

Supporting Information

Silver-Catalyzed Regioselective 1,6-Hydroarylation of *para*-Quinone Methides with Anilines and Phenols

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Reaction optimization:

Table S1. Optimization of solvent. ^a

Entry	Solvent	Catalyst	Temp.	Yield ^b
1	DCE	AgBF₄	80 °C	87%
2	CH ₂ Cl ₂	AgBF ₄	80 °C	77%
3	DMF	AgBF ₄	80 °C	44%
4	CH ₃ CN	AgBF ₄	80 °C	trace
5	1,4-Dioxane	AgBF ₄	80 °C	55%
6	H ₂ O	AgBF ₄	80 °C	N.D.
7	Toluene	AgBF ₄	80 °C	33%
8	THF	AgBF ₄	80 °C	31%

^a Reactions were operated with aniline (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), catalyst (10 mol%) and solvent (1.0 mL) under N₂ conditions. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard.

Table S2. Optimization of additive. ^a

Entry	Solvent	Additive	Catalyst	Temp.	Yield ^b
1	DCE	-	AgBF₄	80 °C	87%
2	DCE	Et ₃ N	AgBF ₄	80 °C	trace
3	DCE	K ₃ PO ₄	AgBF ₄	80 °C	13%
4	DCE	NaOAc	AgBF ₄	80 °C	17%
5	DCE	PhCOOH	AgBF ₄	80 °C	73%
6	DCE	Ph ₂ P(O)OH	AgBF ₄	80 °C	79%

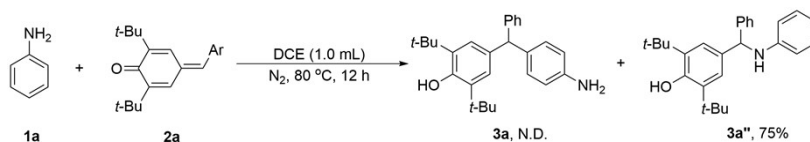
^a Reactions were operated with aniline (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), additive (10 mol%), catalyst (10 mol%) and solvent (1.0 mL) under N₂ conditions. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard.

Table S3. Optimization of the amount of AgBF₄. ^a

Entry	Solvent	Catalyst	Temp.	Yield ^b
1	DCE	AgBF ₄	80 °C	11% ^c

2	DCE	AgBF ₄	80 °C	33% ^d
3	DCE	AgBF ₄	80 °C	66% ^e
4	DCE	AgBF₄	80 °C	87%
5	DCE	AgBF ₄	80 °C	88% ^f

^a Reactions were operated with aniline (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), catalyst (10 mol%) and solvent (1.0 mL) under N₂ conditions. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard. ^c AgBF₄ (1 mol%) ^d AgBF₄ (2 mol%), ^e AgBF₄ (5 mol%), ^f AgBF₄ (20 mol%).



Scheme S1. 1,6-conjugated addition of **1a** to **2a**.

Table S4. Optimization of additive. ^a

Entry	Solvent	Additive	Catalyst	Temp.	Yield ^b
1	DCE	E ₃ N	AgBF ₄	R.T.	trace
2	DCE	DIEA	AgBF ₄	R.T.	trace
3	DCE	DBU	AgBF ₄	R.T.	trace
4	DCE	K ₃ PO ₄	AgBF ₄	R.T.	38%
5	DCE	NaOAc	AgBF ₄	R.T.	35%
6	DCE	PhCOOH	AgBF ₄	R.T.	84%
7	DCE	Ph₂P(O)OH	AgBF₄	R.T.	91%
8	DCE	TfOH	AgBF ₄	R.T.	81%
9	DCE	HBf ₄	AgBF ₄	R.T.	86%
10	DCE	TsOH	AgBF ₄	R.T.	83%
22	DCE	CF ₃ COOH	AgBF ₄	R.T.	84%

^a Reactions were operated with phenol (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), additive (10 mol%), catalyst (10 mol%) and solvent (1.0 mL) under N₂ atmosphere. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard.

Table S5. Optimization of the amount of Ph₂P(O)OH. ^a

Entry	Solvent	Additive	Catalyst	Temp.	Yield ^b
1	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	86% ^c
2	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	88% ^d
3	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	89% ^e
4	DCE	Ph₂P(O)OH	AgBF₄	R.T.	91%

5	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	90% ^f
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^a Reactions were operated with phenol (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), additive (10 mol%), catalyst (10 mol%) and solvent (1.0 mL) under N₂ atmosphere. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard. ^c Ph₂P(O)OH (1 mol%), ^d Ph₂P(O)OH (2 mol%), ^e Ph₂P(O)OH (5 mol%), ^f Ph₂P(O)OH (20 mol%).

Table S6. Optimization of the amount of AgBF₄. ^a

Entry	Solvent	Additive	Catalyst	Temp.	Yield ^b
1	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	24% ^c
2	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	45% ^d
3	DCE	Ph₂P(O)OH	AgBF₄	R.T.	91%
4	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	68% ^e
5	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	90% ^f

^a Reactions were operated with phenol (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), additive (10 mol%), catalyst (10 mol%) and solvent (1.0 mL) under N₂ atmosphere. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard. ^c AgBF₄ (1 mol%), ^d AgBF₄ (2 mol%), ^e AgBF₄ (5 mol%), ^f AgBF₄ (20 mol%).

Table S7. Optimization of solvent. ^a

Entry	Solvent	Additive	Catalyst	Temp.	Yield ^b
1	DCE	Ph₂P(O)OH	AgBF₄	R.T.	91%
2	DCM	Ph ₂ P(O)OH	AgBF ₄	R.T.	81%
3	DMF	Ph ₂ P(O)OH	AgBF ₄	R.T.	76%
4	CH ₃ CN	Ph ₂ P(O)OH	AgBF ₄	R.T.	trace
5	1,4-dioxane	Ph ₂ P(O)OH	AgBF ₄	R.T.	45%
6	H ₂ O	Ph ₂ P(O)OH	AgBF ₄	R.T.	N.D.
7	Toluene	Ph ₂ P(O)OH	AgBF ₄	R.T.	69%
8	THF	Ph ₂ P(O)OH	AgBF ₄	R.T.	67%
9	DME	Ph ₂ P(O)OH	AgBF ₄	R.T.	trace
10	DMSO	Ph ₂ P(O)OH	AgBF ₄	R.T.	N.D.

^a Reactions were operated with phenol (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), additive (10 mol%), catalyst (10 mol%) and solvent (1.0 mL) under N₂ atmosphere. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard.

Table S8. Optimization of atmosphere. ^a

Entry	Solvent	Additive	Catalyst	Temp.	Yield ^b
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1	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	91%
2	DCE	Ph ₂ P(O)OH	AgBF ₄	R.T.	66% ^c

^a Reactions were operated with phenol (0.2 mmol), 4-benzylidene-2,6-di-*tert*-butylcyclohexa-2,5-dienone (0.2 mmol), additive (10 mol%), catalyst (10 mol%) and solvent (1.0 mL) under N₂ atmosphere. ^b Yield was determined by GC analysis, and dodecane was used as the internal standard. ^c air atmosphere.

General considerations:

All solvents used in the reactions were freshly distilled. The other reagents were recrystallized or distilled as necessary. All reactions were performed under an atmosphere of N₂ unless specified otherwise. ¹H (400 MHz), ¹³C (100 MHz), ³¹P (160 MHz) and ¹⁹F (376 MHz) NMR spectra were recorded on a 400 MHz spectrometer in CDCl₃. ¹H NMR chemical shifts were reported using TMS as internal standard while ¹³C NMR chemical shifts were reported relative to CDCl₃. The electron ionization method was used for HRMS measurements, and the mass analyzer type was double-focusing.

General procedure for 1,6-hydroarylation of *p*-QMs with anilines:

A mixture of anilines (0.2 mmol), *p*-QMs compounds (0.2 mmol) and AgBF₄ (10 mol %) were dissolved in DCE (1.0 mL) under a N₂ atmosphere and the mixture was stirred for 12.0 h at 80 °C. Upon completion of the reaction, the mixture was concentrated under vacuum. Removal of the solvent under a reduced pressure gave the crude product; pure product was obtained by passing the crude product through a short silica gel column using hexane/EtOAc (10:1-2:1) as eluent.

General procedure for 1,6-hydroarylation of *p*-QMs with phenols:

A mixture of phenols (0.2 mmol), *p*-QMs compounds (0.2 mmol), Ph₂P(O)OH (10 mol %) and AgBF₄ (10 mol %) were dissolved in DCE (1.0 mL) under a N₂ atmosphere and the mixture was stirred for 12.0 h at room temperature. Upon completion of the reaction, the mixture was concentrated under vacuum. Removal of the solvent under a reduced pressure gave the crude product; pure product was obtained by passing the crude product through a short silica gel column using hexane/EtOAc (10:1-2:1) as eluent.

Crystallographic data

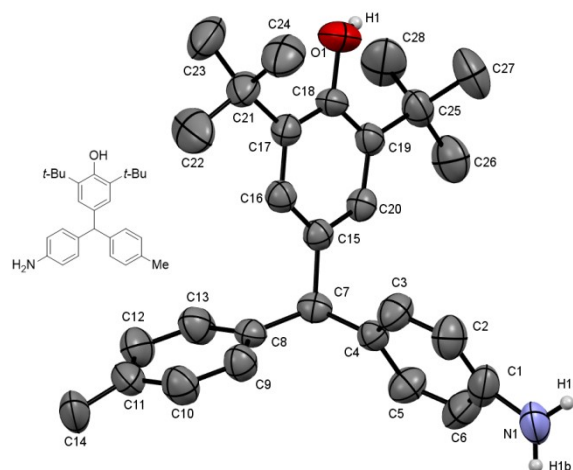


Figure S1. ORTEP drawing (a thermal ellipsoid plot) of compound **4a**.

Hydrogen atoms are omitted for clarity; ellipsoids are drawn at 50% probability.

Table S9. Crystal data and structure refinement for **4a**

Identification code	4a
CCDC	2144010
Empirical formula	C ₂₈ H ₃₅ NO
Formula weight	401.57
Temperature/K	298(2)
Crystal description	block
Space group	P-1
a/Å	9.8041(11)
b/Å	10.9905(12)
c/Å	11.7996(14)
α/°	83.266(2)
β/°	79.515(1)
γ/°	87.061(3)
Volume/Å ³	1241.0(2)
Z	2
ρ calcd/cm ³	1.075
μ/mm ⁻¹	0.064
F(000)	436.168
Crystal size/mm ³	0.40 × 0.30 × 0.21
Radiation	MoKα (λ = 0.71073)
Index ranges	-11 ≤ h ≤ 11, -12 ≤ k ≤ 13, 0 ≤ l ≤ 13

Reflections collected	7762
Independent reflections	4276 [$R_{\text{int}} = 0.0421$, $R_{\text{sigma}} = 0.0768$]
Data/restraints/parameters	4276/0/282
Goodness-of-fit on F_2	1.058
Final R indexes [$I \geq 2\sigma(I)$]	$R_1 = 0.0603$, $wR_2 = 0.1507$
Final R indexes [all data]	$R_1 = 0.1073$, $wR_2 = 0.1784$

Experimental:

Single crystals of $C_{28}H_{35}NO$ (**4a**) were obtained by slow evaporation from its saturated solution of dichloromethane/n-hexane at room temperature. Suitable single crystals were selected for indexing, and intensity data were measured on a Siemens Smart CCD diffractometer with graphite-monochromated $MoK\alpha$ radiation ($\lambda = 0.71073 \text{ \AA}$) at 298(2) K. The raw data frames were integrated into SHELX format reflection files and corrected using SAINT program. Absorption corrections based on multiscan were obtained by the SADABS program. The structure was solved by direct methods and refined with fullmatrix least-squares technique using the ShelXL (Sheldrick, 2015) and ShelXT (Sheldrick, 2015) programs, respectively.

