

Supporting Information

**Ni-Catalyzed Asymmetric Reductive Allylation of  
Aldehydes with Allylic Carbonates**

Zhuozhen Tan, Xiaolong Wan, Zhenhua Zang, Qun Qian,\* Wei Deng,\* and  
Hegui Gong\*

*Department of Chemistry, Innovative Drug Research Center and School of  
Materials Science and Engineering, Shanghai University, 99 Shang-Da Road,  
Shanghai 200444, China*

[Hegui\\_gong@shu.edu.cn](mailto:Hegui_gong@shu.edu.cn)

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**General Methods**

All reactions were carried out under an atmosphere of nitrogen unless otherwise indicated. Anhydrous THF was distilled from sodium/benzophenone ketyl prior to use. All other solvents were technical grade unless noted. Anhydrous DMF (Acros) and other commercially available reagents were used without further purification unless noted. Column chromatography was performed using silica gel 300-400 mesh (purchased from Qingdao-Haiyang Co. China) as the solid support. All NMR spectra were recorded on Bruker Avance 500 MHz spectrometer at STP. <sup>1</sup>H NMR and <sup>13</sup>C NMR chemical shifts are reported in δ units, parts per million (ppm) relative to the chemical shift of residual solvent. Reference peaks for chloroform in <sup>1</sup>H NMR and <sup>13</sup>C NMR spectra were set at 7.26 ppm and 77.0 ppm, respectively. High-resolution mass spectra (HRMS) were obtained using Waters Micromass GCT Premier. Melting point was recorded on a micro melting point apparatus (X-4, YUHUA Co., Ltd, Gongyi, China). Enantiomer ratios were determined by chiral HPLC analysis in comparison with authentic racemic materials. Optical rotations were taken on Autopol IV-T.

## (1) Typical Nickel-Catalyzed Carbonyl Allylation Procedure

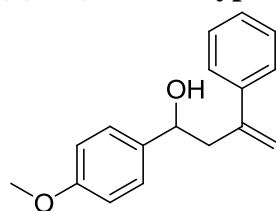
*General Procedure A:* To a flame-dried Schlenk tube equipped with a magnetic stir bar was loaded (*S*)-*t*Bu-Pybox (**2c**) (7.40 mg, 0.023 mmol, 15.0 mol%) and zinc powder (29.3 mg, 0.450 mmol, 300 mol%). The tube was moved into a dry glove box, at which point Ni(ClO<sub>4</sub>)<sub>2</sub> · 6H<sub>2</sub>O (5.50 mg, 0.015 mmol, 10.0 mol%) was added. The tube was capped with a rubber septum before being moved out of the glove box. DMF (1.0 mL), aldehyde (0.150 mmol, 100 mol%) and allylic carbonate (0.225 mmol, 150 mol%) were added via syringe subsequently. The solution was stirred at room temperature overnight. The reaction mixture was directly purified by flash-column chromatography on silica gel with petroleum ether and EtOAc as eluents to give pure products. The enantioselectivity was determined by HPLC. The ratio of anti/syn were determined by <sup>1</sup>H NMR.

*General Procedure B:* To a flame-dried Schlenk tube equipped with a magnetic stir bar was loaded (*S*)-*t*Bu-Pybox (**2c**) (7.40 mg, 0.023 mmol, 15.0 mol%) and zinc powder (29.3 mg, 0.450 mmol, 300 mol%). The tube was moved into a dry glove box, at which point Ni(ClO<sub>4</sub>)<sub>2</sub> · 6H<sub>2</sub>O (5.50 mg, 0.015 mmol, 10.0 mol%) and CsI (19.5 mg, 0.075 mmol, 50.0 mol%) was added. The tube was capped with a rubber septum before being moved out of the glove box. DMF (1.0 mL), aldehyde (0.150 mmol, 100 mol%) and allylic carbonate (0.225 mmol, 150 mol%) were added via syringe subsequently. The reaction mixture was allowed to stir 12 hours under N<sub>2</sub> atmosphere at -15 °C/-25 °C.

*Procedure for preparing compound 5:* To a flame-dried Schlenk tube equipped with a magnetic stir bar was loaded (*S*)-*t*Bu-Pybox (**2c**) (7.40 mg, 0.023 mmol, 15.0 mol%) and zinc powder (29.3 mg, 0.450 mmol, 300 mol%). The tube was moved into a dry glove box, at which point Ni(COD)<sub>2</sub> (4.20 mg, 0.015 mmol, 10.0 mol%) and CuI (14.3mg, 0.075 mmol, 50.0 mol%) was added. The tube was capped with a rubber septum, and it was moved out of the glove box. DMF (1.0 mL), 4-Anisaldehyde (22.4 mg, 0.150 mmol, 100 mol%) and allyl methyl carbonate (26.1 mg, 0.225 mmol, 150 mol%) were added via syringe subsequently. After the reaction mixture was allowed to stir 12 hours under N<sub>2</sub> atmosphere at 0 °C, it was directly loaded onto a silica column without work-up. Flash column chromatography (SiO<sub>2</sub>: 8% ethyl acetate in petroleum ether) gave **5** as a colorless oil.

## (2) Analytical Data for Homoallylic Alcohols

### (+)-1-(4-Methoxyphenyl)-3-phenylbut-3-en-1-ol (**1**)



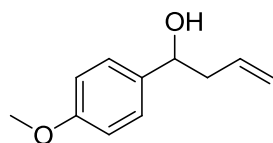
Colorless oil, 95% yield (36.2 mg), 91% ee; [ $\alpha$ ]<sub>D</sub><sup>22</sup> = +34.3 (*c* = 1.00, CHCl<sub>3</sub>).

<sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>):  $\delta$  2.06 (s, 1H), 2.90 (dd, *J* = 8.9, 14.3 Hz, 1H), 3.00 (ddd, *J* = 0.9, 4.5, 14.3 Hz, 1H), 3.83 (s, 3H), 4.72 (m, 1H), 5.18 (d, *J* = 0.8 Hz, 1H), 5.43 (d, *J* = 1.3 Hz, 1H), 6.93-6.88 (m, 2H), 7.27-7.50 (m, 7H).

<sup>13</sup>C NMR (125 MHz, CDCl<sub>3</sub>):  $\delta$  45.9, 55.3, 71.7, 113.8, 115.7, 126.3, 127.1, 127.8, 128.5, 136.1,

145.2, 159.1. HRMS (EI) calcd for C<sub>17</sub>H<sub>18</sub>O<sub>2</sub>: 254.1307, found 254.1304. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 85/15, flow rate 0.8 mL/min, UV 214 nm, *t<sub>R</sub>* = 9.7 min (minor), *t<sub>R</sub>* = 11.5 min (major).

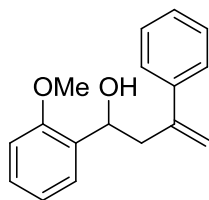
**(S)-(-)-1-(4-Methoxyphenyl)but-3-en-1-ol (5)<sup>1</sup>**



Colorless oil, 90% yield (24.0 mg), 66% ee;  $[\alpha]_D^{23} = -38.5$  (*c* = 0.40, CHCl<sub>3</sub>).

<sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>): δ 2.10 (s, 1H), 2.49 (t, *J* = 6.8 Hz, 2H), 3.80 (s, 3H), 4.67 (t, *J* = 6.5 Hz, 1H 2H), 5.10-5.18 (m, 2H), 5.82-5.73 (m, 1H), 6.86-6.90 (m, 2H), 7.326-7.29 (m, 2H). HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 92/8, flow rate 0.6 mL/min, UV 220 nm, *t<sub>R</sub>* = 15.4 min (minor), *t<sub>R</sub>* = 16.8 min (major).

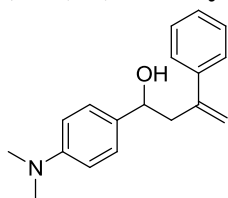
**(+)-1-(2-Methoxyphenyl)-3-phenylbut-3-en-1-ol (6)**



Colorless oil, 96% yield (36.6 mg), 88% ee;  $[\alpha]_D^{26} = +24.8$  (*c* = 1.00, CHCl<sub>3</sub>). <sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>): δ 2.48 (d, *J* = 4.8 Hz, 1H), 2.77 (dd, *J* = 9.0, 14.3 Hz, 1H), 3.17 (ddd, *J* = 0.9, 4.5, 14.3 Hz, 1H), 3.86 (s, 3H), 5.00 (m, 1H), 5.18 (s, 1H), 5.44 (d, *J* = 1.2 Hz, 1H), 6.86-6.90 (m, 1H), 6.94-6.99 (m, 1H), 7.22-7.27 (m, 1H), 7.28-7.32 (m, 1H), 7.34-7.39 (m, 3H), 7.51-7.55 (m, 2H). <sup>13</sup>C

NMR (125 MHz, CDCl<sub>3</sub>): δ 43.9, 55.2, 68.5, 110.3, 115.1, 120.7, 126.3, 126.6, 127.6, 128.2, 128.3, 132.0, 140.3, 145.3, 156.2. HRMS (EI) calcd for C<sub>17</sub>H<sub>18</sub>O<sub>2</sub>: 254.1307, found 254.1305. HPLC condition: Chiralcel OD-H, *n*-hexane/2-propanol 80/20, flow rate 0.7 mL/min, UV 214 nm, *t<sub>R</sub>* = 7.4 min (major), *t<sub>R</sub>* = 7.9 min (minor).

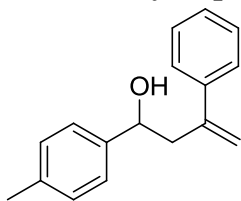
**(+)-1-(4-(Dimethylamino)phenyl)-3-phenylbut-3-en-1-ol (7)**



Colorless oil, 95% yield (38.1 mg), 85% ee;  $[\alpha]_D^{22} = +41.6$  (*c* = 1.03, CHCl<sub>3</sub>). <sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>): δ 1.99 (s, 1H), 2.91 (dd, *J* = 8.9, 14.3 Hz, 1H), 2.96 (s, 6H), 2.98 (dd, *J* = 4.5, 14.3 Hz, 1H), 4.67 (m, 1H), 5.18 (s, 1H), 5.41 (d, *J* = 1.2 Hz, 1H), 6.71-6.76 (m, 2H), 7.21-7.26 (m, 2H), 7.29-7.34 (m, 1H), 7.35-7.40 (m, 2H), 7.44-7.49 (m, 2H). <sup>13</sup>C NMR (125 MHz, CDCl<sub>3</sub>): δ 40.7, 45.6, 71.9, 112.6, 115.4, 126.4, 126.9, 127.7, 128.5, 131.9, 140.6, 145.4, 150.3. HRMS

calcd for C<sub>18</sub>H<sub>21</sub>NO (ED): 267.1623, found 267.1622. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 85/15, flow rate 0.8 mL/min, UV 214 nm, *t<sub>R</sub>* = 13.2 min (minor), *t<sub>R</sub>* = 16.8 min (major).

**(+)-3-Phenyl-1-*p*-tolylbut-3-en-1-ol (8)**



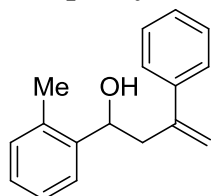
Colorless oil, 95% yield (33.9 mg), 88% ee;  $[\alpha]_D^{24} = +33.4$  (*c* = 0.98, CHCl<sub>3</sub>).

<sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>): δ 2.07 (s, 1H), 2.39 (s, 3H), 2.90 (dd, *J* = 9.0, 14.3 Hz, 1H), 3.02 (dd, *J* = 3.9, 14.3 Hz, 1H), 4.74

<sup>1</sup> V. De Sio, A. Massa, A. Scettri, *Org. Biomol. Chem.* **2010**, *8*, 3055-3059.

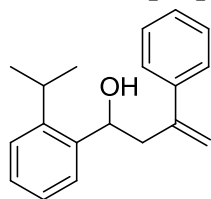
(m, 1H), 5.20 (s, 1H), 5.44 (s, 1H), 7.16-7.22 (m, 2H), 7.26-7.30 (m, 2H), 7.32-7.38 (m, 1H), 7.38-7.42 (m, 2H), 7.44-7.51 (m, 2H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  21.1, 45.9, 71.9, 115.7, 125.8, 126.3, 127.8, 128.5, 129.1, 137.2, 140.4, 141.0, 145.2. HRMS (EI) calcd for  $\text{C}_{17}\text{H}_{18}\text{O}$ : 238.1358, found 238.1356. HPLC condition: Chiralcel AD-H, *n*-hexane/ 2-propanol 92/8, flow rate 0.6 mL/min, UV 214 nm,  $t_{\text{R}}$  = 16.3 min (minor),  $t_{\text{R}}$  = 20.0 min (major).

**(-)-3-phenyl-1-*o*-tolylbut-3-en-1-ol (9)**



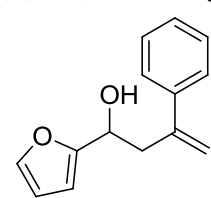
Colorless oil, 85% yield (30.3 mg), 86% ee;  $[\alpha]_{\text{D}}^{13} = -4.4$  ( $c = 0.98$ ,  $\text{CHCl}_3$ ).  
 $^1\text{H}$  NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.01 (s, 1H), 2.25(s, 3H), 2.77 (dd,  $J = 9.5, 14.4$  Hz, 1H), 3.02 (ddd,  $J = 0.9, 3.3, 14.5$  Hz, 1H), 4.95 (m, 1H), 5.24 (s, 1H), 5.45 (d,  $J = 1.2$  Hz, 1H), 7.11-7.14 (m, 1H), 7.15-7.20 (m, 1H), 7.22-7.27 (m, 1H), 7.30-7.34 (m, 1H), 7.35-7.40 (m, 2H), 7.45-7.49 (m, 2H), 7.54-7.57 (m, 1H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  19.0, 44.7, 68.4, 115.7, 125.3, 126.3, 127.2, 127.8, 128.4, 130.3, 134.3, 140.3, 141.9, 145.4. HRMS (EI) calcd for  $\text{C}_{17}\text{H}_{18}\text{O}$ : 238.1358, found 238.1362. HPLC condition: Chiralpak ID-3, *n*-hexane/ 2-propanol 90/10, flow rate 0.7 mL/min, UV 214 nm,  $t_{\text{R}}$  = 5.0 min (minor),  $t_{\text{R}}$  = 5.7 min (major).

**(+)-1-(2-isopropylphenyl)-3-phenylbut-3-en-1-ol (10)**



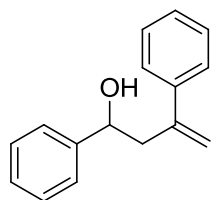
Colorless oil, 48% yield (19.1 mg), 85% ee;  $[\alpha]_{\text{D}}^{15} = +2.8$  ( $c = 0.98$ ,  $\text{CHCl}_3$ ).  
 $^1\text{H}$  NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.18 (d,  $J = 6.7$  Hz, 3H), 1.25 (d,  $J = 6.7$  Hz, 3H), 2.04 (s, 1H), 2.85 (dd,  $J = 9.5, 14.5$  Hz, 1H), 3.02 (ddd,  $J = 1.1, 3.4, 14.5$  Hz, 1H), 3.10 (m, 1H), 5.09 (dd,  $J = 2.7, 9.54$  Hz, 1H), 5.26 (d,  $J = 0.5$  Hz, 1H), 5.45 (d,  $J = 1.3$ , 1H), 7.23-7.36 (m, 4H), 7.36-7.42 (m, 2H), 7.46-7.49 (m, 2H), 7.55-7.59 (m, 1H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  23.8, 24.4, 27.9, 45.6, 67.8, 115.6, 125.2, 125.5, 126.0, 126.3, 127.6, 127.8, 128.4, 140.4, 145.4, 145.8. HRMS (EI) calcd for  $\text{C}_{19}\text{H}_{22}\text{O}$ : 238.1671, found 238.1669. HPLC condition: Lux Amylose-2, *n*-hexane/ 2-propanol 90/10, flow rate 0.7 mL/min, UV 214 nm,  $t_{\text{R}}$  = 7.0 min (major),  $t_{\text{R}}$  = 7.5 min (minor).

**(+)-1-(Furan-2-yl)-3-phenylbut-3-en-1-ol (11)**



Colorless oil, 95% yield (30.5 mg), 80% ee;  $[\alpha]_{\text{D}}^{26} = +26.2$  ( $c = 1.03$ ,  $\text{CHCl}_3$ ).  
 $^1\text{H}$  NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.09 (s, 1H), 3.00 (dd,  $J = 8.8, 14.3$  Hz, 1H), 3.15 (dd,  $J = 4.9, 14.3$  Hz, 1H), 4.77 (m, 1H), 5.20 (s, 1H), 5.41 (d,  $J = 1.0$  Hz, 1H), 6.22 (d,  $J = 3.0$  Hz, 1H), 6.32 (dd,  $J = 1.8, 3.0$  Hz, 1H), 7.27-7.32 (m, 1H), 7.32-7.39 (m, 3H), 7.40-7.45 (m, 2H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  41.9, 65.9, 106.3, 110.1, 115.9, 126.3, 127.8, 128.5, 140.2, 142.0, 144.4, 155.8. HRMS (EI) calcd for  $\text{C}_{14}\text{H}_{14}\text{O}_2$ : 214.0994, found 214.0995. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_{\text{R}}$  = 15.2 min (minor),  $t_{\text{R}}$  = 17.0 min (major).

**(S)-(+)-1,3-Diphenylbut-3-en-1-ol (12)<sup>2</sup>**

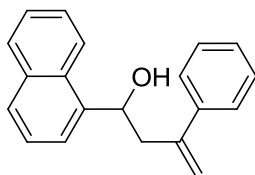


Colorless oil, 95% yield (32.0 mg), 87% ee;  $[\alpha]_D^{23} = +28.5$  ( $c = 1.09$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.09 (s, 1H), 2.88 (dd,  $J = 9.2, 14.3$  Hz, 1H), 3.02 (dd,  $J = 3.6, 14.3$  Hz, 1H), 4.73 (m, 1H), 5.18 (s, 1H), 5.43 (s, 1H), 7.29-7.47 (m, 10H). HPLC condition: Chiralcel

AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_R = 12.9$  min (minor),  $t_R = 14.4$  min (major). (lit:  $t_R = 34.1$  min (*R*)-isomer,  $t_R = 38.2$  min (*S*)-isomer (Daicel Chiralpak AD-H, flow rate 0.5 mL/min, *n*-hexane/2-propanol 30/1)).

**(-)-1-(naphthalen-2-yl)-3-phenylbut-3-en-1-ol (13)**

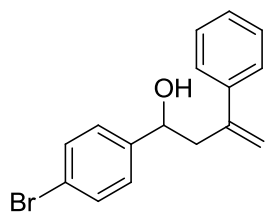


Colorless oil, 95% yield (39.0 mg), 73% ee;  $[\alpha]_D^{26} = -0.9$  ( $c = 1.05$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.21 (d,  $J = 1.9$  Hz, 1H), 2.91 (dd,  $J = 9.6, 14.6$  Hz, 1H), 3.26 (ddd,  $J = 1.0, 3.2, 14.6$  Hz, 1H), 5.28 (d,  $J = 0.5$  Hz, 1H), 5.45-5.51 (m, 2H), 7.23-7.37 (m,

1H), 7.38-7.44 (m, 2H), 7.46-7.53 (m, 5H), 7.71 (d,  $J = 7.2$  Hz, 1H), 7.78 (d,  $J = 8.2$  Hz, 1H), 7.85-7.92 (m, 1H), 7.95-8.00 (m, 1H).  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  45.2, 68.9, 115.8, 122.9, 123.0, 125.4, 125.5, 126.5, 127.8, 127.9, 128.5, 128.9, 130.3, 133.8, 139.6, 140.5, 145.7. HRMS (EI) calcd for  $\text{C}_{20}\text{H}_{18}\text{O}$ : 274.1358, found 274.1362. HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 70/30, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 6.8$  min (major),  $t_R = 7.4$  min (minor).

**(+)-1-(4-Bromophenyl)-3-phenylbut-3-en-1-ol (14)**



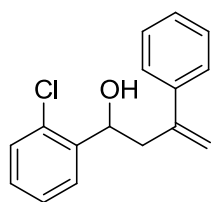
This compound was prepared according to the *General Procedure A* at 25 °C. Colorless oil, 95% yield (43.2 mg), 77% ee;  $[\alpha]_D^{27} = +36.8$  ( $c = 1.05$ ,  $\text{CHCl}_3$ ).

This compound was also prepared according to the *General Procedure B* at -25 °C. 77% yield (34.8 mg), 85% ee;  $[\alpha]_D^{25} = +42.0$  ( $c = 1.00$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.11 (s, 1H), 2.82 (ddd,  $J = 0.6, 8.9, 14.2$  Hz, 1H), 2.96 (ddd,  $J = 1.1, 4.4, 14.2$  Hz, 1H), 4.68 (q,  $J = 4.3$  Hz, 1H), 5.14 (d,  $J = 1.0$  Hz, 1H), 5.41 (d,  $J = 1.3$  Hz), 7.19-7.23 (m, 2H), 7.30-7.34 (m, 1H), 7.35-7.39 (m, 2H), 7.40-7.48 (m, 4H).  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  46.0, 71.4, 116.1, 121.2, 126.2, 127.5, 127.9, 128.6, 131.4, 140.0, 142.8, 144.6. HRMS (EI) calcd for  $\text{C}_{16}\text{H}_{15}\text{OBr}$ : 302.0306, found 302.0307. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_R = 15.2$  min (minor),  $t_R = 17.1$  min (major).

<sup>2</sup> S. Kotani, S. Hashimoto, M. Nakajima, *Tetrahedron*, **2007**, *63*, 3122–3132.

**(-)-1-(2-Chlorophenyl)-3-phenylbut-3-en-1-ol (15)**

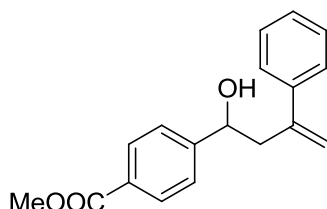


Colorless solid, M.p: 56-58 °C, 94% yield (36.5 mg), 77% ee;  $[\alpha]_D^{27} = -12.6$  ( $c = 1.05$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.28 (s, 1H), 2.84 (ddd,  $J = 0.4, 9.8, 14.3$  Hz, 1H), 3.00 (ddd,  $J = 1.2, 3.1, 14.3$  Hz, 1H), 5.14 (m, 1H), 5.26 (s, 1H), 5.51 (s, 1H), 7.18-7.24 (td,  $J = 1.7, 7.7$  Hz, 1H),

7.27-7.35 (m, 3H), 7.35-7.40 (m, 2H), 7.52-7.56 (m, 2H), 7.62-7.62 (dd,  $J = 1.5, 7.8$  Hz, 1H).  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  44.01, 68.6, 116.0, 126.3, 127.0, 127.1, 127.9, 128.3, 128.4, 129.3, 131.4, 139.8, 141.3, 144.9. HRMS (EI) calcd for  $\text{C}_{16}\text{H}_{15}\text{OCl}$ : 258.0811, found 258.0804. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 11.0$  min (minor),  $t_R = 11.7$  min (major).

**(+)-Methyl-4-(1-hydroxy-3-phenylbut-3-enyl)benzoate (16)**

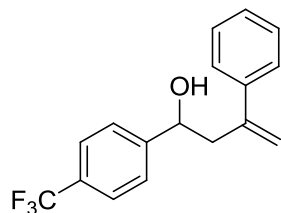


Colorless oil, 95% yield (40.2 mg), 71% ee;  $[\alpha]_D^{27} = +40.6$  ( $c = 1.02$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.28 (s, 1H), 2.84 (dd,  $J = 9.0, 14.3$  Hz, 1H), 3.00 (dd,  $J = 4.4, 14.3$  Hz, 1H), 3.91 (s, 3H), 4.77 (m, 1H), 5.14 (s, 1H), 5.41 (s, 1H), 7.29-7.34 (m, 1H), 7.34-7.41 (m, 4H), 7.42-7.46 (m, 2H), 7.97-8.02

(m, 2H).  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  45.9, 52.0, 71.6, 116.1, 125.7, 126.2, 127.9, 128.6, 129.2, 129.7, 140.0, 144.5, 149.0, 166.9. HRMS (EI) calcd for  $\text{C}_{18}\text{H}_{18}\text{O}_3$ : 282.1256, found 282.1255. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_R = 20.3$  min (minor),  $t_R = 23.4$  min (major).

**(+)-3-Phenyl-1-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (17)**

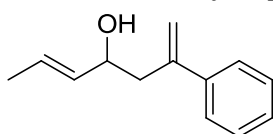


Colorless oil, 95% yield (41.6 mg), 69% ee;  $[\alpha]_D^{23} = +16.6$  ( $c = 1.05$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.28 (s, 1H), 2.84 (dd,  $J = 9.1, 14.3$  Hz, 1H), 3.00 (ddd,  $J = 0.9, 4.5, 14.3$  Hz, 1H), 4.77 (m, 1H), 5.16 (d,  $J = 0.7$  Hz, 1H), 5.43 (d,  $J = 1.1$  Hz, 1H), 7.31-7.36 (m, 1H), 7.36-7.41 (m, 2H), 7.42-7.45 (m, 4H),

7.57-7.62 (d,  $J = 8.1$  Hz, 2H).  $^{13}\text{C NMR}$  (125MHz,  $\text{CDCl}_3$ ):  $\delta$  46.0, 71.4, 116.2, 124.1 (q,  $J = 272.4$  Hz,  $\text{CF}_3$ ), 125.3 (q,  $J = 3.8$  Hz), 126.1, 126.2, 127.9, 128.6, 129.7 (q,  $J = 32.2$  Hz), 140.0, 144.5, 147.8. HRMS (EI) calcd for  $\text{C}_{17}\text{H}_{15}\text{F}_3\text{O}$ : 292.1075, found 292.1080. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_R = 11.1$  min (minor),  $t_R = 12.1$  min (major).

**(+)-(E)-2-Phenylhepta-1,5-dien-4-ol (18)**

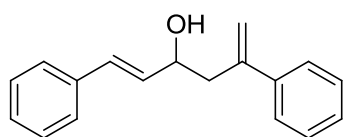


Colorless oil, 91% yield (26.5 mg), 85% ee;  $[\alpha]_D^{25} = +90.3$  ( $c = +0.40$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.67-1.68 (m, 4 H), 2.66 (dd,  $J = 4.6, 14.1$  Hz, 1H), 2.92 (dd,  $J = 5.0, 14.1$  Hz, 1H), 4.12 (m,

1H), 5.17 (s, 1H), 5.40 (d,  $J = 1.1$  Hz, 1H), 5.50 (ddd,  $J = 1.3, 6.8, 15.6$  Hz, 1H), 6.55 (dq,  $J = 6.4, 15.6$  Hz, 1H), 7.24-7.31 (m, 1H), 7.31-7.37 (m, 2H), 7.38-7.44 (m, 2H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  17.6, 43.8, 70.6, 115.4, 126.3, 127.0, 127.7, 128.4, 133.2, 140.6, 145.0. HRMS (EI) calcd for  $\text{C}_{13}\text{H}_{16}\text{O}$ : 188.1201, found 188.1201. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_{\text{R}} = 10.3$  min (minor),  $t_{\text{R}} = 11.1$  min (major).

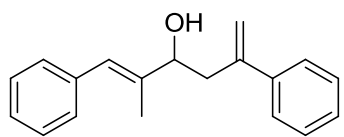
**(+)-(E)-1,5-Diphenylhexa-1,5-dien-3-ol (19)<sup>3</sup>**



Colorless oil, 94% yield (35.3 mg), 85% ee;  $[\alpha]_{\text{D}}^{25} = +24.6$  ( $c = 1.00$ ,  $\text{CHCl}_3$ ).

$^1\text{H}$  NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.90 (s, 1H), 2.81 (dd,  $J = 8.3, 14.1$  Hz, 1H), 2.92 (dd,  $J = 5.0, 14.1$  Hz, 1H), 4.37 (m, 1H), 5.23 (s, 1H), 5.44 (d,  $J = 1.2$  Hz, 1H), 6.22 (dd,  $J = 6.4, 15.9$  Hz, 1H), 6.55 (d,  $J = 15.9$  Hz, 1H), 7.22-7.27 (m, 1H), 7.28-7.40 (m, 7H), 7.43-7.48 (m, 2H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  43.9, 70.6, 115.8, 126.3, 126.4, 127.6, 127.7, 128.4, 128.5, 130.2, 131.5, 136.7, 140.5, 144.7. HRMS (EI) calcd for  $\text{C}_{18}\text{H}_{18}\text{O}$ : 250.1358, found 250.1357. HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_{\text{R}} = 19.5$  min (minor),  $t_{\text{R}} = 22.4$  min (major).

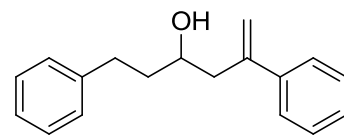
**(+)-(E)-2-methyl-1,5-diphenylhexa-1,5-dien-3-ol (20)**



Colorless oil, 87% yield (35.3 mg), 91% ee;  $[\alpha]_{\text{D}}^{16} = +78.5$  ( $c = 1.01$ ,  $\text{CHCl}_3$ ).

$^1\text{H}$  NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.87 (d,  $J = 2.5$  Hz, 1H), 1.90 (d,  $J = 1.5$  Hz, 3H), 2.79 (ddd,  $J = 1.0, 8.3, 14.2$  Hz, 1H), 2.96 (ddd,  $J = 1.0, 4.8, 14.2$  Hz, 1H), 4.25 (m, 1H), 5.23 (d,  $J = 0.9$  Hz, 1H), 5.42 (d,  $J = 1.4$  Hz, 1H), 6.43 (s, 1H), 7.20-7.23 (m, 3H), 7.23-7.32 (m, 5H), 7.32-7.35 (m, 2H).  $^{13}\text{C}$  NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  13.4, 42.0, 75.6, 115.6, 126.0, 126.4, 126.5, 127.7, 128.1, 128.5, 129.0, 137.5, 139.3, 140.6, 145.3. HRMS (EI) calcd for  $\text{C}_{19}\text{H}_{20}\text{O}$ : 264.1508, found 264.1510. HPLC condition: Chiralpak ID-3, *n*-hexane/2-propanol 95/5, flow rate 0.7 mL/min, UV 214 nm,  $t_{\text{R}} = 6.9$  min (major),  $t_{\text{R}} = 7.9$  min (minor).

**(+)-1,5-Diphenylhex-5-en-3-ol (21)<sup>4</sup>**



This compound was prepared according to the *General Procedure A* at 25 °C. Colorless oil, 95% yield (35.9 mg), 68% ee;  $[\alpha]_{\text{D}}^{27} = +41.4$  ( $c = 0.98$ ,  $\text{CHCl}_3$ ).

This compound was also prepared according to the *General Procedure B* at -25 °C. 52% yield (19.5 mg), 80% ee;  $[\alpha]_{\text{D}}^{25} = +48.0$  ( $c = 1.00$ ,  $\text{CHCl}_3$ ).

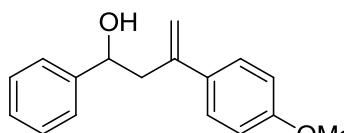
$^1\text{H}$  NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.76 (s, 1H), 1.80-1.88 (m, 2H), 2.60 (dd,  $J = 8.8, 14.2$

<sup>3</sup> T. Kamei, K. Fujita, K. Itami, J.-I. Yoshida, *Org. Lett.* **2005**, 7, 4725-4728.

<sup>4</sup> S.-F. Zhu, X.-C. Qiao, Y.-Z. Zhang, L.-X. Wang, Q.-L. Zhou, *Chem. Sci.* **2011**, 2, 1135-1140.

Hz, 1H), 2.79-2.87 (m, 2H), 3.73 (m, 1H), 5.19 (s, 1H), 5.43 (d,  $J = 1.2$  Hz, 1H), 7.16-7.23 (m, 3H), 7.27-7.33 (m, 3H), 7.33-7.38 (m, 2H), 7.39-7.44 (m, 2H). HPLC condition: Chiralcel AD-H, *n*-hexane/2-propanol 90/10, flow rate 0.6 mL/min, UV 214 nm,  $t_R = 12.0$  min (minor),  $t_R = 13.6$  min (major).

**(+)-3-(4-Methoxyphenyl)-1-phenylbut-3-en-1-ol (22)**

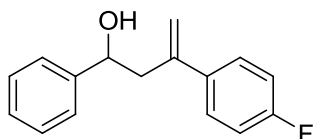


This compound was prepared according to the *General Procedure A* at 25 °C. Colorless oil, 95% yield (36.1 mg), 72% ee;  $[\alpha]_D^{27} = +37.3$  ( $c = 1.00$ , CHCl<sub>3</sub>).

This compound was also prepared according to the *General Procedure B* at -15 °C. 60% yield (22.8 mg), 81% ee;  $[\alpha]_D^{25} = +41.0$  ( $c = 1.00$ , CHCl<sub>3</sub>).

<sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>):  $\delta$  2.18 (d,  $J = 2.2$  Hz, 1H), 2.81 (dd,  $J = 9.2, 14.3$  Hz, 1H), 2.99 (ddd,  $J = 0.9, 4.2, 14.3$  Hz, 1H), 3.83 (s, 3H), 4.72 (dq,  $J = 2.2, 9.2$  Hz, 1H), 5.09 (d,  $J = 0.7$  Hz, 1H), 5.36 (d,  $J = 1.3$  Hz, 1H), 6.88-6.93 (m, 2H), 7.27-7.32 (m, 1H), 7.33-7.38 (m, 4H), 7.39-7.43 (m, 2H). <sup>13</sup>C NMR (125 MHz, CDCl<sub>3</sub>):  $\delta$  46.0, 55.2, 71.9, 113.8, 114.1, 125.7, 127.3, 127.4, 128.3, 132.5, 143.9, 144.2, 159.3. HRMS (EI) calcd for C<sub>17</sub>H<sub>18</sub>O<sub>2</sub>: 254.1307, found 254.2310. HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 90/10, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 14.4$  min (major),  $t_R = 15.4$  min (minor).

