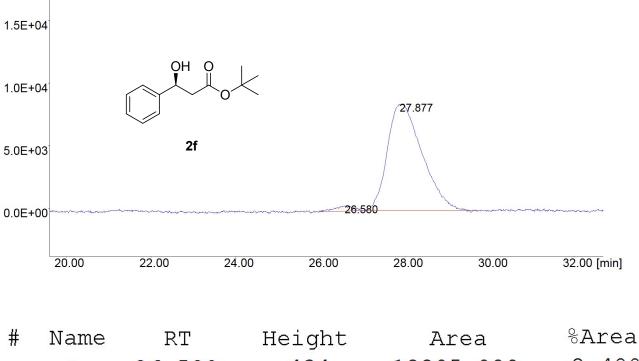




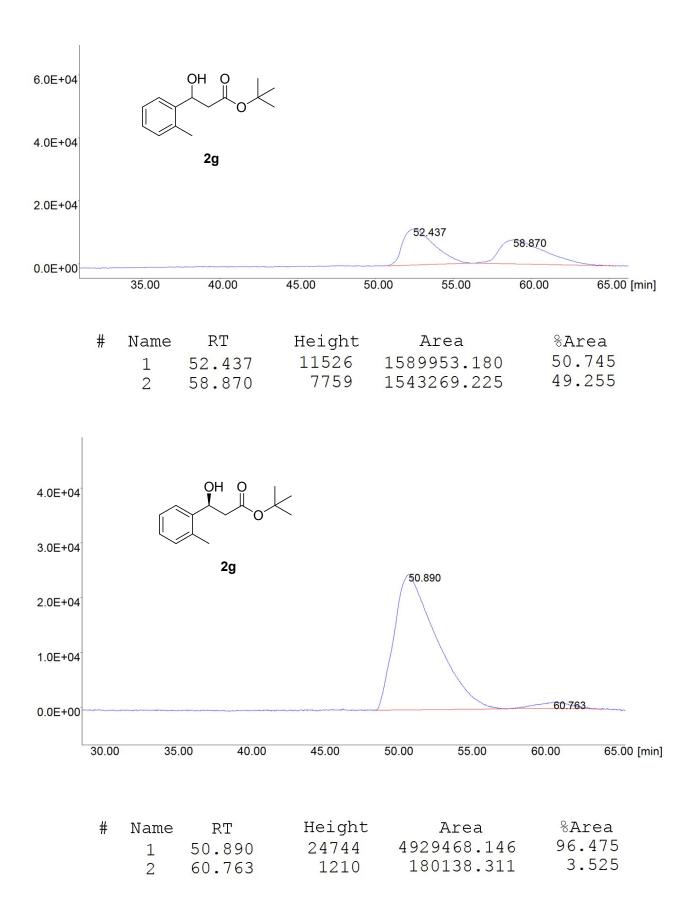


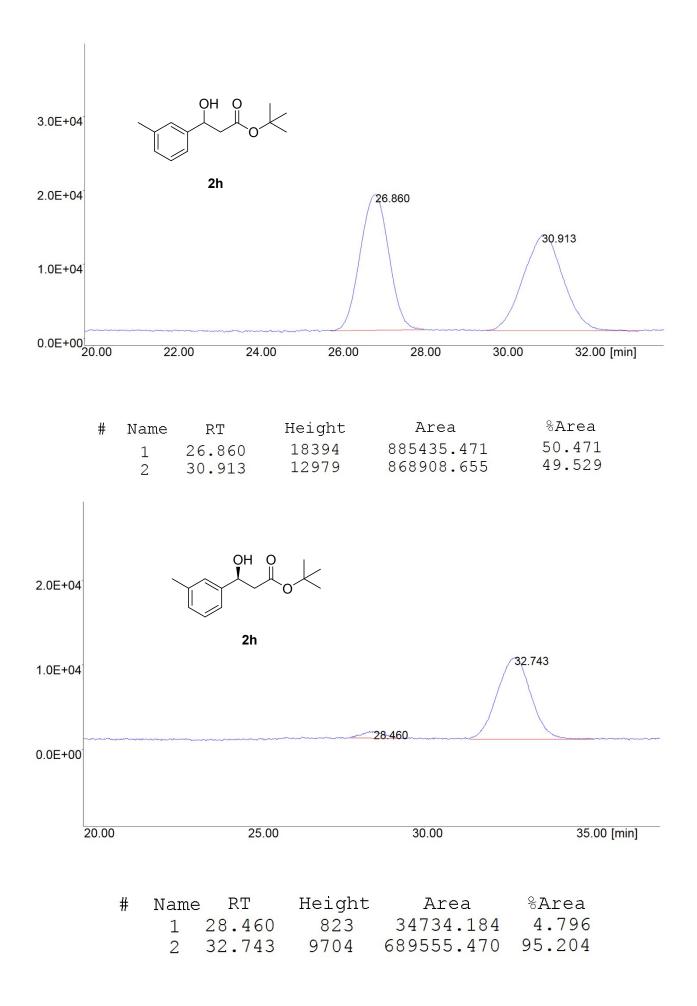