Crystal structure determination of **4a**:

Crystal Data for $C_{28}H_{35}NO$ ($M = 401.57 \text{ g/mol}$): triclinic, space group P-1, $a = 9.8041(11) \text{ \AA}$, $b = 10.9905(12) \text{ \AA}$, $c = 11.7996(14) \text{ \AA}$, $V = 1241.0(2) \text{ \AA}^3$, $Z = 2$, $T = 298(2) \text{ K}$, $\mu(MoK\alpha) = 0.71073 \text{ mm}^{-1}$, $D_{\text{calc}} = 1.075 \text{ g/cm}^3$, 7762 reflections measured, 4276 unique ($R_{\text{int}} = 0.0421$, $R_{\text{sigma}} = 0.0768$) which were used in all calculations. The final R_1 was 0.0603 ($I > 2\sigma(I)$) and wR_2 was 0.1784 (all data).

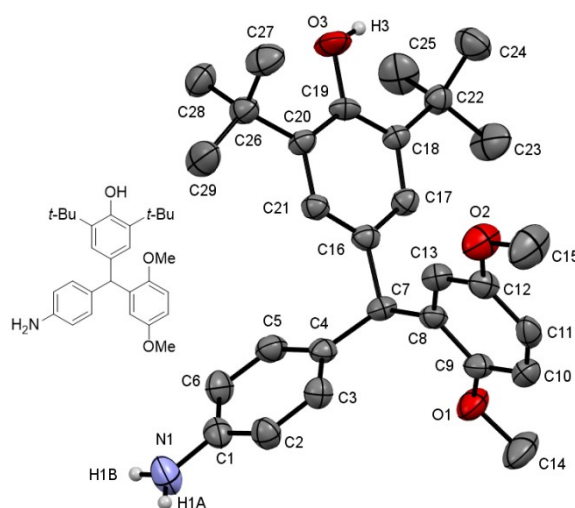


Figure S2. ORTEP drawing (a thermal ellipsoid plot) of compound **4j**.

Hydrogen atoms are omitted for clarity; ellipsoids are drawn at 50% probability.

Table S10. Crystal data and structure refinement for **4j**

Identification code	4j
CCDC	2144013
Empirical formula	C ₂₉ H ₃₇ NO ₃
Formula weight	447.60
Temperature/K	298(2)
Crystal description	block
Space group	P2(1)/c
a/Å	9.9675(11)
b/Å	12.6997(13)
c/Å	20.234(2)
α/°	90.00
β/°	101.754(3)
γ/°	90.00
Volume/Å ³	2507.6(5)
Z	4
ρ calcd/cm ³	1.186
μ/mm ⁻¹	0.076
F(000)	968
Crystal size/mm ³	0.25 * 0.20 * 0.12
Radiation	MoKα (λ = 0.71073)
Index ranges	-11 ≤ h ≤ 8, -15 ≤ k ≤ 11, -24 ≤ l ≤ 21
Reflections collected	11864
Independent reflections	4418 [R _{int} = 0.0940, R _{sigma} = 0.1251]
Data/restraints/parameters	4418/0/308
Goodness-of-fit on F ₂	1.091
Final R indexes [I ≥ 2σ (I)]	R ₁ = 0.0747, wR ₂ = 0.1607
Final R indexes [all data]	R ₁ = 0.1518, wR ₂ = 0.1828

Experimental:

Single crystals of C₂₉H₃₇NO₃ (**4j**) were obtained by slow evaporation from its saturated solution of dichloromethane/n-hexane at room temperature. Suitable single crystals were selected for indexing, and intensity data were measured on a Siemens Smart CCD diffractometer with graphite-monochromated MoKα radiation (λ = 0.71073 Å) at 298(2) K. The raw data frames were integrated into SHELX format reflection files and corrected using SAINT program. Absorption corrections based on multiscan were obtained by the SADABS program. The structure was solved by direct methods and refined with fullmatrix least-squares technique using the ShelXL (Sheldrick, 2015) and ShelXT (Sheldrick, 2015) programs, respectively.

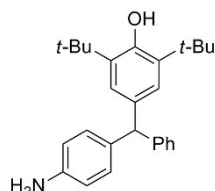
Crystal structure determination of **4j**:

Crystal Data for C₂₉H₃₇NO₃ (M = 447.60 g/mol): monoclinic, space group P2(1)/c, a = 9.9675(11) Å, b = 12.6997(13) Å, c = 20.234(2) Å, V = 2507.6(5) Å³, Z = 4, T =

298(2) K, $\mu(\text{MoK}\alpha) = 0.71073 \text{ mm}^{-1}$, $D_{\text{calc}} = 1.186 \text{ g/cm}^3$, 11864 reflections measured, 4418 unique ($R_{\text{int}} = 0.0940$, $R_{\text{sigma}} = 0.1251$) which were used in all calculations. The final R_1 was 0.0747 ($I > 2\sigma(I)$) and wR_2 was 0.1828 (all data).

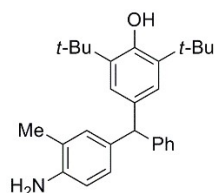
Analytical data for the compounds

4-((4-Aminophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3a)



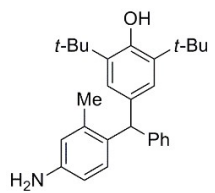
According to the general procedure, work-up and flash column chromatography (Hexane/EtOAc: 10:1) gave product **3a** (68.9 mg, 0.17 mmol, 85%) as a yellow solid. mp: 108.1-108.5 °C. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): $\delta = 7.23\text{--}7.27$ (m, 2H), 7.10-7.18 (m, 3H), 6.87-6.90 (m, 4H), 6.59-6.61 (m, 2H), 5.33 (s, 1H), 5.05 (s, 1H), 3.56 (s, 2H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): $\delta = 152.0$ (s), 145.5 (s), 144.3 (s), 135.4 (s), 135.2 (s), 134.7 (s), 130.2 (s), 129.4 (s), 128.1 (s), 126.0 (s), 125.8 (s), 115.1 (s), 56.1 (s), 36.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{34}\text{NO}$ $[\text{M}+\text{H}]^+$: 388.2635, found: 388.2629.

4-((4-Amino-3-methylphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3b)



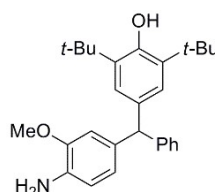
According to the general procedure, work-up and flash column chromatography (Hexane/EtOAc: 10:1) gave product **3b** (66.7 mg, 0.17 mmol, 83%) as a yellow solid. mp: 147.3-147.5 °C. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): $\delta = 7.22\text{--}7.26$ (m, 2H), 7.10-7.17 (m, 3H), 6.91-6.92 (m, 2H), 6.82-6.83 (m, 1H), 6.72-6.75 (m, 1H), 6.57-6.59 (m, 1H), 5.31 (s, 1H), 5.05 (s, 1H), 3.49 (s, 2H), 2.10 (s, 3H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): $\delta = 152.0$ (s), 145.6 (s), 142.5 (s), 135.3 (s), 135.2 (s), 134.8 (s), 131.5 (s), 129.4 (s), 128.0 (s), 127.9 (s), 126.0 (s), 125.8 (s), 122.2 (s), 114.9 (s), 56.2 (s), 34.4 (s), 30.4 (s), 17.5 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{28}\text{H}_{36}\text{NO}$ $[\text{M}+\text{H}]^+$: 402.2792, found: 402.2788.

4-((4-Amino-2-methylphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3c)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3c** (68.3 mg, 0.17 mmol, 85%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): $\delta = 7.22\text{--}7.25$ (m, 2H), 7.14-7.17 (m, 1H), 7.05-7.06 (m, 2H), 6.82-6.83 (m, 2H), 6.58-6.60 (m, 1H), 6.50-6.51 (m, 1H), 6.42-6.44 (m, 1H), 5.45 (s, 1H), 5.04 (s, 1H), 3.50 (s, 2H), 2.12 (s, 3H), 1.34 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): $\delta = 151.9$ (s), 144.8 (s), 144.3 (s), 137.5 (s), 135.3 (s), 134.3 (s), 133.8 (s), 130.3 (s), 129.5 (s), 128.0 (s), 126.2 (s), 125.8 (s), 117.4 (s), 122.4 (s), 52.8 (s), 34.3 (s), 30.4 (s), 20.1 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{28}\text{H}_{36}\text{NO}$ $[\text{M}+\text{H}]^+$: 402.2792, found: 402.2786.

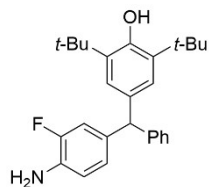
4-((4-Amino-3-methoxyphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3d)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3d** (70.3 mg, 0.17 mmol, 84%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): $\delta = 7.22\text{--}7.26$ (m, 2H), 7.10-7.17 (m, 3H), 6.92-6.93 (m, 2H), 6.57-6.62 (m, 2H), 6.48-6.51 (m, 1H), 5.34 (s, 1H), 5.06 (s, 1H), 3.71 (s,

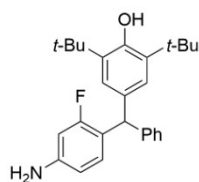
3H), 3.36 (s, 2H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): δ = 152.1 (s), 147.3 (s), 145.5 (s), 135.4 (s), 135.3 (s), 134.7 (s), 134.0 (s), 129.4 (s), 128.1 (s), 126.0 (s), 125.9 (s), 122.0 (s), 114.8 (s), 112.0 (s), 56.5 (s), 55.5 (s), 34.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{28}\text{H}_{36}\text{NO}_2$ $[\text{M}+\text{H}]^+$: 418.2741, found: 418.2735.

4-((4-Amino-3-fluorophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3e)



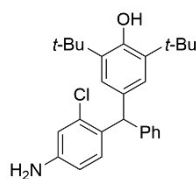
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3e** (66.6 mg, 0.16 mmol, 82%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): δ = 7.24-7.28 (m, 2H), 7.16-7.19 (m, 1H), 7.09-7.11 (m, 2H), 6.88-6.89 (m, 2H), 6.65-6.74 (m, 3H), 5.32 (s, 1H), 5.07 (s, 1H), 3.61 (s, 2H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): δ = 151.6 (d, $J(\text{C},\text{F})$ = 237.0 Hz), 152.1 (s), 144.8 (s), 136.1 (d, $J(\text{C},\text{F})$ = 5.5 Hz), 135.5 (s), 134.2 (s), 132.2 (d, $J(\text{C},\text{F})$ = 13.2 Hz), 129.3 (s), 128.2 (s), 126.1 (s), 125.9 (s), 125.2 (d, $J(\text{C},\text{F})$ = 3.1 Hz), 116.7 (d, $J(\text{C},\text{F})$ = 3.7 Hz), 116.2 (d, $J(\text{C},\text{F})$ = 19.1 Hz), 55.9 (s), 34.4 (s), 30.4 (s); ^{19}F NMR (376 MHz, CDCl_3 , 25 °C, TMS): δ = -135.2 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{33}\text{FNO}$ $[\text{M}+\text{H}]^+$: 406.2541, found: 406.2534.