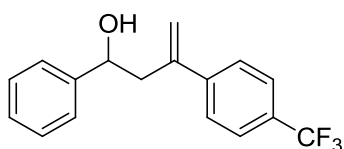
**(+)-3-(4-fluorophenyl)-1-phenylbut-3-en-1-ol (23)**



Colorless oil, 60% yield (21.8 mg), 81% ee;  $[\alpha]_D^{25} = +7.8$  ( $c = 1.10$ , CHCl<sub>3</sub>).

<sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>):  $\delta$  2.08 (s, 1H), 2.86 (dd,  $J = 8.9, 14.3$  Hz, 1H), 2.99 (dd,  $J = 4.3, 14.3$  Hz, 1H), 4.71 (m, 1H), 5.15 (s, 1H), 5.36 (s, 1H), 7.02-7.07 (m, 2H), 7.25-7.30 (m, 1H), 7.32-7.37 (m, 4H), 7.38-7.42 (m, 2H). <sup>13</sup>C NMR (125 MHz, CDCl<sub>3</sub>):  $\delta$  45.9, 72.1, 115.2, 115.3, 115.6, 125.7, 127.6, 127.8, 127.9, 128.7, 136.4 (d,  $J = 3.4$  Hz), 143.8 (d,  $J = 29.4$  Hz), 162.4 (d,  $J = 246.9$  Hz). HRMS (EI) calcd for C<sub>16</sub>H<sub>15</sub>FO: 242.1107, found 242.1104. HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 80/20, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 6.6$  min (minor),  $t_R = 7.0$  min (major).

**(+)-1-phenyl-3-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (24)**



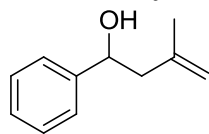
Colorless oil, 90% yield (39.4 mg), 82% ee;  $[\alpha]_D^{16} = +17.0$  ( $c = 1.00$ , CHCl<sub>3</sub>).

<sup>1</sup>H NMR (500 MHz, CDCl<sub>3</sub>):  $\delta$  2.14 (s, 1H), 2.93 (ddd,  $J = 0.8, 8.5, 14.6$  Hz, 1H), 2.98 (ddd,  $J = 1.1, 5.1, 14.6$  Hz, 1H), 4.71 (dd,  $J = 5.1, 8.3$  Hz, 1H), 5.25 (d,  $J = 1.0$ , Hz, 1H), 5.45 (d,  $J = 0.8$  Hz, 1H), 7.25-7.37 (m, 5H), 7.50-7.54 (m, 2H), 7.60-7.63 (m, 2H). <sup>13</sup>C NMR (125 MHz, CDCl<sub>3</sub>):  $\delta$  45.4, 72.3, 117.5, 124.1 (q,  $J = 272.1$  Hz), 125.3 (q,  $J = 3.8$  Hz), 125.8, 127.7, 128.4, 129.6 (q,  $J = 32.5$  Hz), 143.6, 143.9, 144.2. HRMS (EI) calcd for C<sub>17</sub>H<sub>15</sub>F<sub>3</sub>O: 292.1075, found 292.1073. HPLC condition:



Chiralpak ID-3, *n*-hexane/2-propanol 95/5, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 5.0$  min (minor),  $t_R = 5.3$  min (major).

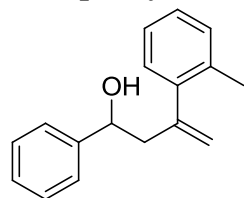
**(S)-3-Methyl-1-phenylbut-3-en-1-ol (25)<sup>5</sup>**



Colorless oil, 96% yield (23.3 mg), 60% ee;  $[\alpha]_D^{27} = -36.8$  ( $c = 1.00$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.81 (s, 3H), 2.14 (d,  $J = 1.8$  Hz, 1H), 2.43 (d,  $J = 7.1$  Hz, 2H), 4.83 (t,  $J = 6.5$  Hz, 1H), 4.86 (d,  $J = 0.9$  Hz, 1H), 4.93 (m, 1H), 7.25-7.31 (m, 1H), 7.32-7.42 (m, 4H).  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  22.4, 48.4, 71.4, 114.1, 125.7, 127.5, 128.4, 142.4, 144.1. HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 90/10, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 6.9$  min (major),  $t_R = 7.5$  min (minor).

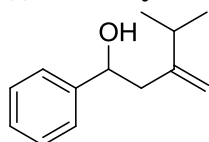
**(+)-1-phenyl-3-*o*-tolylbut-3-en-1-ol (26)**



Colorless oil, 53% yield (18.9 mg), 41% ee;  $[\alpha]_D^{13} = +5.4$  ( $c = 0.95$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.14 (d,  $J = 1.7$  Hz, 1H), 2.33 (s, 3H), 2.77 (dd,  $J = 9.5, 14.4$  Hz, 1H), 2.83 (ddd,  $J = 0.7, 4.4, 14.4$  Hz, 1H), 4.62 (m, 1H), 5.07 (s, 1H), 5.36 (d,  $J = 0.6$  Hz, 1H), 7.13-7.16 (m, 1H), 7.16-7.23 (m, 3H), 7.25-7.30 (m, 1H), 7.30-7.36 (m, 4H).  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  19.9, 47.9, 71.6, 117.5, 125.6, 125.8, 127.1, 127.5, 128.3, 128.4, 130.3, 134.9, 141.7, 143.8, 146.6. HRMS (EI) calcd for  $\text{C}_{17}\text{H}_{18}\text{O}$ : 238.1358, found 238.1357. HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 80/20, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 6.3$  min (minor),  $t_R = 6.8$  min (major).

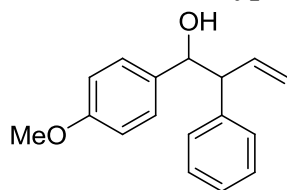
**(-)-4-methyl-3-methylene-1-phenylpentan-1-ol (27)**



Colorless oil, 45% yield (12.9 mg), 51% ee;  $[\alpha]_D^{14} = -35.4$  ( $c = 0.90$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  1.06 (d,  $J = 6.9$  Hz, 3H), 1.08 (d,  $J = 6.9$  Hz, 3H), 2.20 (d,  $J = 2.2$  Hz, 3H), 2.29 (m, 1H), 2.39 (ddd,  $J = 0.5, 9.7, 14.5$  Hz, 1H), 2.43 (ddd,  $J = 1.0, 3.8, 14.5$  Hz, 1H), 4.81 (m, 1H), 4.90 (d,  $J = 0.9$  Hz, 1H), 4.93 (m, 1H), 7.25-7.30 (m, 1H), 7.33-7.41 (m, 4H). HRMS (EI) calcd for  $\text{C}_{13}\text{H}_{18}\text{O}$ : 190.1358, found 90.1359.  $^{13}\text{C NMR}$  (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  21.5, 21.9, 45.4, 71.8, 110.2, 125.7, 127.4, 128.4, 144.1, 152.5. HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 80/20, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 5.3$  min (major),  $t_R = 5.6$  min (minor).

***anti*-1-(4-Methoxyphenyl)-2-phenylbut-3-en-1-ol (28)<sup>2</sup>**

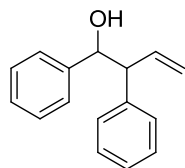


Colorless oil, 95% yield (36.2 mg), *anti:syn* = 99:1, 64% ee;  $[\alpha]_D^{26} = -3.2$  ( $c = 1.00$ ,  $\text{CHCl}_3$ ).

$^1\text{H NMR}$  (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.30 (d,  $J = 2.1$  Hz, 1H), 3.53 (t,  $J = 8.4$  Hz, 1H), 3.75 (s, 3H), 4.80 (dd,  $J = 1.2, 8.3$  Hz, 1H), 5.21-5.29 (m, 2H), 6.25 (ddd,  $J = 7.9, 10.2, 18.1$  Hz, 1H), 6.72-6.76 (m, 2H), 7.03-7.08 (m, 4H), 7.12-7.16 (m, 1H), 7.18-7.23 (m, 2H).

HPLC condition: Lux Amylose-2, *n*-hexane/2-propanol 90/10, flow rate 0.7 mL/min, UV 214 nm,  $t_R = 14.1$  min (major),  $t_R = 16.0$  min (minor).

**(1*S*,2*R*)-(-)-1,2-Diphenylbut-3-en-1-ol (29)<sup>5</sup>**

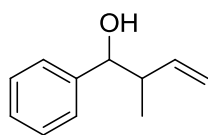


Colorless oil, 95% yield (32.0 mg), *anti:syn* = 97:3, 50% ee;  $[\alpha]_D^{26} = -6.7$  ( $c = 1.10$ ,  $\text{CHCl}_3$ ).

This compound was prepared according to the *General Procedure A* at  $-25$  °C. 67% yield (21.9 mg), *anti:syn* = 97:3, 64% ee;

<sup>1</sup>H NMR (500 MHz,  $\text{CDCl}_3$ ):  $\delta$  2.34 (d,  $J = 1.3$ , 1H), 3.56 (t,  $J = 8.3$  Hz, 1H), 4.85 (d,  $J = 7.7$  Hz, 1H), 5.21-5.29 (m, 2H), 6.26 (ddd,  $J = 8.7, 10.2, 16.8$  Hz, 1H), 7.04-7.09 (m, 2H), 7.13-7.24 (m, 8H). <sup>13</sup>C NMR (125 MHz,  $\text{CDCl}_3$ ):  $\delta$  59.1, 77.2, 118.4, 126.5, 126.6, 127.4, 127.9, 128.2, 128.3, 137.8, 140.6, 141.8. HPLC chiral condition: Chiralcel AD-H, *n*-hexane/2-propanol 98/2, flow rate 0.9 mL/min, UV 214 nm,  $t_R = 24.0$  min (major),  $t_R = 26.2$  min (minor).

**2-Methyl-1-phenylbut-3-en-1-ol (30)<sup>5</sup>**



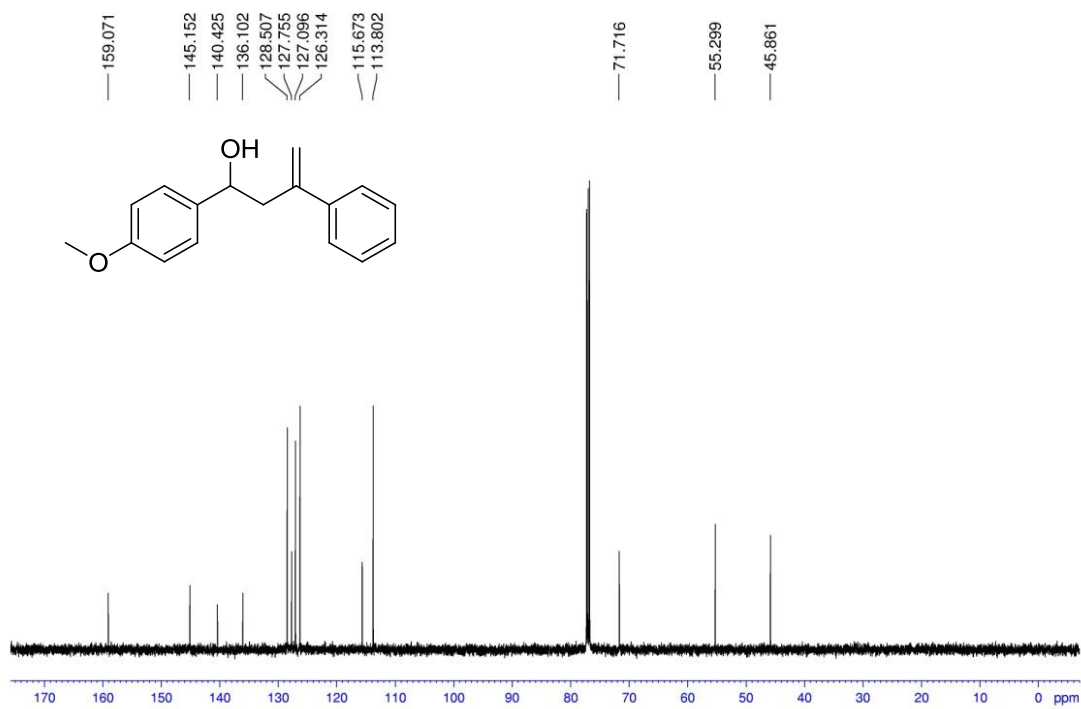
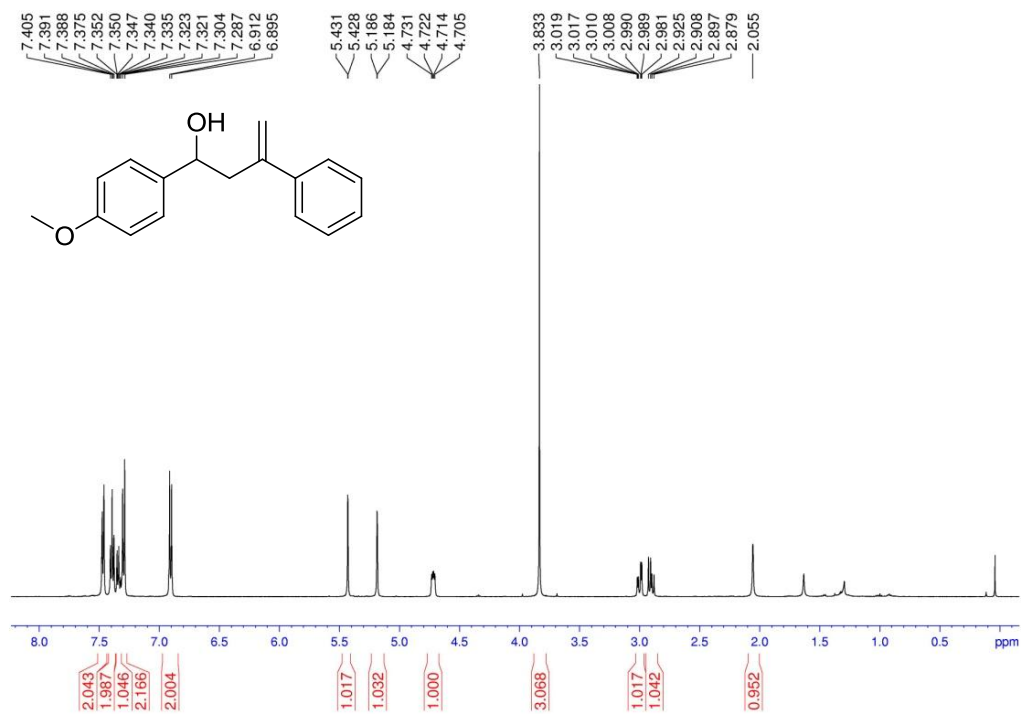
Colorless oil obtained from table 3, entry 4: 95% yield (23.1 mg), *anti:syn* = 59:41; 60% ee (for *anti*), 60% ee (for *syn*);  $[\alpha]_D^{27} = -49.1$  ( $c = 1.00$ ,  $\text{CHCl}_3$ ); colorless oil obtained from table 3, entry 5: 95% yield (23.3 mg), *anti:syn* = 59:41, 66% ee (for *anti*), 66% ee (for *syn*);  $[\alpha]_D^{27} = -55.0$  ( $c = 1.01$ ,  $\text{CHCl}_3$ ).

<sup>1</sup>H NMR (500 MHz,  $\text{CDCl}_3$ ): (a) *anti*-isomer:  $\delta$  0.88 (d,  $J = 6.8$  Hz, 3H), 2.16 (d,  $J = 2.3$  Hz, 1H), 2.49 (dq,  $J = 7.3, 15.3$  Hz, 1H), 4.36 (dd,  $J = 1.4, 7.9$  Hz, 1H), 5.16-5.24 (m, 2H), 5.81 (ddd,  $J = 7.6, 10.3, 17.2$  Hz, 1H), 7.24-7.38 (m, 5H); (b) *syn*-isomer:  $\delta$  1.02 (d,  $J = 6.8$  Hz, 3H), 1.96 (d,  $J = 2.3$  Hz, 1H), 2.59 (m, 1H), 4.61 (m, 1H), 5.04 (m, 1H), 5.07 (m, 1H), 5.76 (m, 1H), 7.24-7.38 (m, 5H). HPLC condition: Chiralcel IC, *n*-hexane/2-propanol 98/2, flow rate 0.7 mL/min, UV 214 nm, *syn*:  $t_R = 10.2$  min (minor)(*1S,2R*),  $t_R = 11.2$  min (major) (*1R,2S*); *anti*:  $t_R = 11.9$  min (minor) (*1R,2R*),  $t_R = 12.9$  min (major) (*1S,2S*).

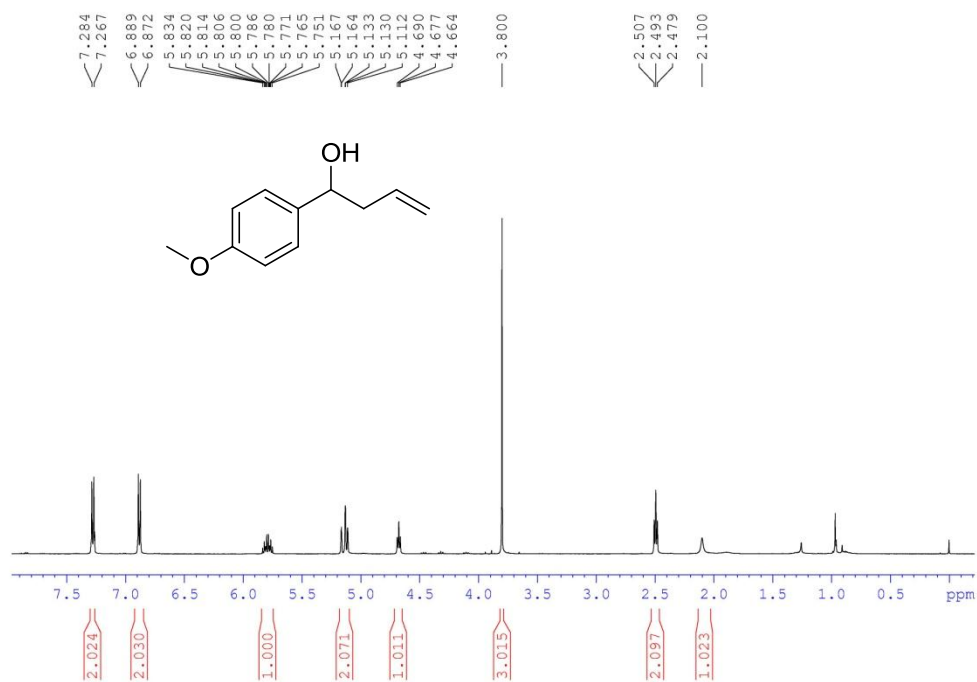
<sup>5</sup> J.-Z. Chen, D.-L. Liu, D.-Y. Fan, Y.-G. Liu, W.-B. Zhang, *Tetrahedron*, **2013**, *69*, 8161–8168.

### (3) NMR Spectra for Homoallylic Alcohols

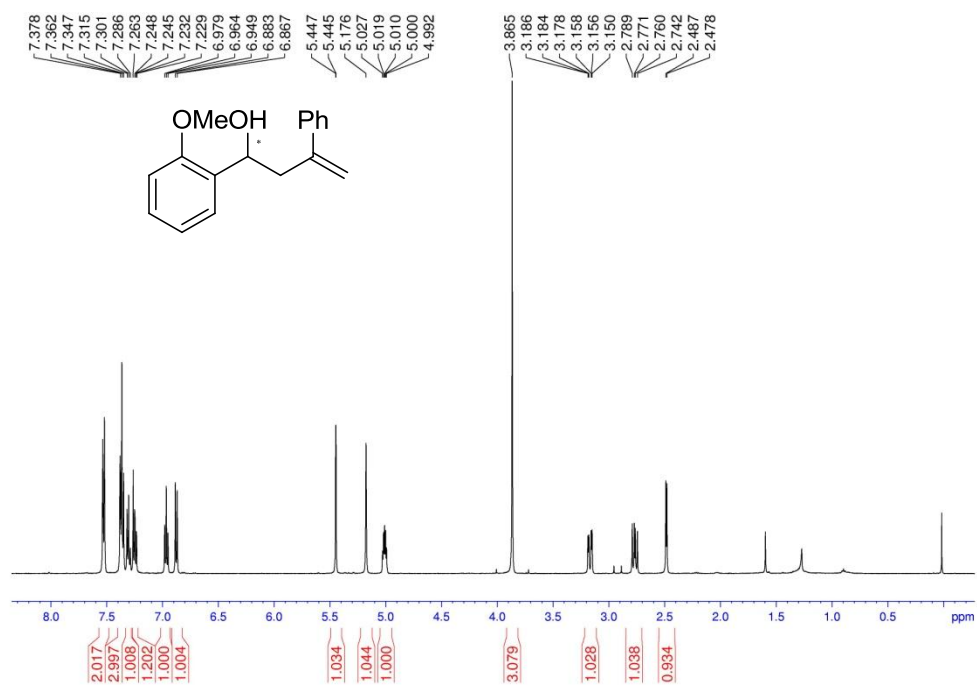
#### 1-(4-Methoxyphenyl)-3-phenylbut-3-en-1-ol (1)

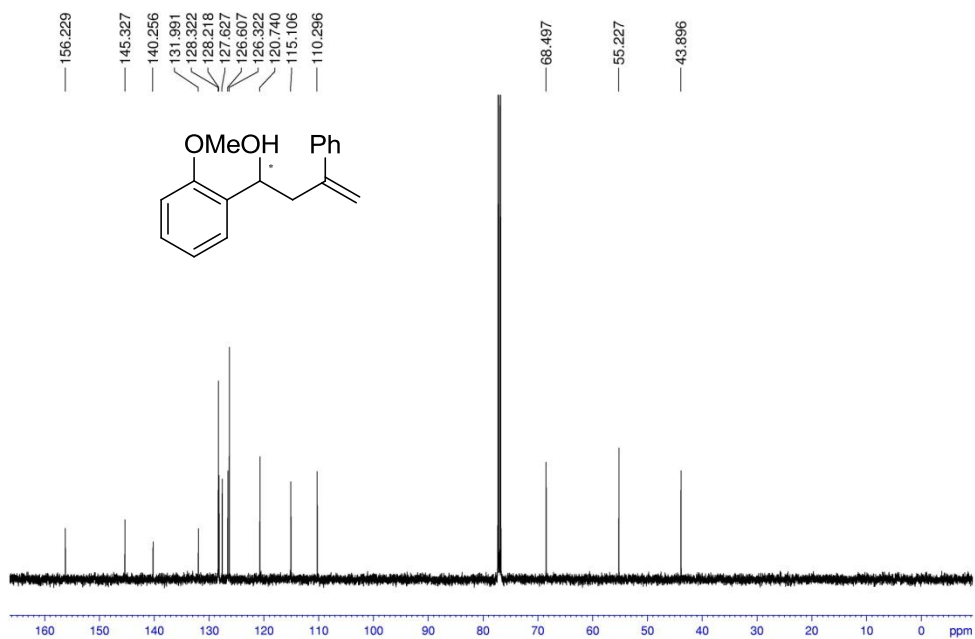


### 1-(4-Methoxyphenyl)but-3-en-1-ol (5)

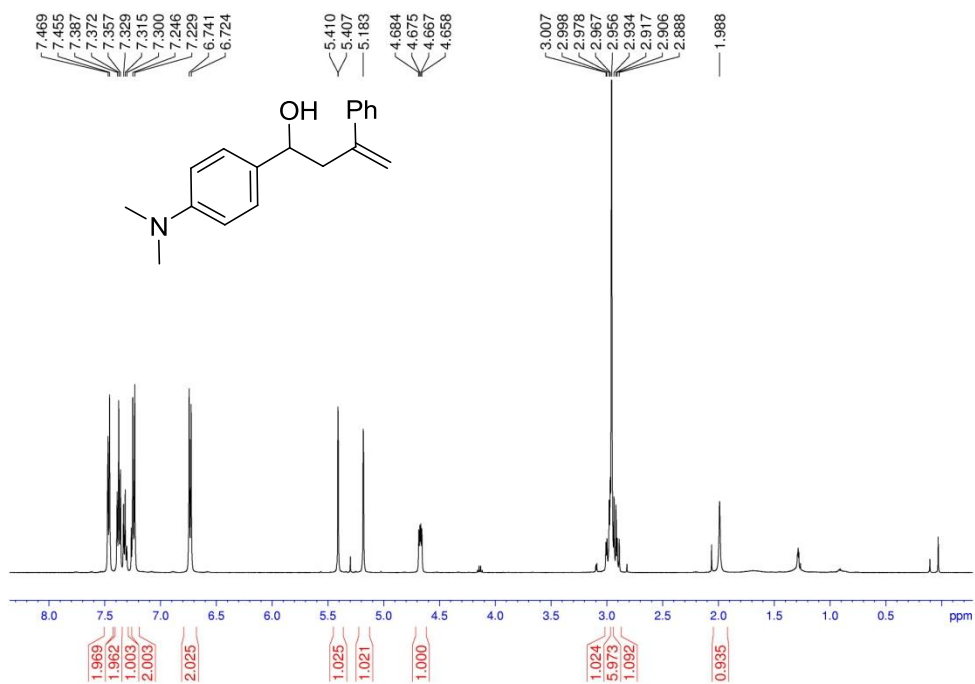


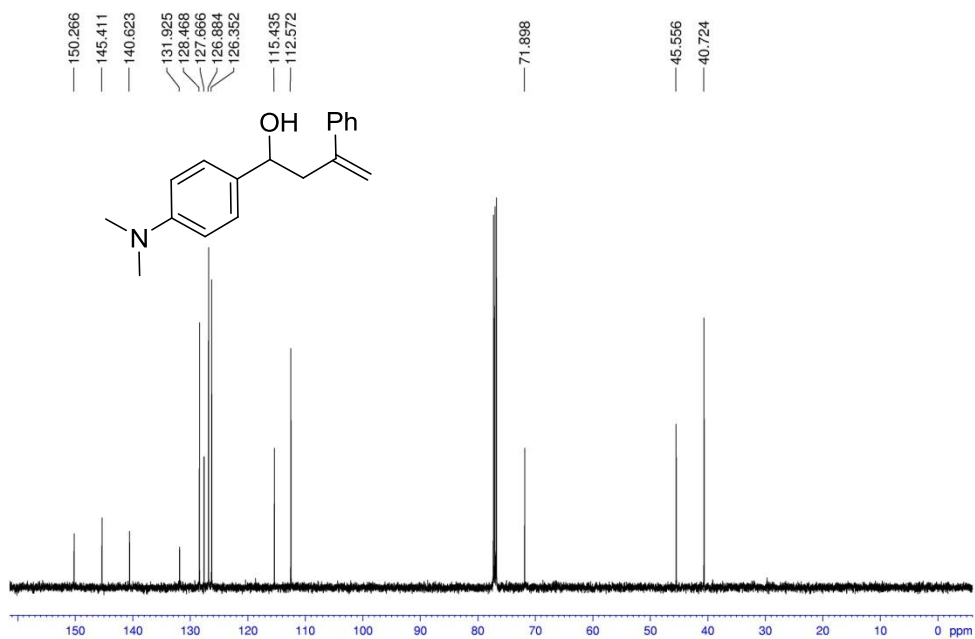
### 1-(2-Methoxyphenyl)-3-phenylbut-3-en-1-ol (6)



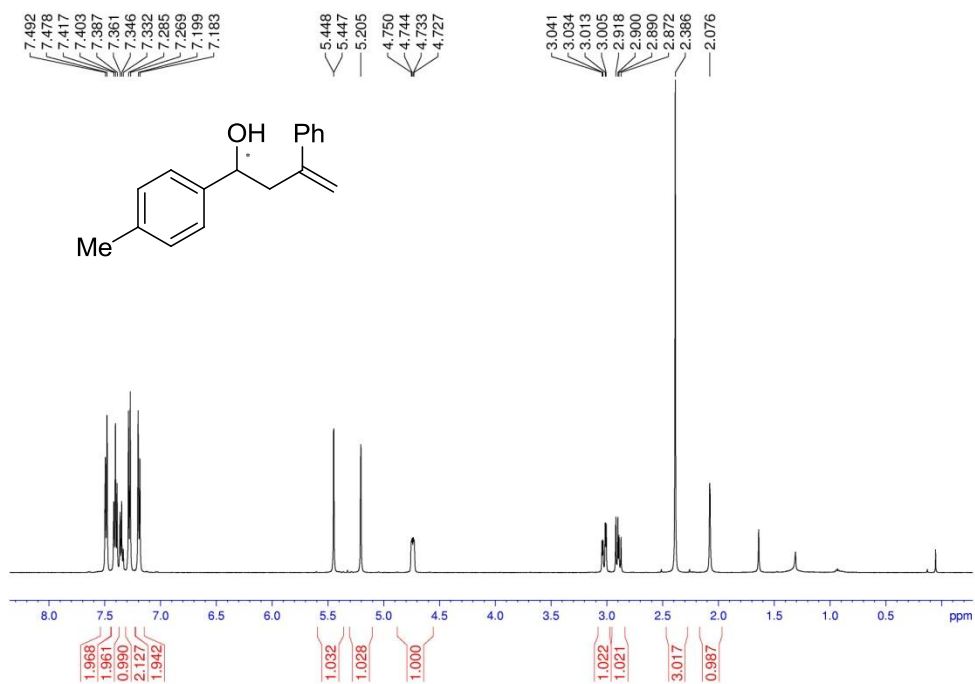


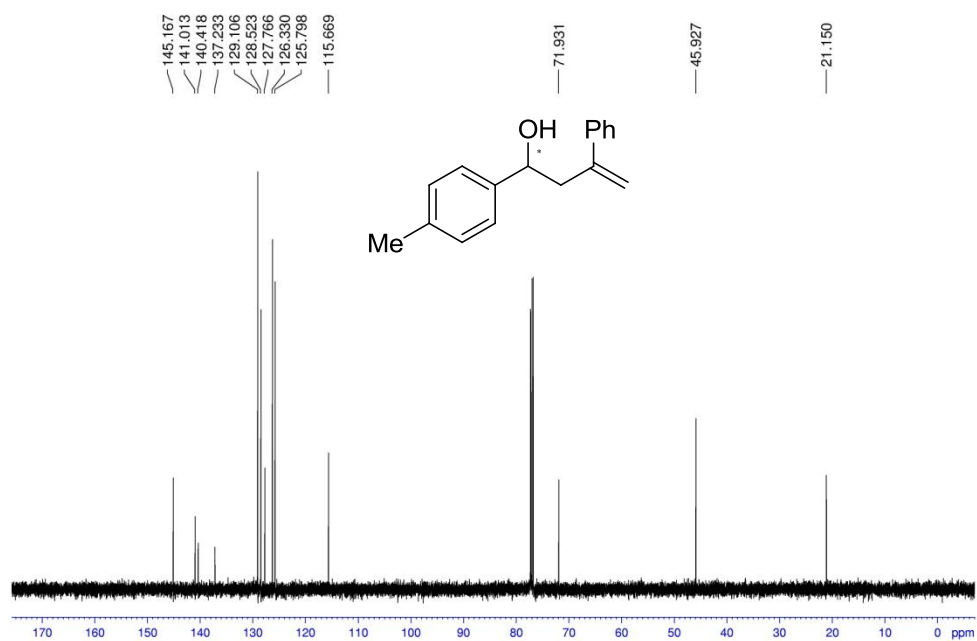
**1-(4-(Dimethylamino)phenyl)-3-phenylbut-3-en-1-ol (7)**



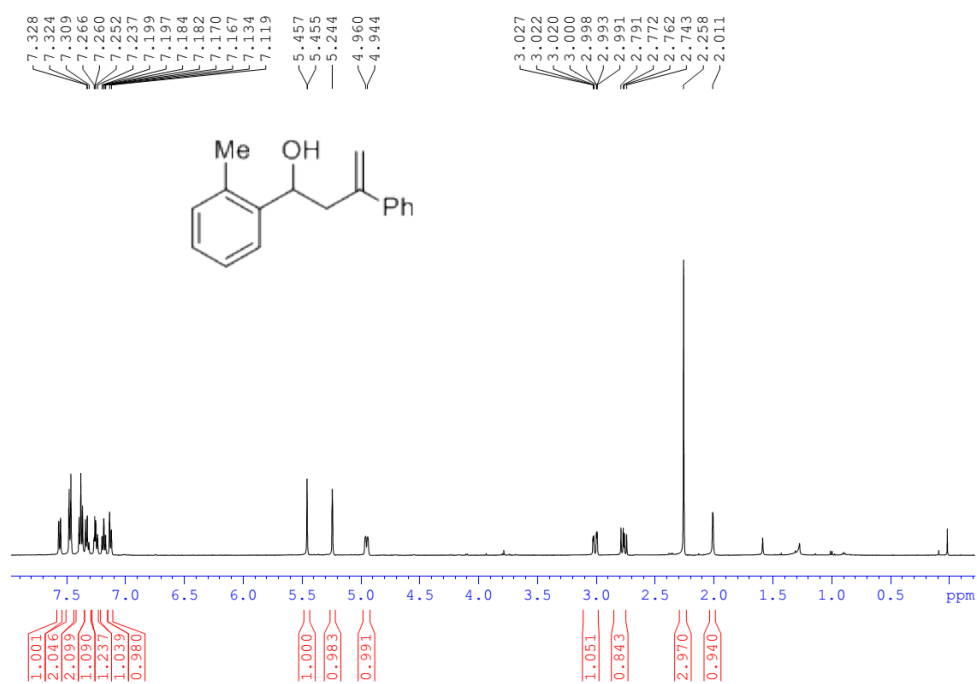


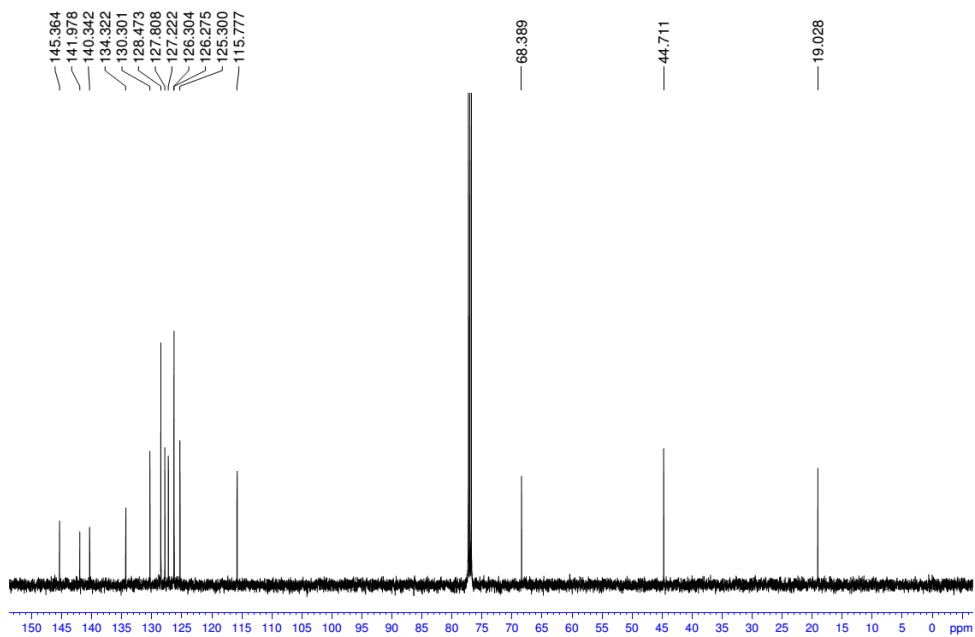
### 3-Phenyl-1-p-tolylbut-3-en-1-ol (8)