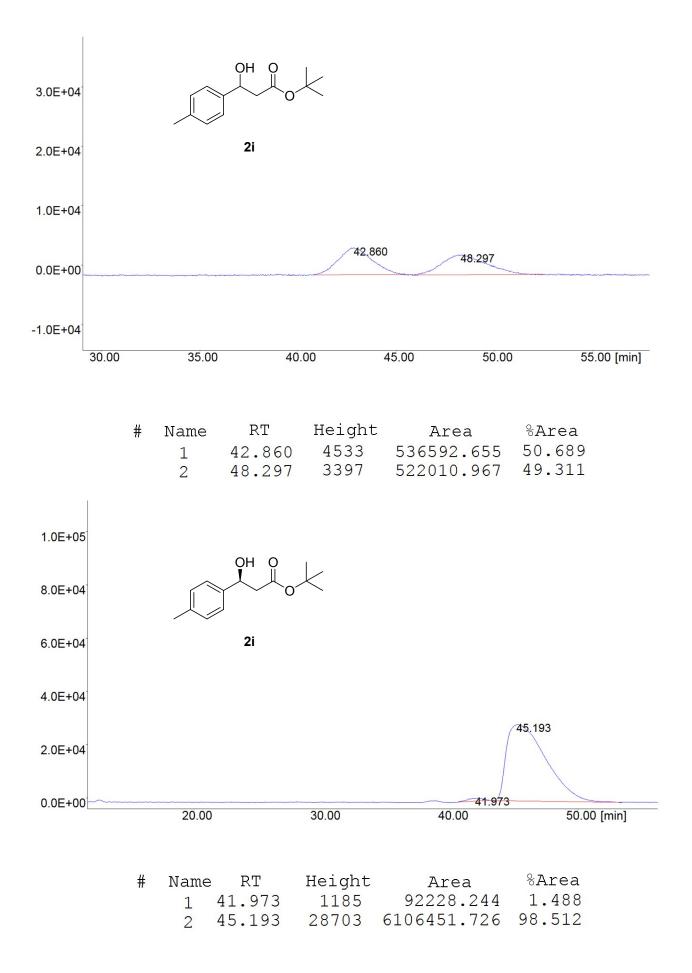
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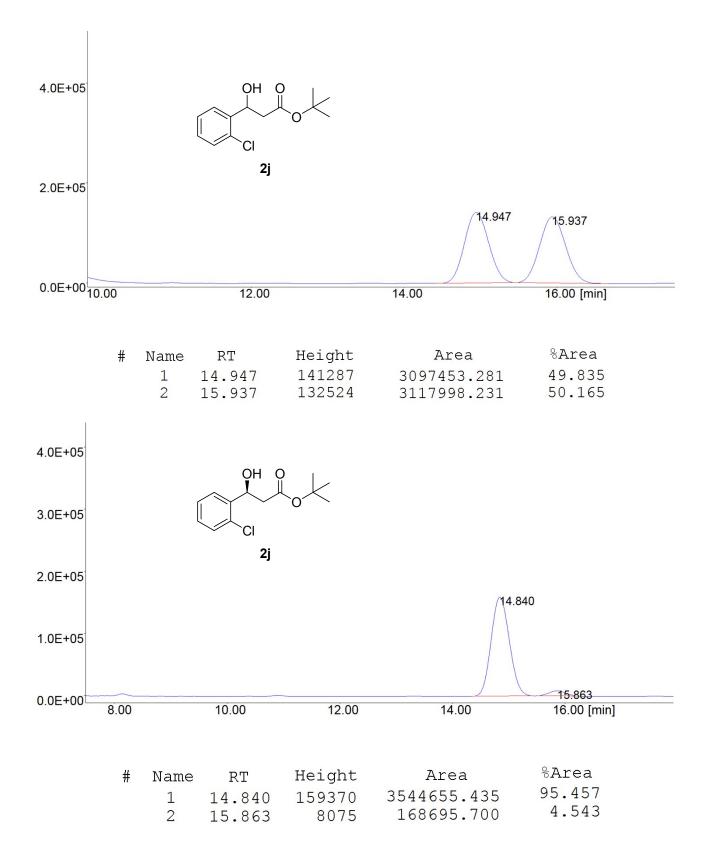