4-((4-Amino-2-fluorophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3f)



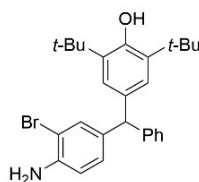
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3f** (65.6 mg, 0.16 mmol, 81%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): δ = 7.23-7.27 (m, 2H), 7.15-7.19 (m, 1H), 7.09-7.11 (m, 2H), 6.88-6.89 (m, 2H), 6.67-6.71 (m, 1H), 6.34-6.36 (m, 2H), 5.60 (s, 1H), 5.02 (s, 1H), 3.63 (s, 2H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): δ = 161.3 (d, $J(\text{C},\text{F})$ = 243.2 Hz), 152.1 (s), 146.4 (d, $J(\text{C},\text{F})$ = 10.9 Hz), 144.2 (s), 135.4 (s), 133.4 (s), 131.4 (d, $J(\text{C},\text{F})$ = 6.0 Hz), 129.1 (s), 128.1 (s), 126.0 (s), 125.9 (s), 121.7 (d, $J(\text{C},\text{F})$ = 15.1 Hz), 110.5 (d, $J(\text{C},\text{F})$ = 2.7 Hz), 102.2 (d, $J(\text{C},\text{F})$ = 25.6 Hz), 48.7 (d, $J(\text{C},\text{F})$ = 2.6 Hz), 34.4 (s), 30.4 (s); ^{19}F NMR (376 MHz, CDCl_3 , 25 °C, TMS): δ = -116.2 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{33}\text{FNO}$ $[\text{M}+\text{H}]^+$: 406.2541, found: 406.2535.

4-((4-Amino-2-chlorophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3g)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3g** (68.4 mg, 0.16 mmol, 81%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): δ = 7.24-7.27 (m, 2H), 7.18-7.19 (m, 1H), 7.06-7.08 (m, 2H), 6.84-6.85 (m, 2H), 6.70-6.72 (m, 2H), 6.47-6.49 (m, 1H), 5.73 (s, 1H), 5.07 (s, 1H), 3.62 (s, 2H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): δ = 152.0 (s), 145.5 (s), 144.0 (s), 135.3 (s), 135.3 (s), 134.8 (s), 133.5 (s), 132.5 (s), 131.6 (s), 129.4 (s), 128.1 (s), 126.2 (s), 126.0 (s), 115.8 (s), 52.4 (s), 34.3 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{33}\text{ClNO}$ $[\text{M}+\text{H}]^+$: 422.2246, found: 422.2241.

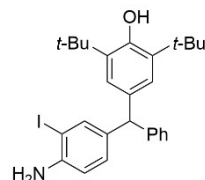
4-((4-Amino-3-bromophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3h)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3h** (77.2 mg, 0.17 mmol, 83%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 °C, TMS): δ = 7.22-7.27 (m, 2H), 7.15-7.19 (m, 2H), 7.08-7.10 (m, 2H), 6.87-6.88 (m, 2H), 6.82-6.84 (m, 1H), 6.64-6.66 (m, 1H), 5.30 (s, 1H), 5.08 (s, 1H), 3.95 (s, 2H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 °C, TMS): δ = 152.2 (s), 144.8 (s), 142.1 (s), 136.4 (s), 135.5 (s), 134.1 (s), 133.2 (s), 129.4 (s), 129.3 (s), 128.2 (s), 126.1 (s), 125.9 (s),

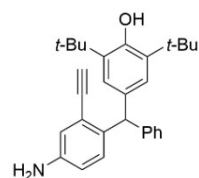
115.6 (s), 109.3 (s), 55.7 (s), 34.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $C_{27}H_{33}BrNO$ $[M+H]^+$: 466.1741, found: 466.1735.

4-((4-Amino-3-iodophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3i)



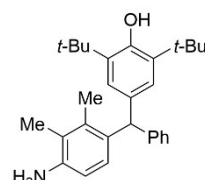
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3i** (81.1 mg, 0.16 mmol, 79%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.41-7.42 (m, 1H), 7.24-7.28 (m, 2H), 7.16-7.20 (m, 1H), 7.08-7.10 (m, 2H), 6.84-6.88 (m, 3H), 6.64-6.66 (m, 1H), 5.28 (s, 1H), 5.08 (s, 1H), 3.98 (s, 2H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 152.1 (s), 144.8 (s), 139.5 (s), 136.8 (s), 135.5 (s), 134.0 (s), 130.4 (s), 129.3 (s), 129.2 (s), 128.2 (s), 126.1 (s), 125.9 (s), 114.5 (s), 84.4 (s), 55.5 (s), 34.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $C_{27}H_{33}INO$ $[M+H]^+$: 514.1602, found: 514.1595.

4-((4-Amino-2-ethynylphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3j)



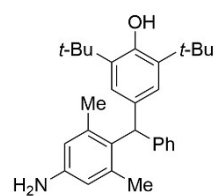
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **3j** (65.9 mg, 0.16 mmol, 80%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.22-7.26 (m, 2H), 7.16-7.18 (m, 1H), 7.09-7.11 (m, 2H), 6.89-6.90 (m, 2H), 6.79-6.83 (m, 2H), 6.59-6.61 (m, 1H), 5.86 (s, 1H), 5.05 (s, 1H), 3.59 (s, 2H), 3.15 (s, 1H), 1.35 (m, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 151.9 (s), 144.6 (s), 144.1 (s), 138.0 (s), 135.2 (s), 134.0 (s), 130.2 (s), 129.4 (s), 127.9 (s), 126.2 (s), 125.8 (s), 122.7 (s), 119.0 (s), 116.1 (s), 82.7 (s), 80.9 (s), 53.3 (s), 34.3 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $C_{29}H_{34}NO$ $[M+H]^+$: 412.2635, found: 412.2631. HRMS (ESI) m/z : calcd. for $C_{29}H_{34}NO$ $[M+H]^+$: 412.2635, found: 412.2630.

4-((4-Amino-2,3-dimethylphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3k)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **3k** (66.7 mg, 0.16 mmol, 81%) as a white solid. mp: 153.2-153.5 °C. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.21-7.25 (m, 2H), 7.13-7.17 (m, 1H), 7.04-7.05 (m, 2H), 6.82-6.83 (m, 2H), 6.44-6.49 (m, 2H), 5.54 (s, 1H), 5.04 (s, 1H), 3.50 (s, 2H), 2.12 (s, 3H), 2.09 (s, 3H), 1.34 (m, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 151.8 (s), 145.2 (s), 142.5 (s), 135.6 (s), 135.2 (s), 134.6 (s), 134.0 (s), 129.5 (s), 128.0 (s), 127.6 (s), 126.4 (s), 125.7 (s), 121.4 (s), 112.5 (s), 53.7 (s), 34.3 (s), 30.4 (s), 16.1 (s), 13.5 (s). HRMS (ESI) m/z : calcd. for $C_{29}H_{38}NO$ $[M+H]^+$: 416.2948, found: 416.2940.

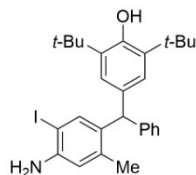
4-((4-Amino-2,6-dimethylphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3l)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3l** (58.2 mg, 0.14 mmol, 68%) as a white solid. mp: 143.3-143.7 °C. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.17-7.28 (m, 5H), 7.04-7.05 (m, 2H), 6.45 (s, 1H), 6.30 (s, 1H), 5.70 (s, 1H), 5.09 (s, 1H), 3.30 (s, 2H), 2.22 (s, 3H), 2.20 (s, 3H), 1.35 (m, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 152.2 (s), 145.5 (s), 143.4 (s), 137.9 (s), 136.7 (s), 135.7 (s), 132.2 (s), 128.9 (s), 125.8 (s), 124.6 (s), 122.2 (s), 116.4 (s),

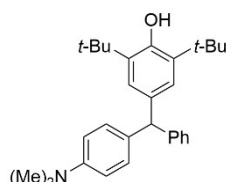
49.8 (s), 34.4 (s), 30.4 (s), 21.5 (s), 21.0 (s). HRMS (ESI) m/z : calcd. for $C_{29}H_{38}NO$ $[M+H]^+$: 416.2948, found: 416.2942.

4-((4-Amino-5-iodo-2-methylphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3m)



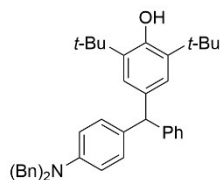
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **3m** (87.6 mg, 0.17 mmol, 83%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.23-7.27 (m, 2H), 7.17-7.18 (m, 1H), 7.03-7.05 (m, 3H), 6.82 (s, 2H), 6.55 (s, 1H), 5.38 (s, 1H), 5.07 (s, 1H), 3.90 (s, 2H), 2.07 (s, 3 H), 1.36 (m, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 152.1 (s), 144.8 (s), 144.1 (s), 139.5 (s), 138.2 (s), 135.7 (s), 135.5 (s), 133.6 (s), 129.3 (s), 128.2 (s), 126.1 (s), 126.0 (s), 116.8 (s), 80.9 (s), 52.6 (s), 34.4 (s), 30.4 (s), 19.9 (s). HRMS (ESI) m/z : calcd. for $C_{28}H_{35}INO$ $[M+H]^+$: 528.1758, found: 528.1753.

2,6-Di-*tert*-butyl-4-((4-(dimethylamino)phenyl)(phenyl)methyl)phenol (3n)



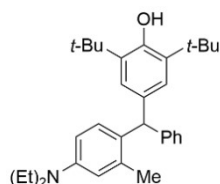
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3n** (71.6 mg, 0.17 mmol, 86%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.23-7.27 (m, 2H), 7.16-7.18 (m, 3H), 6.92-6.98 (m, 4H), 6.66-6.68 (m, 2H), 5.35 (s, 1H), 5.05 (s, 1H), 2.91 (s, 6H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 151.9 (s), 148.9 (s), 145.7 (s), 135.3 (s), 134.8 (s), 133.1 (s), 130.0 (s), 129.4 (s), 128.0 (s), 126.0 (s), 125.8 (s), 112.6 (s), 56.0 (s), 40.8 (s), 34.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $C_{29}H_{38}NO$ $[M+H]^+$: 416.2948, found: 416.2942.

2,6-Di-*tert*-butyl-4-((4-(dibenzylamino)phenyl)(phenyl)methyl)phenol (3o)



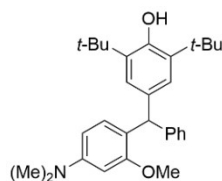
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3o** (88.6 mg, 0.16 mmol, 78%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.11-7.22 (m, 12H), 7.02-7.08 (m, 3H), 6.79-6.82 (m, 4H), 6.55-6.57 (m, 2H), 5.23 (s, 1H), 4.94 (s, 1H), 4.51 (s, 4H), 1.26 (s, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 152.0 (s), 147.5 (s), 145.7 (s), 138.9 (s), 135.3 (s), 134.9 (s), 133.1 (s), 130.1 (s), 129.5 (s), 128.7 (s), 128.1 (s), 126.9 (s), 126.8 (s), 126.1 (s), 125.9 (s), 112.5 (s), 56.0 (s), 54.5 (s), 34.4 (s), 30.5 (s). HRMS (ESI) m/z : calcd. for $C_{41}H_{46}NO$ $[M+H]^+$: 568.3574, found: 568.3570.

2,6-Di-*tert*-butyl-4-((4-(diethylamino)-2-methylphenyl)(phenyl)methyl)phenol (3p)



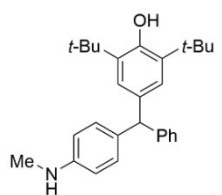
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3p** (65.1 mg, 0.14 mmol, 74%) as a red oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.22-7.26 (m, 2H), 7.12-7.17 (m, 1H), 7.07-7.09 (m, 2H), 6.86 (s, 2H), 6.62-6.64 (m, 1H), 6.43-6.50 (m, 2H), 5.44 (s, 1H), 5.03 (s, 1H), 3.28-3.33 (m, 4H), 2.15 (s, 3H), 1.35 (s, 18H), 1.13 (t, J = 7.0 Hz, 6H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 146.5 (s), 140.9 (s), 139.9 (s), 131.9 (s), 129.9 (s), 129.3 (s), 125.3 (s), 124.8 (s), 124.1 (s), 122.6 (s), 121.0 (s), 120.3 (s), 108.7 (s), 104.0 (s), 47.51 (s), 38.96 (s), 29.02 (s), 25.07 (s), 15.36 (s), 7.40 (s). HRMS (ESI) m/z : calcd. for $C_{32}H_{44}NO$ $[M+H]^+$: 458.3418, found: 458.3413.