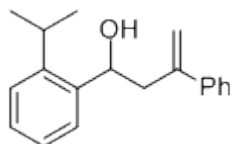
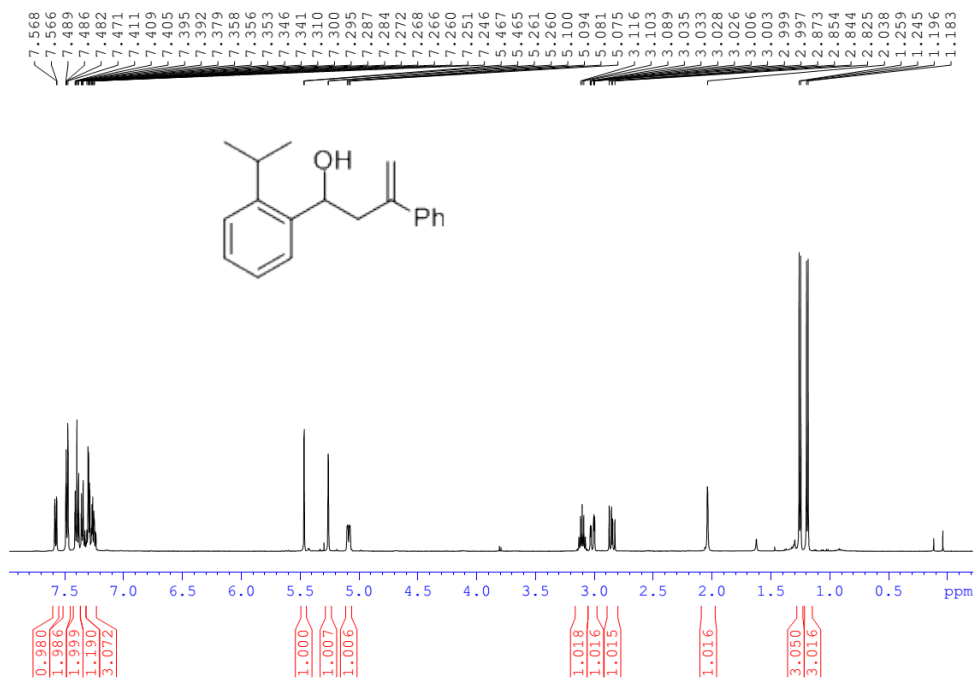


### 3-phenyl-1-o-tolylbut-3-en-1-ol (9)

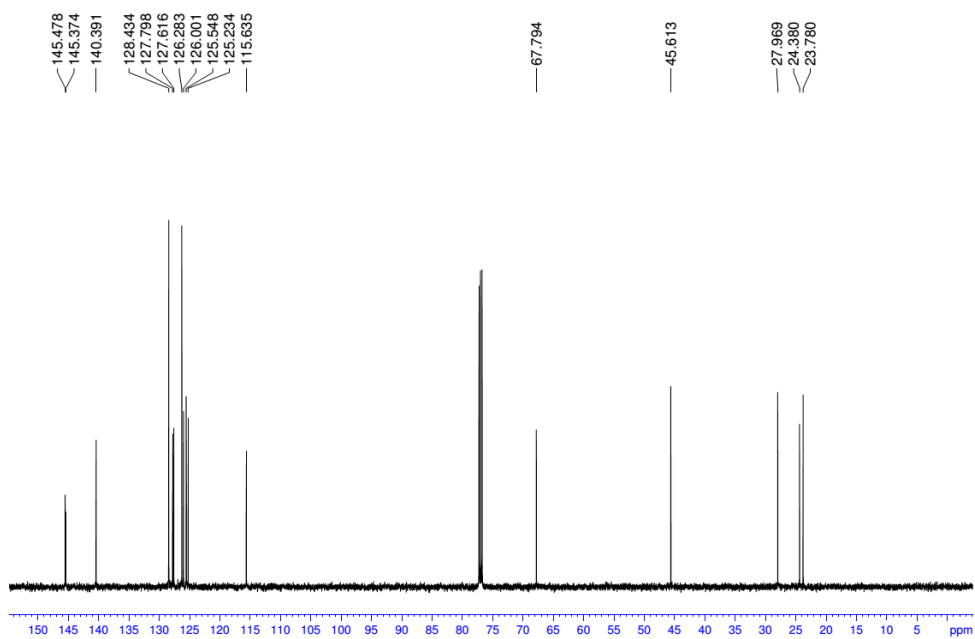




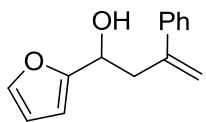
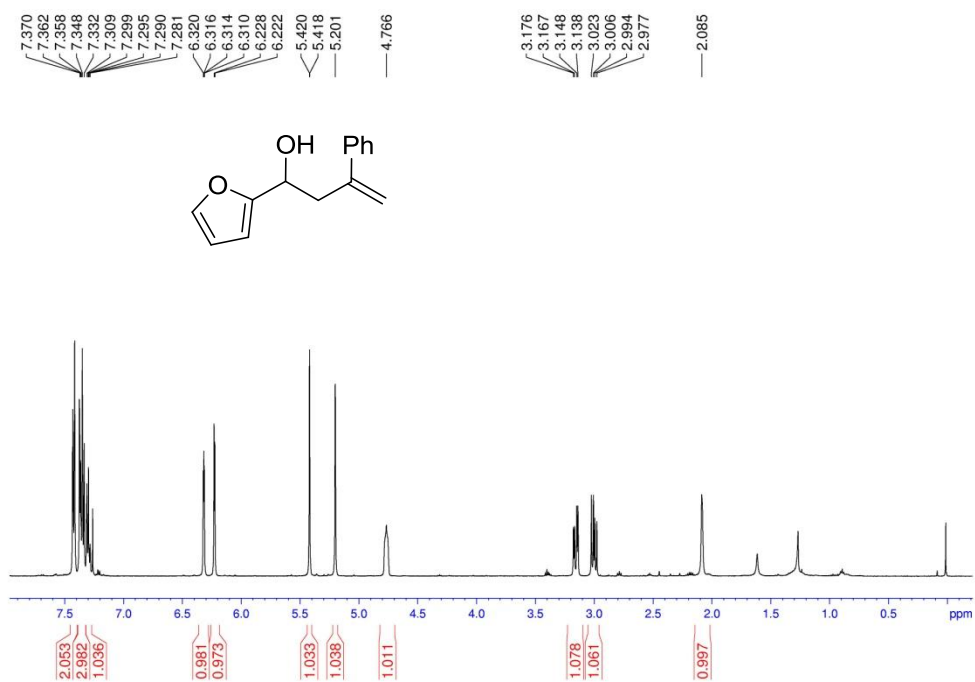
**1-(2-isopropylphenyl)-3-phenylbut-3-en-1-ol (10)**

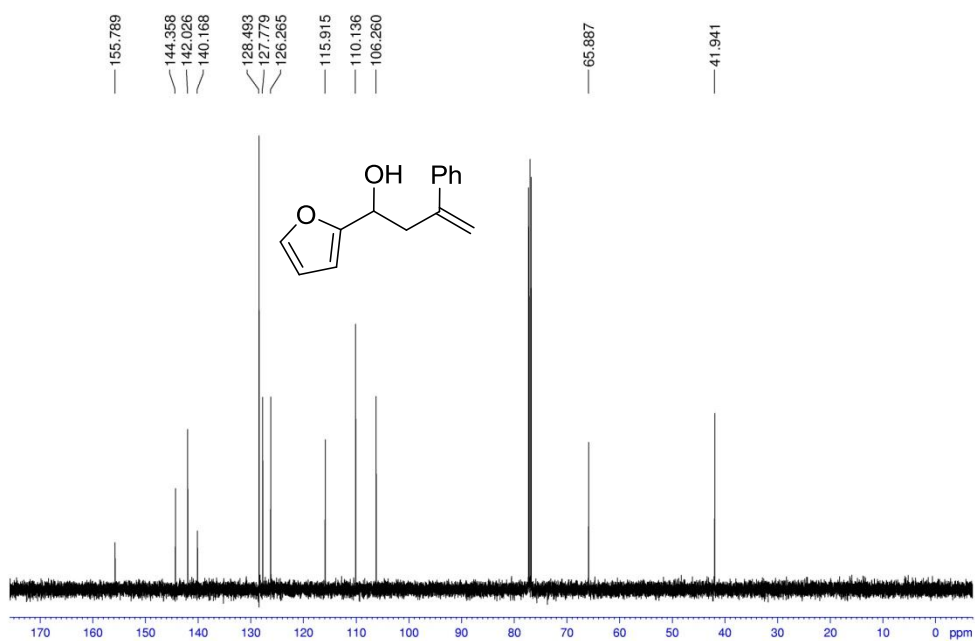




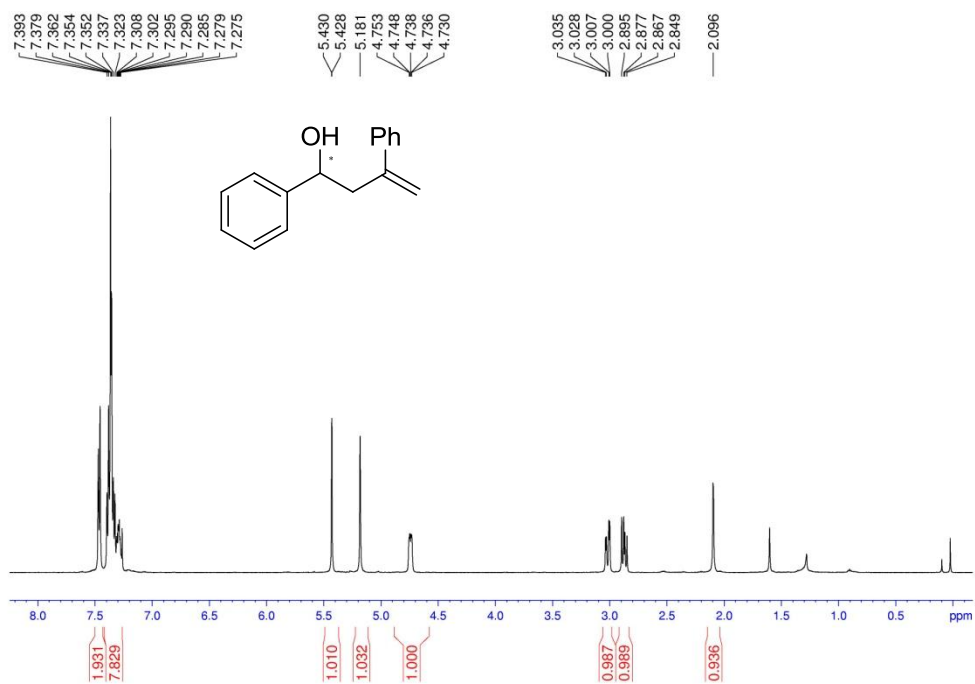


**1-(Furan-2-yl)-3-phenylbut-3-en-1-ol (11)**

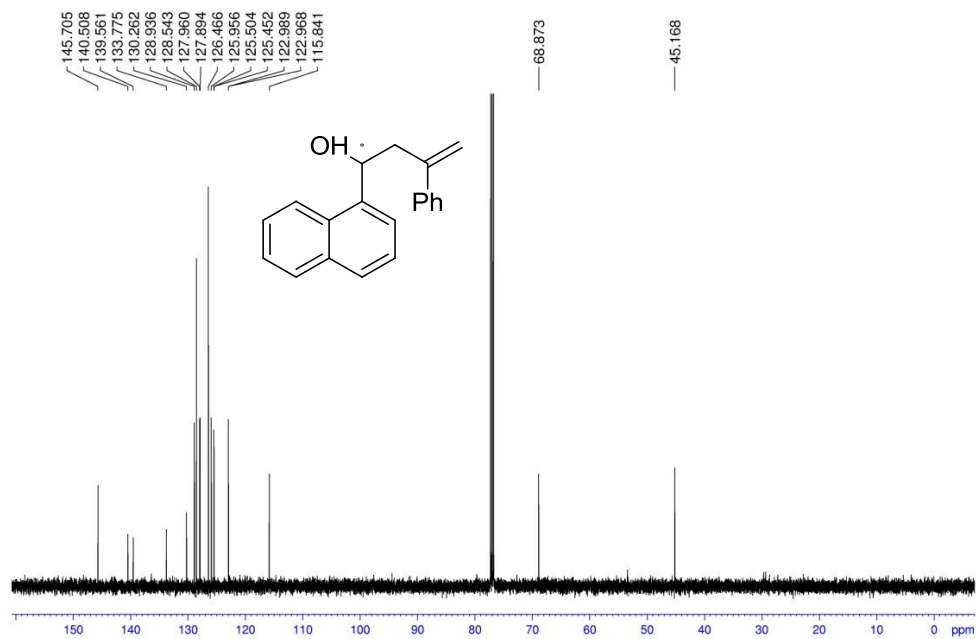
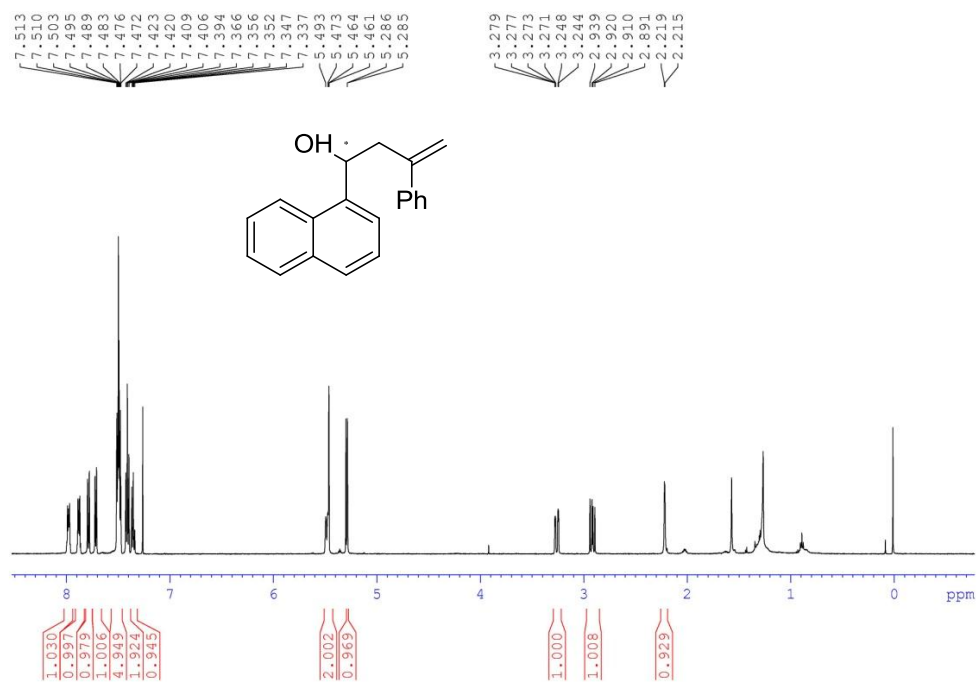




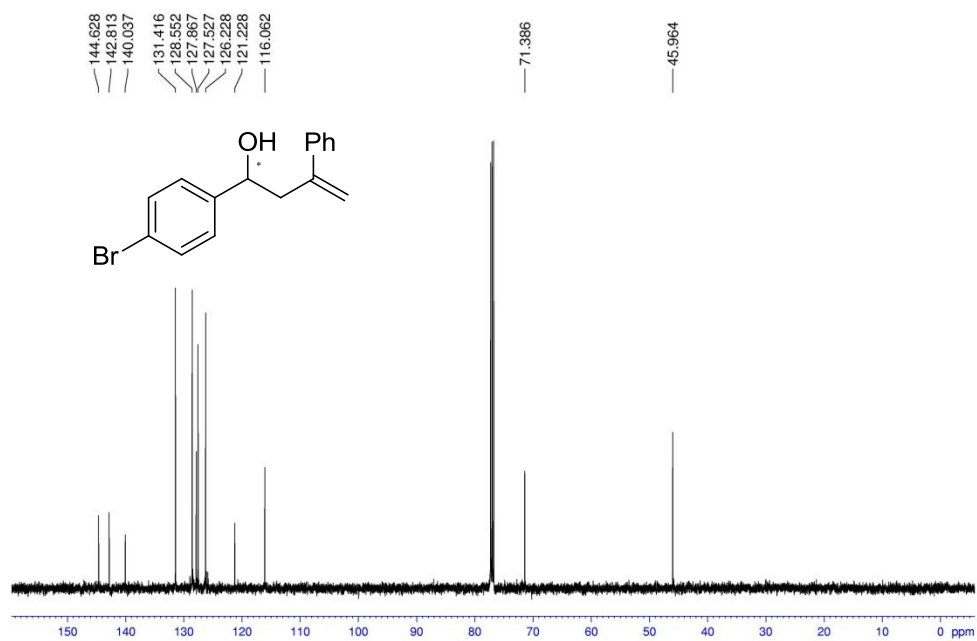
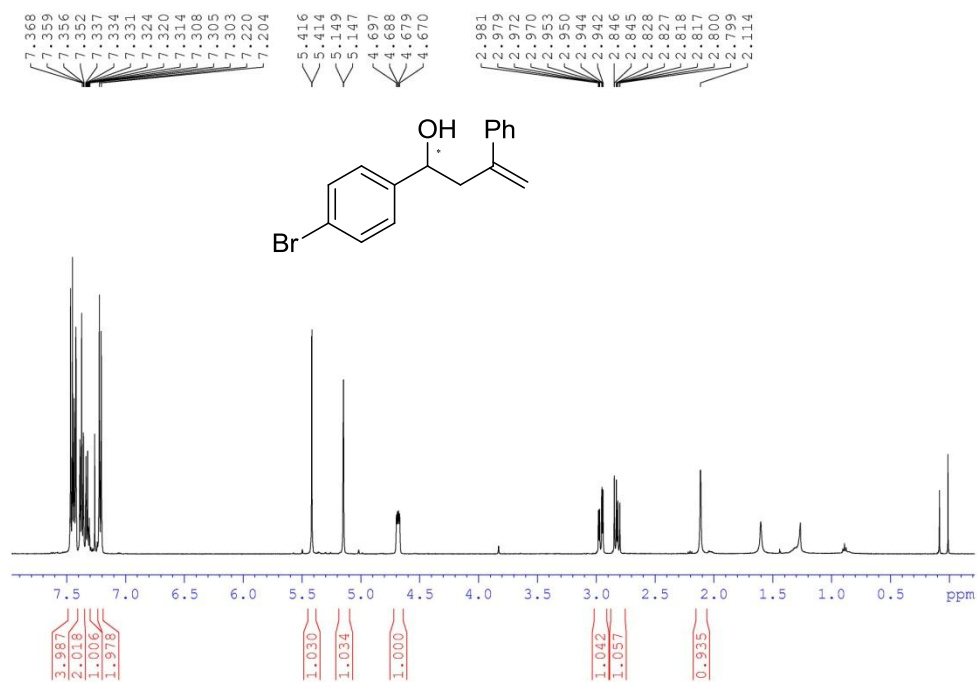
**(+)-1,3-Diphenylbut-3-en-1-ol (12)**



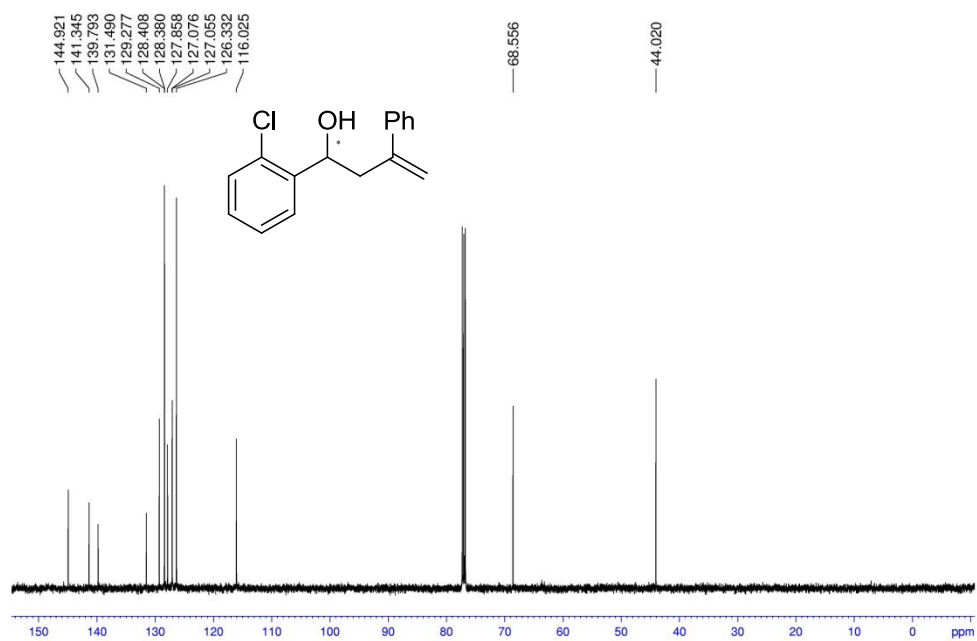
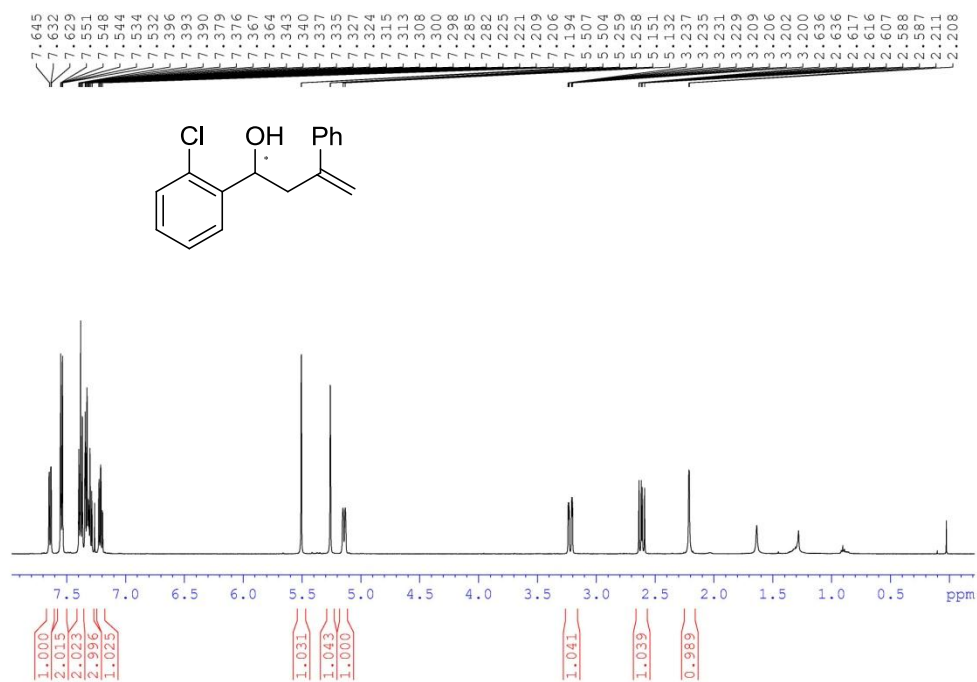
# 1-(Naphthalen-2-yl)-3-phenylbut-3-en-1-ol (13)



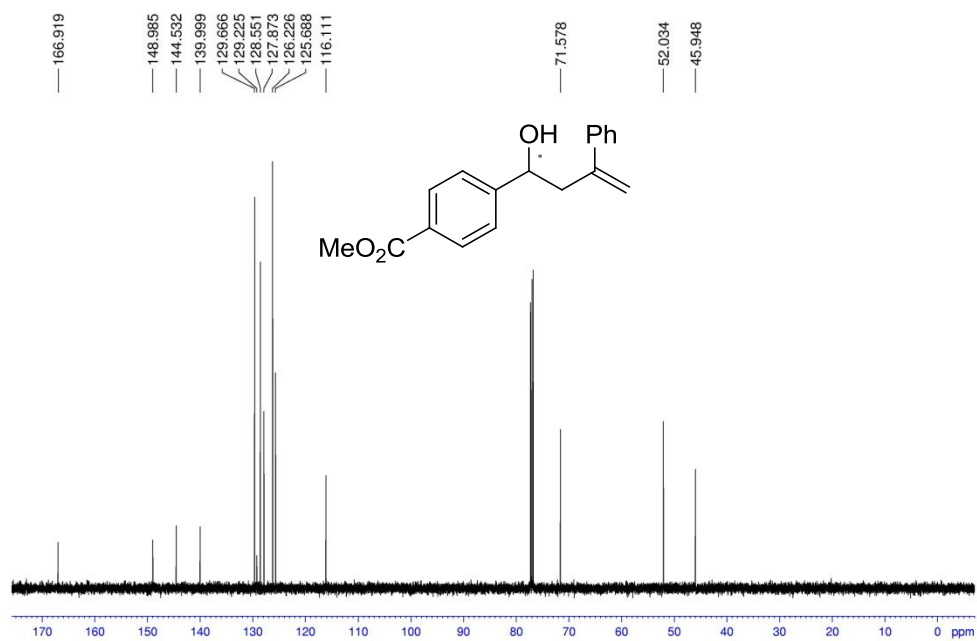
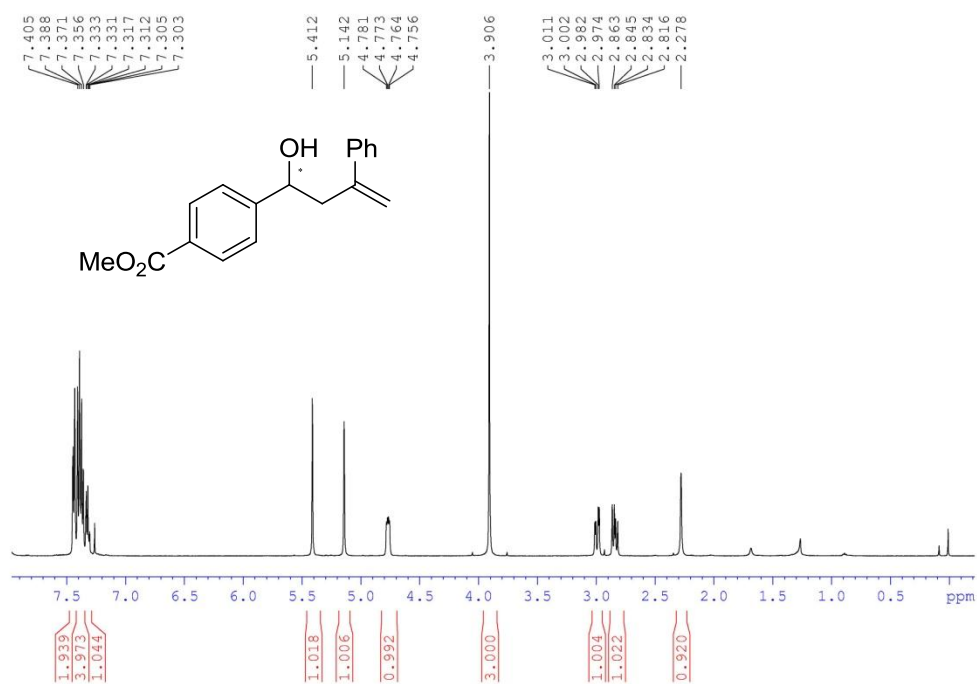
**(+)-1-(4-Bromophenyl)-3-phenylbut-3-en-1-ol (14)**



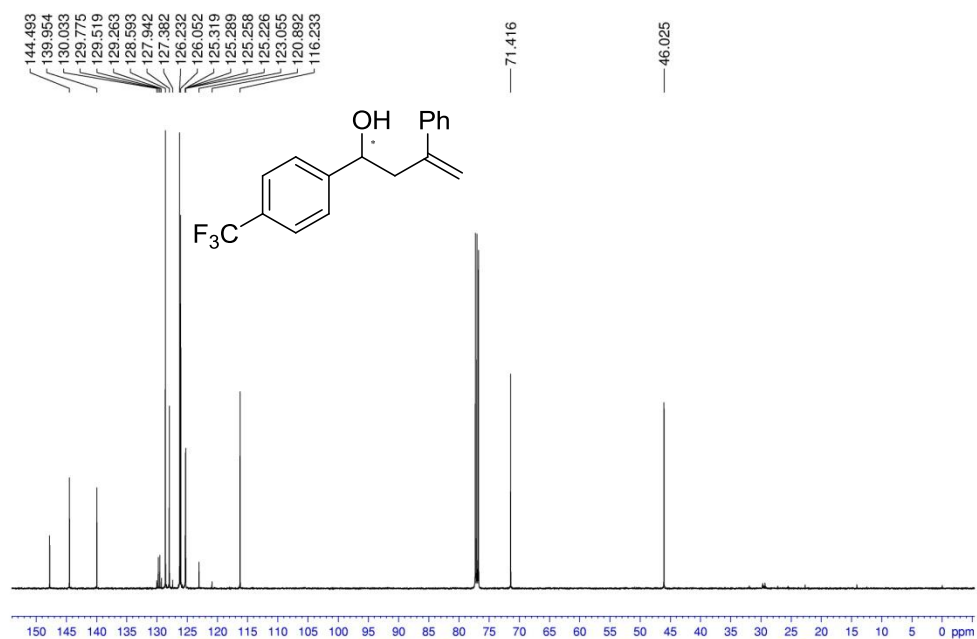
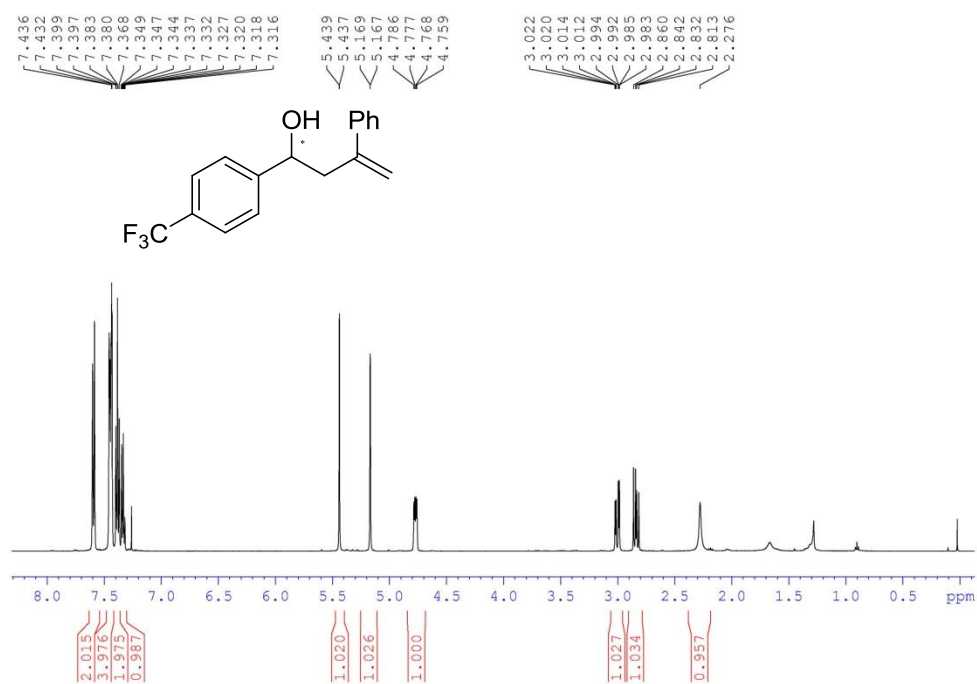
**(-)-1-(2-Chlorophenyl)-3-phenylbut-3-en-1-ol (15)**