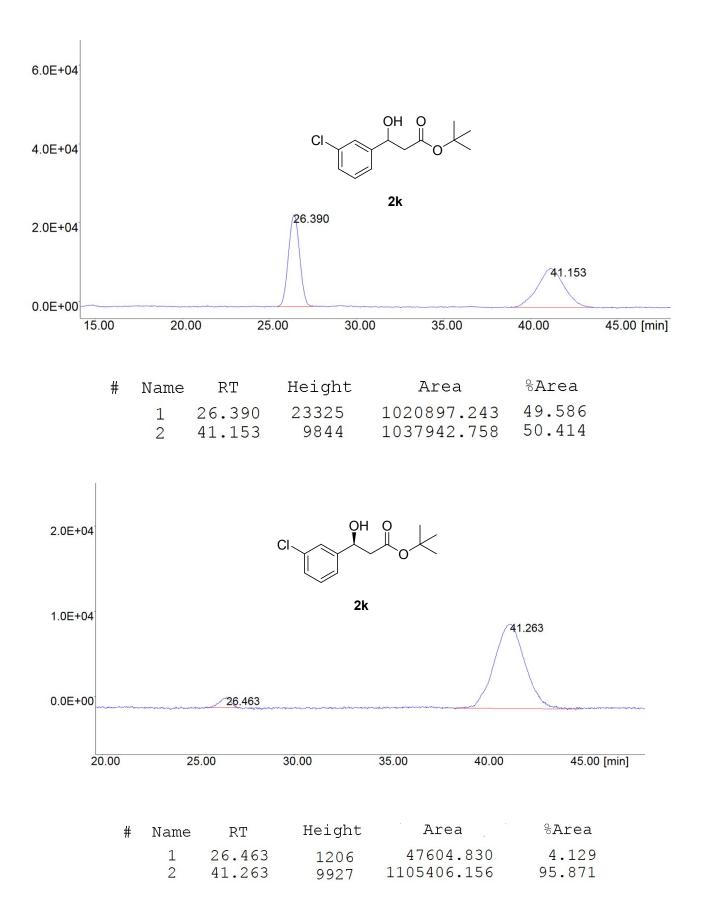
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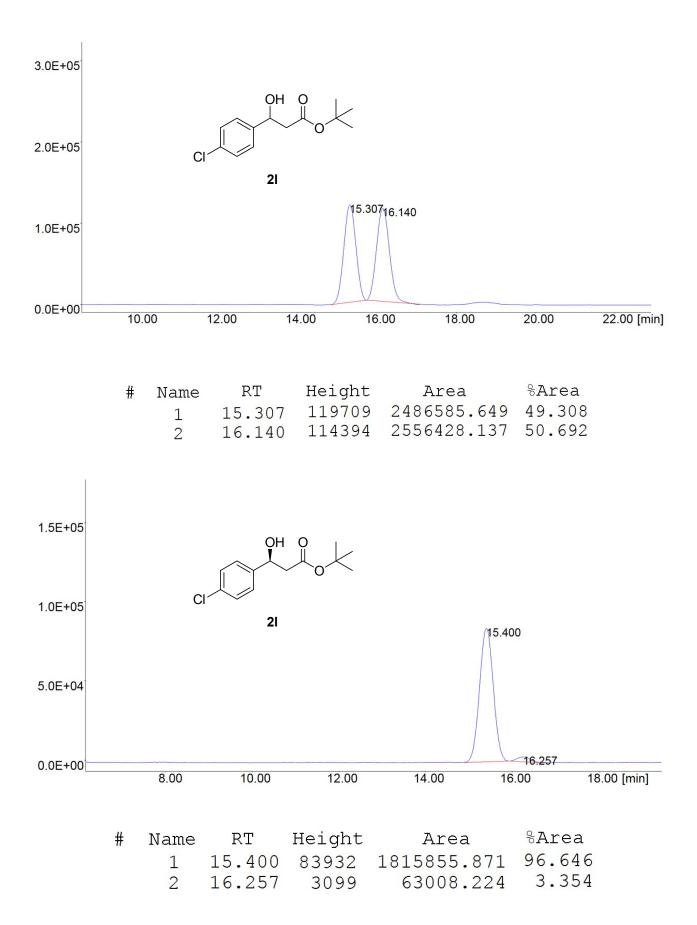