2,6-Di-*tert*-butyl-4-((4-(dimethylamino)-2-methoxyphenyl)(phenyl)methyl)phenol (**3q**)



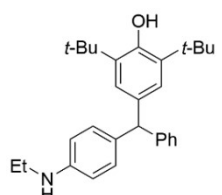
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3q** (66.2 mg, 0.14 mmol, 71%) as a yellow solid. mp: 136.1-136.3 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.11-7.14 (m, 2H), 7.00-7.05 (m, 3H), 6.82-6.83 (m, 2H), 6.63-6.65 (m, 1H), 6.16-6.18 (m, 2H), 5.61 (s, 1H), 4.92 (s, 1H), 3.60 (s, 3H), 2.83 (s, 6H), 1.27 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 157.9 (s), 151.8 (s), 150.5 (s), 145.7 (s), 135.1 (s), 134.8 (s), 130.6 (s), 129.3 (s), 127.8 (s), 126.3 (s), 125.5 (s), 122.3 (s), 104.6 (s), 96.8 (s), 55.7 (s), 49.0 (s), 40.9 (s), 34.4 (s), 30.5 (s). HRMS (ESI) *m/z*: calcd. for C₃₀H₄₀NO₂ [M+H]⁺: 446.3054, found: 446.3050.

2,6-Di-*tert*-butyl-4-((4-(methylamino)phenyl)(phenyl)methyl)phenol (**3r**)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 30:1) gave product **3r** (67.5 mg, 0.17 mmol, 84%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.26 (m, 2H), 7.11-7.17 (m, 3H), 6.92-6.93 (m, 4H), 6.52-6.54 (m, 2H), 5.34 (s, 1H), 5.05 (s, 1H), 3.60 (s, 1H), 2.80 (s, 3H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 147.5 (s), 145.7 (s), 135.3 (s), 134.8 (s), 133.8 (s), 130.1 (s), 129.4 (s), 128.1 (s), 126.0 (s), 125.8 (s), 112.3 (s), 56.1 (s), 34.4 (s), 31.0 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₆NO [M+H]⁺: 402.2792, found: 402.2787.

2,6-Di-*tert*-butyl-4-((4-(ethylamino)phenyl)(phenyl)methyl)phenol (**3s**)

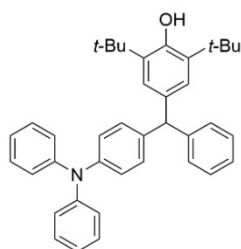


According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 30:1) gave product **3s** (68.2 mg, 0.16 mmol, 82%) as a red oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.22-7.26 (m, 2H), 7.10-7.17 (m, 3H), 6.90-6.91 (m, 4H), 6.52-6.54 (m, 2H), 5.33 (s, 1H), 5.04 (s, 1H), 3.42 (s, 1H), 3.10-3.15 (m, 2H), 1.35 (s, 18H), 1.23 (t, *J* = 7.1 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 146.6 (s), 145.7 (s), 135.3 (s), 134.9 (s), 133.8 (s), 130.2 (s), 129.4 (s), 128.0 (s), 126.0 (s), 125.8 (s), 112.6 (s), 56.1 (s), 38.7 (s), 34.4 (s), 30.4 (s), 15.0 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₈NO [M+H]⁺: 416.2948, found: 416.2944.

2,6-Di-*tert*-butyl-4-((4-(naphthalen-2-ylamino)phenyl)(phenyl)methyl)phenol (**3t**)

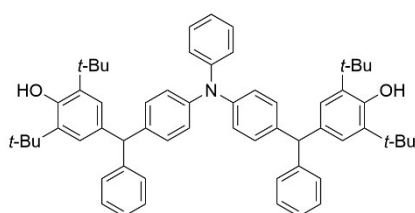
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3t** (75.0 mg, 0.15 mmol, 73%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.69-7.71 (m, 2H), 7.59-7.61 (m, 1H), 7.35-7.39 (m, 2H), 7.24-7.30 (m, 3H), 7.15-7.22 (m, 4H), 7.04-7.09 (m, 4H), 6.94 (s, 2H), 5.80 (s, 1H), 5.42 (s, 1H), 5.07 (s, 1H), 1.37 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 145.1 (s), 141.3 (s), 140.7 (s), 138.2 (s), 135.5 (s), 134.7 (s), 134.4 (s), 130.3 (s), 129.4 (s), 129.1 (s), 129.0 (s), 128.2 (s), 127.7 (s), 126.5 (s), 126.4 (s), 126.1 (s), 126.0 (s), 123.3 (s), 119.8 (s), 118.5 (s), 110.9 (s), 56.3 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₇H₄₀NO [M+H]⁺: 514.3105, found: 514.3100.

2,6-Di-*tert*-butyl-4-((4-(diphenylamino)phenyl)(phenyl)methyl)phenol (3u)



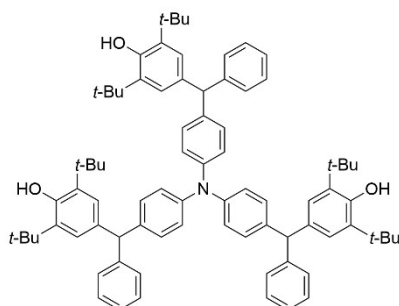
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3u** as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.18-7.23 (m, 2H), 7.10-7.16 (m, 5H), 7.06-7.08 (m, 2H), 6.98-7.00 (m, 4H), 6.88-6.92 (m, 6H), 6.83 (m, 2H), 5.32 (s, 1H), 5.00 (s, 1H), 1.30 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.0 (s), 146.9 (s), 144.7 (s), 143.9 (s), 138.4 (s), 134.4 (s), 133.3 (s), 129.1 (s), 128.4 (s), 128.2 (s), 128.1 (s), 127.1 (s), 125.0 (s), 123.2 (s), 122.8 (s), 121.4 (s), 55.2 (s), 33.3 (s), 29.3 (s). HRMS (ESI) *m/z*: calcd. for C₃₉H₄₂NO [M+H]⁺: 540.3261, found: 540.3254.

4,4'-(((Phenylazanediy)bis(4,1-phenylene))bis(phenylmethylene))bis(2,6-di-*tert*-butylphenol) (3u')



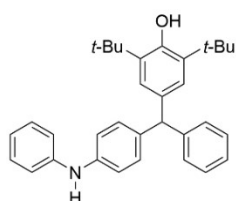
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3u'** as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.17-7.22 (m, 4H), 7.09-7.13 (m, 4H), 7.05-7.07 (m, 4H), 6.96-6.98 (m, 2H), 6.86-6.90 (m, 9H), 6.82 (m, 4H), 5.31 (s, 2H), 4.99 (s, 2H), 1.28 (s, 36H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.0 (s), 146.9 (s), 144.7 (s), 143.9 (s), 138.1 (s), 134.3 (s), 133.4 (s), 129.1 (s), 128.4 (s), 128.0 (s), 127.1 (s), 125.0 (s), 122.9 (s), 122.8 (s), 122.4 (s), 121.0 (s), 55.2 (s), 33.3 (s), 29.3 (s). HRMS (ESI) *m/z*: calcd. for C₆₀H₆₈NO₂ [M+H]⁺: 834.5245, found: 834.5241.

4,4',4''-((Nitrilotris(benzene-4,1-diyl))tris(phenylmethylene))tris(2,6-di-*tert*-butylphenol) (3u'')



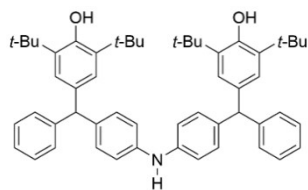
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **3u''** as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.28 (m, 6H), 7.17-7.19 (m, 2H), 7.11-7.13 (m, 6H), 6.92-6.98 (m, 11H), 6.89 (m, 6H), 5.37 (s, 3H), 5.05 (s, 2H), 1.35 (s, 54H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 145.9 (s), 145.05 (s), 138.8 (s), 135.4 (s), 134.5 (s), 130.0 (s), 129.4 (s), 128.1 (s), 126.1 (s), 126.0 (s), 123.7 (s), 56.2 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₈₁H₉₄NO₃ [M+H]⁺: 1128.7229, found: 1128.7223.

2,6-Di-*tert*-butyl-4-(phenyl(4-(phenylamino)phenyl)methyl)phenol (3v)



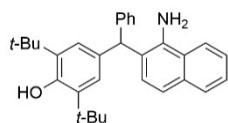
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3v** as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.05-7.22 (m, 7H), 6.90-6.96 (m, 6H), 6.79-6.85 (m, 3H), 5.57 (s, 1H), 5.31 (s, 1H), 4.99 (s, 1H), 1.29 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.0 (s), 144.1 (s), 142.5 (s), 139.9 (s), 136.7 (s), 134.4 (s), 133.4 (s), 129.2 (s), 128.3 (s), 128.2 (s), 127.1 (s), 124.9 (s), 124.8 (s), 119.5 (s), 117.0 (s), 116.3 (s), 55.2 (s), 33.3 (s), 29.3 (s). HRMS (ESI) *m/z*: calcd. for C₃₃H₃₈NO [M+H]⁺: 464.2948, found: 464.2944.

4,4'-((Azanediylbis(4,1-phenylene))bis(phenylmethylene))bis(2,6-di-*tert*-butylphenol) (**3v'**)



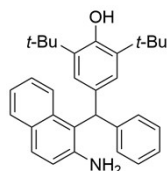
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3v'** as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.24-7.28 (m, 4H), 7.15-7.19 (m, 2H), 7.11-7.13 (m, 4H), 6.94-6.99 (m, 8H), 6.91 (m, 4H), 5.59 (s, 1H), 5.37 (s, 2H), 5.06 (s, 2H), 1.36 (s, 36H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.0 (s), 145.2 (s), 141.3 (s), 137.3 (s), 135.4 (s), 134.5 (s), 130.2 (s), 129.4 (s), 128.1 (s), 126.0 (s), 125.9 (s), 117.5 (s), 56.2 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₅₄H₆₄NO₂ [M+H]⁺: 758.4932, found: 758.4928.

4-((1-Aminonaphthalen-2-yl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (**3w**)



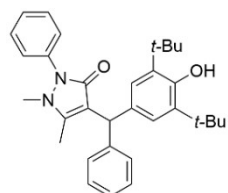
According to the general procedure, work-up and flash column chromatography (Hexane/EtOAc: 10:1) gave product **3w** (49.1 mg, 0.11 mmol, 56%) as a colorless oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.99-8.01 (m, 1H), 7.84-7.86 (m, 1H), 7.36-7.46 (m, 2H), 7.21-7.25 (m, 2H), 7.10-7.17 (m, 3H), 6.88 (s, 2H), 6.69-6.75 (m, 2H), 6.06 (s, 1H), 5.05 (s, 1H), 4.22 (s, 2H), 1.33 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.0 (s), 145.0 (s), 140.5 (s), 135.4 (s), 134.5 (s), 134.1 (s), 132.7 (s), 131.9 (s), 129.5 (s), 128.1 (s), 127.9 (s), 126.4 (s), 125.9 (s), 125.2 (s), 124.4 (s), 124.3 (s), 121.3 (s), 109.5 (s), 52.7 (s), 34.3 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₃₆NO [M+H]⁺: 438.2792, found: 438.2786.