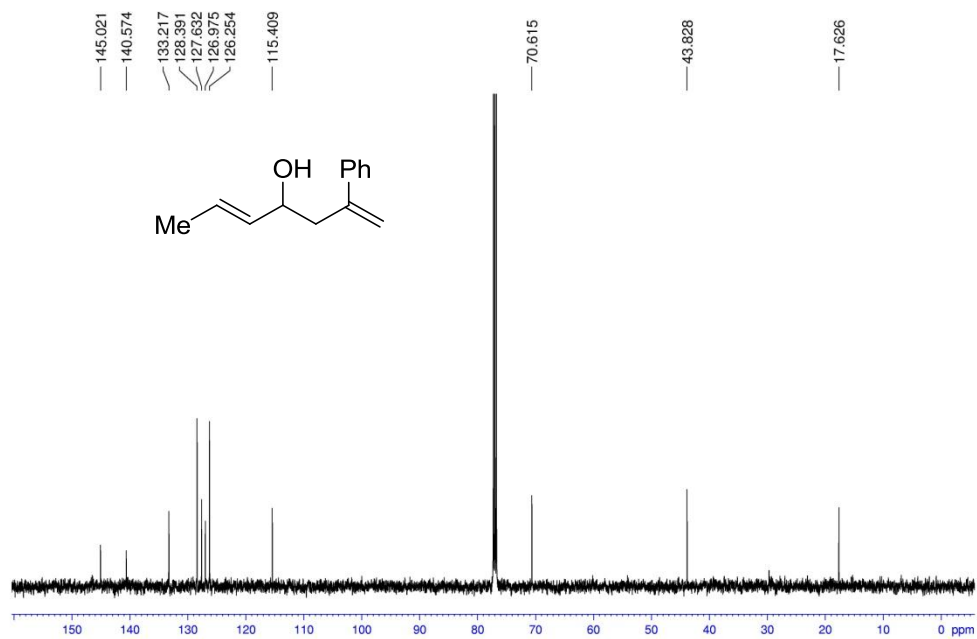
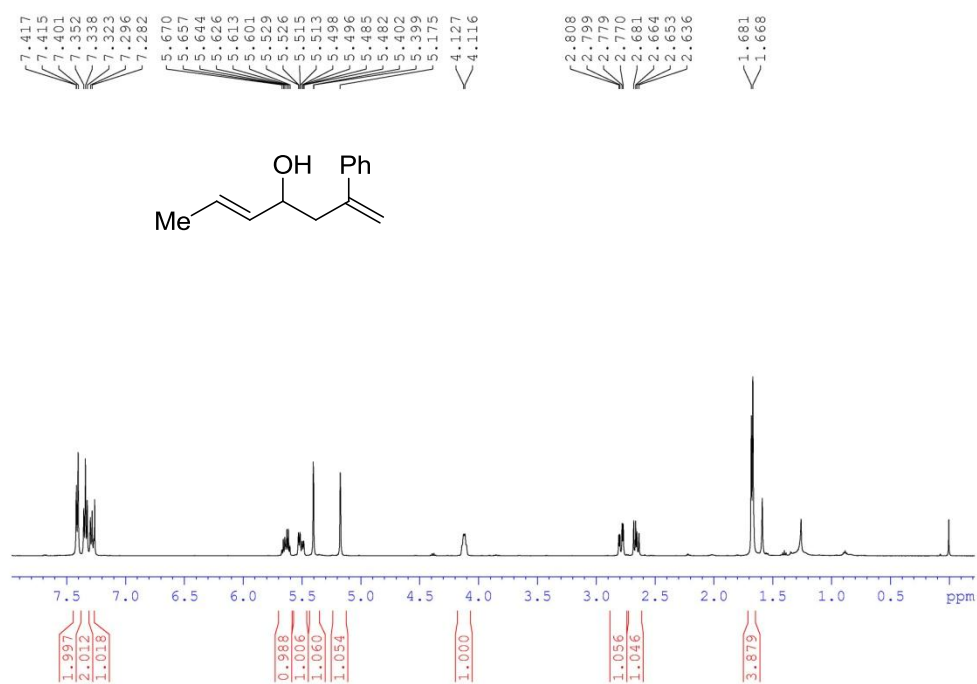
# Methyl-4-(1-hydroxy-3-phenylbut-3-enyl)benzoate (16)



### 3-Phenyl-1-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (17)

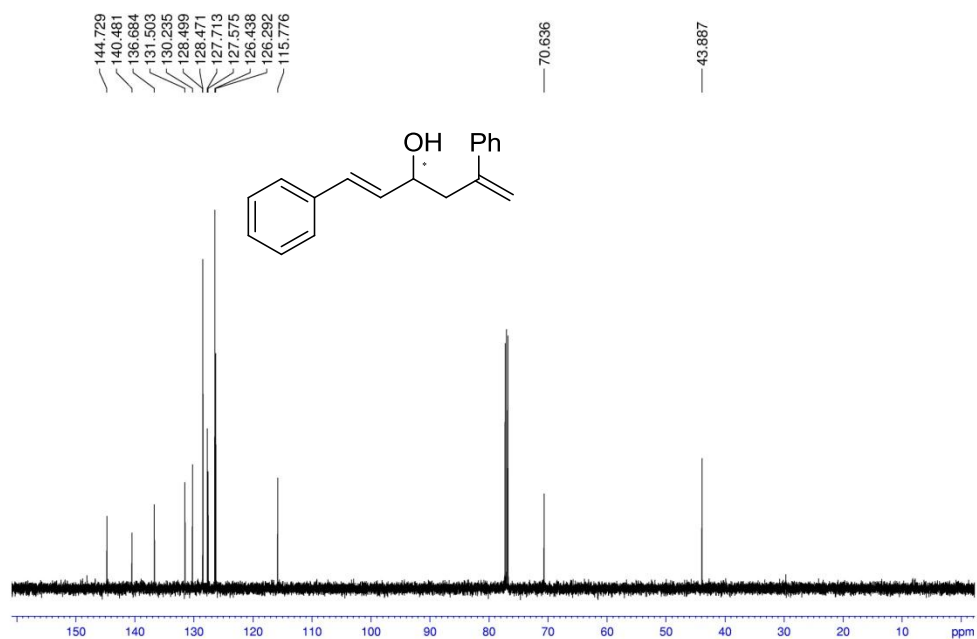
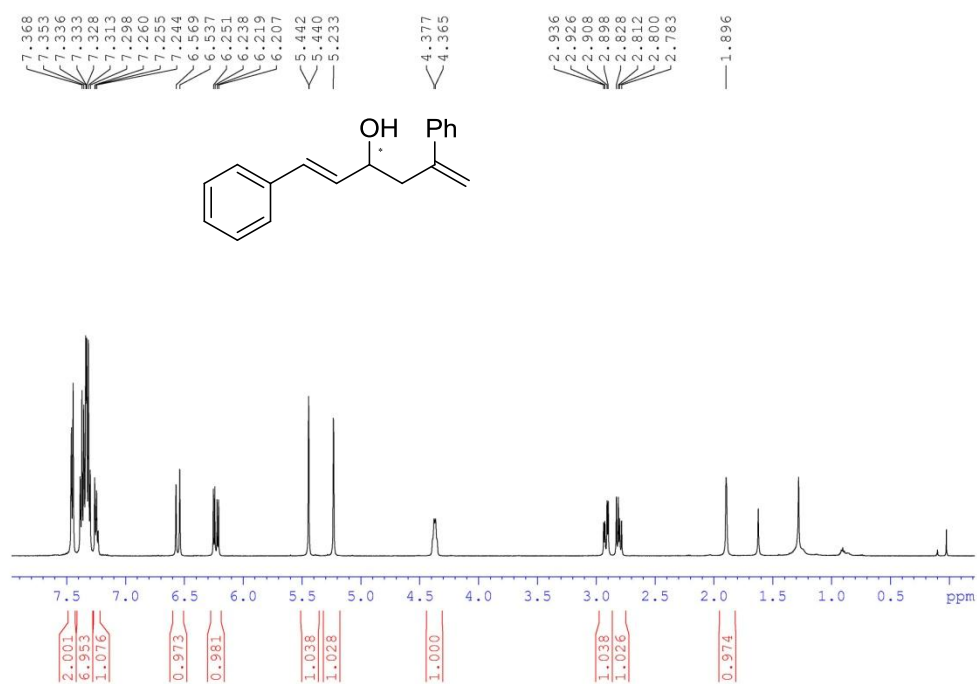


**(E)-2-Phenylhepta-1,5-dien-4-ol (18)**

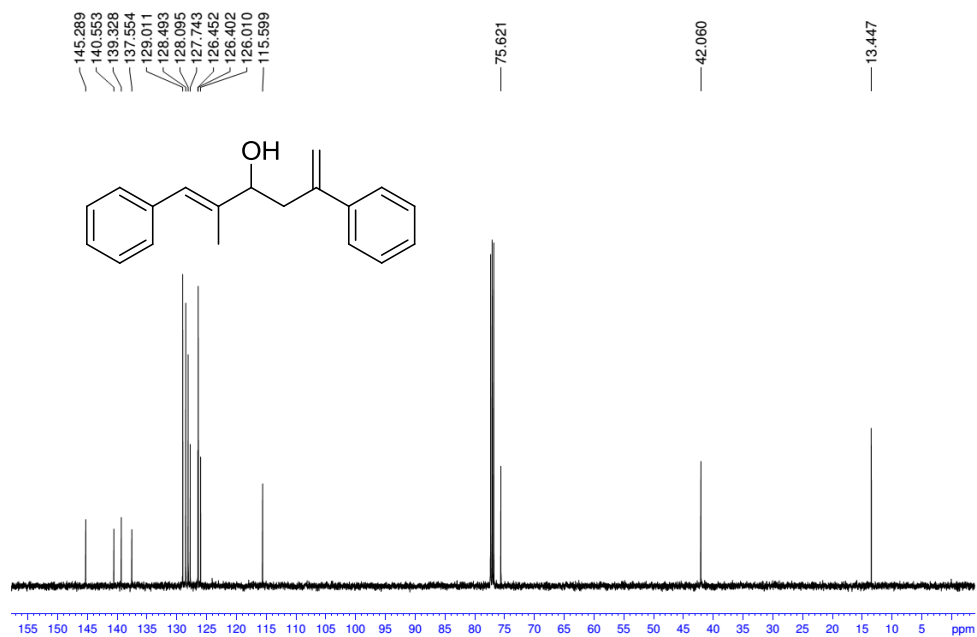
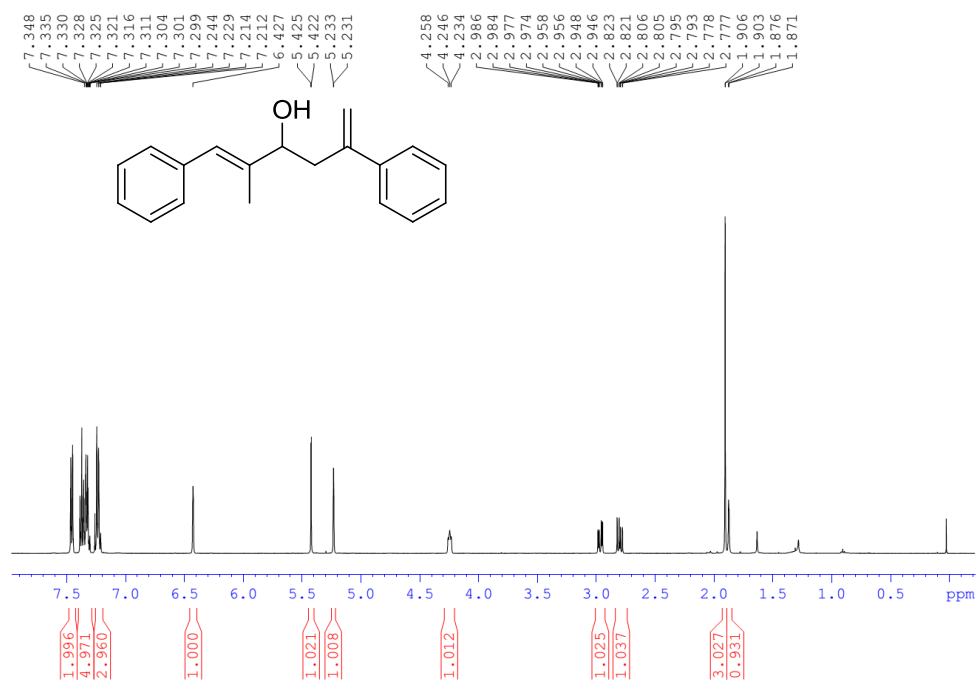




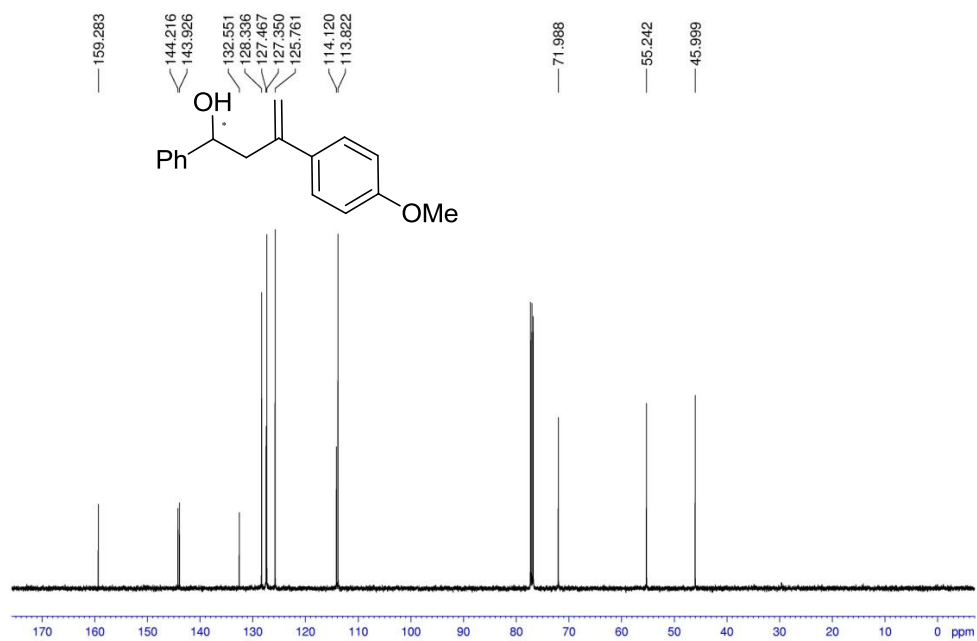
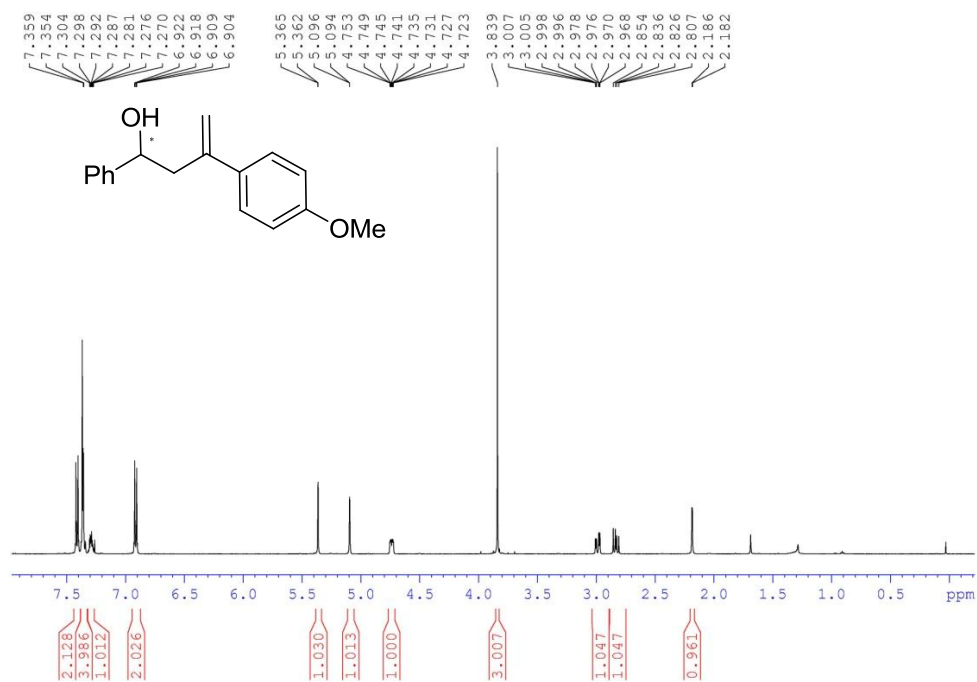
**(E)-1,5-Diphenylhexa-1,5-dien-3-ol (19)**



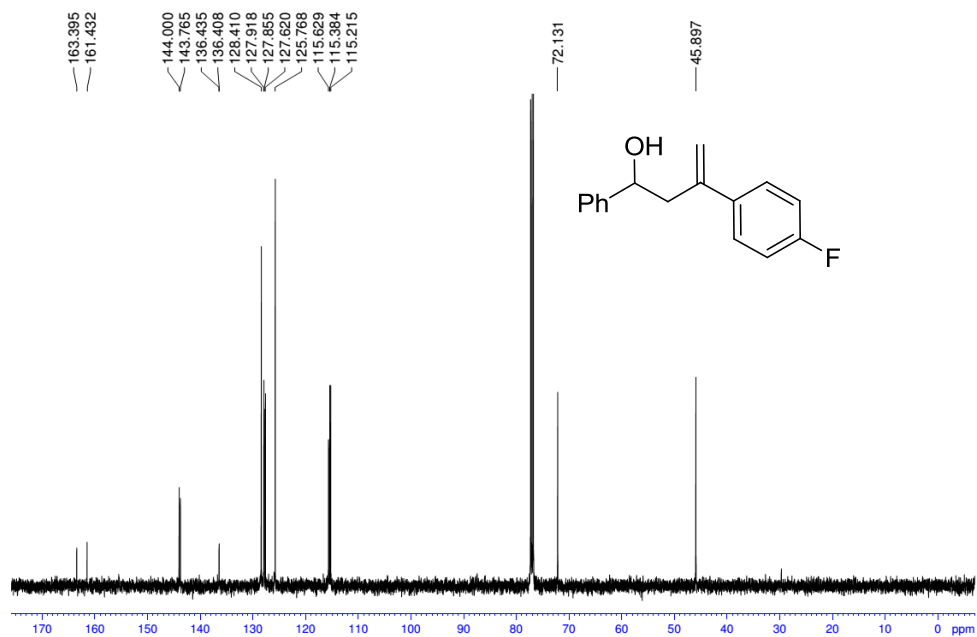
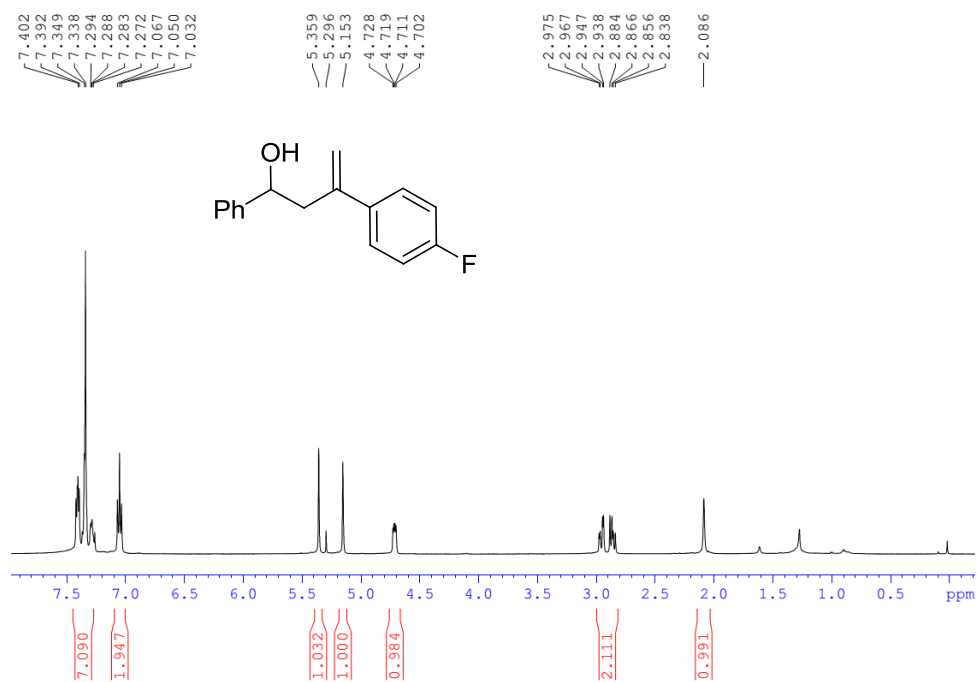
**(E)-2-methyl-1,5-diphenylhexa-1,5-dien-3-ol (20)**



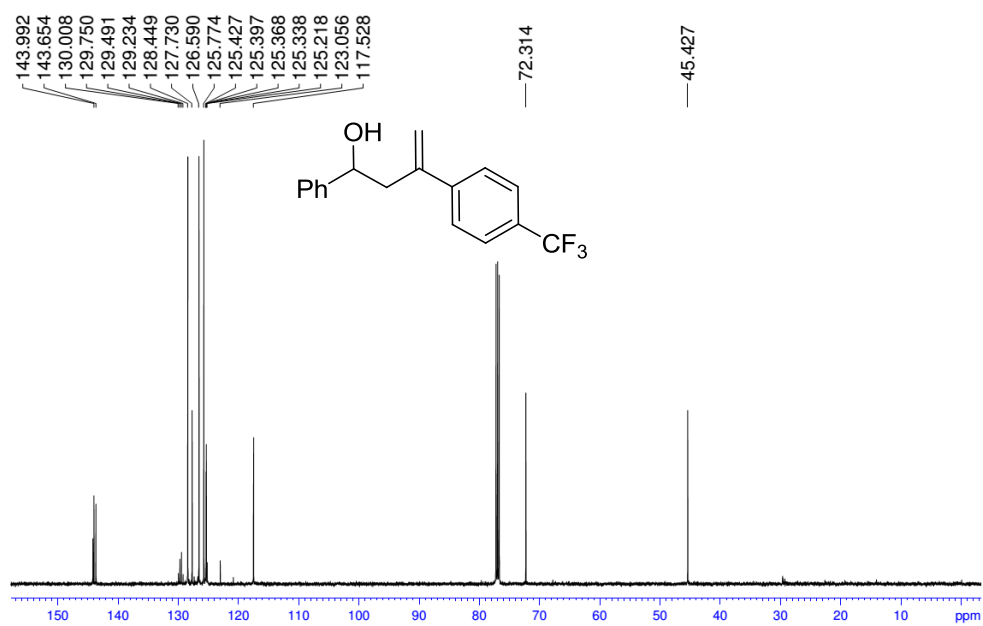
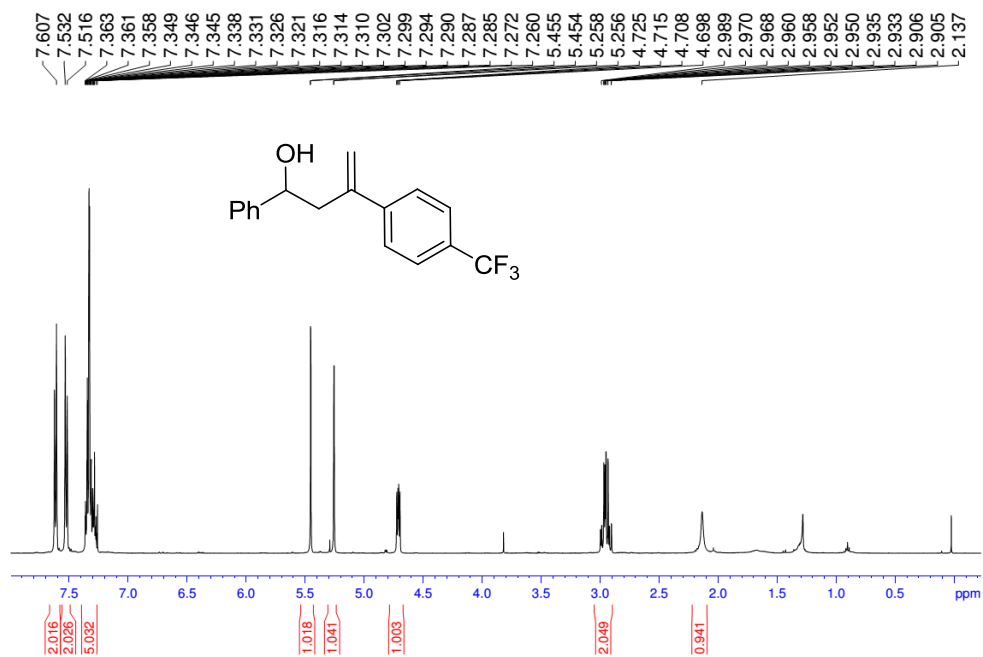
### 3-(4-Methoxyphenyl)-1-phenylbut-3-en-1-ol (22)



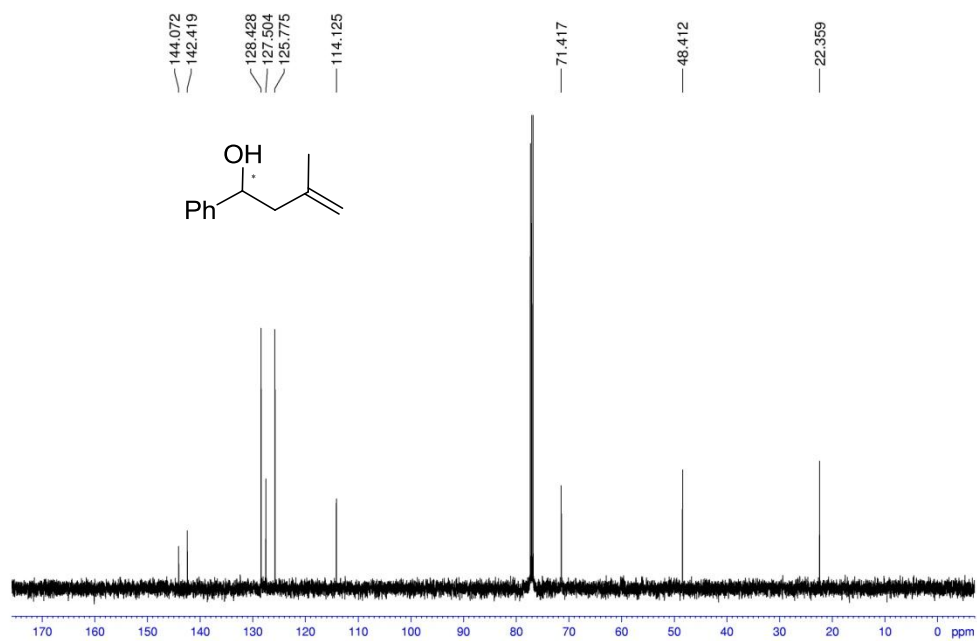
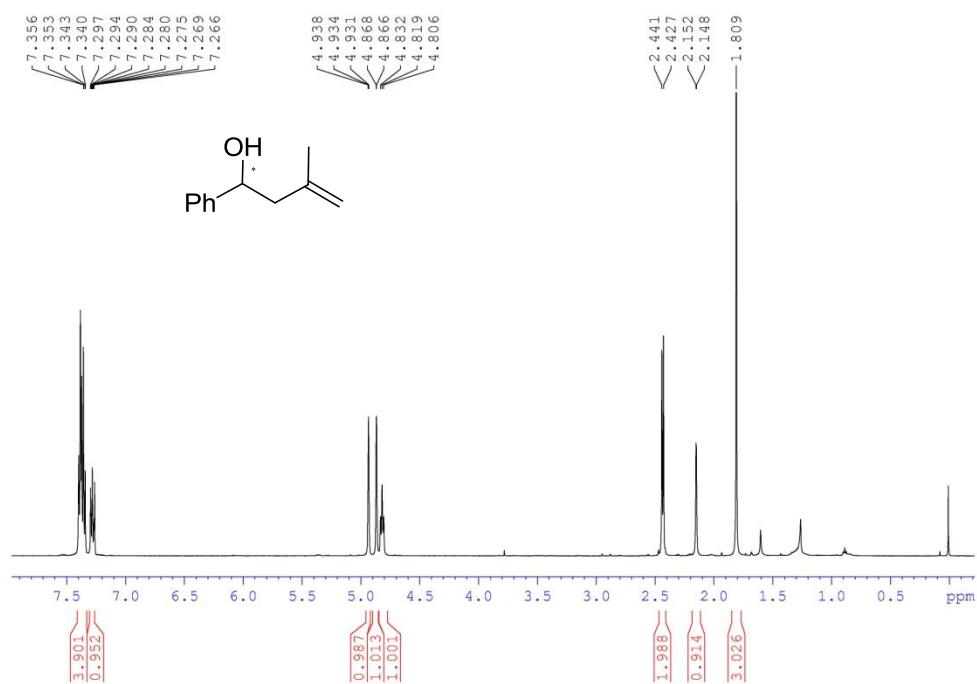
### 3-(4-fluorophenyl)-1-phenylbut-3-en-1-ol (23)



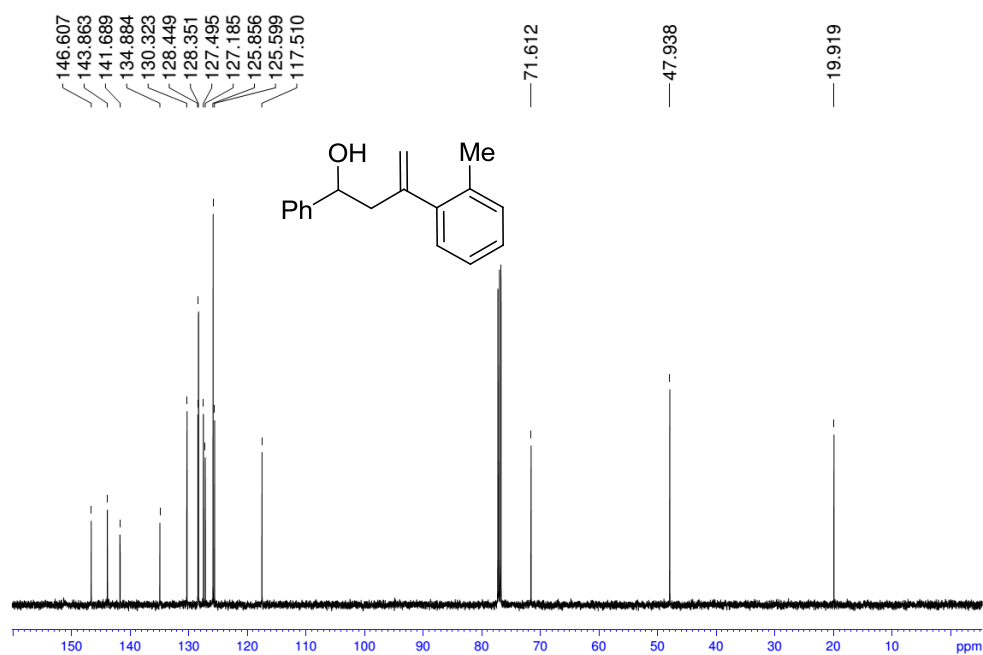
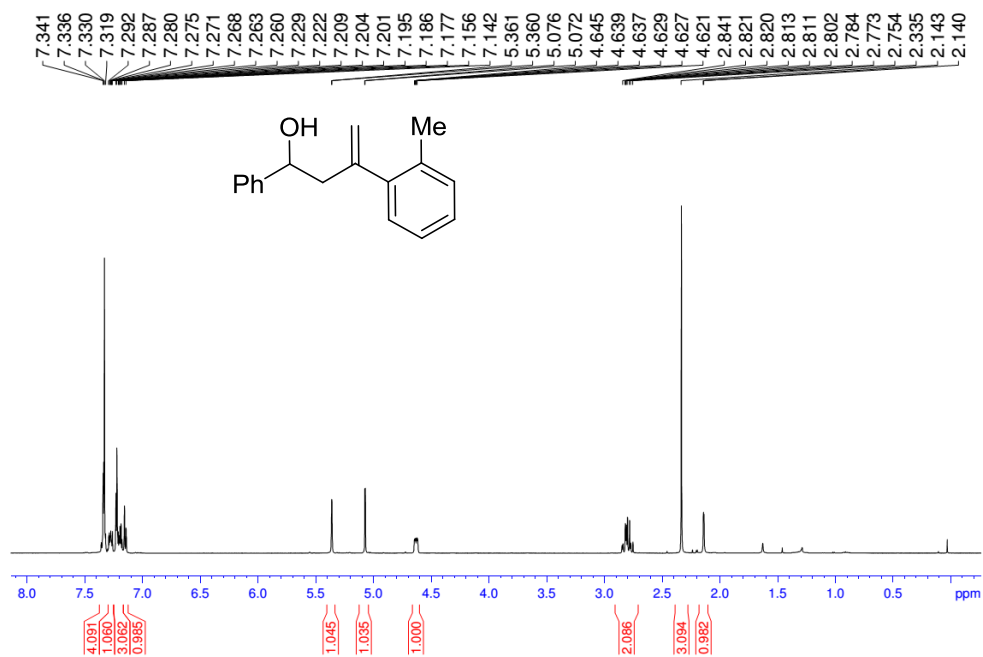
# 1-phenyl-3-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (24)