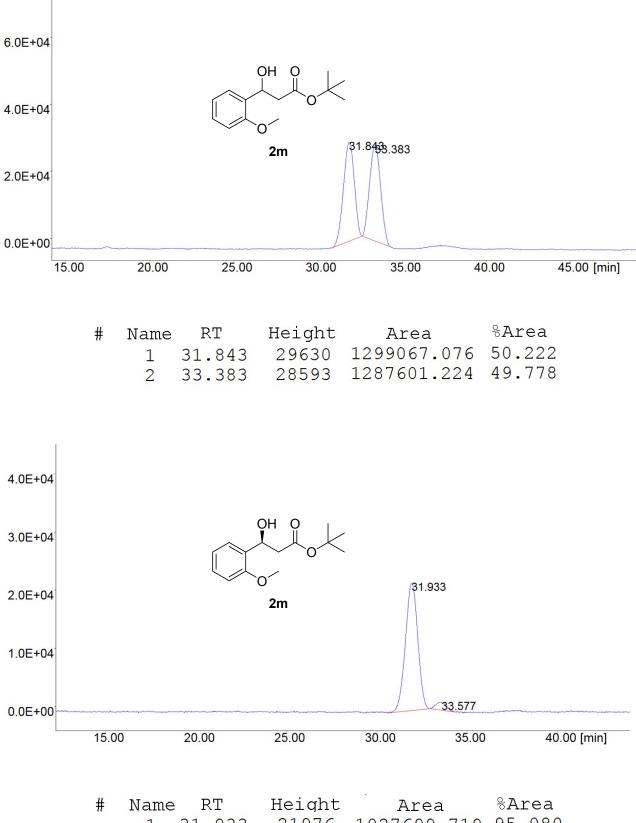




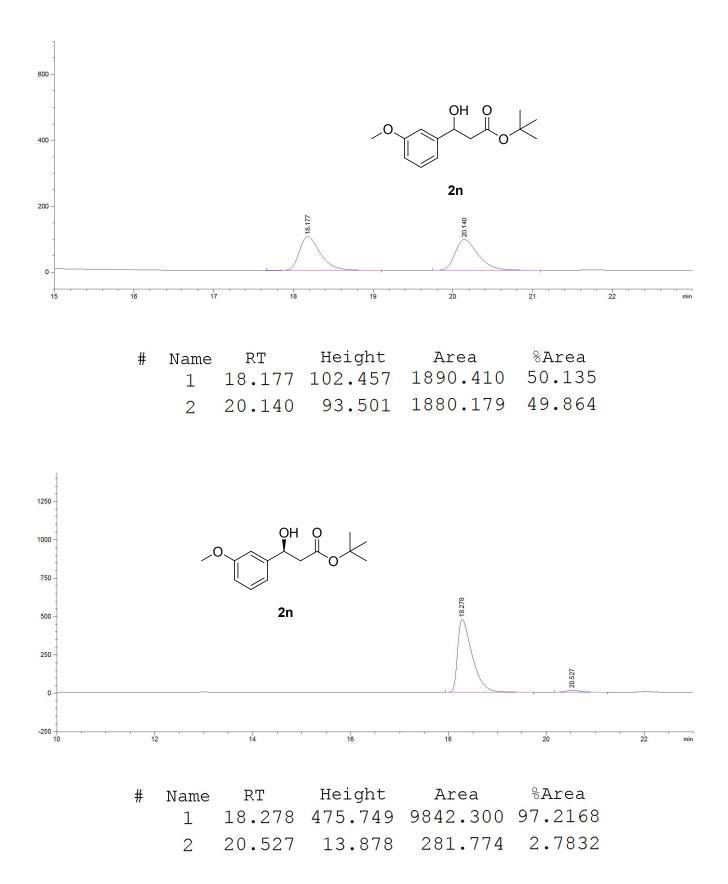


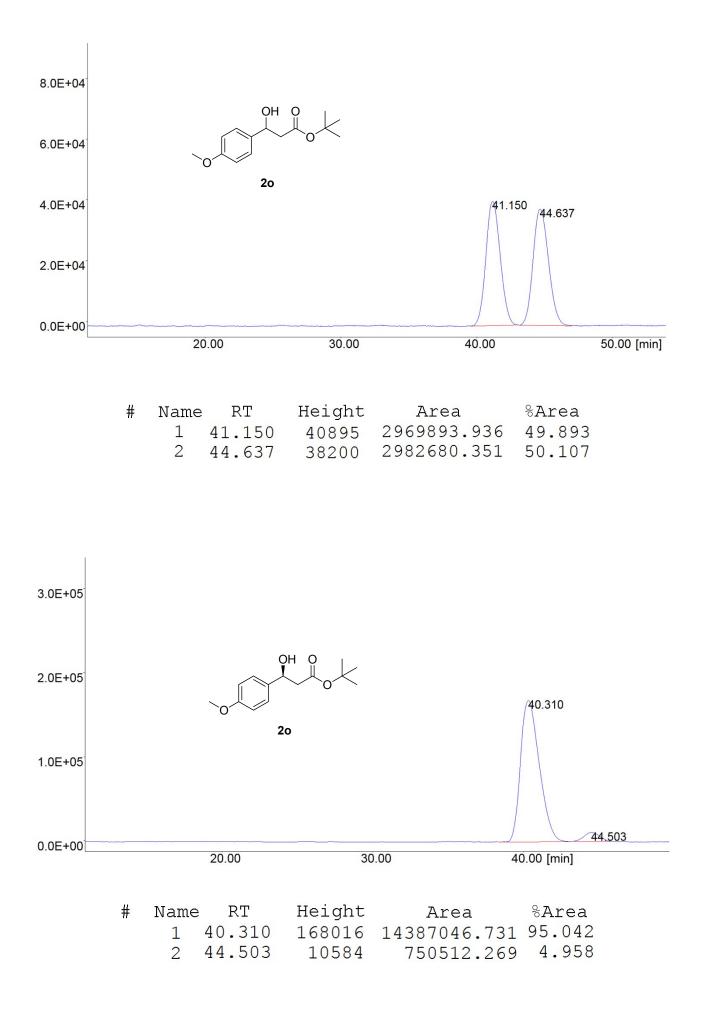


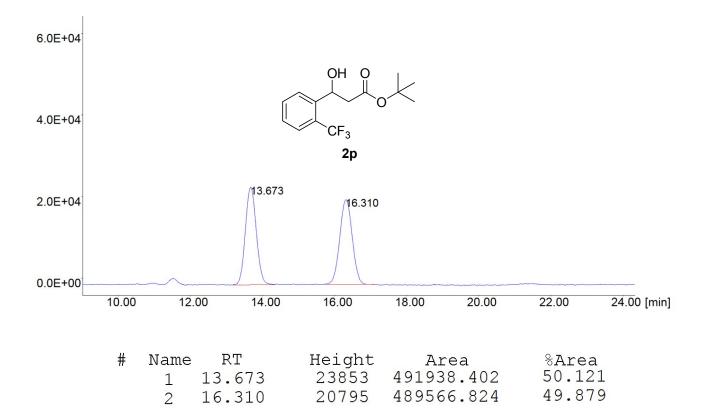