4-((2-Aminonaphthalen-1-yl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (**3x**)



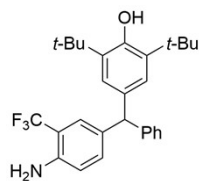
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3x** (64.8 mg, 0.15 mmol, 74%) as a colorless oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.95-7.97 (m, 1H), 7.69-7.71 (m, 1H), 7.60-7.62 (m, 1H), 7.32-7.36 (m, 1H), 7.17-7.28 (m, 6H), 7.07 (s, 2H), 6.84-6.86 (m, 1H), 6.28 (s, 1H), 5.10 (s, 1H), 3.57 (s, 2H), 1.33 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.3 (s), 143.6 (s), 143.0 (s), 135.9 (s), 134.3 (s), 132.1 (s), 129.0 (s), 128.8 (s), 128.7 (s), 128.6 (s), 128.5 (s), 126.5 (s), 126.3 (s), 125.9 (s), 122.7 (s), 121.8 (s), 120.5 (s), 118.8 (s), 48.8 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₃₆NO [M+H]⁺: 438.2792, found: 438.2788.

4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-1,5-dimethyl-2-phenyl-1,2-dihydro-3H-pyrazol-3-one (**3y**)



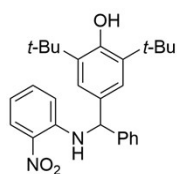
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3y** (55.1 mg, 0.11 mmol, 57%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.39-7.46 (m, 4H), 7.16-7.28 (m, 6H), 7.03 (s, 2H), 5.42 (s, 1H), 5.08 (s, 1H), 2.99 (s, 3H), 1.86 (s, 3H), 1.38 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 165.7 (s), 154.1 (s), 152.2 (s), 143.4 (s), 135.6 (s), 135.5 (s), 132.1 (s), 128.9 (s), 128.8 (s), 128.1 (s), 125.9 (s), 125.8 (s), 125.6 (s), 123.2 (s), 114.9 (s), 45.1 (s), 36.6 (s), 34.4 (s), 30.4 (s), 12.0 (s). HRMS (ESI) *m/z*: calcd. for C₃₂H₃₉N₂O₂ [M+H]⁺: 483.3007, found: 483.3001.

4-((4-Amino-3-(trifluoromethyl)phenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3z)



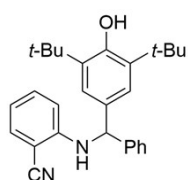
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **3z** (72.1 mg, 0.16 mmol, 79%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.17-7.28 (m, 4H), 7.02-7.10 (m, 3H), 6.88 (s, 2H), 6.64-6.66 (m, 1H), 5.34 (s, 1H), 5.09 (s, 1H), 4.05 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.2 (s), 144.7 (s), 142.6 (s), 135.6 (s), 134.4 (s), 133.9 (s), 133.8 (s), 129.2 (s), 128.3 (s), 127.2 (d, *J*(C,F) = 5.1 Hz), 126.2 (s), 125.9 (s), 125.1 (d, *J*(C,F) = 270.7 Hz), 117.2 (s), 113.6 (d, *J*(C,F) = 29.5 Hz), 55.8 (s), 34.4 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -62.5 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₃F₃NO [M+H]⁺: 456.2509, found: 456.2505.

2,6-Di-*tert*-butyl-4-(((2-nitrophenyl)amino)(phenyl)methyl)phenol (3aa')



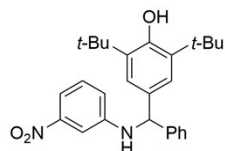
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 30:1) gave product **3aa'** (75.3 mg, 0.17 mmol, 87%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 8.63-8.64 (m, 1H), 8.18-8.20 (m, 1H), 7.24-7.39 (m, 6H), 7.09 (s, 2H), 6.72-6.74 (m, 1H), 6.60-6.65 (m, 1H), 5.66 (d, *J* = 5.6 Hz, 1H), 5.21 (s, 1H), 1.39 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.4 (s), 144.4 (s), 141.7 (s), 136.3 (s), 136.1 (s), 132.3 (s), 132.2 (s), 128.9 (s), 127.5 (s), 126.9 (s), 126.8 (s), 124.1 (s), 115.7 (s), 115.4 (s), 62.1 (s), 34.4 (s), 30.2 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃N₂O₃ [M+H]⁺: 433.2486, found: 433.2481.

2-(((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)amino)benzonitrile (3ab')



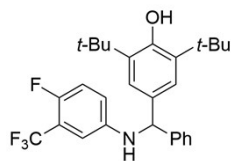
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 20:1) gave product **3ab'** (66.9 mg, 0.16 mmol, 81%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.31-7.41 (m, 5H), 7.20-7.28 (m, 2H), 7.04 (s, 2H), 6.64-6.66 (m, 1H), 6.49-6.51 (m, 1H), 5.53 (d, *J* = 4.8 Hz, 1H), 5.21 (s, 1H), 5.11 (d, *J* = 4.7 Hz, 1H), 1.38 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.5 (s), 149.4 (s), 141.6 (s), 136.3 (s), 134.2 (s), 132.8 (s), 132.3 (s), 128.8 (s), 127.5 (s), 127.1 (s), 124.5 (s), 118.1 (s), 116.9 (s), 112.3 (s), 96.0 (s), 62.4 (s), 34.4 (s), 30.2 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₃N₂O [M+H]⁺: 413.2588, found: 413.2584.

2,6-Di-*tert*-butyl-4-(((3-nitrophenyl)amino)(phenyl)methyl)phenol (3ac')



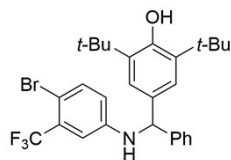
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 30:1) gave product **3ac'** (73.6 mg, 0.17 mmol, 85%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.46-7.48 (m, 1H), 7.32-7.40 (m, 5H), 7.17-7.27 (m, 2H), 7.08 (s, 2H), 6.77-6.80 (m, 1H), 5.48 (d, *J* = 4.4 Hz, 1H), 5.21 (s, 1H), 4.61 (d, *J* = 4.3 Hz, 1H), 1.39 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.4 (s), 149.2 (s), 148.1 (s), 141.8 (s), 136.3 (s), 132.8 (s), 129.6 (s), 128.9 (s), 127.5 (s), 127.1 (s), 124.5 (s), 119.0 (s), 112.0 (s), 107.6 (s), 63.0 (s), 34.5 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃N₂O₃ [M+H]⁺: 433.2486, found: 433.2480.

2,6-Di-*tert*-butyl-4-(((4-fluoro-3-(trifluoromethyl)phenyl)amino)(phenyl)methyl)phenol (**3ad'**)



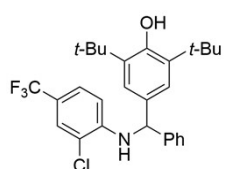
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 40:1) gave product **3ad'** (76.7 mg, 0.17 mmol, 83%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.30-7.39 (m, 4H), 7.21-7.26 (m, 1H), 7.08 (s, 2H), 6.86-6.90 (m, 1H), 6.71-6.73 (m, 1H), 6.55-6.59 (m, 1H), 5.35 (s, 1H), 5.19 (s, 1H), 4.31 (d, *J* = 2.1 Hz, 1H), 1.38 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.3 (s), 152.1 (d, *J*(C,F) = 241.9 Hz), 143.7 (d, *J*(C,F) = 2.1 Hz), 142.2 (s), 136.2 (s), 133.2 (s), 127.4 (s), 127.0 (s), 124.5 (s), 122.8 (d, *J*(C,F) = 269.7 Hz), 117.3 (d, *J*(C,F) = 11.1 Hz), 117.2 (d, *J*(C,F) = 2.6 Hz), 111.2 (d, *J*(C,F) = 4.6 Hz), 63.6 (s), 34.5 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -61.5 (d, *J* = 12.8 Hz), -130.7 (q, *J* = 12.7 Hz). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₂F₄NO [M+H]⁺: 474.2415, found: 474.2411.

4-(((4-Bromo-3-(trifluoromethyl)phenyl)amino)(phenyl)methyl)-2,6-di-*tert*-butylphenol (**3ae'**)



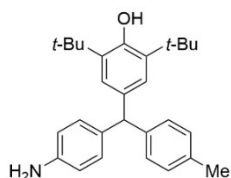
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 40:1) gave product **3ae'** (82.2 mg, 0.15 mmol, 77%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.30-7.37 (m, 5H), 7.20-7.26 (m, 1H), 7.06 (s, 2H), 6.87-6.88 (m, 1H), 6.44-6.47 (m, 1H), 5.38 (d, *J* = 4.1 Hz, 1H), 5.20 (s, 1H), 4.47 (d, *J* = 4.1 Hz, 1H), 1.38 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.4 (s), 146.4 (s), 141.8 (s), 136.3 (s), 135.3 (s), 132.9 (s), 127.5 (s), 127.0 (s), 124.5 (s), 123.0 (d, *J*(C,F) = 271.6 Hz), 116.9 (s), 113.1 (d, *J*(C,F) = 5.7 Hz), 105.4 (s), 63.2 (s), 34.5 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -62.7. HRMS (ESI) *m/z*: calcd. for C₂₈H₃₂BrF₃NO [M+H]⁺: 534.1614, found: 534.1607.

2,6-Di-*tert*-butyl-4-(((2-chloro-4-(trifluoromethyl)phenyl)amino)(phenyl)methyl)phenol (**3af'**)



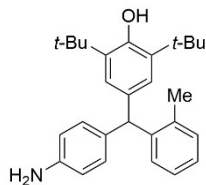
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 40:1) gave product **3af'** (76.4 mg, 0.16 mmol, 78%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.51-7.52 (m, 1H), 7.32-7.35 (m, 4H), 7.22-7.29 (m, 2H), 7.06 (s, 2H), 6.49-6.51 (m, 1H), 5.52 (d, *J* = 4.9 Hz, 1H), 5.22 (s), 5.21 (s), 1.38 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.4 (s), 145.6 (s), 141.6 (s), 136.2 (s), 132.5 (s), 128.9 (s), 127.5 (s), 127.1 (s), 126.2 (d, *J*(C,F) = 3.8 Hz), 125.0 (d, *J*(C,F) = 3.7 Hz), 124.4 (s), 121.0 (d, *J*(C,F) = 365.8 Hz), 118.9 (s), 118.7 (s), 111.8 (s), 62.4 (s), 34.4 (s), 30.2 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -61.1. HRMS (ESI) *m/z*: calcd. for C₂₈H₃₂ClF₃NO [M+H]⁺: 490.2120, found: 490.2115.

4-((4-Aminophenyl)(*p*-tolyl)methyl)-2,6-di-*tert*-butylphenol (**4a**)



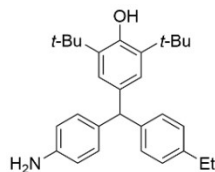
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4a** (65.1 mg, 0.16 mmol, 81%) as a yellow solid. mp: 147.3-147.5 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.98-7.06 (m, 4H), 6.87-6.91 (m, 4H), 6.57-6.60 (m, 2H), 5.28 (s, 1H), 5.04 (s, 1H), 3.43 (s, 2H), 2.30 (s, 3H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 144.2 (s), 142.5 (s), 135.4 (s), 135.3 (s), 135.2 (s), 134.9 (s), 130.2 (s), 129.2 (s), 128.8 (s), 126.0 (s), 115.0 (s), 55.7 (s), 34.4 (s), 30.4 (s), 21.1 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₆NO [M+H]⁺: 402.2792, found: 402.2786.