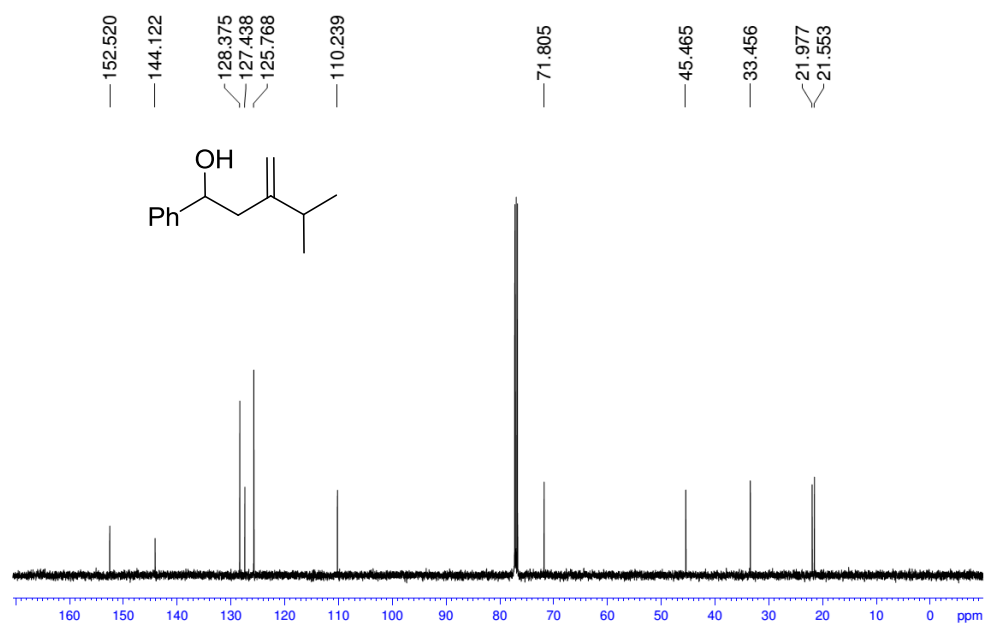
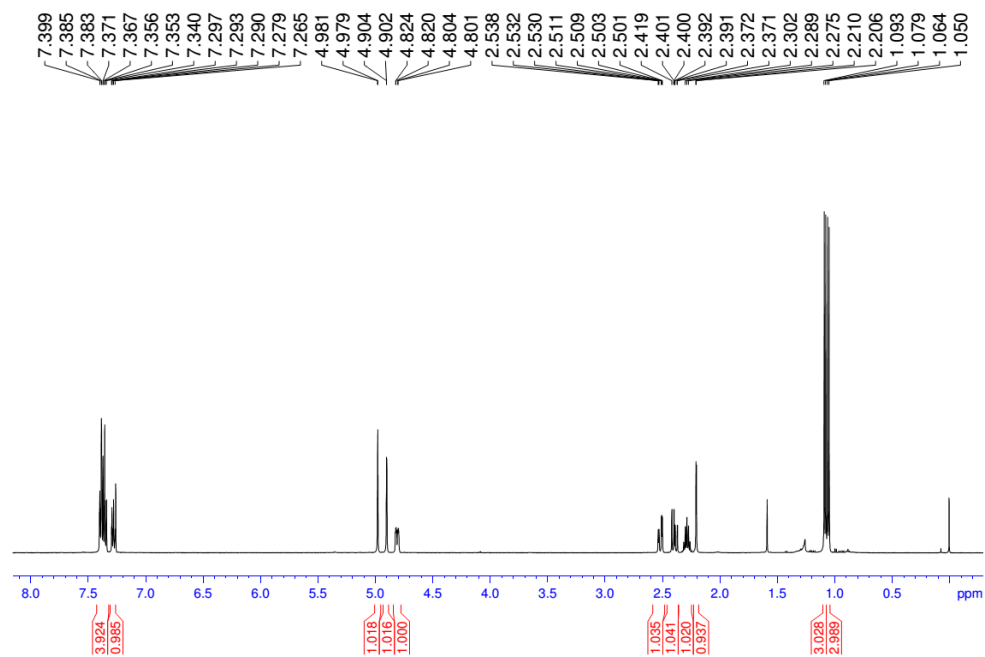
**(S)-3-Methyl-1-phenylbut-3-en-1-ol (25)**



# 1-phenyl-3-o-tolylbut-3-en-1-ol (26)

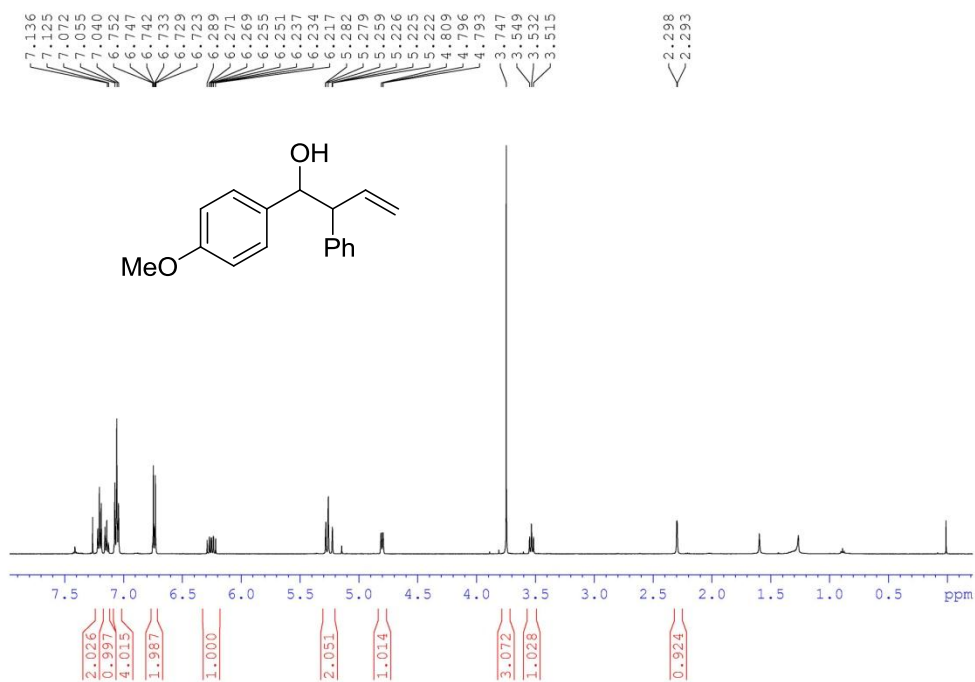


**4-methyl-3-methylene-1-phenylpentan-1-ol (27)**

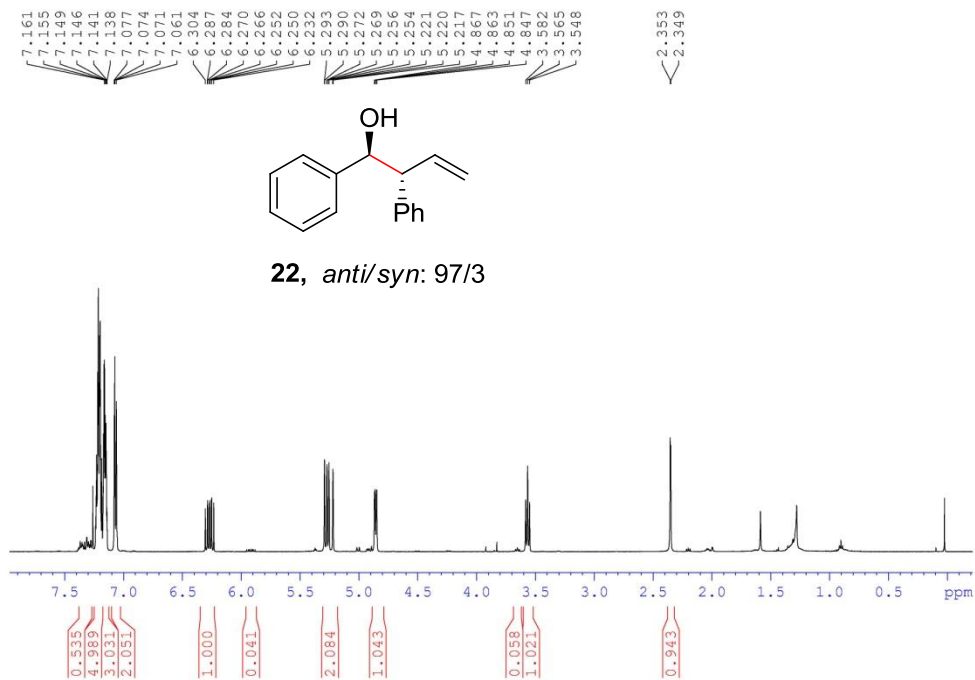




***anti*-1-(4-Methoxyphenyl)-2-phenylbut-3-en-1-ol (28)**

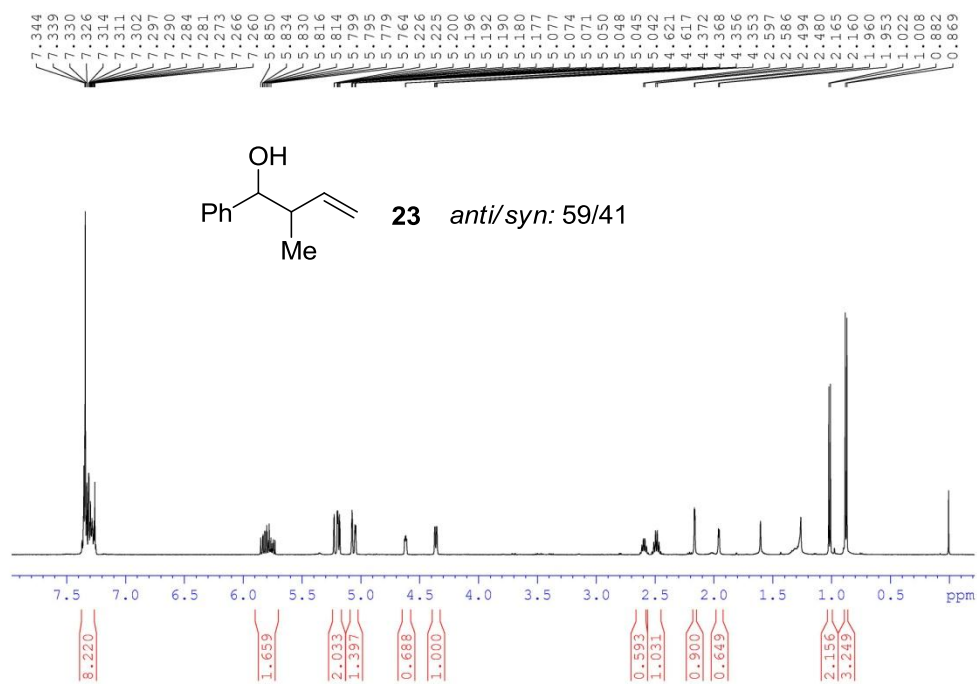


**(1*S*,2*R*)-(-)-1,2-Diphenylbut-3-en-1-ol (29)**

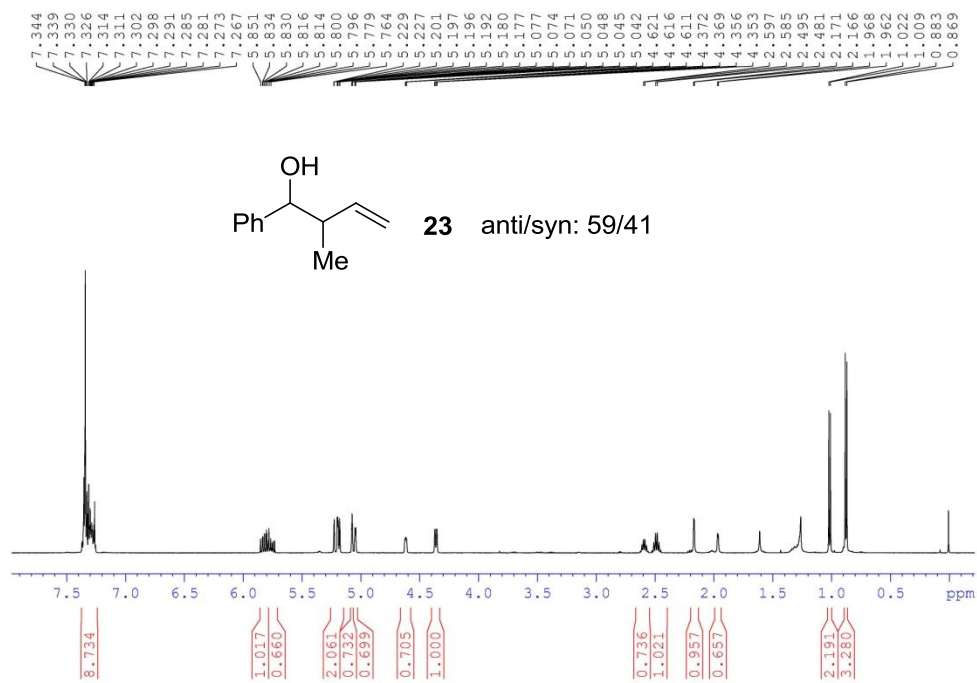


## 2-Methyl-1-phenylbut-3-en-1-ol (30)

For (table 3, entry 4)



For (table 3, entry 5)

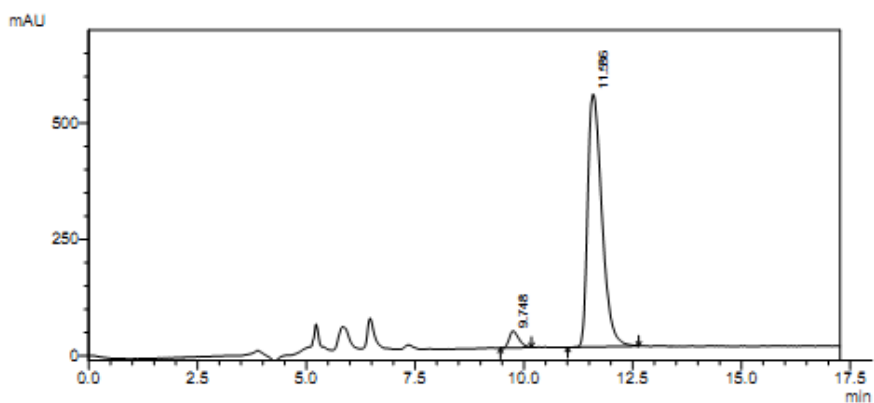
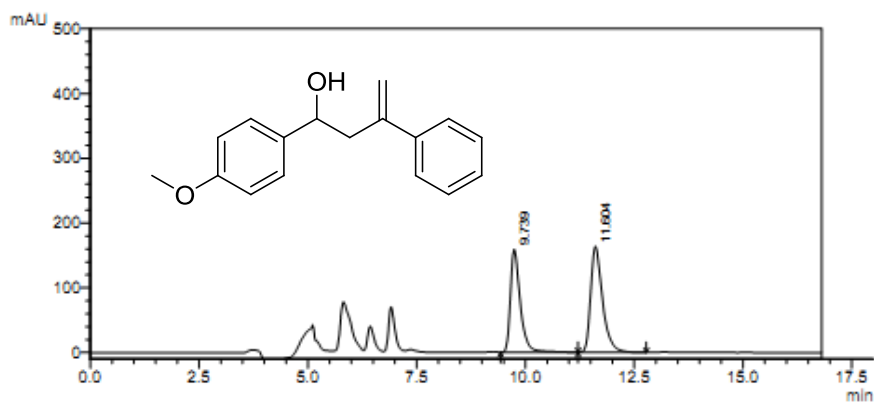


## (4) HPLC Charts for Homoallylic Alcohols

### (+)-1-(4-Methoxyphenyl)-3-phenylbut-3-en-1-ol (1)

2013-10-13 15:09:51 1 / 1

#### ==== Shimadzu LcSolution Report====



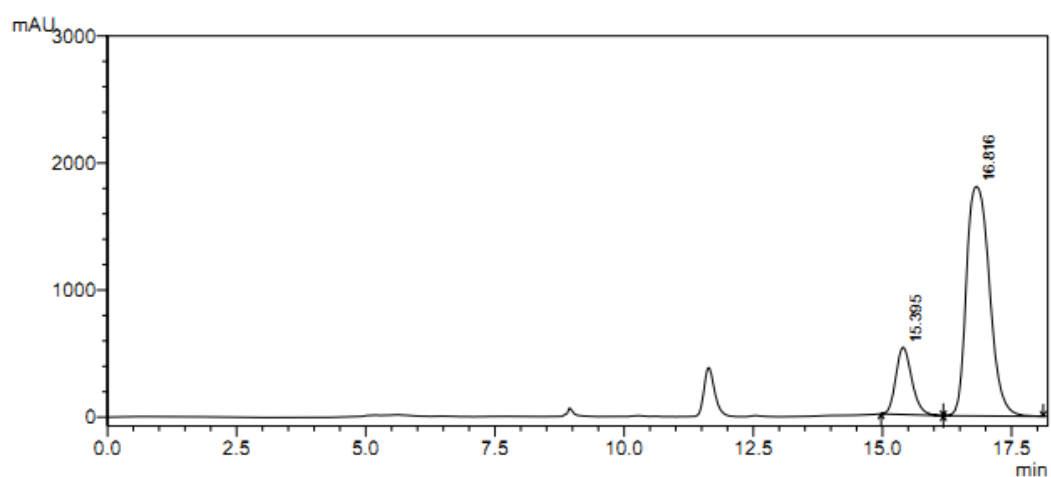
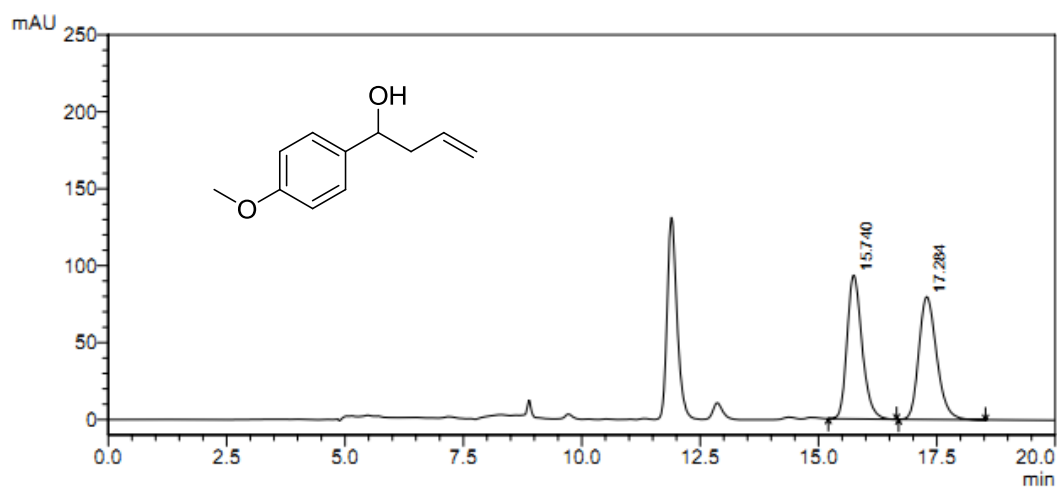
PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	9.748	552331	36456	4.275	6.290
2	11.586	12369144	543121	95.725	93.710
总计		12921475	579577	100.000	100.000

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(S)-1-(4-Methoxyphenyl)but-3-en-1-ol (5)

==== Shimadzu LCsolution Report====

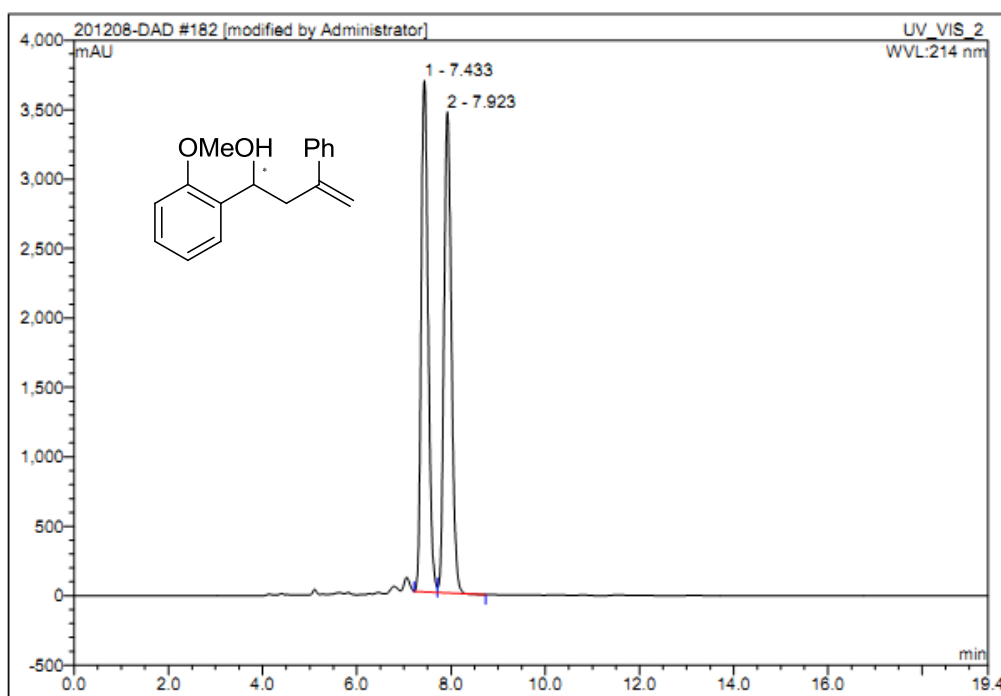


1 PDA 多色谱图 1/220nm 4nm

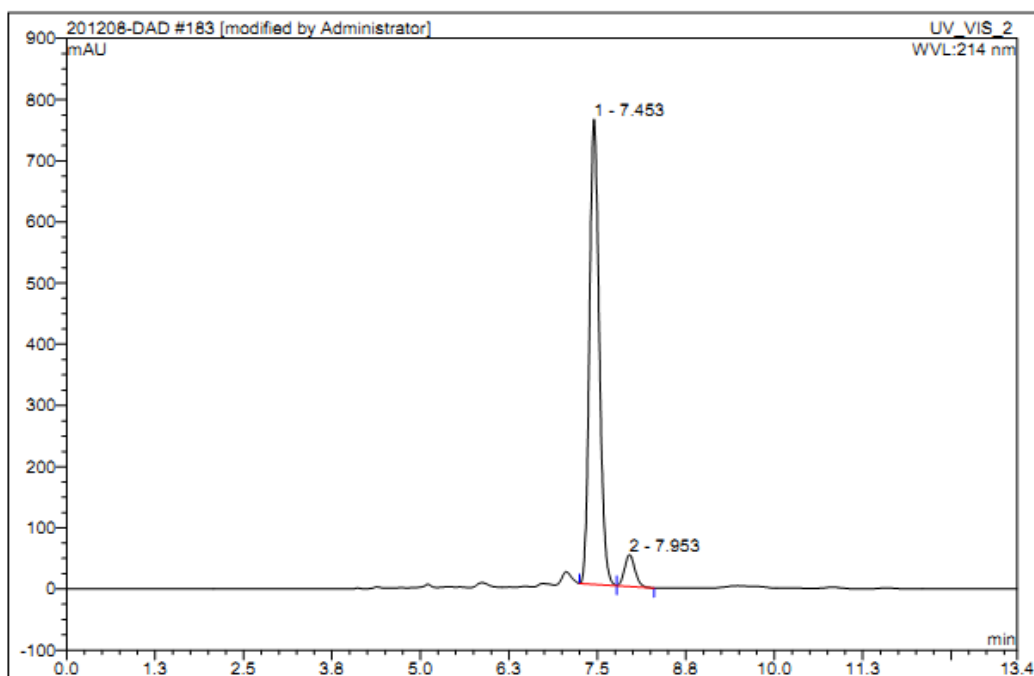
PDA Ch1 220nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	15.395	11346787	527809	16.845	22.613
2	16.816	56012294	1806253	83.155	77.387
总计		67359081	2334062	100.000	100.000

**(+)-1-(2-Methoxyphenyl)-3-phenylbut-3-en-1-ol (6)**



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	7.43	n.a.	3683.258	630.886	49.94	n.a.	BM *
2	7.92	n.a.	3466.734	632.403	50.06	n.a.	MB*
<b>Total:</b>			7149.992	1263.289	100.00	0.000	

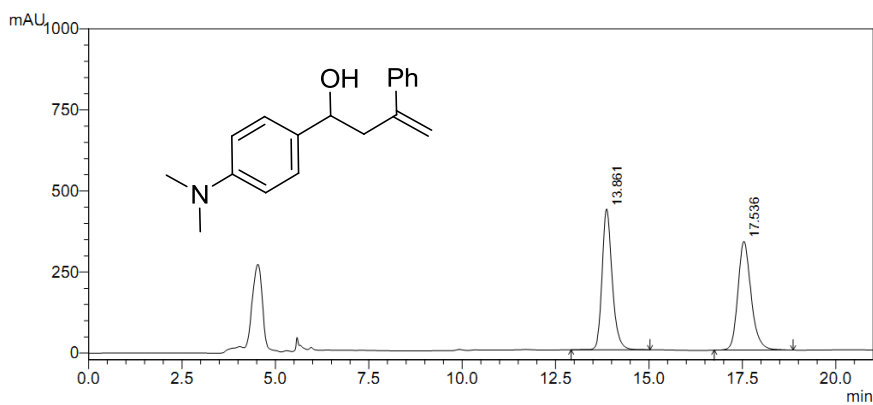


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	7.45	n.a.	760.628	121.461	93.25	n.a.	BM *
2	7.95	n.a.	51.723	8.796	6.75	n.a.	MB*
<b>Total:</b>			812.351	130.256	100.00	0.000	

**(+)-1-(4-(Dimethylamino)phenyl)-3-phenylbut-3-en-1-ol (7)**

2013-8-9 16:15:21 1 / 1

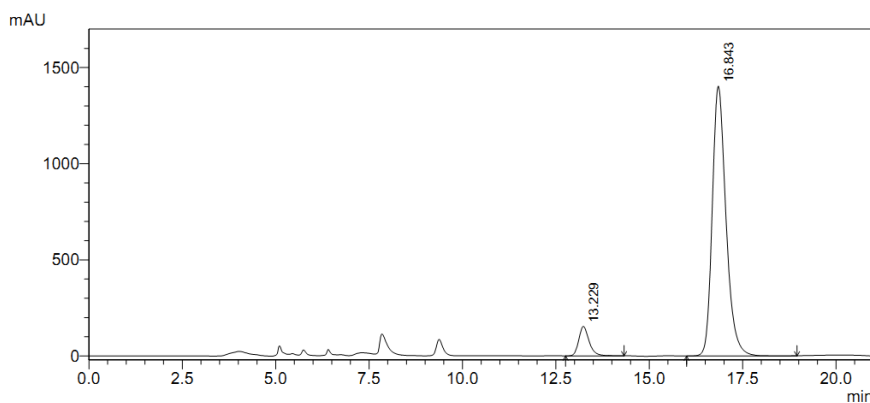
**==== Shimadzu LCsolution Report====**



1 PDA 多色谱图 1/214nm 4nm

**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	13.861	8170530	434380	49.898	56.463
2	17.536	8204019	334941	50.102	43.537
总计		16374549	769321	100.000	100.000



1 PDA 多色谱图 1/214nm 4nm

**PDA Ch1 214nm 4nm**

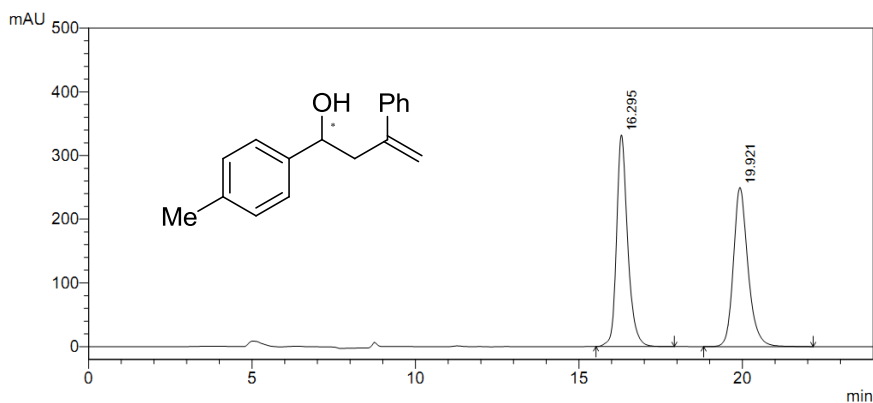
Peak#	Resolution Time	Area	Height	Area %	Height %
1	13.229	2967293	153486	7.631	9.860
2	16.843	35916152	1403177	92.369	90.140
总计		38883445	1556663	100.000	100.000

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(+)-3-Phenyl-1-p-tolylbut-3-en-1-ol (8)

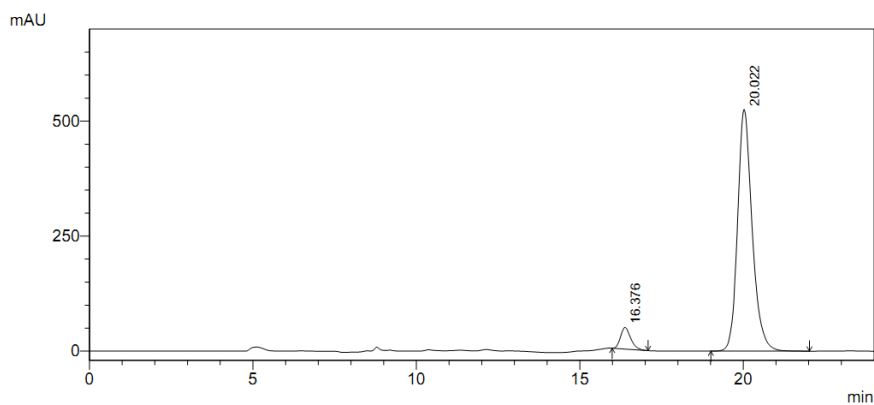
2013-8-9 16:02:12 1 / 1

==== Shimadzu LCsolution Report====



PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	16.295	7832214	332353	50.163	57.080
2	19.921	7781255	249904	49.837	42.920
总计		15613469	582256	100.000	100.000

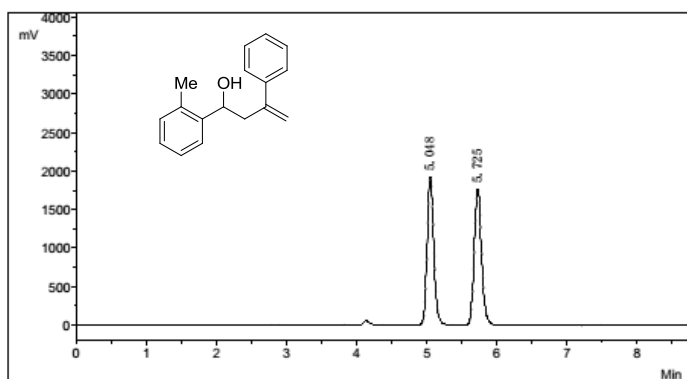


PDA Ch1 214nm 4nm

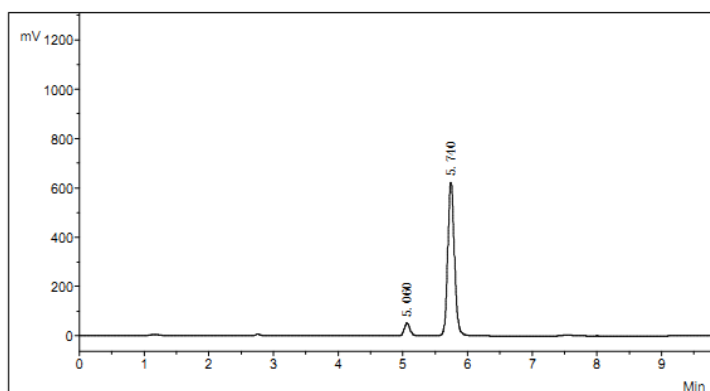
Peak#	Resolution Time	Area	Height	Area %	Height %
1	16.376	1008733	47047	5.822	8.220
2	20.022	16316561	525274	94.178	91.780
总计		17325294	572321	100.000	100.000

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**(-)-3-phenyl-1-o-tolylbut-3-en-1-ol (9)**



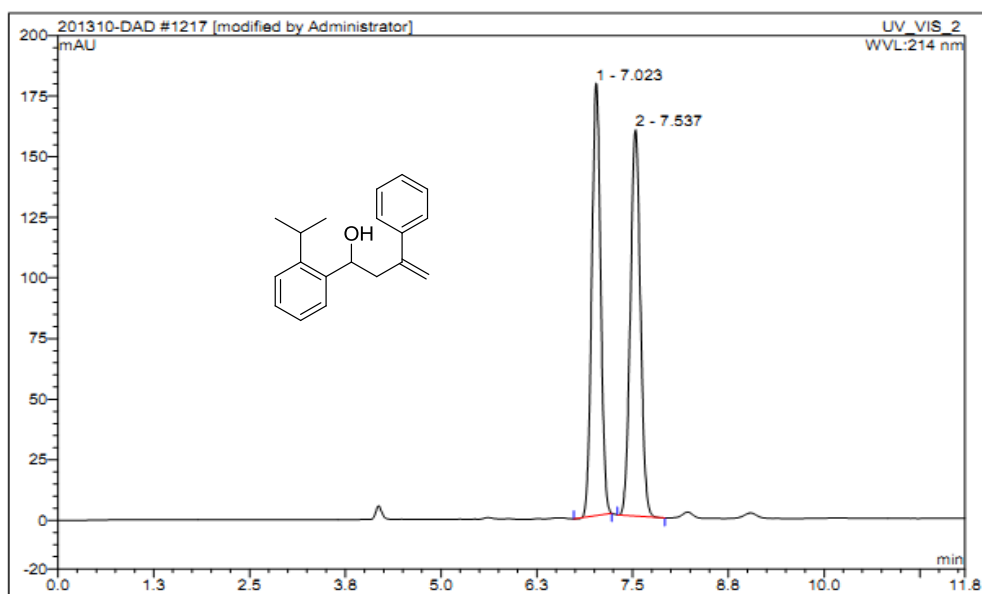
No.	PeakNo	ID. Name	R. Time	PeakHeight	PeakArea	PerCent
1	1		5.048	1923843.3	12669561.1	49.5999
2	2		5.725	1766288.1	12873952.8	50.4001
Total				3690131.4	25543513.9	100.0000