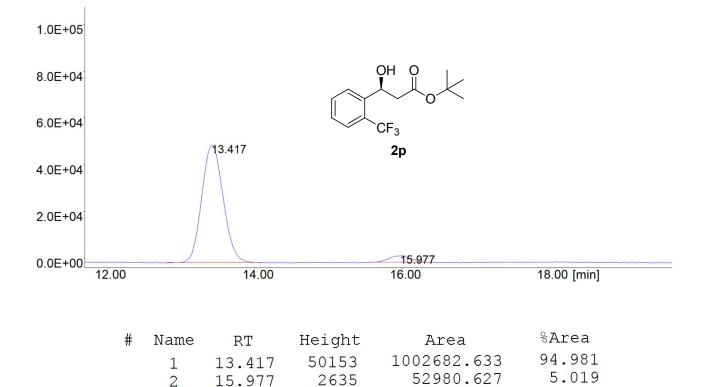


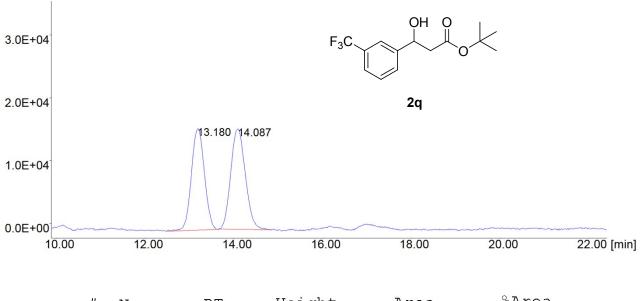
L	31.933	21976	102/699./19	95.080
2	33.577	1381	53183.528	4.920