4-((4-Aminophenyl)(*o*-tolyl)methyl)-2,6-di-*tert*-butylphenol (4b)



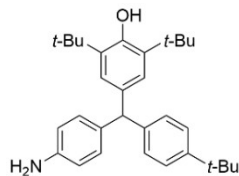
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4b** (68.3 mg, 0.17 mmol, 83%) as a yellow solid. mp: 147.3-147.5 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.06-7.13 (m, 3H), 6.82-6.84 (m, 5H), 6.57-6.61 (m, 2H), 5.44 (s, 1H), 5.04 (s, 1H), 3.29 (s, 2H), 2.21 (s, 3H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 144.1 (s), 143.8 (s), 136.6 (s), 135.3 (s), 134.4 (s), 134.2 (s), 130.3 (s), 130.2 (s), 129.2 (s), 126.2 (s), 125.9 (s), 125.5 (s), 115.1 (s), 52.6 (s), 34.3 (s), 30.4 (s), 20.0 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₆NO [M+H]⁺: 402.2792, found: 402.2787.

4-((4-Aminophenyl)(4-ethylphenyl)methyl)-2,6-di-*tert*-butylphenol (4c)



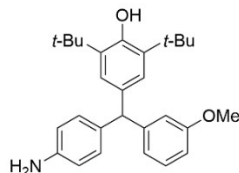
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4c** (69.9 mg, 0.41 mmol, 84%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.07-7.09 (m, 2H), 7.00-7.02 (m, 2H), 6.88-6.91 (m, 4H), 6.58-6.60 (m, 2H), 5.29 (s, 1H), 5.04 (s, 1H), 3.54 (s, 2H), 2.58-2.63 (m, 2H), 1.35 (s, 18H), 1.21 (t, *J* = 7.6 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 144.2 (s), 142.7 (s), 141.6 (s), 135.5 (s), 135.3 (s), 134.9 (s), 130.2 (s), 129.2 (s), 127.5 (s), 126.0 (s), 115.1 (s), 55.8 (s), 34.4 (s), 30.4 (s), 28.5 (s), 15.6 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₈NO [M+H]⁺: 416.2948, found: 416.2941.

4-((4-Aminophenyl)(4-(*tert*-butyl)phenyl)methyl)-2,6-di-*tert*-butylphenol (4d)



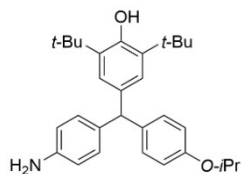
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4d** (71.0 mg, 0.16 mmol, 80%) as a yellow solid. mp: 145.2-145.5 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.22-7.27 (m, 2H), 7.02-7.03 (m, 2H), 6.89-6.91 (m, 4H), 6.57-6.60 (m, 2H), 5.28 (s, 1H), 5.03 (s, 1H), 3.53 (s, 2H), 1.36 (s, 18H), 1.29 (s, 9H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 148.5 (s), 144.3 (s), 142.3 (s), 135.5 (s), 135.3 (s), 134.9 (s), 130.2 (s), 128.9 (s), 126.0 (s), 124.9 (s), 115.0 (s), 55.7 (s), 34.4 (s), 31.5 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₄₂NO [M+H]⁺: 444.3261, found: 444.3255.

4-((4-Aminophenyl)(3-methoxyphenyl)methyl)-2,6-di-*tert*-butylphenol (4e)



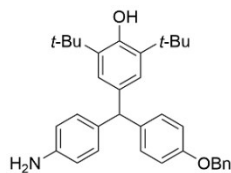
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4e** (71.9 mg, 0.17 mmol, 83%) as a yellow solid. mp: 108.1-108.5 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.15-7.19 (m, 1H), 6.88-6.91 (m, 4H), 6.67-6.72 (m, 3H), 6.58-6.60 (m, 2H), 5.29 (s, 1H), 5.05 (s, 1H), 3.73 (s, 3H), 3.47 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 159.4 (s), 152.0 (s), 147.1 (s), 144.3 (s), 135.3 (s), 135.0 (s), 134.5 (s), 130.2 (s), 128.9 (s), 126.0 (s), 122.0 (s), 115.3 (s), 115.1 (s), 111.1 (s), 56.1 (s), 55.1 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₆NO₂ [M+H]⁺: 418.2741, found: 418.2734.

4-((4-Aminophenyl)(4-isopropoxyphenyl)methyl)-2,6-di-*tert*-butylphenol (4f)



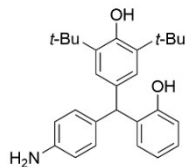
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4f** (73.1 mg, 0.16 mmol, 82%) as a yellow solid. mp: 120.6-120.8 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.98-7.00 (m, 2H), 6.88-6.90 (m, 4H), 6.76-6.79 (m, 2H), 6.59-6.61 (m, 2H), 5.26 (s, 1H), 5.04 (s, 1H), 6.45-6.54 (m, 1H), 3.55 (s, 2H), 1.35 (s, 18H), 1.31 (d, *J* = 6.1 Hz, 6H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 156.0 (s), 151.9 (s), 144.2 (s), 137.5 (s), 135.6 (s), 135.3 (s), 135.1 (s), 130.2 (s), 130.1 (s), 125.9 (s), 115.5 (s), 115.0 (s), 69.9 (s), 55.2 (s), 34.3 (s), 30.4 (s), 22.1 (s). HRMS (ESI) *m/z*: calcd. for C₃₀H₄₀NO₂ [M+H]⁺: 446.3054, found: 446.3050.

4-((4-Aminophenyl)(4-(benzyloxy)phenyl)methyl)-2,6-di-*tert*-butylphenol (4g)



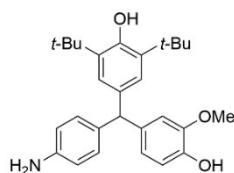
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4g** (82.1 mg, 0.17 mmol, 83%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.41-7.43 (m, 2H), 7.34-7.38 (m, 2H), 7.28-7.32 (m, 1H), 7.00-7.02 (m, 2H), 6.86-6.89 (m, 6H), 6.58-6.60 (m, 2H), 5.28 (s, 1H), 5.04 (s, 1H), 5.02 (s, 2H), 3.54 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 157.0 (s), 151.9 (s), 144.3 (s), 138.0 (s), 137.3 (s), 135.5 (s), 135.3 (s), 135.0 (s), 130.3 (s), 130.1 (s), 128.6 (s), 127.9 (s), 127.6 (s), 125.9 (s), 115.0 (s), 114.4 (s), 70.1 (s), 55.3 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₄H₄₀NO₂ [M+H]⁺: 494.3054, found: 494.3049.

4-((4-Aminophenyl)(2-hydroxyphenyl)methyl)-2,6-di-*tert*-butylphenol (4h)



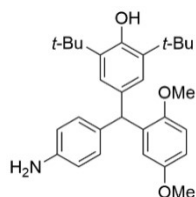
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 2:1) gave product **4h** (58.2 mg, 0.14 mmol, 72%) as a yellow oil. ¹H NMR (400 MHz, DMSO-*d*₆, 25 °C, TMS): δ = 9.21 (s, 1H), 6.95-6.99 (m, 1H), 6.81-6.82 (m, 2H), 6.74-6.77 (m, 2H), 6.67-6.72 (m, 4H), 6.44-6.46 (m, 2H), 5.52 (s, 1H), 4.83 (s, 2H), 1.29 (s, 18H); ¹³C NMR (100 MHz, DMSO-*d*₆, 25 °C, TMS): δ = 154.9 (s), 152.0 (s), 146.8 (s), 138.9 (s), 135.9 (s), 132.2 (s), 132.1 (s), 130.1 (s), 129.9 (s), 127.0 (s), 125.6 (s), 118.9 (s), 115.4 (s), 114.1 (s), 48.4 (s), 34.9 (s), 30.9 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₄NO₂ [M+H]⁺: 404.2585, found: 404.2578.

4-((4-Aminophenyl)(4-hydroxy-3-methoxyphenyl)methyl)-2,6-di-*tert*-butylphenol (4i)



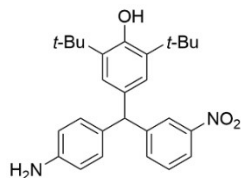
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 2:1) gave product **4i** (66.1 mg, 0.15 mmol, 76%) as a yellow solid. mp: 114.2-114.5 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.79-6.91 (m, 5H), 6.47-6.74 (m, 5H), 5.25 (s, 1H), 5.05 (s, 1H), 3.76 (s, 3H), 3.69 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 146.2 (s), 144.2 (s), 143.7 (s), 137.5 (s), 135.5 (s), 135.3 (s), 134.9 (s), 130.1 (s), 125.9 (s), 122.1 (s), 115.1 (s), 113.9 (s), 112.1 (s), 55.9 (s), 55.7 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₆NO₃ [M+H]⁺: 434.2690, found: 434.2685.

4-((4-Aminophenyl)(2,5-dimethoxyphenyl)methyl)-2,6-di-*tert*-butylphenol (4j)



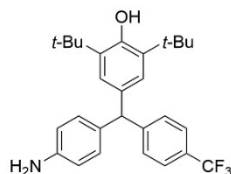
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4j** (79.3 mg, 0.17 mmol, 83%) as a yellow solid. mp: 126.8-127.2 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.86-6.89 (m, 4H), 6.76-6.78 (m, 1H), 6.67-6.70 (m, 1H), 6.56-6.58 (m, 2H), 6.49-6.50 (m, 1H), 5.67 (s, 1H), 5.03 (s, 1H), 3.66 (s, 3H), 3.63 (s, 3H), 3.53 (s, 3H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.3 (s), 151.9 (s), 151.6 (s), 144.1 (s), 135.7 (s), 135.2 (s), 134.7 (s), 134.4 (s), 130.1 (s), 126.1 (s), 117.0 (s), 115.0 (s), 112.0 (s), 110.9 (s), 56.6 (s), 55.6 (s), 48.8 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₈NO₃ [M+H]⁺: 478.2847, found: 478.2840.

4-((4-Aminophenyl)(3-nitrophenyl)methyl)-2,6-di-*tert*-butylphenol (4k)



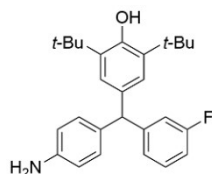
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4k** (50.2 mg, 0.12 mmol, 58%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 8.01-8.06 (m, 2H), 7.40-7.47 (m, 2H), 6.86-6.88 (m, 4H), 6.62-6.64 (m, 2H), 5.42 (s, 1H), 5.12 (s, 1H), 3.64 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.4 (s), 147.8 (s), 144.9 (s), 144.6 (s), 135.8 (s), 135.5 (s), 133.4 (s), 133.2 (s), 130.1 (s), 128.9 (s), 125.8 (s), 124.2 (s), 121.2 (s), 115.2 (s), 55.7 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃N₂O₃ [M+H]⁺: 433.2486, found: 433.2482.

4-((4-Aminophenyl)(4-(trifluoromethyl)phenyl)methyl)-2,6-di-*tert*-butylphenol (4l)



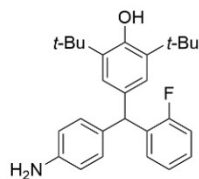
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4l** (55.6 mg, 0.12 mmol, 61%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.49-7.51 (m, 2H), 7.21-7.23 (m, 2H), 6.85-6.87 (m, 4H), 6.61-6.63 (m, 2H), 5.37 (s, 1H), 5.09 (s, 1H), 3.61 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.2 (s), 149.7 (s), 144.7 (s), 135.6 (s), 134.0 (s), 133.8 (s), 130.2 (s), 129.6 (s), 127.1 (d, *J*(C,F) = 239.4 Hz), 125.9 (s), 125.0 (d, *J*(C,F) = 11.1 Hz), 124.9 (d, *J*(C,F) = 3.8 Hz), 115.1 (s), 55.9 (s), 34.4 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -62.2 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₃F₃NO [M+H]⁺: 456.2509, found: 456.2503.