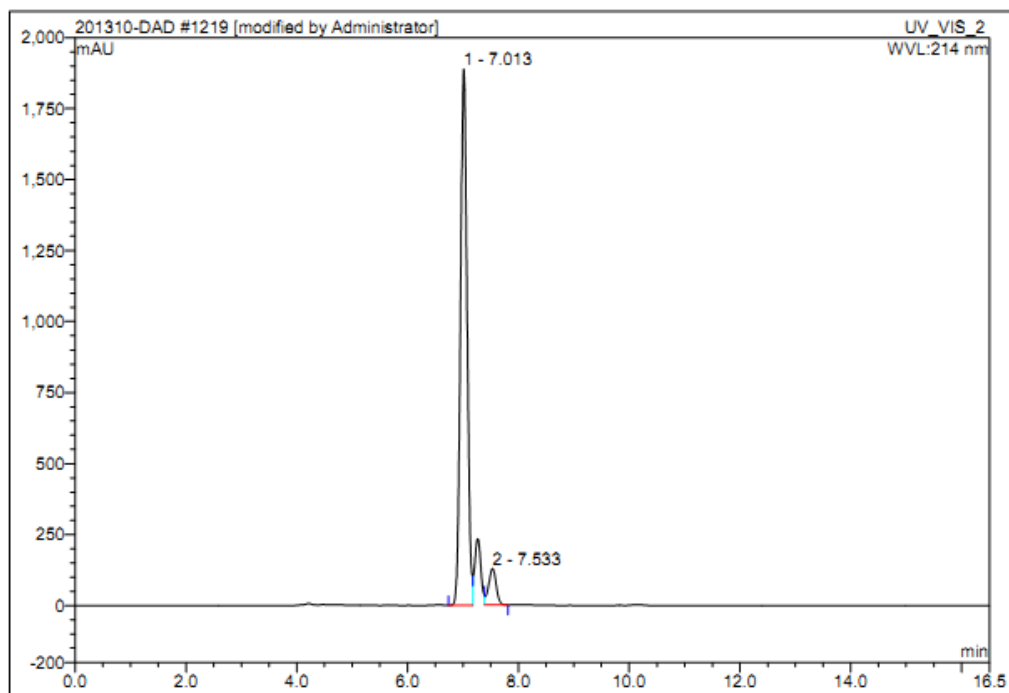
No.	PeakNo	ID. Name	R. Time	PeakHeight	PeakArea	PerCent
1	1		5.060	51758.4	330863.6	6.8633
2	2		5.740	613868.4	4489893.7	93.1367
Total				665626.8	4820757.3	100.0000



**(+)-1-(2-isopropylphenyl)-3-phenylbut-3-en-1-ol (10)**



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	7.02	n.a.	178.394	23.918	49.61	n.a.	BMB*
2	7.54	n.a.	159.435	24.292	50.39	n.a.	BMB*
<b>Total:</b>			337.829	48.210	100.00	0.000	

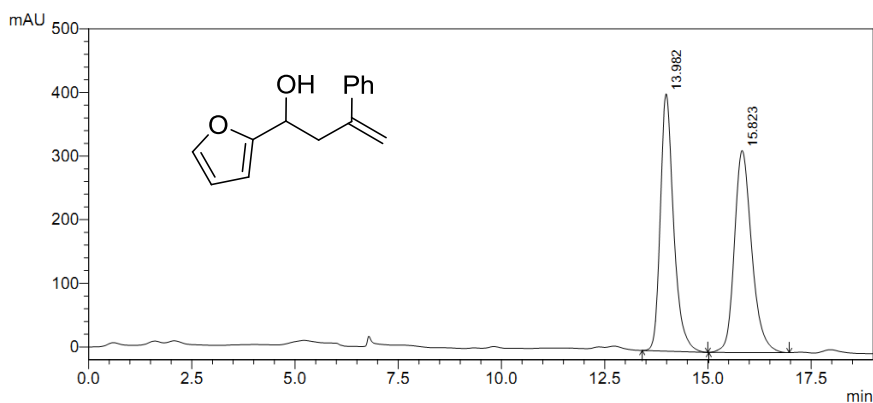


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	7.01	n.a.	1888.505	261.852	92.67	n.a.	BM *
2	7.53	n.a.	128.936	20.727	7.33	n.a.	MB*
<b>Total:</b>			2017.441	282.579	100.00	0.000	

**(+)-1-(Furan-2-yl)-3-phenylbut-3-en-1-ol (11)**

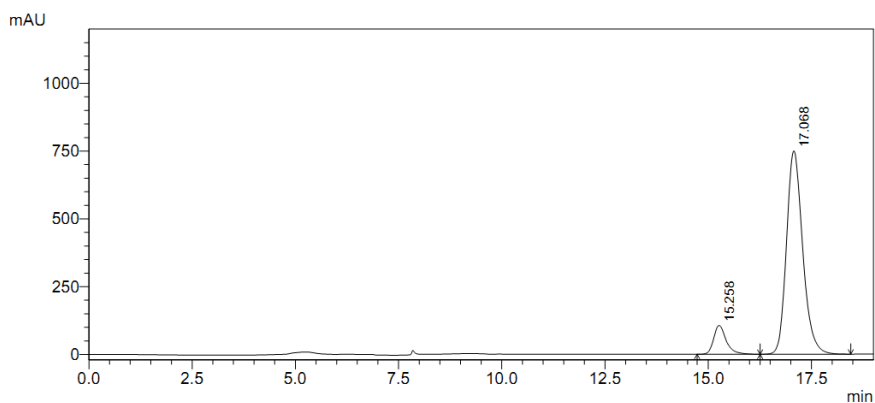
2013-8-9 17:27:49 1 / 1

**==== Shimadzu LCsolution Report====**



**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	13.982	9031417	404626	49.356	56.020
2	15.823	9267121	317658	50.644	43.980
总计		18298537	722283	100.000	100.000



**PDA Ch1 214nm 4nm**

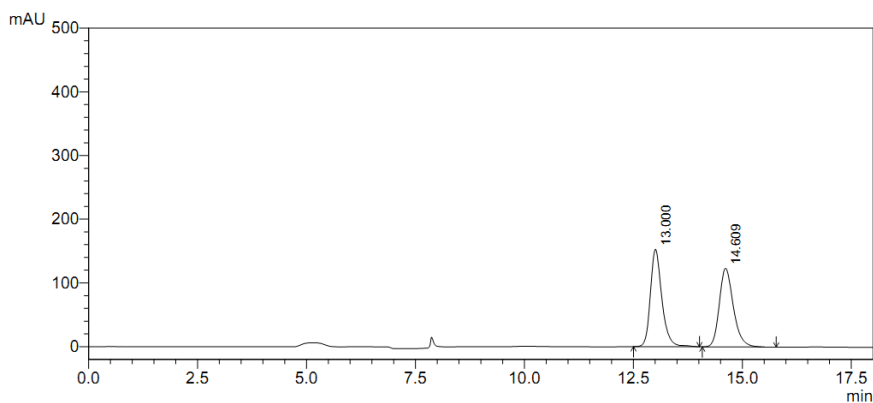
Peak#	Resolution Time	Area	Height	Area %	Height %
1	15.258	2235000	106626	9.841	12.448
2	17.068	20475586	749943	90.159	87.552
总计		22710586	856568	100.000	100.000

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# (S)-(+)-1,3-Diphenylbut-3-en-1-ol (12)

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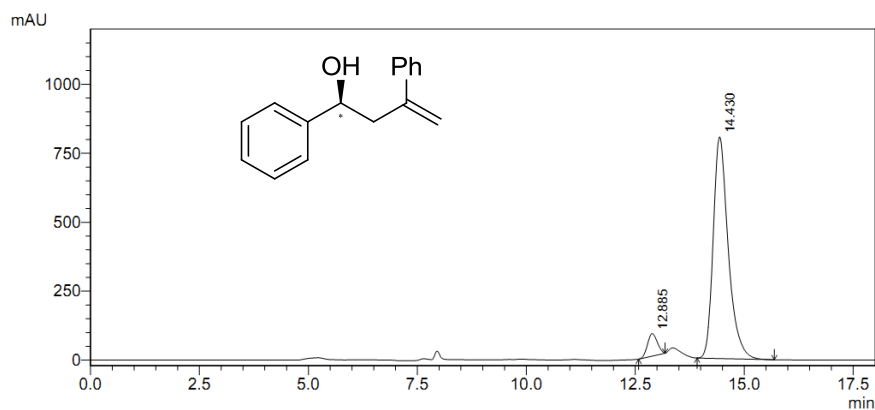
## ==== Shimadzu LCsolution Report====



1 PDA 多色谱图 1/214nm 4nm

### PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	13.000	2796315	152937	50.247	55.322
2	14.609	2768824	123510	49.753	44.678
总计		5565139	276447	100.000	100.000



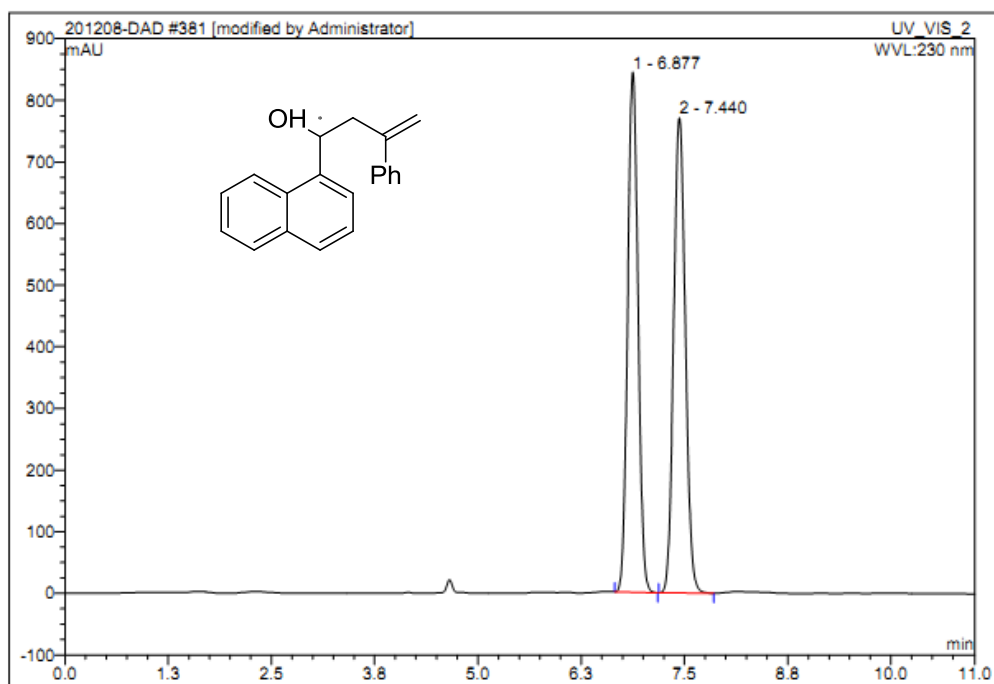
1 PDA 多色谱图 1/214nm 4nm

### PDA Ch1 214nm 4nm

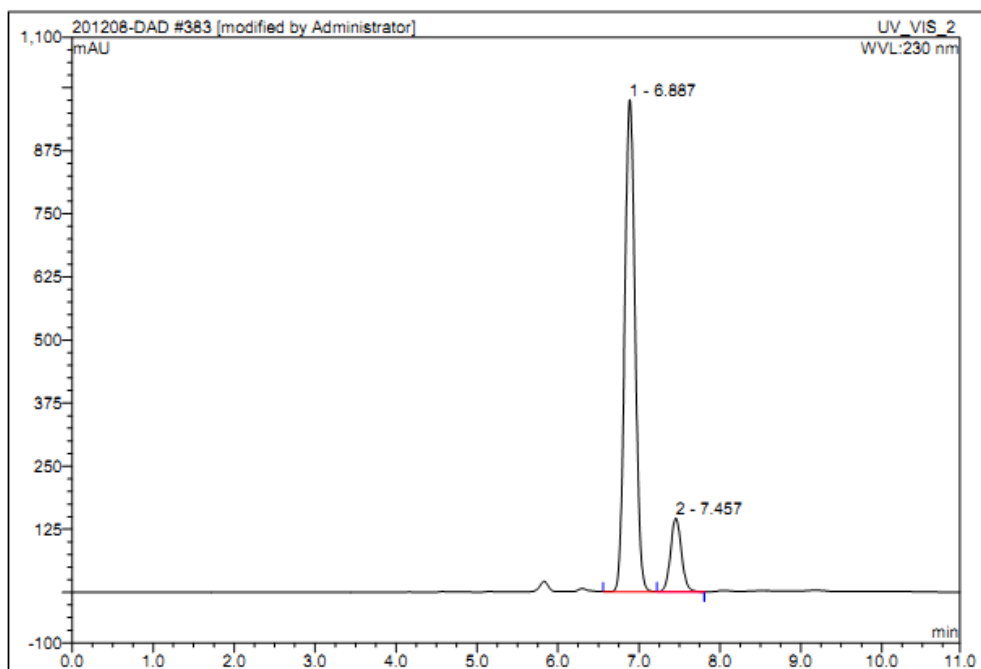
Peak#	Resolution Time	Area	Height	Area %	Height %
1	12.885	1273583	81670	6.213	9.225
2	14.430	19226592	803680	93.787	90.775
总计		20500176	885350	100.000	100.000

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**(-)-1-(Naphthalen-2-yl)-3-phenylbut-3-en-1-ol (13)**



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.88	n.a.	844.134	123.975	49.50	n.a.	BMB
2	7.44	n.a.	770.729	126.476	50.50	n.a.	BMB
<b>Total:</b>			1614.863	250.452	100.00	0.000	

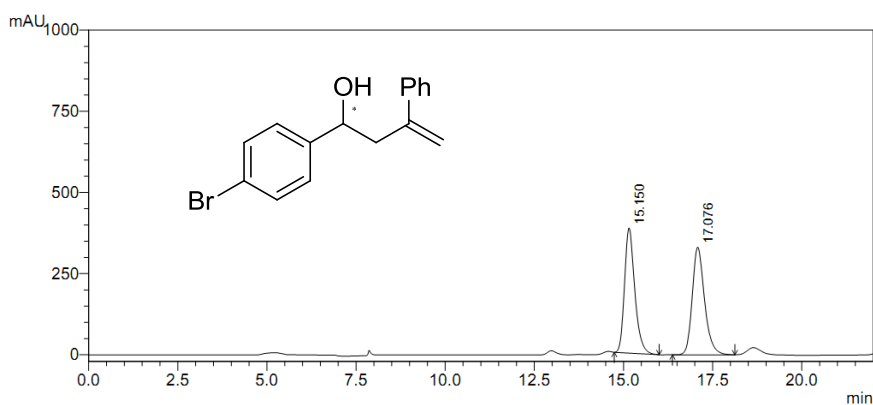


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.89	n.a.	975.580	145.048	86.36	n.a.	BMB
2	7.46	n.a.	145.911	22.915	13.64	n.a.	Rd
<b>Total:</b>			1121.491	167.963	100.00	0.000	

# (+)-1-(4-Bromophenyl)-3-phenylbut-3-en-1-ol (14)

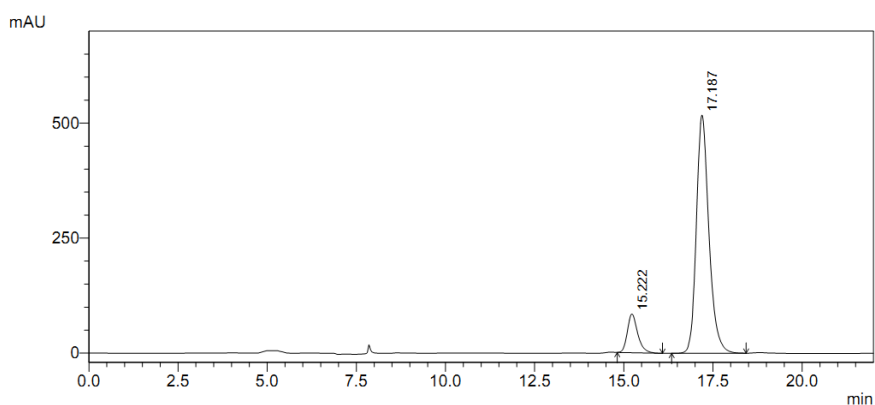
2013-8-9 16:43:40 1 / 1

## ==== Shimadzu LCsolution Report====



PDA Ch1 214nm 4nm

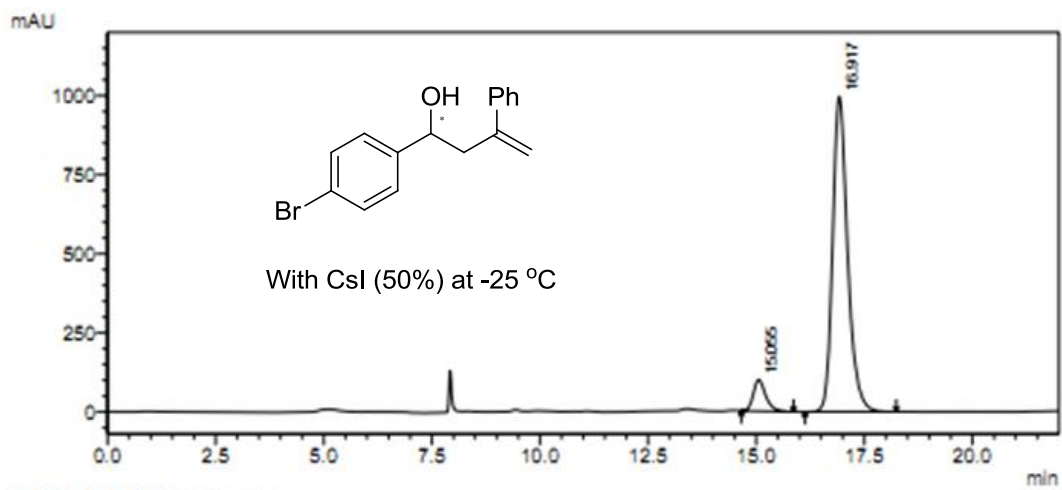
Peak#	Resolution Time	Area	Height	Area %	Height %
1	15.150	7677411	385580	49.041	53.764
2	17.076	7977672	331591	50.959	46.236
总计		15655083	717171	100.000	100.000



PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	15.222	1693613	84257	11.774	13.993
2	17.187	12691050	517877	88.226	86.007
总计		14384663	602134	100.000	100.000

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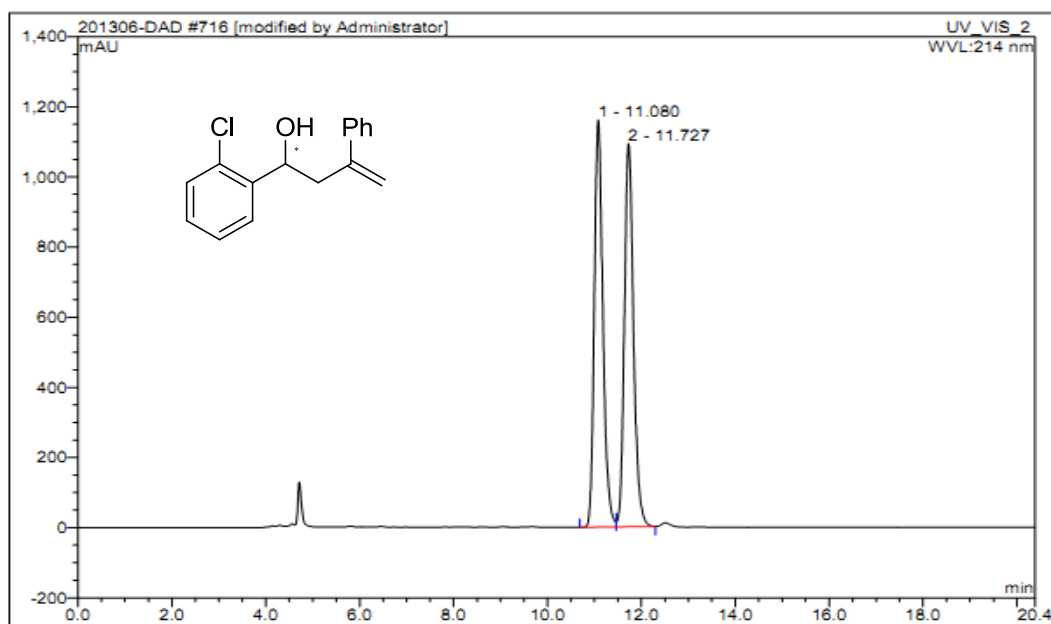


1 PDA 多色谱图 1/214nm 4nm

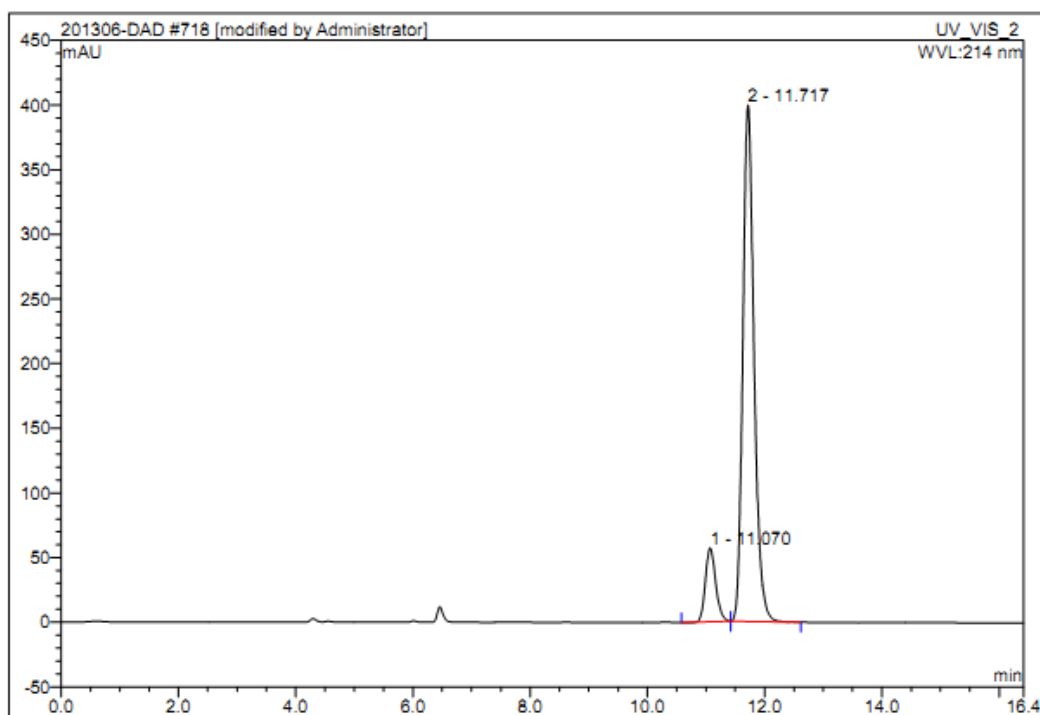
PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	15.055	1957688	99268	7.564	9.054
2	16.917	23923433	997114	92.436	90.946
总计		25881121	1096382	100.000	100.000

**(-)-1-(2-Chlorophenyl)-3-phenylbut-3-en-1-ol (15)**



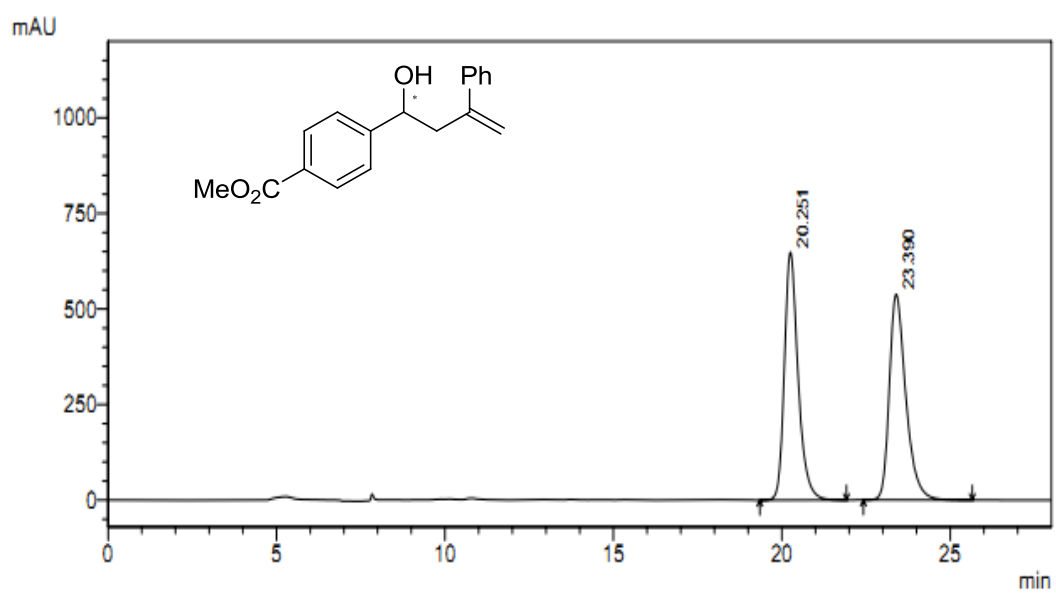
No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	11.08	n.a.	1161.416	246.889	50.09	n.a.	BM *
2	11.73	n.a.	1092.674	245.961	49.91	n.a.	MB*
<b>Total:</b>			2254.090	492.850	100.00	0.000	



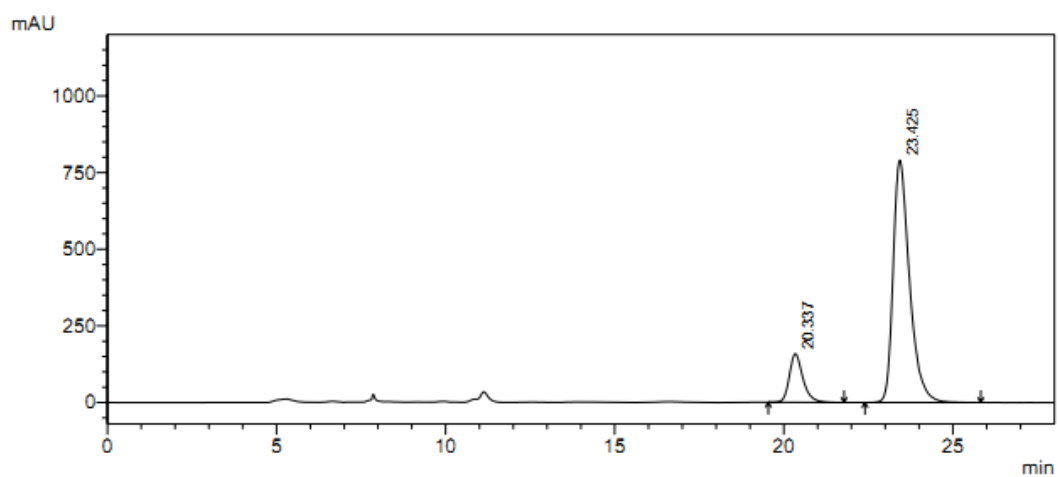
No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	11.07	n.a.	57.322	11.617	11.59	n.a.	BMb*
2	11.72	n.a.	399.575	88.615	88.41	n.a.	bMB*
<b>Total:</b>			456.897	100.231	100.00	0.000	

**(+)-Methyl-4-(1-hydroxy-3-phenylbut-3-enyl)benzoate (16)**

**==== Shimadzu LCsolution Report====**







1 PDA 多色谱图 1/214nm 4nm

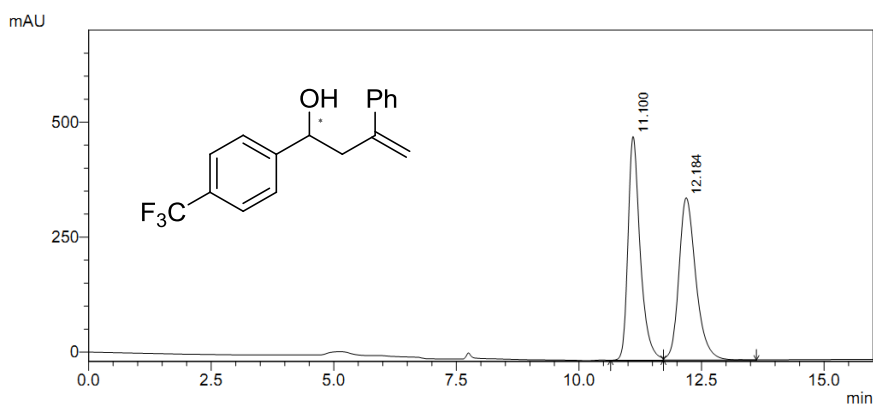
**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	20.337	4407872	158499	14.196	16.703
2	23.425	26642303	790406	85.804	83.297
总计		31050176	948904	100.000	100.000

# (+)-3-Phenyl-1-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (17)

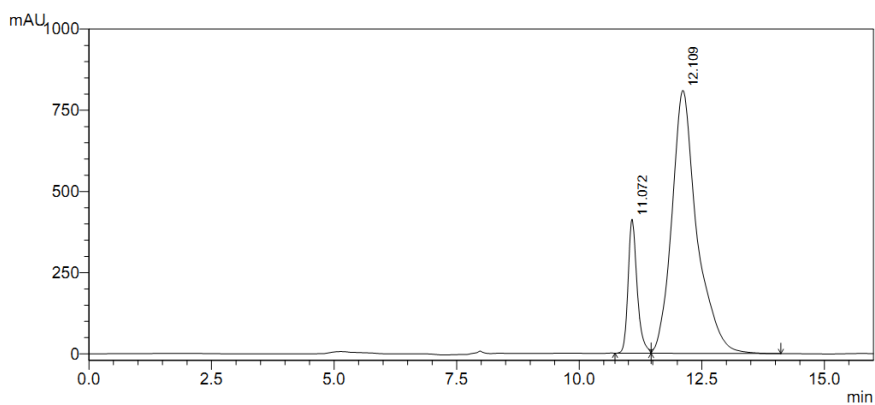
2013-8-9 17:31:27 1 / 1

## ==== Shimadzu LCsolution Report====



PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	11.100	8395465	486622	49.717	57.956
2	12.184	8491182	353023	50.283	42.044
总计		16886646	839646	100.000	100.000



PDA Ch1 214nm 4nm

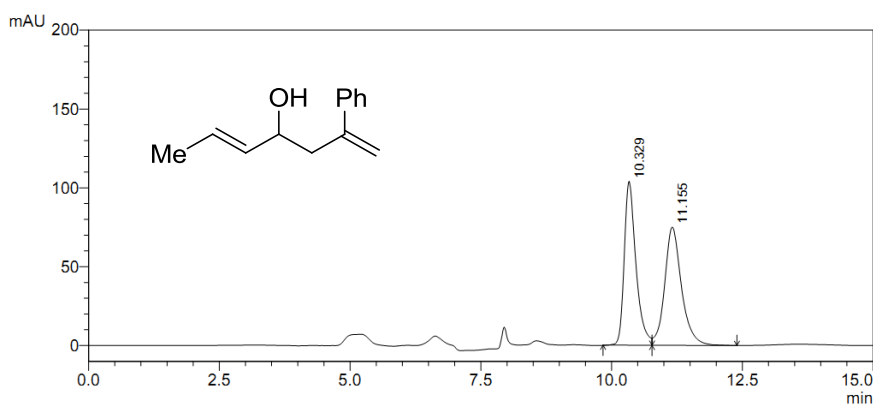
Peak#	Resolution Time	Area	Height	Area %	Height %
1	11.072	5275339	412584	15.495	33.753
2	12.109	28770701	809790	84.505	66.247
总计		34046039	1222374	100.000	100.000

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**(+)-(E)-2-Phenylhepta-1,5-dien-4-ol (18)**

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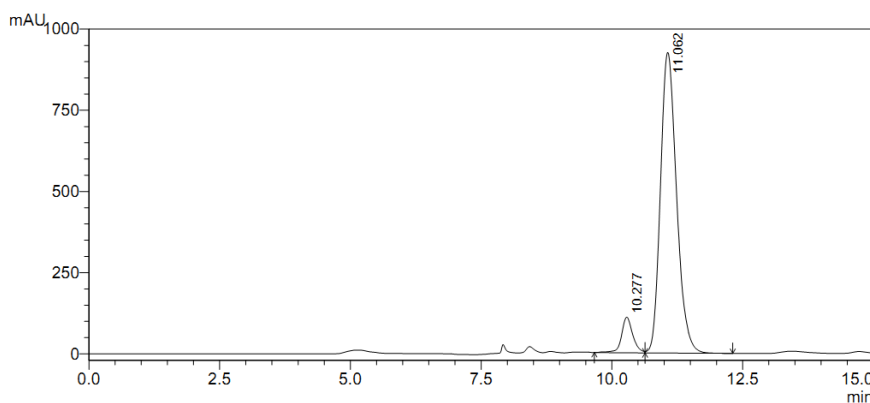
**==== Shimadzu LCsolution Report====**



1 PDA 多色谱图 1/214nm 4nm

**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	10.329	1623358	104004	49.260	58.117
2	11.155	1672129	74951	50.740	41.883
总计		3295487	178956	100.000	100.000



1 PDA 多色谱图 1/214nm 4nm

**PDA Ch1 214nm 4nm**

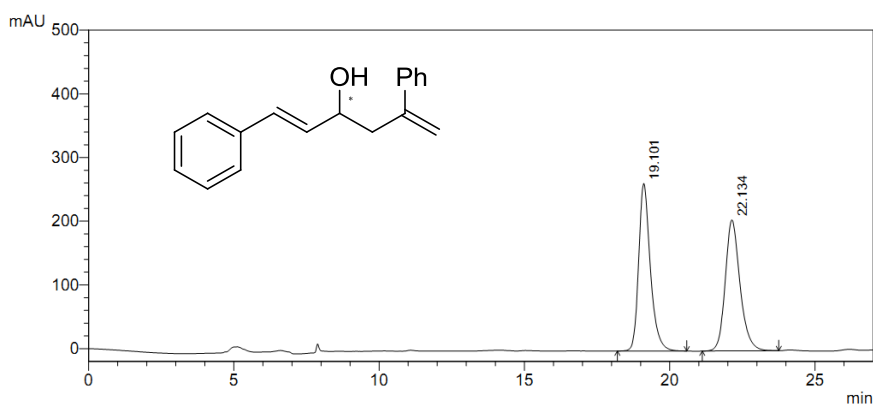
Peak#	Resolution Time	Area	Height	Area %	Height %
1	10.277	1641517	109735	7.660	10.598
2	11.062	19788754	925731	92.340	89.402
总计		21430271	1035466	100.000	100.000