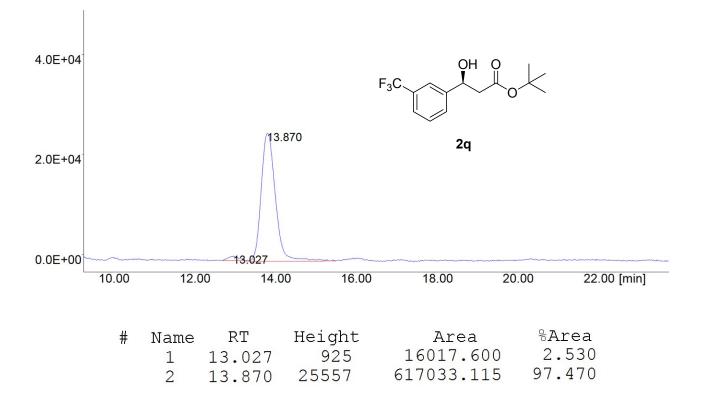


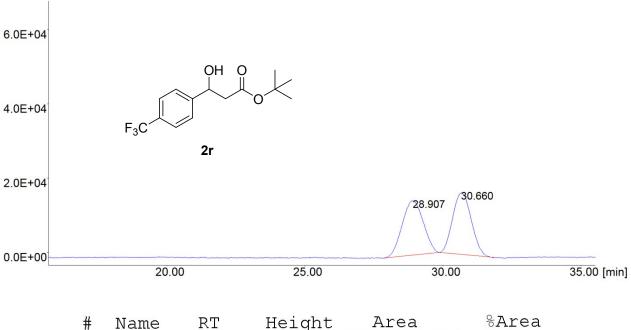




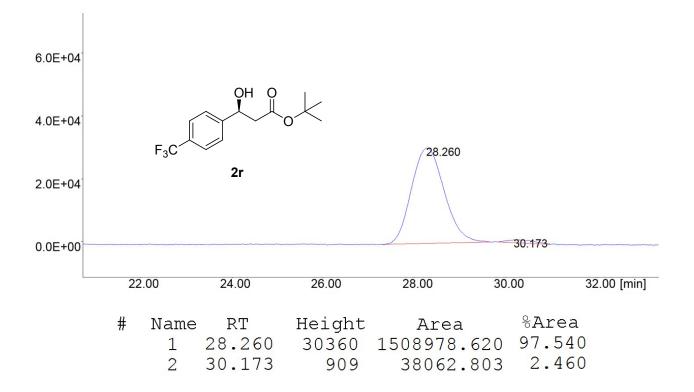


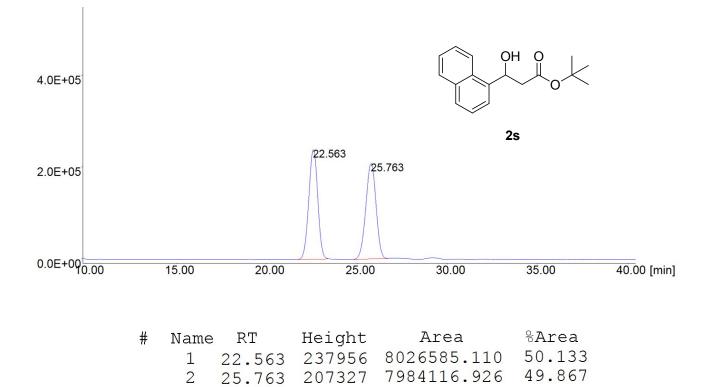
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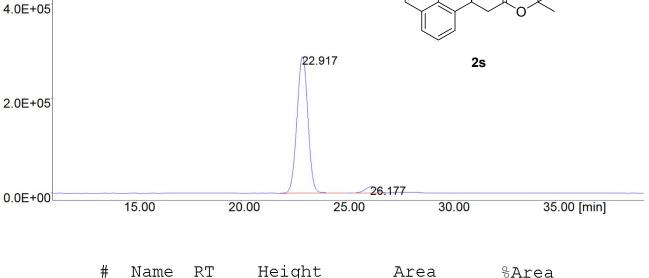


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2	30.660	16647	773290.619	49.990