4-((4-Aminophenyl)(3-fluorophenyl)methyl)-2,6-di-*tert*-butylphenol (4m)



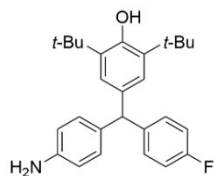
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4m** (57.7 mg, 0.14 mmol, 71%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.18-7.23 (m, 2H), 6.78-6.91 (m, 7H), 6.60-6.62 (m, 2H), 5.32 (s, 1H), 5.07 (s, 1H), 3.58 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 162.8 (d, *J*(C,F) = 243.4 Hz), 152.1 (s), 148.3 (d, *J*(C,F) = 6.5 Hz), 135.5 (s), 134.1 (s), 130.1 (s), 129.3 (d, *J*(C,F) = 8.3 Hz), 125.9 (s), 125.0 (d, *J*(C,F) = 2.7 Hz), 116.3 (s), 116.1 (s), 115.1 (s), 112.8 (s), 112.6 (s), 55.8 (s), 34.4 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -113.9 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃FNO [M+H]⁺: 406.2541, found: 406.2537.

4-((4-Aminophenyl)(2-fluorophenyl)methyl)-2,6-di-*tert*-butylphenol (4n)



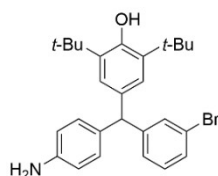
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4n** (60.1 mg, 0.15 mmol, 74%) as a yellow solid. mp: 132.2-132.4 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.14-7.19 (m, 1H), 6.93-7.04 (m, 3H), 6.87-6.89 (m, 4H), 6.59-6.62 (m, 2H), 5.63 (s, 1H), 5.06 (s, 1H), 3.57 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 160.7 (d, *J*(C,F) = 244.5 Hz), 144.5 (s), 135.4 (s), 133.7 (s), 133.3 (s), 132.5 (d, *J*(C,F) = 14.3 Hz), 130.8 (d, *J*(C,F) = 4.2 Hz), 130.0 (s), 127.6 (d, *J*(C,F) = 8.1 Hz), 125.9 (s), 123.6 (d, *J*(C,F) = 3.5 Hz), 115.3 (s), 115.1 (s), 115.0 (s), 48.3 (s), 34.3 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -116.8 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃FNO [M+H]⁺: 406.2541, found: 406.2535.

4-((4-Aminophenyl)(4-fluorophenyl)methyl)-2,6-di-*tert*-butylphenol (4o)



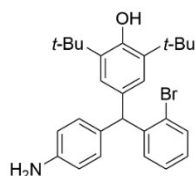
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4o** (58.8 mg, 0.14 mmol, 72%) as a white solid. mp: 126.6-126.9 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.03-7.07 (m, 2H), 6.90-6.96 (m, 2H), 6.85-6.87 (m, 4H), 6.59-6.62 (m, 2H), 5.31 (s, 1H), 5.06 (s, 1H), 3.58 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 161.2 (d, *J*(C,F) = 242.4 Hz), 144.4 (s), 141.2 (s), 135.4 (s), 134.9 (s), 134.6 (s), 130.7 (d, *J*(C,F) = 7.8 Hz), 130.1 (s), 125.9 (s), 115.1 (s), 114.9 (s), 114.7 (s), 55.3 (s), 34.4 (s), 30.4 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -117.8 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃FNO [M+H]⁺: 406.2541, found: 406.2533.

4-((4-Aminophenyl)(3-bromophenyl)methyl)-2,6-di-*tert*-butylphenol (4p)



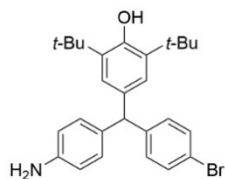
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4p** (75.5 mg, 0.16 mmol, 81%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.28-7.31 (m, 2H), 7.10-7.14 (m, 1H), 7.02-7.03 (m, 1H), 6.85-6.87 (m, 4H), 6.60-6.62 (m, 2H), 5.28 (s, 1H), 5.08 (s, 1H), 3.60 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.2 (s), 147.9 (s), 144.6 (s), 135.5 (s), 134.2 (s), 133.9 (s), 132.4 (s), 130.1 (s), 129.6 (s), 129.0 (s), 128.0 (s), 125.9 (s), 122.3 (s), 115.1 (s), 55.7 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃BrNO [M+H]⁺: 466.1741, found: 466.1735.

4-((4-Aminophenyl)(2-bromophenyl)methyl)-2,6-di-*tert*-butylphenol (4q)



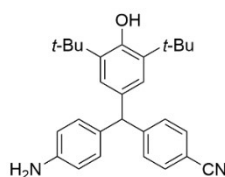
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4q** (74.6 mg, 0.16 mmol, 80%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.52-7.55 (m, 1H), 7.14-7.21 (m, 1H), 7.02-7.06 (m, 1H), 6.96-6.98 (m, 1H), 6.84-6.86 (m, 3H), 6.68-6.78 (m, 1H), 6.60-6.62 (m, 2H), 5.72 (s, 1H), 5.02 (s, 1H), 3.59 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 135.4 (s), 133.5 (s), 131.2 (s), 130.4 (s), 127.5 (s), 127.0 (s), 126.2 (s), 125.5 (s), 125.3 (s), 115.06 (s), 55.0 (s), 34.3 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃BrNO [M+H]⁺: 466.1741, found: 466.1737.

4-((4-Aminophenyl)(4-bromophenyl)methyl)-2,6-di-*tert*-butylphenol (4r)



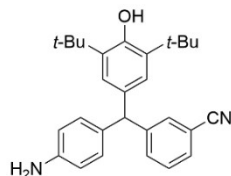
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4r** (77.4 mg, 0.17 mmol, 83%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.35-7.38 (m, 2H), 6.96-6.98 (m, 2H), 6.86-6.88 (m, 4H), 6.65-6.67 (m, 2H), 5.27 (s, 1H), 5.08 (s, 1H), 3.66 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 144.5 (s), 135.5 (s), 134.0 (s), 131.2 (s), 131.1 (s), 130.16 (s), 125.9 (s), 119.7 (s), 115.8 (s), 55.5 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃BrNO [M+H]⁺: 466.1741, found: 466.1733.

4-((4-Aminophenyl)(3,5-di-*tert*-butyl-4-hydroxyphenyl)methyl)benzonitrile (4s)



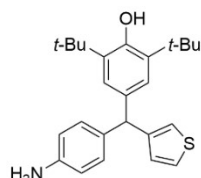
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4s** (55.3 mg, 0.13 mmol, 67%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.45-7.49 (m, 1H), 7.33-7.39 (m, 3H), 6.83-6.85 (m, 4H), 6.62-6.64 (m, 2H), 5.35 (s, 1H), 5.12 (s, 1H), 3.66 (s, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.4 (s), 147.2 (s), 144.8 (s), 135.7 (s), 133.9 (s), 133.3 (s), 132.8 (s), 130.1 (s), 129.8 (s), 128.8 (s), 125.8 (s), 115.2 (s), 112.0 (s), 55.6 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₃N₂O [M+H]⁺: 413.2588, found: 413.2583.

3-((4-Aminophenyl)(3,5-di-*tert*-butyl-4-hydroxyphenyl)methyl)benzonitrile (4t)



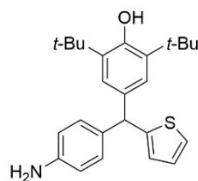
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4t** (57.1 mg, 0.14 mmol, 69%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.53-7.56 (m, 2H), 7.20-7.22 (m, 2H), 6.83-6.85 (m, 4H), 6.61-6.63 (m, 2H), 5.36 (s, 1H), 5.11 (s, 1H), 3.65 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.3 (s), 151.3 (s), 144.8 (s), 135.7 (s), 133.4 (s), 133.3 (s), 132.0 (s), 131.9 (s), 130.2 (s), 130.1 (s), 125.8 (s), 125.7 (s), 119.2 (s), 115.2 (s), 109.7 (s), 56.1 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₃N₂O [M+H]⁺: 413.2588, found: 413.2581.

4-((4-Aminophenyl)(thiophen-3-yl)methyl)-2,6-di-*tert*-butylphenol (4u)



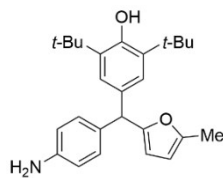
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4u** (60.2 mg, 0.15 mmol, 76%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.21-7.23 (m, 1H), 6.92-6.94 (m, 4H), 6.85-6.86 (m, 1H), 6.70-6.71 (m, 1H), 6.60-6.62 (m, 2H), 5.28 (s, 1H), 5.05 (s, 1H), 3.57 (s, 2H), 1.37 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 146.6 (s), 144.4 (s), 135.4 (s), 135.0 (s), 134.8 (s), 129.7 (s), 129.0 (s), 125.5 (s), 125.0 (s), 122.1 (s), 115.1 (s), 51.8 (s), 34.3 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₅H₃₂NOS [M+H]⁺: 394.2200, found: 394.2196.

4-((4-Aminophenyl)(thiophen-2-yl)methyl)-2,6-di-*tert*-butylphenol (4v)



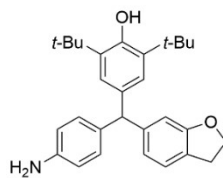
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4v** (56.7 mg, 0.14 mmol, 72%) as a yellow solid. mp: 104.5-104.8 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.15-7.16 (m, 1H), 6.98-7.01 (m, 4H), 6.89-6.91 (m, 1H), 6.66-6.67 (m, 1H), 6.60-6.62 (m, 2H), 5.45 (s, 1H), 5.08 (s, 1H), 3.58 (s, 2H), 1.37 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.3 (s), 150.1 (s), 144.7 (s), 135.4 (s), 135.1 (s), 134.7 (s), 129.6 (s), 126.4 (s), 125.8 (s), 125.4 (s), 124.1 (s), 115.1 (s), 51.5 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₅H₃₂NOS [M+H]⁺: 394.2200, found: 394.2193.

4-((4-Aminophenyl)(5-methylfuran-2-yl)methyl)-2,6-di-*tert*-butylphenol (4w)



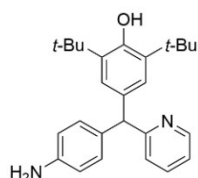
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **4w** (54.1 mg, 0.14 mmol, 69%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.94-6.97 (m, 4H), 6.59-6.61 (m, 2H), 5.84-5.85 (m, 1H), 5.70-5.71 (m, 1H), 5.17 (s, 1H), 3.42 (s, 2H), 2.23 (s, 3H), 1.37 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 156.4 (s), 152.3 (s), 151.0 (s), 144.6 (s), 135.4 (s), 133.1 (s), 133.0 (s), 129.6 (s), 125.3 (s), 115.2 (s), 108.5 (s), 105.8 (s), 50.1 (s), 34.4 (s), 30.4 (s), 13.7 (s). HRMS (ESI) *m/z*: calcd. for C₂₆H₃₄NO₂ [M+H]⁺: 392.2585, found: 392.2579.

4-((4-Aminophenyl)(2,3-dihydrobenzofuran-6-yl)methyl)-2,6-di-*tert*-butylphenol (4x)



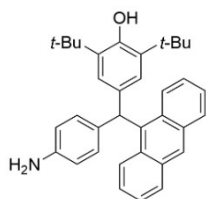
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4x** (62.8 mg, 0.15 mmol, 73%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.87-6.94 (m, 5H), 6.81-6.83 (m, 1H), 6.65-6.67 (m, 1H), 6.58-6.60 (m, 2H), 5.26 (s, 1H), 5.04 (s, 1H), 4.52 (t, *J* = 8.7, 2H), 3.54 (s, 2H), 3.12 (t, *J* = 8.6, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 158.2 (s), 151.9 (s), 144.2 (s), 137.7 (s), 135.7 (s), 135.3 (s), 135.2 (s), 130.1 (s), 128.8 (s), 126.7 (s), 125.9 (s), 125.8 (s), 115.0 (s), 108.6 (s), 71.2 (s), 55.5 (s), 34.4 (s), 30.4 (s), 29.9 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₆NO₂ [M+H]⁺: 430.2741, found: 430.2735.