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(+)-(E)-1,5-Diphenylhexa-1,5-dien-3-ol (19)

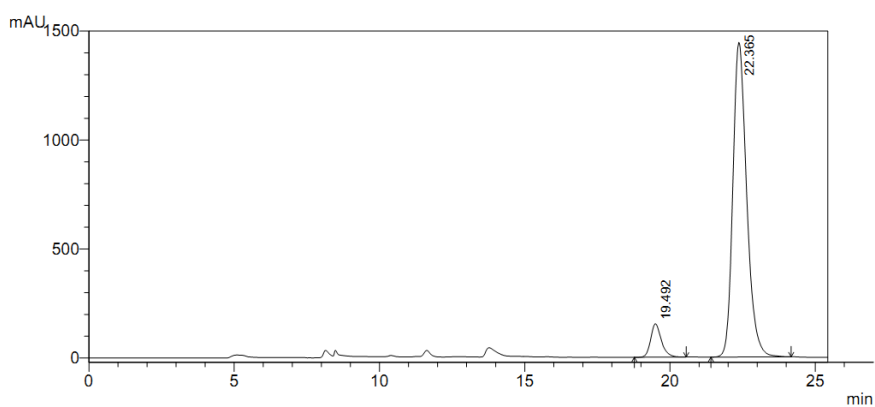
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==== Shimadzu LCsolution Report====



PDA Ch1 214nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	19.101	7339174	263096	50.275	56.142
2	22.134	7258777	205527	49.725	43.858
总计		14597950	468623	100.000	100.000

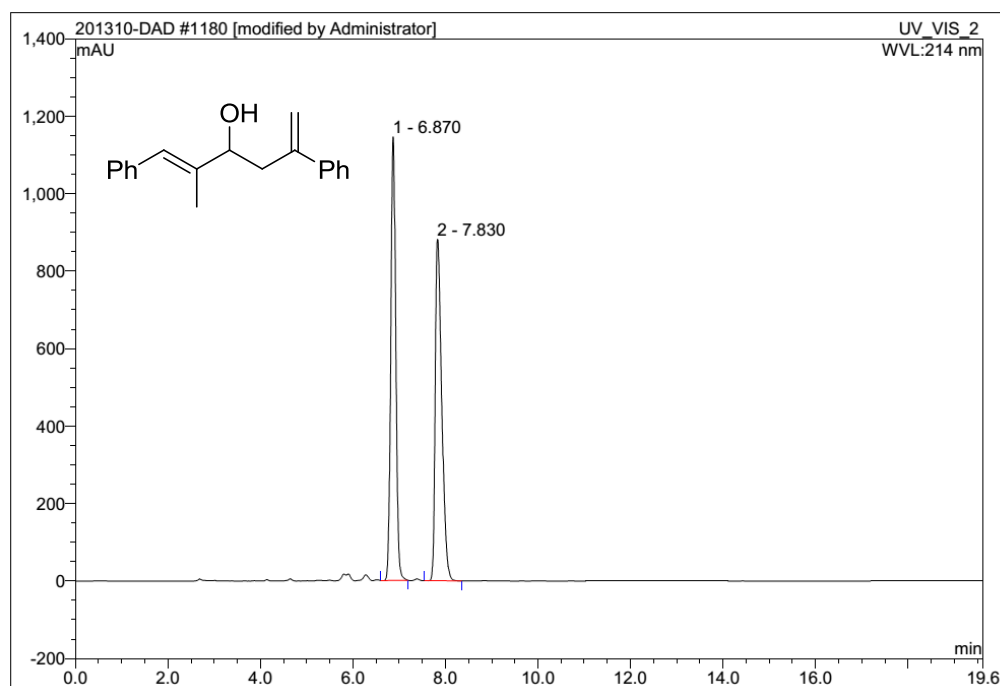


PDA Ch1 214nm 4nm

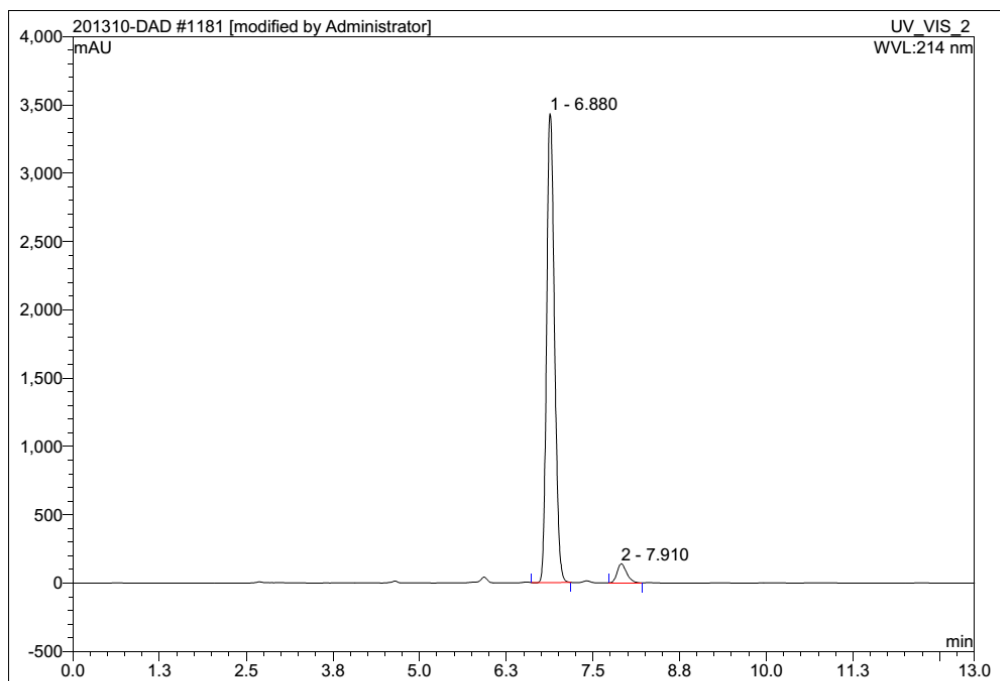
Peak#	Resolution Time	Area	Height	Area %	Height %
1	19.492	3910126	153269	7.570	9.593
2	22.365	47743356	1444413	92.430	90.407
总计		51653482	1597682	100.000	100.000

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**(+)-(E)-2-methyl-1,5-diphenylhexa-1,5-dien-3-ol (20)**



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.87	n.a.	1145.460	144.246	50.04	n.a.	BMB*
2	7.83	n.a.	881.695	144.028	49.96	n.a.	BMB*
<b>Total:</b>			2027.155	288.274	100.00	0.000	

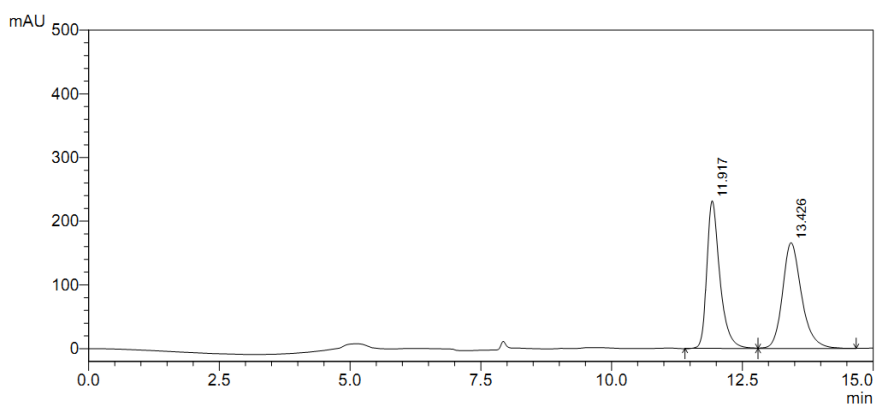


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.88	n.a.	3434.469	452.596	95.25	n.a.	BMB*
2	7.91	n.a.	140.001	22.551	4.75	n.a.	BMB*
<b>Total:</b>			3574.470	475.146	100.00	0.000	

**(+)-1,5-Diphenylhex-5-en-3-ol (21)**

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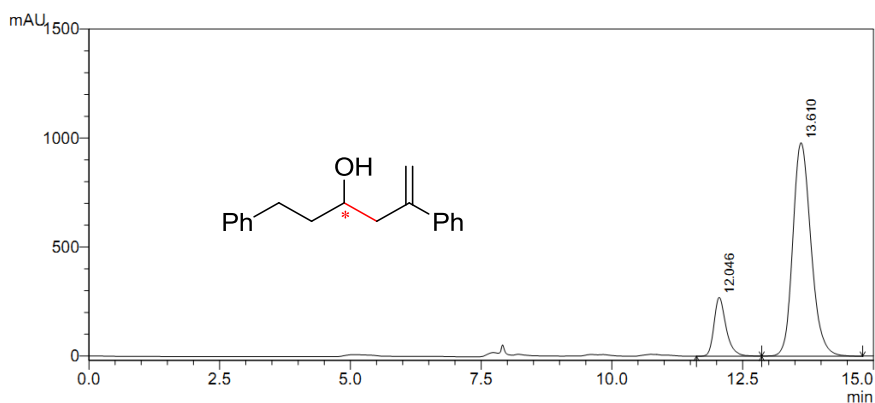
**==== Shimadzu LCsolution Report====**



1 PDA 多色谱图 1/214nm 4nm

**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	11.917	4001984	231795	48.802	58.280
2	13.426	4198449	165928	51.198	41.720
总计		8200433	397723	100.000	100.000



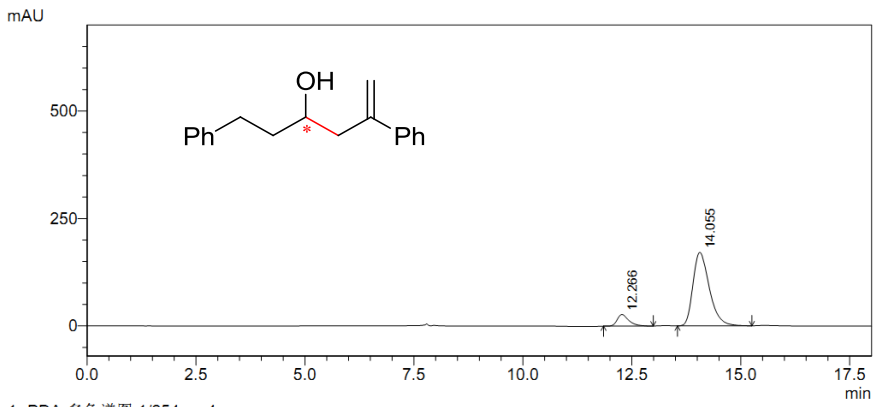
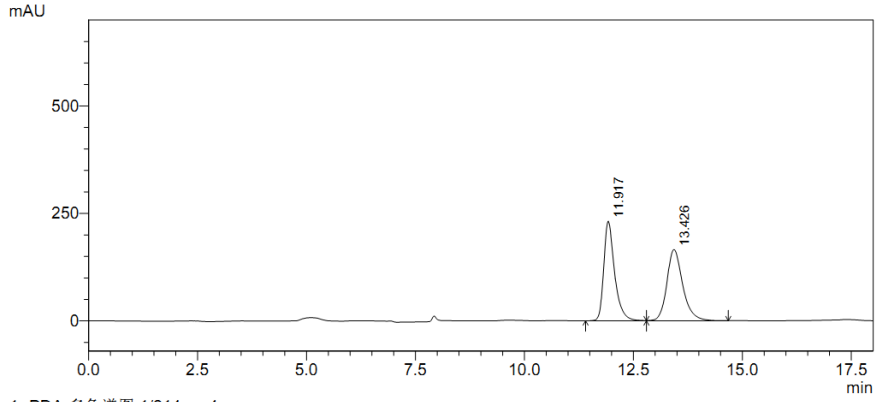
1 PDA 多色谱图 1/214nm 4nm

**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	12.046	4338400	270258	15.504	21.624
2	13.610	23643659	979539	84.496	78.376
总计		27982059	1249797	100.000	100.000

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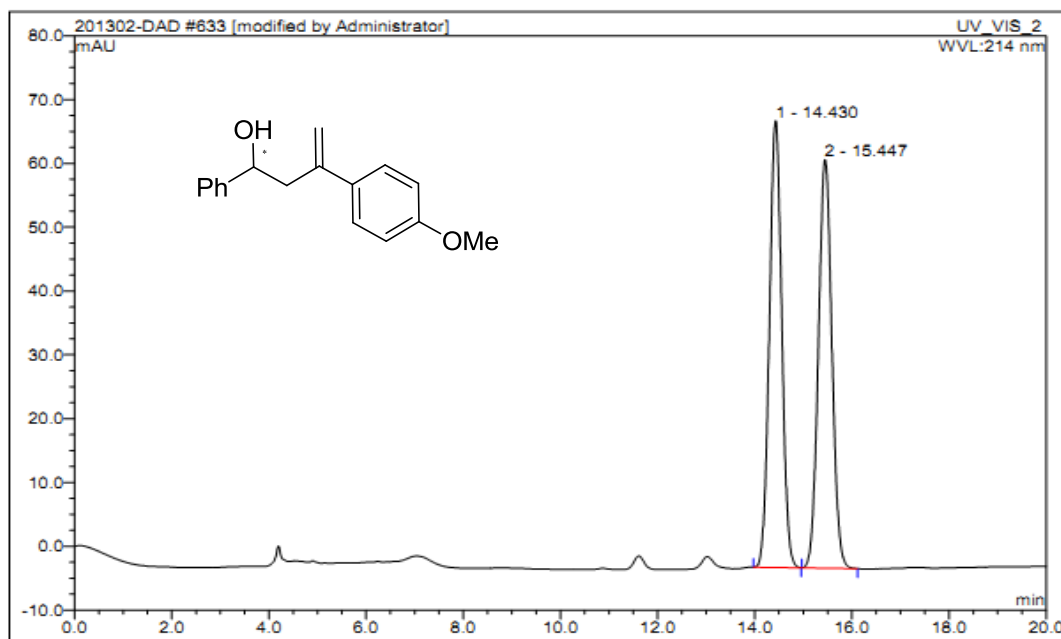
==== Shimadzu LCsolution Report====



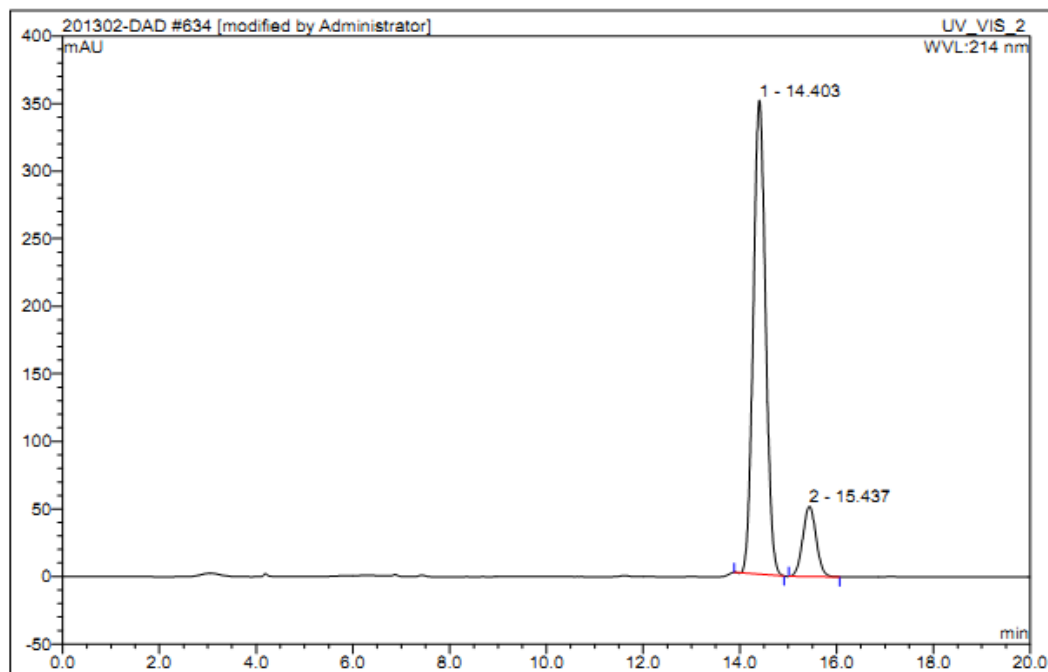
PDA Ch1 254nm 4nm

Peak#	Resolution Time	Area	Height	Area %	Height %
1	12.266	491310	26982	9.897	13.626
2	14.055	4472812	171032	90.103	86.374
总计		4964122	198015	100.000	100.000

**(+)-3-(4-Methoxyphenyl)-1-phenylbut-3-en-1-ol (22)**

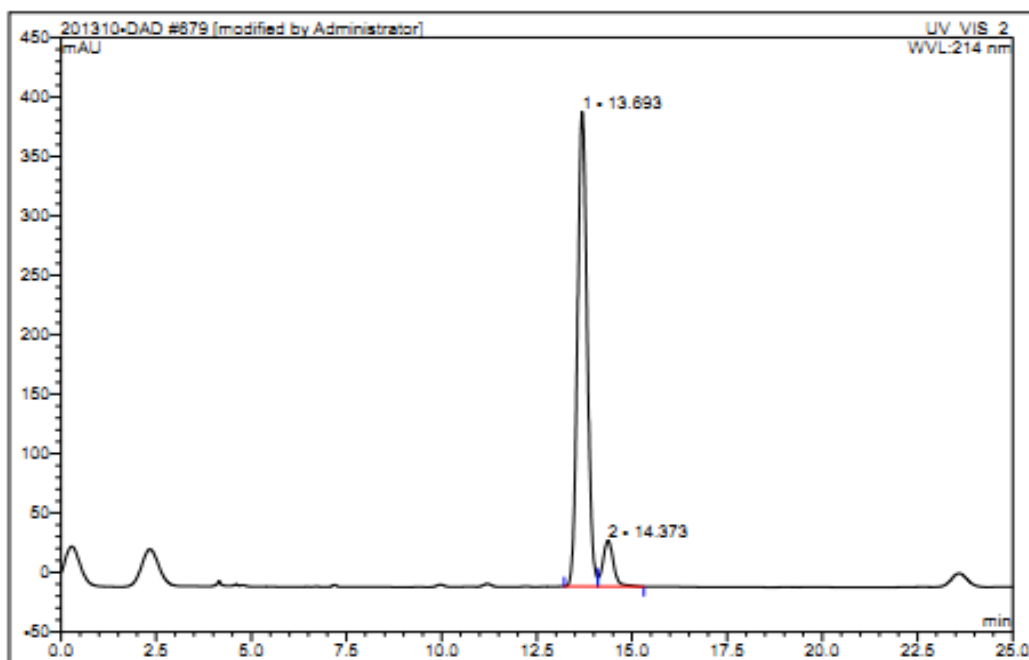


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	14.43	n.a.	70.053	20.640	50.02	n.a.	BMB
2	15.45	n.a.	64.062	20.619	49.98	n.a.	BMB
<b>Total:</b>			134.115	41.259	100.00	0.000	



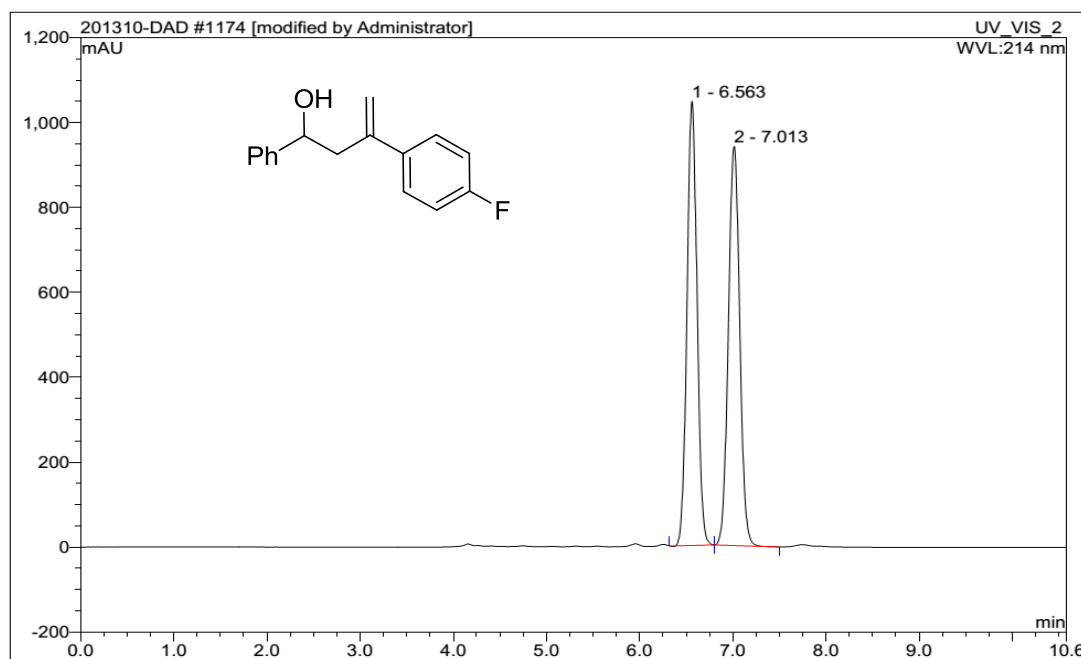
No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	14.40	n.a.	350.868	103.261	86.09	n.a.	BMB*
2	15.44	n.a.	52.123	16.686	13.91	n.a.	BMB*
<b>Total:</b>			402.991	119.947	100.00	0.000	



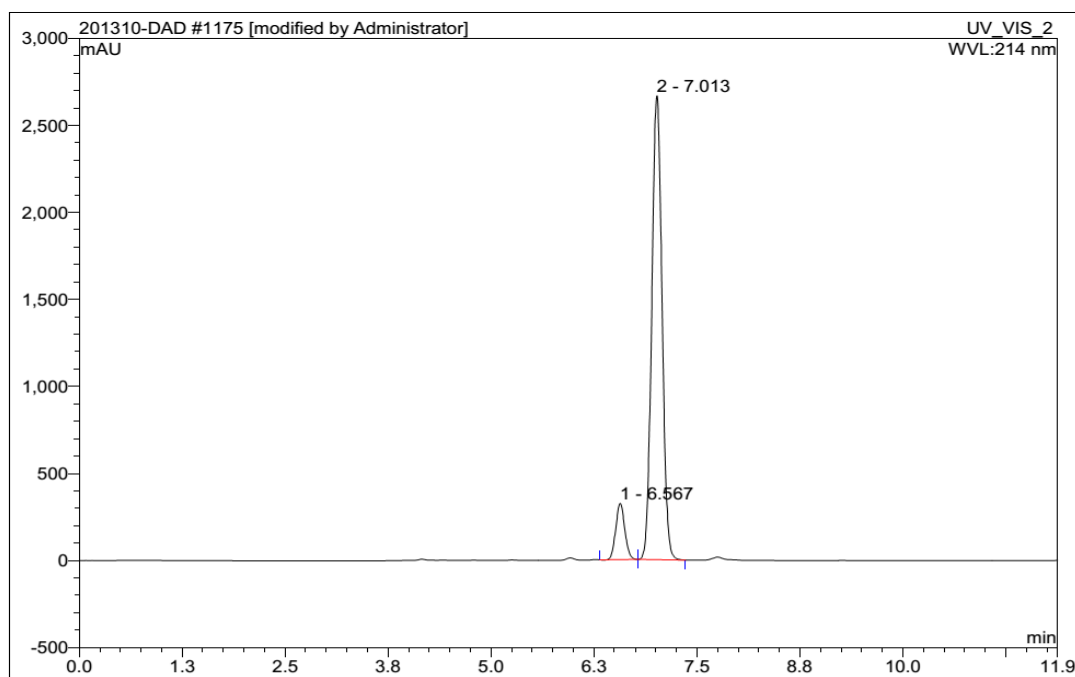


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	13.69	n.a.	399.649	117.478	90.35	n.a.	BM
2	14.37	n.a.	38.810	12.554	9.65	n.a.	MB
<b>Total:</b>			<b>438.459</b>	<b>130.032</b>	<b>100.00</b>	<b>0.000</b>	

**(+)-3-(4-fluorophenyl)-1-phenylbut-3-en-1-ol (23)**

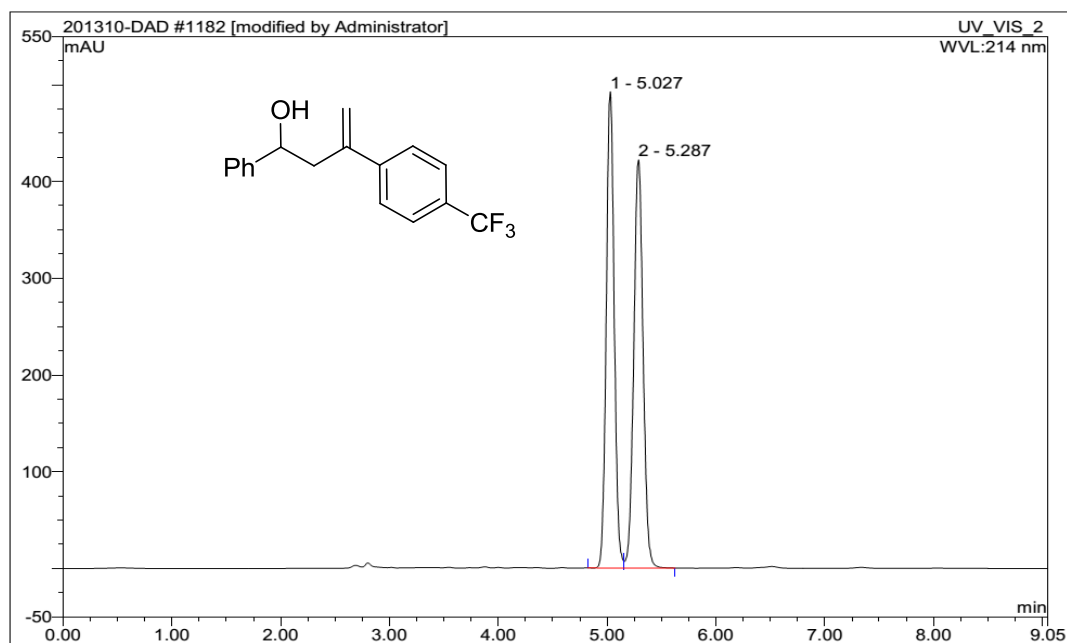


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.56	n.a.	1045.955	131.028	49.67	n.a.	BMb*
2	7.01	n.a.	940.762	132.752	50.33	n.a.	bMB*
<b>Total:</b>			<b>1986.717</b>	<b>263.780</b>	<b>100.00</b>	<b>0.000</b>	

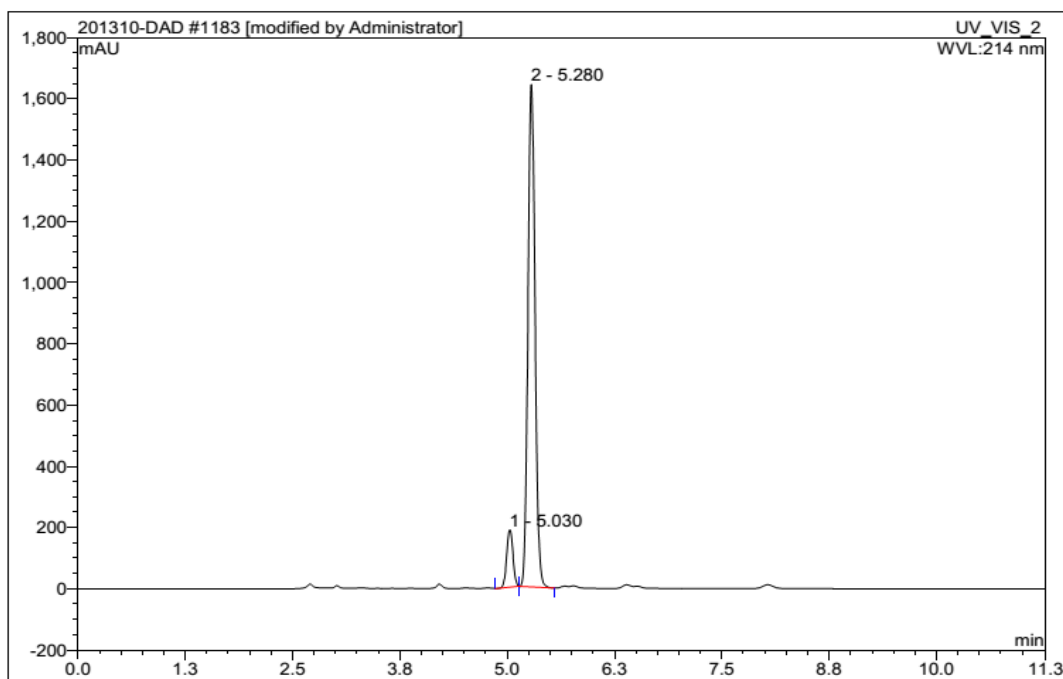


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.57	n.a.	324.151	41.182	9.65	n.a.	BMb*
2	7.01	n.a.	2666.026	385.444	90.35	n.a.	bMB*
<b>Total:</b>			2990.177	426.626	100.00	0.000	

**(+)-1-phenyl-3-(4-(trifluoromethyl)phenyl)but-3-en-1-ol (24)**

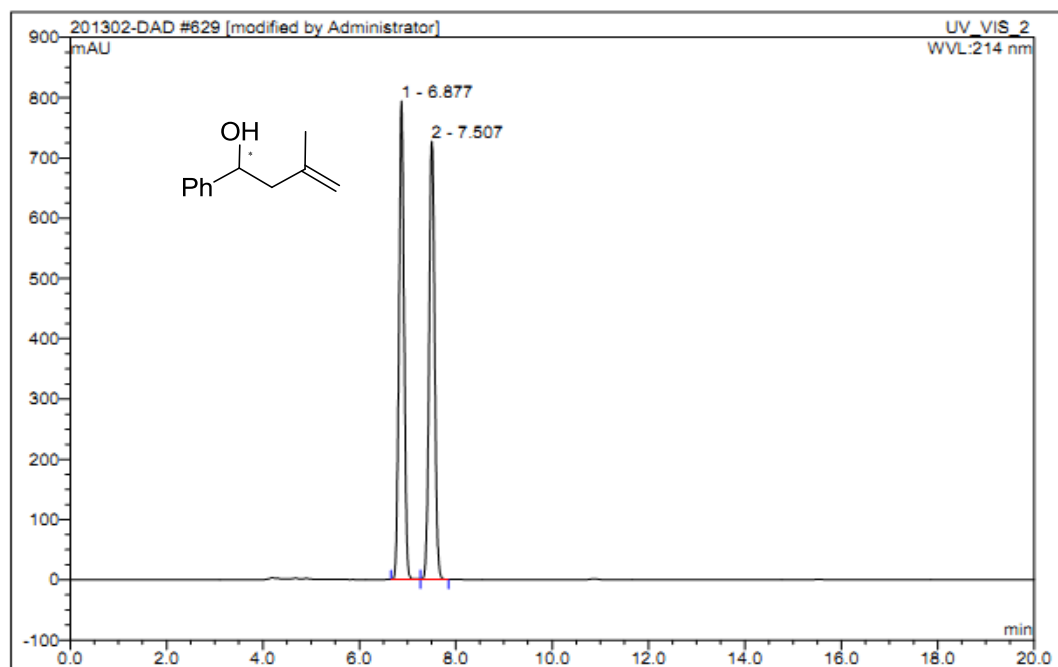


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	5.03	n.a.	492.522	41.575	49.71	n.a.	BM *
2	5.29	n.a.	422.321	42.053	50.29	n.a.	MB*
<b>Total:</b>			914.844	83.628	100.00	0.000	

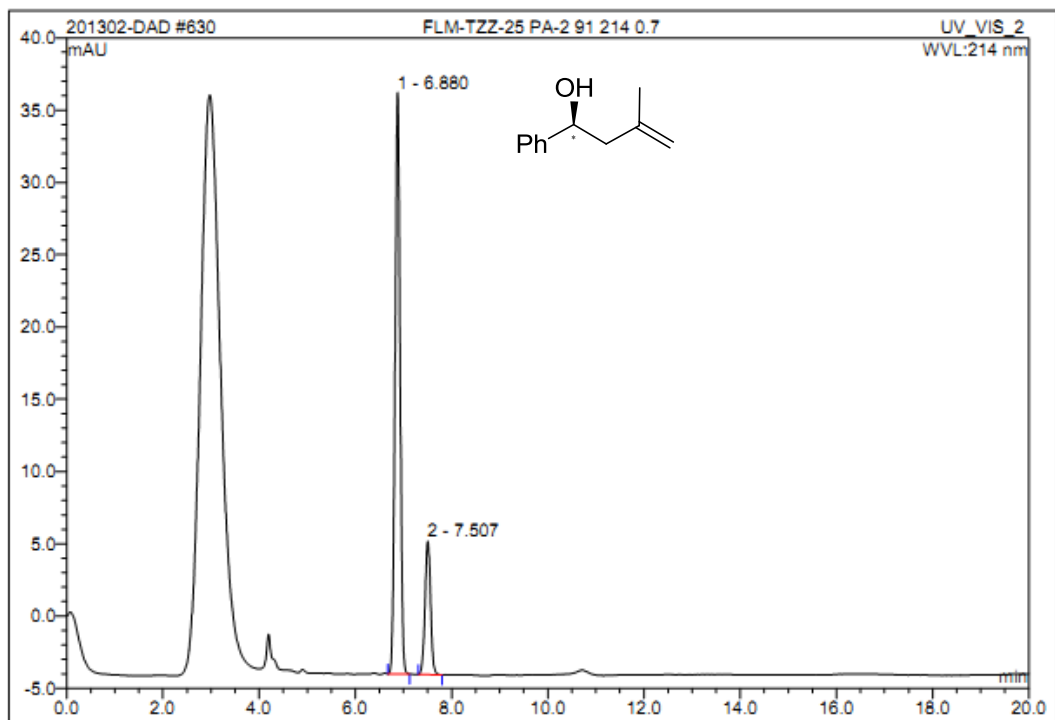


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	5.03	n.a.	186.836	15.710	9.08	n.a.	BMb*
2	5.28	n.a.	1641.379	157.310	90.92	n.a.	bMB*
<b>Total:</b>			1828.215	173.020	100.00	0.000	