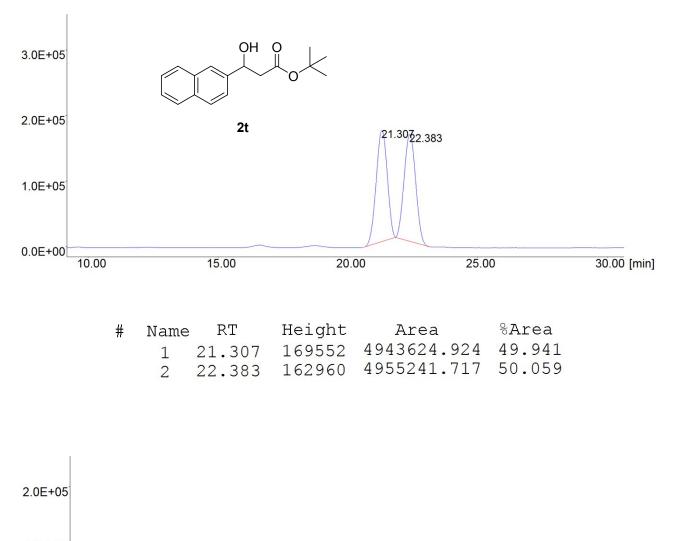


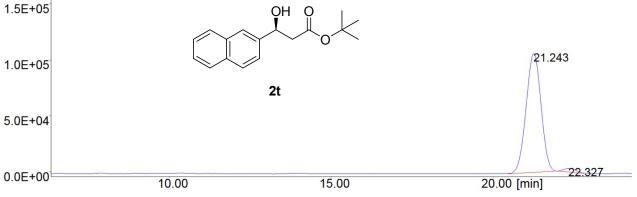


)E+05		OH O U O
	22.917	2s

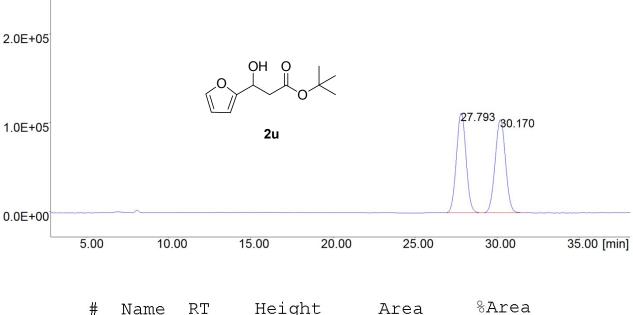


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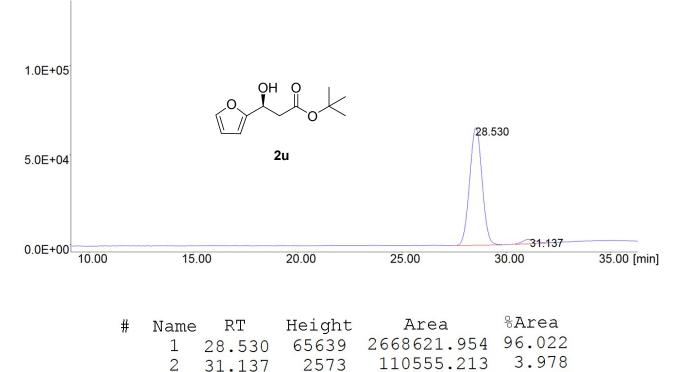