**(S)-3-Methyl-1-phenylbut-3-en-1-ol (25)**

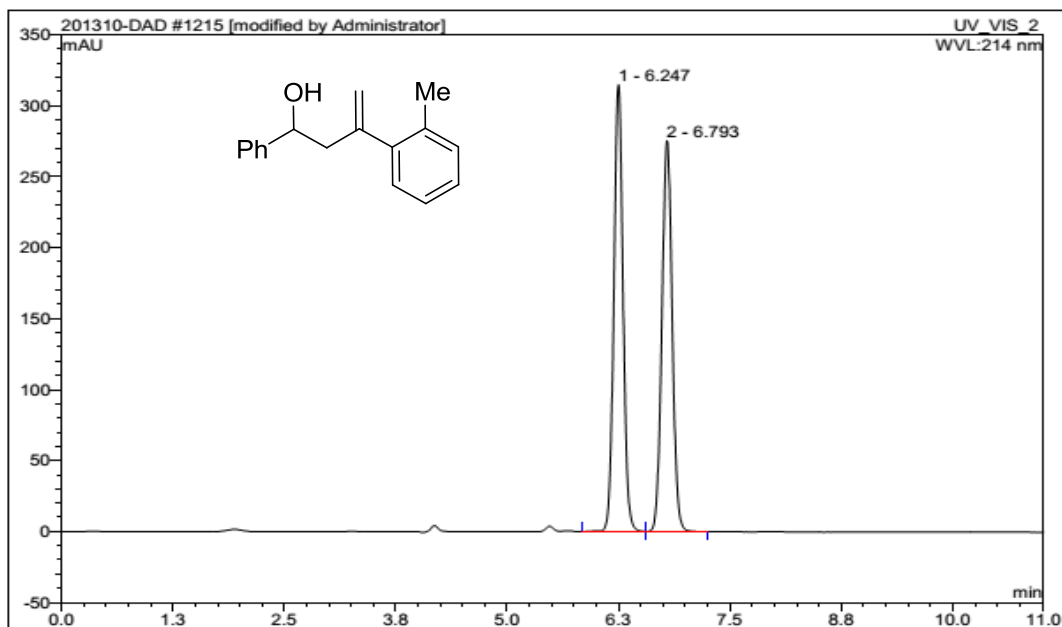


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.88	n.a.	793.794	97.261	49.95	n.a.	BMb*
2	7.51	n.a.	726.983	97.445	50.05	n.a.	bMB
<b>Total:</b>			1520.777	194.706	100.00	0.000	

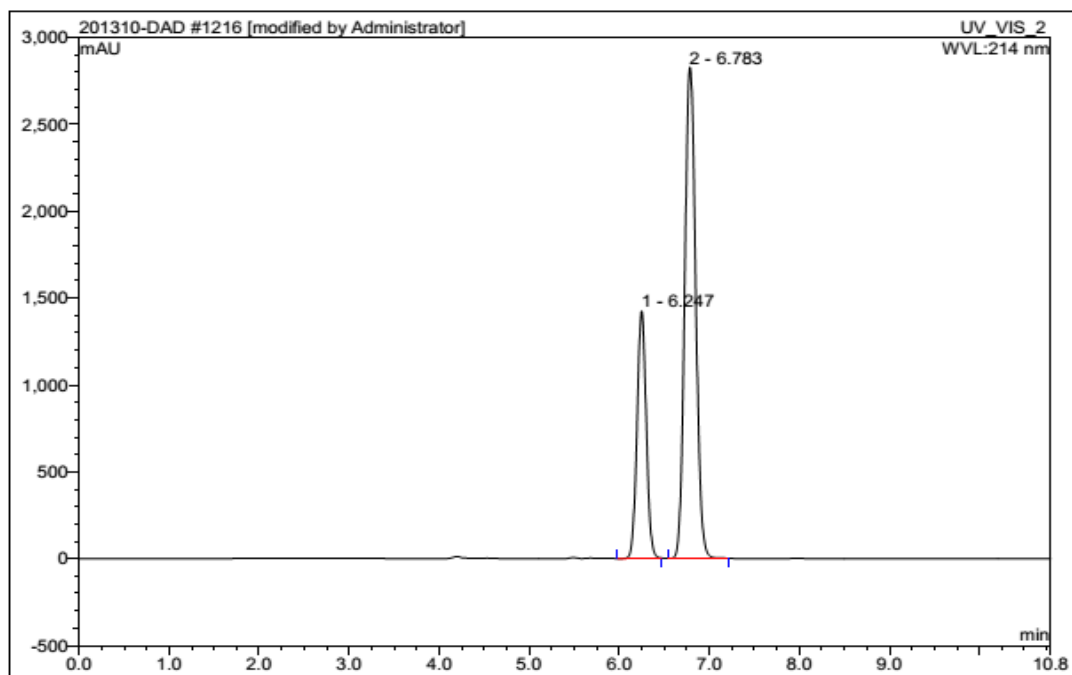


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.88	n.a.	40.260	4.890	79.75	n.a.	BMB
2	7.51	n.a.	9.242	1.242	20.25	n.a.	BMB
<b>Total:</b>			49.502	6.131	100.00	0.000	

**(+)-1-phenyl-3-o-tolylbut-3-en-1-ol (26)**

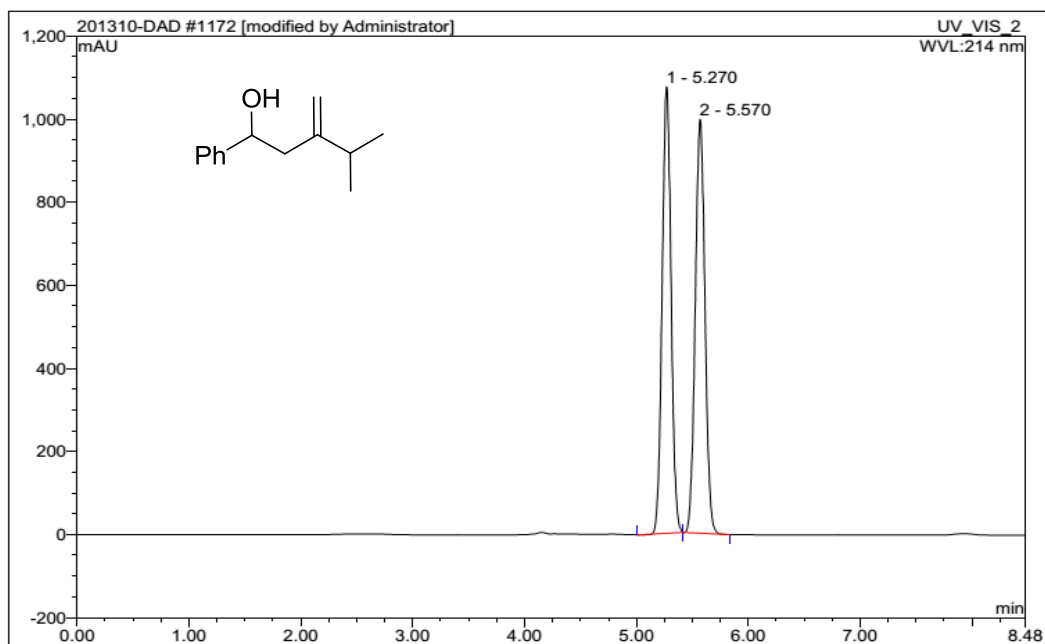


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.25	n.a.	314.888	37.695	50.60	n.a.	BMB*
2	6.79	n.a.	275.564	36.801	49.40	n.a.	bMB*
<b>Total:</b>			590.452	74.496	100.00	0.000	

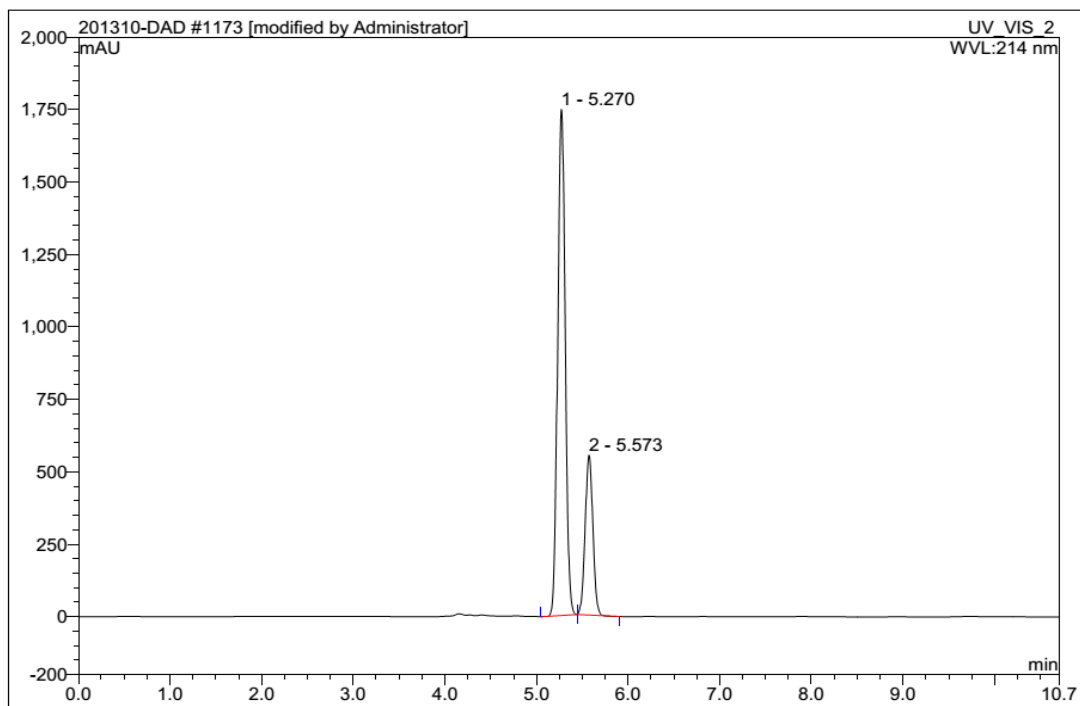


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	6.25	n.a.	1423.438	169.708	29.65	n.a.	BMB*
2	6.78	n.a.	2827.475	402.592	70.35	n.a.	BMB*
<b>Total:</b>			4250.913	572.300	100.00	0.000	

**(-)-4-methyl-3-methylene-1-phenylpentan-1-ol (27)**

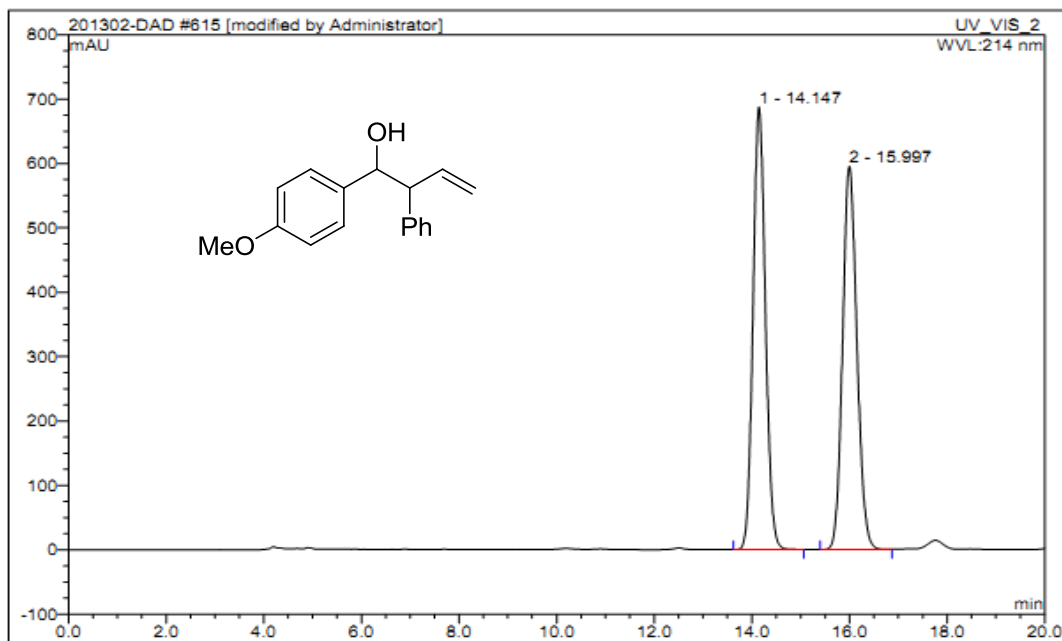


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	5.27	n.a.	1075.557	99.499	49.92	n.a.	BMB*
2	5.57	n.a.	997.549	99.823	50.08	n.a.	bMB*
<b>Total:</b>			2073.106	199.322	100.00	0.000	

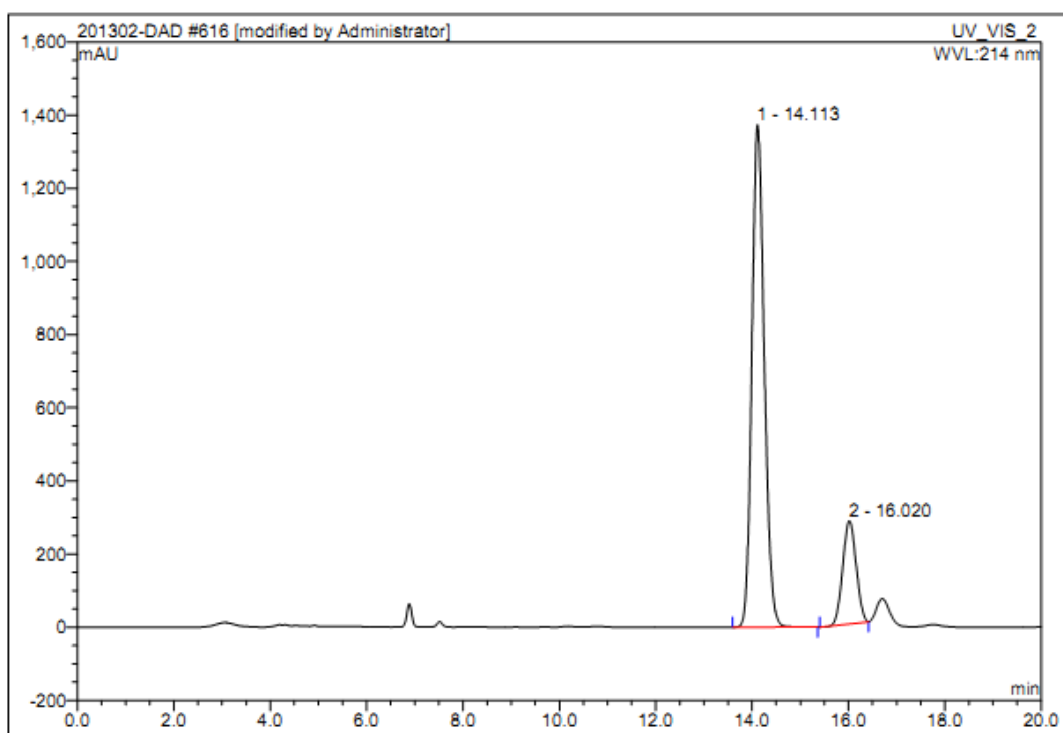


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	5.27	n.a.	1747.876	166.810	75.39	n.a.	BMB*
2	5.57	n.a.	552.778	54.457	24.61	n.a.	bMB*
<b>Total:</b>			2300.654	221.267	100.00	0.000	

***anti*-1-(4-Methoxyphenyl)-2-phenylbut-3-en-1-ol (28)**

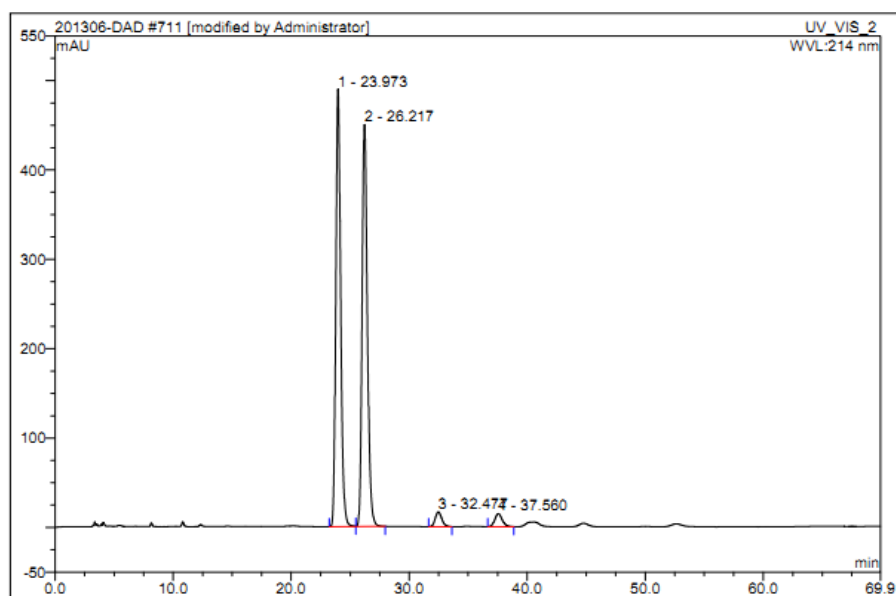


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	14.15	n.a.	687.538	202.371	50.03	n.a.	BMB
2	16.00	n.a.	595.411	202.159	49.97	n.a.	BMB*
<b>Total:</b>			1282.949	404.530	100.00	0.000	

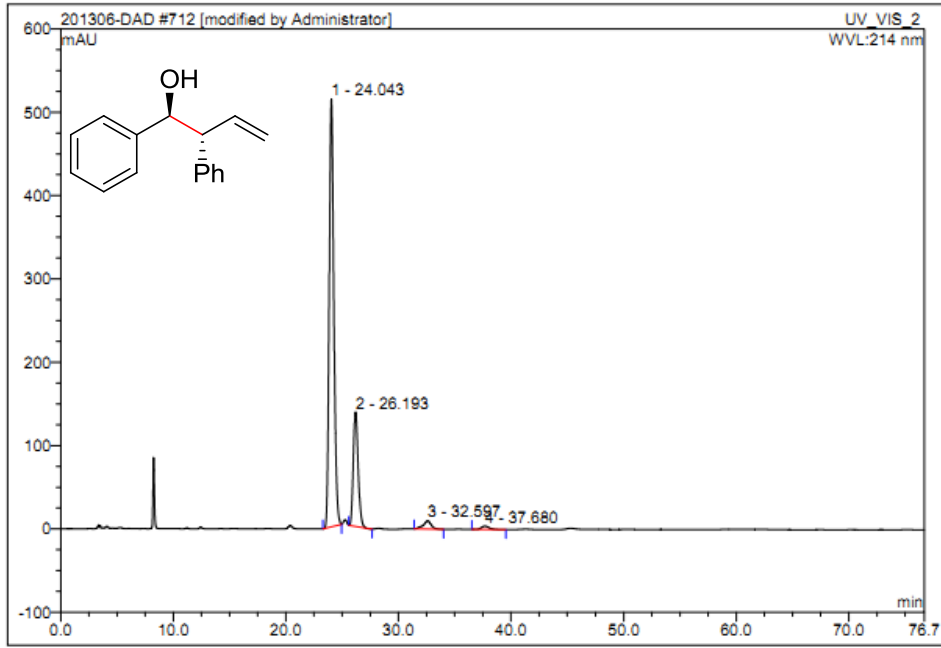


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	14.11	n.a.	1374.186	408.941	81.85	n.a.	BMB
2	16.02	n.a.	282.014	90.693	18.15	n.a.	BMB*
<b>Total:</b>			1656.200	499.635	100.00	0.000	

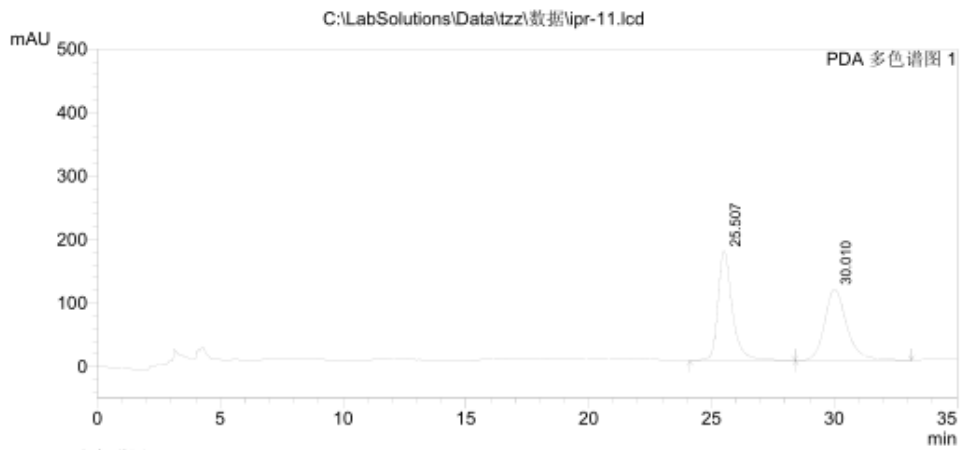
**(1S,2R)-(-)-1,2-Diphenylbut-3-en-1-ol (29)**



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	23.97	n.a.	489.449	233.362	47.81	n.a.	BM
2	26.22	n.a.	449.714	233.497	47.83	n.a.	MB
3	32.48	n.a.	16.715	10.662	2.18	n.a.	BMB
4	37.56	n.a.	14.452	10.616	2.17	n.a.	BMB
<b>Total:</b>			970.329	488.137	100.00	0.000	



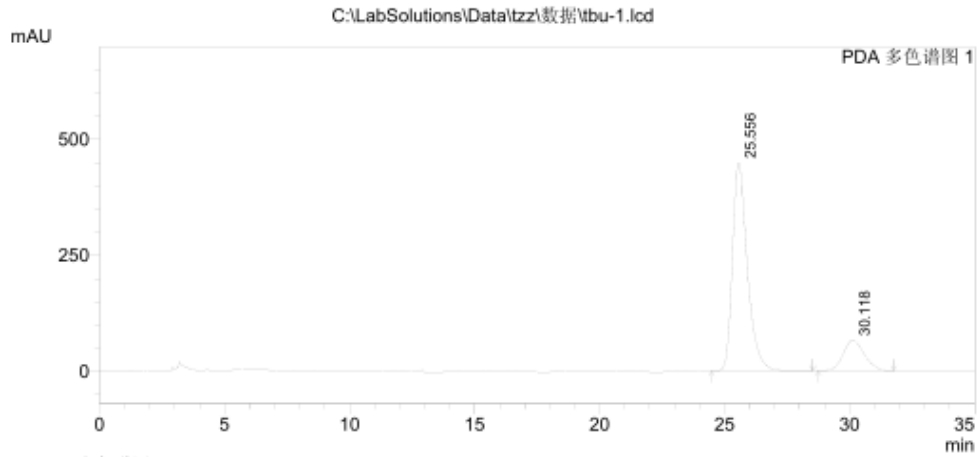
No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	24.04	n.a.	513.854	245.129	75.39	n.a.	BMB*
2	26.19	n.a.	137.104	68.880	21.18	n.a.	BMB*
3	32.60	n.a.	10.008	8.058	2.48	n.a.	BMB*
4	37.68	n.a.	4.050	3.077	0.95	n.a.	BMB*
<b>Total:</b>			665.016	325.143	100.00	0.000	



**PDA Ch1 214nm 4nm**

Peak#	Resolution Time	Area	Height	Area %	Height %
1	25.507	7092318	172969	49.949	60.543
2	30.010	7106896	112729	50.051	39.457
总计		14199215	285698	100.000	100.000



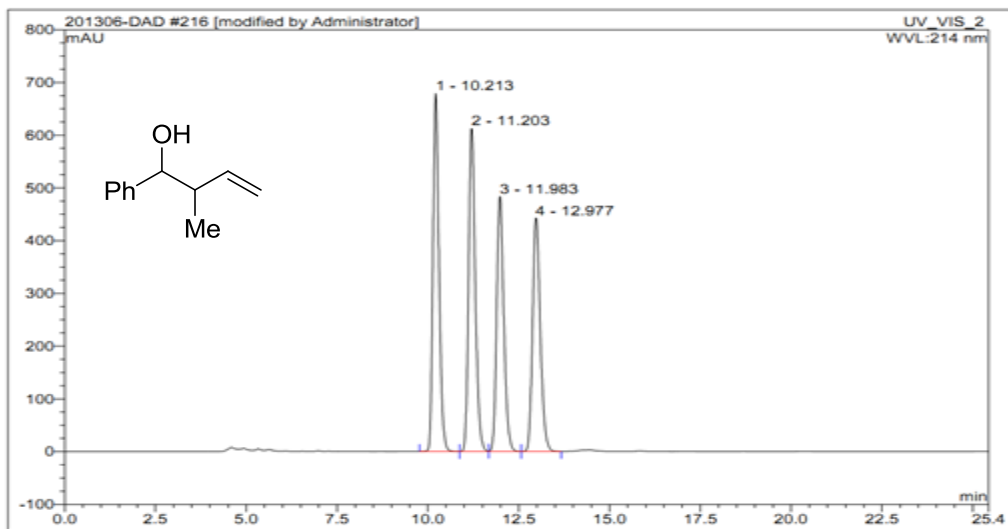


**PDA Ch1 214nm 4nm**

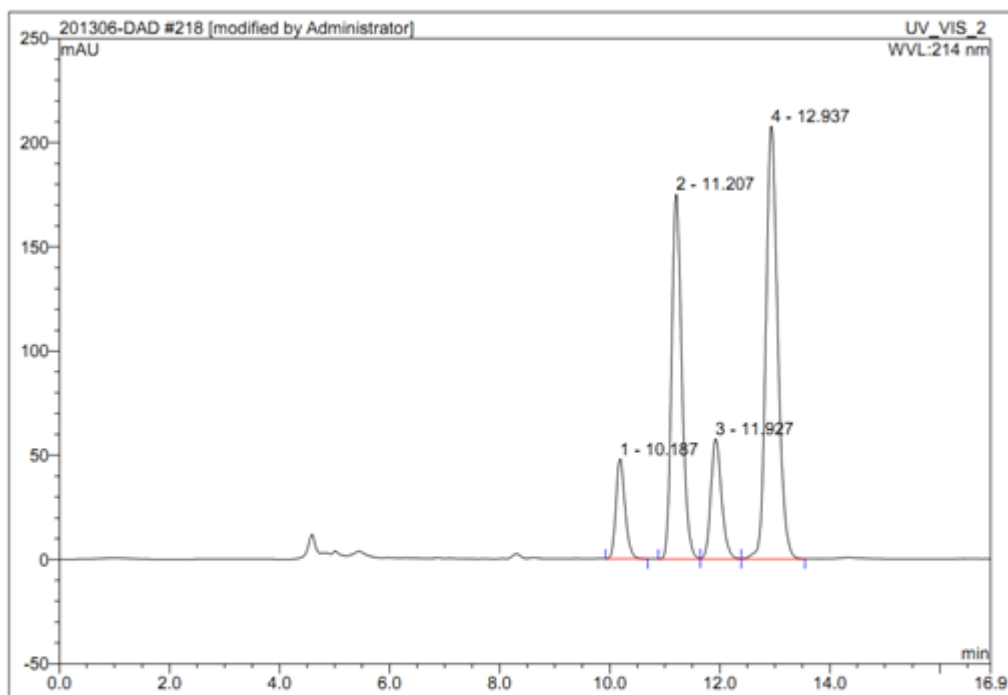
Peak#	Resolution Time	Area	Height	Area %	Height %
1	25.556	18627474	449591	82.615	87.123
2	30.118	3919789	66449	17.385	12.877
总计		22547262	516040	100.000	100.000

**2-Methyl-1-phenylbut-3-en-1-ol (30)**

For (Table 3, entry 4)

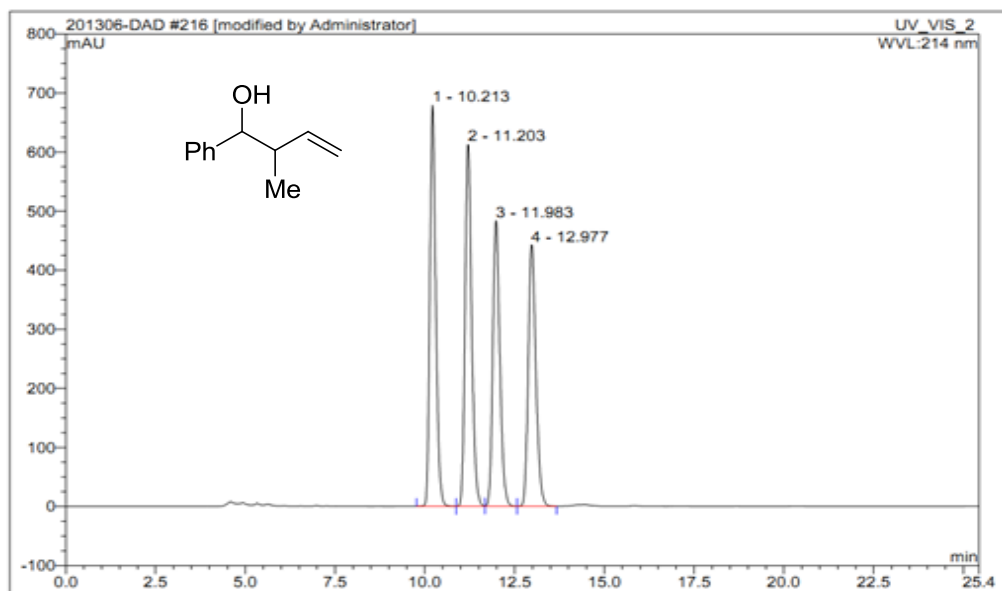


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	10.21	n.a.	679.166	130.605	27.18	n.a.	BM *
2	11.20	n.a.	612.951	130.737	27.21	n.a.	M *
3	11.98	n.a.	483.892	109.886	22.87	n.a.	M *
4	12.98	n.a.	442.822	109.313	22.75	n.a.	MB*
<b>Total:</b>			2218.831	480.540	100.00	0.000	

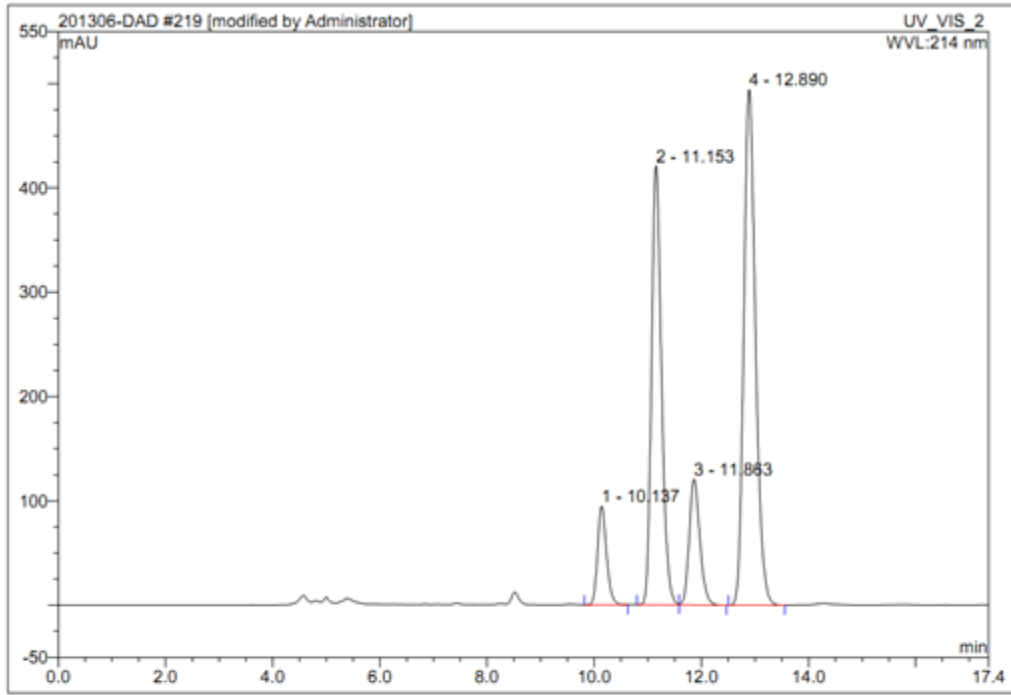


No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	10.19	n.a.	48.033	9.206	8.24	n.a.	BMB
2	11.21	n.a.	175.214	37.567	33.61	n.a.	BM
3	11.93	n.a.	57.838	13.051	11.68	n.a.	M
4	12.94	n.a.	207.803	51.955	46.48	n.a.	MB
<b>Total:</b>			488.889	111.779	100.00	0.000	

For (Table 3, entry 5)



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	10.21	n.a.	679.166	130.605	27.18	n.a.	BM *
2	11.20	n.a.	612.951	130.737	27.21	n.a.	M *
3	11.98	n.a.	483.892	109.886	22.87	n.a.	M *
4	12.98	n.a.	442.822	109.313	22.75	n.a.	MB*
<b>Total:</b>			2218.831	480.540	100.00	0.000	



No.	Ret.Time min	Peak Name	Height mAU	Area mAU*min	Rel.Area %	Amount	Type
1	10.14	n.a.	94.665	18.360	7.01	n.a.	BMB
2	11.15	n.a.	421.204	90.847	34.66	n.a.	BM
3	11.86	n.a.	120.681	27.846	10.62	n.a.	MB
4	12.89	n.a.	494.483	125.031	47.71	n.a.	BMB
<b>Total:</b>			1131.033	262.083	100.00	0.000	