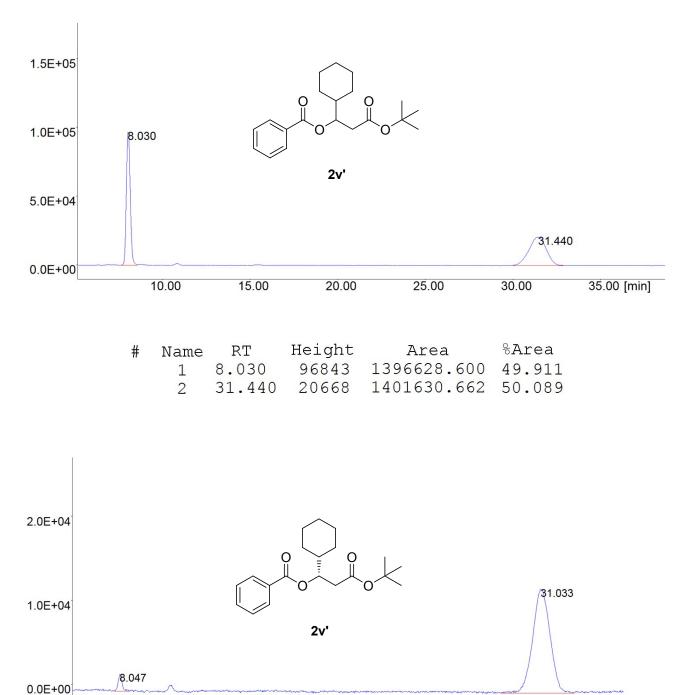


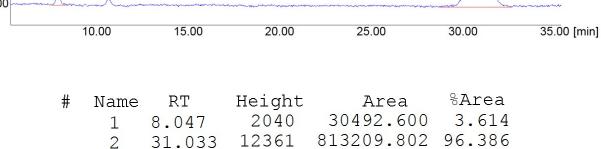
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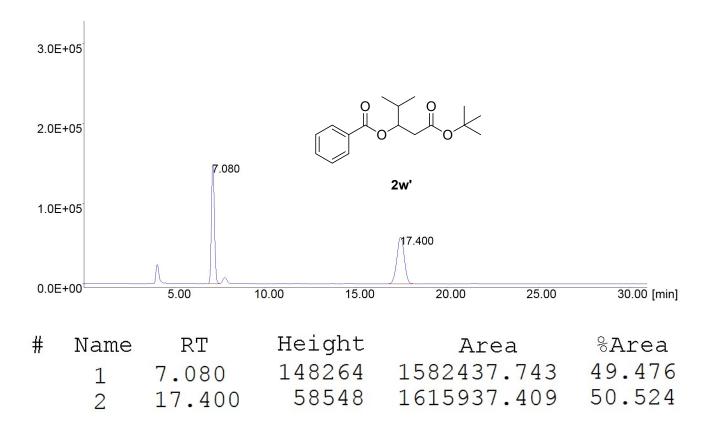


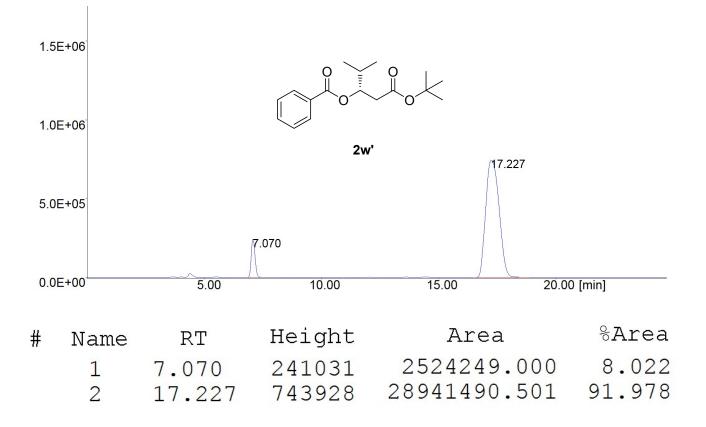
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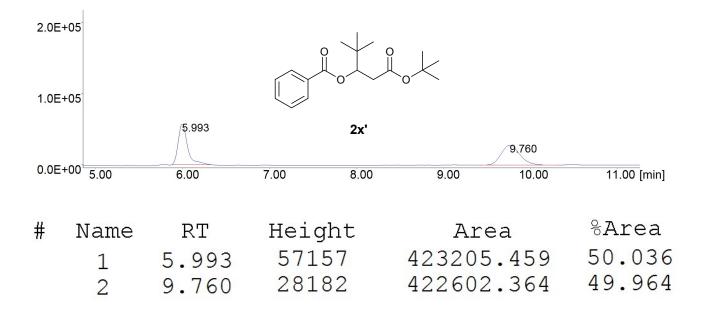


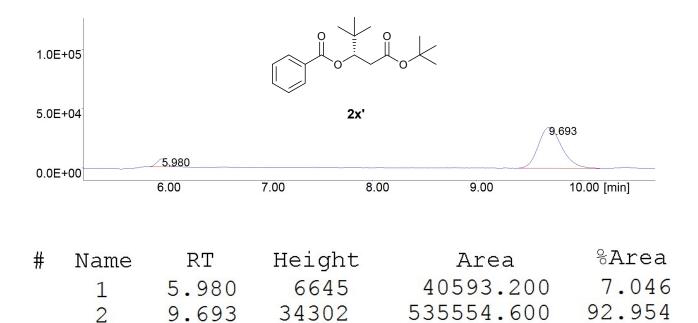












## 3. The comparative kinetic study of the ligands