4-((4-Aminophenyl)(pyridin-2-yl)methyl)-2,6-di-*tert*-butylphenol (4y)



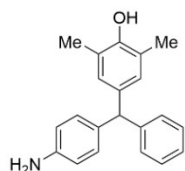
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4y** (58.3 mg, 0.15 mmol, 75%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 8.57-8.58 (m, 1H), 7.55-7.60 (m, 1H), 7.05-7.11 (m, 2H), 6.93-6.95 (m, 4H), 6.61-6.63 (m, 2H), 5.49 (s, 1H), 5.06 (s, 1H), 3.59 (s, 2H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 164.8 (s), 152.2 (s), 149.3 (s), 144.6 (s), 136.2 (s), 135.4 (s), 133.7 (s), 133.5 (s), 130.1 (s), 126.0 (s), 123.5 (s), 121.0 (s), 115.2 (s), 58.6 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₆H₃₃N₂O [M+H]⁺: 389.2588, found: 389.2581.

4-((4-Aminophenyl)(anthracen-9-yl)methyl)-2,6-di-*tert*-butylphenol (4aa)



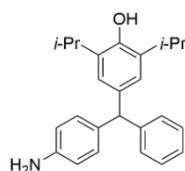
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4aa** (54.7 mg, 0.11 mmol, 56%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.73-7.74 (m, 1H), 7.34-7.42 (m, 3H), 7.15-7.30 (m, 5H), 6.97-7.01 (m, 3H), 6.92 (s, 1H), 6.48-6.50 (m, 2H), 5.19 (s, 1H), 5.11 (s, 1H), 3.49 (s, 2H), 1.32 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.1 (s), 144.5 (s), 141.2 (s), 139.3 (s), 139.1 (s), 135.4 (s), 134.7 (s), 134.2 (s), 128.9 (s), 128.5 (s), 128.4 (s), 127.9 (s), 127.3 (s), 126.6 (s), 126.4 (s), 125.2 (s), 123.7 (s), 115.2 (s), 51.6 (s), 34.3 (s), 30.2 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₆NO₃ [M+H]⁺: 488.2948, found: 488.2943.

4-((4-Aminophenyl)(phenyl)methyl)-2,6-dimethylphenol (4ab)



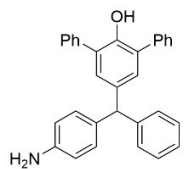
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4ab** (51.7 mg, 0.17 mmol, 85%) as a red oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.22-7.26 (m, 2H), 7.16-7.18 (m, 1H), 7.08-7.10 (m, 2H), 6.86-6.88 (m, 2H), 6.69 (s, 2H), 6.56-6.58 (m, 2H), 5.30 (s, 1H), 3.99 (s, 3H), 2.13 (s, 6H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 150.6 (s), 145.1 (s), 144.4 (s), 136.1 (s), 134.9 (s), 130.3 (s), 129.5 (s), 129.4 (s), 128.3 (s), 126.1 (s), 122.9 (s), 115.3 (s), 55.4 (s), 16.2 (s). HRMS (ESI) *m/z*: calcd. for C₂₁H₂₂NO [M+H]⁺: 304.1696, found: 304.1692.

4-((4-Aminophenyl)(phenyl)methyl)-2,6-diisopropylphenol (4ac)



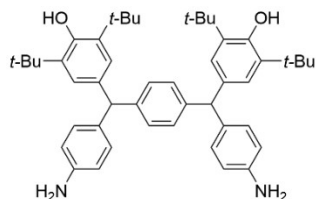
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 2:1) gave product **4ac** (59.8 mg, 0.17 mmol, 83%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.22-7.26 (m, 2H), 7.14-7.18 (m, 1H), 7.09-7.11 (m, 2H), 6.87-6.89 (m, 2H), 6.78 (s, 2H), 6.57-6.61 (m, 2H), 5.36 (s, 1H), 3.92 (s, 3H), 3.03-3.14 (m, 2H), 1.16 (d, *J* = 6.9, 12H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 148.3 (s), 145.3 (s), 144.3 (s), 136.1 (s), 135.1 (s), 133.3 (s), 130.2 (s), 129.4 (s), 128.1 (s), 125.9 (s), 124.6 (s), 115.2 (s), 55.9 (s), 22.8 (s), 22.7 (s). HRMS (ESI) *m/z*: calcd. for C₂₅H₃₀NO [M+H]⁺: 360.2322, found: 360.2316.

5'-((4-Aminophenyl)(phenyl)methyl)-[1,1':3',1''-terphenyl]-2'-ol (4ad)



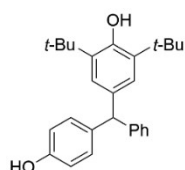
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **4ad** (52.2 mg, 0.12 mmol, 61%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.49-7.50 (m, 4H), 7.35-7.42 (m, 5H), 7.28-7.34 (m, 3H), 7.15-7.20 (m, 3H), 7.04 (s, 2H), 6.92-6.97 (m, 2H), 6.57-6.59 (m, 2H), 5.43 (s, 1H), 5.36 (s, 1H), 3.53 (s, 2H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 147.7 (s), 144.7 (s), 144.6 (s), 137.8 (s), 136.8 (s), 134.3 (s), 130.9 (s), 130.2 (s), 129.4 (s), 129.3 (s), 128.8 (s), 128.5 (s), 128.3 (s), 127.6 (s), 126.2 (s), 115.2 (s), 55.5 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₂₆NO [M+H]⁺: 428.2009, found: 428.2001.

4,4'-(1,4-Phenylenebis((4-Aminophenyl)methylene))bis(2,6-di-*tert*-butylphenol) (4ae)



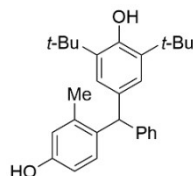
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 2:1) gave product **4ae** (91.9 mg, 0.13 mmol, 66%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 6.99 (s, 4H), 6.86-6.88 (m, 8H), 6.57-6.59 (m, 4H), 5.29 (s, 2H), 5.03 (s, 2H), 3.56 (s, 4H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 151.9 (s), 144.2 (s), 142.8 (s), 142.7 (s), 135.4 (s), 135.2 (s), 135.1 (s), 135.0 (s), 130.2 (s), 129.0 (s), 126.0 (s), 115.0 (s), 55.7 (s), 34.3 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{48}\text{H}_{61}\text{N}_2\text{O}_2$ $[\text{M}+\text{H}]^+$: 697.4728, found: 697.4722.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(phenyl)methyl)phenol (6a)



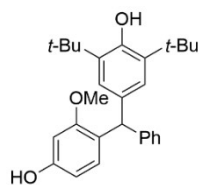
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6a** (69.2 mg, 0.18 mmol, 89%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.23-7.27 (m, 2H), 7.15-7.19 (m, 1H), 7.09-7.11 (m, 2H), 6.95-6.98 (m, 2H), 6.89 (s, 2H), 6.70-6.74 (m, 2H), 5.38 (s, 1H), 5.07 (s, 1H), 5.00 (s, 1H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 153.7 (s), 152.1 (s), 145.1 (s), 137.2 (s), 135.5 (s), 134.5 (s), 130.6 (s), 129.4 (s), 128.3 (s), 128.2 (s), 126.0 (s), 115.0 (s), 56.0 (s), 34.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{33}\text{O}_2$ $[\text{M}+\text{H}]^+$: 389.2476, found: 389.2471.

2,6-Di-*tert*-butyl-4-((4-hydroxy-2-methylphenyl)(phenyl)methyl)phenol (6b)



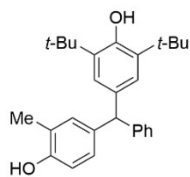
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6b** (72.5 mg, 0.18 mmol, 90%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.23-7.27 (m, 2H), 7.15-7.19 (m, 1H), 7.03-7.05 (m, 2H), 6.81 (s, 2H), 6.63-6.67 (m, 2H), 6.54-6.57 (m, 1H), 5.47 (s, 1H), 5.05 (s, 1H), 4.54 (s, 1H), 2.16 (s, 3H), 1.34 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 153.5 (s), 152.0 (s), 144.5 (s), 138.2 (s), 135.8 (s), 135.4 (s), 134.0 (s), 130.5 (s), 129.4 (s), 128.1 (s), 126.2 (s), 125.9 (s), 117.1 (s), 112.2 (s), 52.8 (s), 34.3 (s), 30.4 (s), 20.1 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{28}\text{H}_{35}\text{O}_2$ $[\text{M}+\text{H}]^+$: 403.2632, found: 403.2625.

2,6-Di-*tert*-butyl-4-((4-hydroxy-2-methoxyphenyl)(phenyl)methyl)phenol (6c)



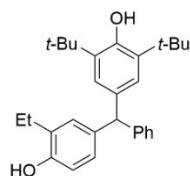
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6c** (74.6 mg, 0.18 mmol, 89%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.21-7.32 (m, 4H), 7.13-7.15 (m, 2H), 6.92 (s, 2H), 6.66-6.68 (m, 1H), 6.40-6.42 (m, 2H), 5.47 (s, 1H), 5.13 (s, 1H), 4.78 (s, 1H), 3.76 (s, 3H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 159.5 (s), 154.6 (s), 152.5 (s), 143.2 (s), 135.9 (s), 132.5 (s), 130.9 (s), 129.2 (s), 128.5 (s), 126.5 (s), 125.9 (s), 123.3 (s), 106.0 (s), 102.3 (s), 55.3 (s), 50.9 (s), 34.4 (s), 30.3 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{28}\text{H}_{35}\text{O}_3$ $[\text{M}+\text{H}]^+$: 419.2581, found: 419.2574.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3-methylphenyl)(phenyl)methyl)phenol (**6d**)



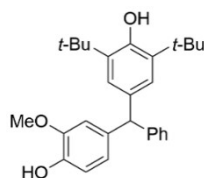
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6d** (70.2 mg, 0.17 mmol, 87%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.27 (m, 2H), 7.15-7.18 (m, 1H), 7.07-7.11 (m, 2H), 6.90 (s, 3H), 6.76-6.78 (m, 1H), 6.64-6.66 (m, 1H), 5.34 (s, 1H), 5.06 (s, 1H), 4.76 (s, 1H), 2.17 (s, 1H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.71 (s), 152.0 (s), 145.3 (s), 137.1 (s), 135.4 (s), 134.5 (s), 132.1 (s), 129.3 (s), 128.1 (s), 127.9 (s), 126.0 (s), 125.9 (s), 123.3 (s), 114.6 (s), 56.0 (s), 34.4 (s), 30.4 (s), 15.9 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₅O₂ [M+H]⁺: 403.2632, found: 403.2627.

2,6-Di-*tert*-butyl-4-((3-ethyl-4-hydroxyphenyl)(phenyl)methyl)phenol (**6e**)



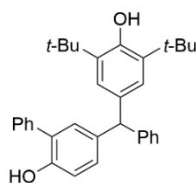
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6e** (71.2 mg, 0.17 mmol, 86%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.22-7.27 (m, 2H), 7.14-7.18 (m, 1H), 7.09-7.11 (m, 2H), 6.90 (s, 2H), 6.76-6.79 (m, 1H), 6.23-6.25 (m, 1H), 5.35 (s, 1H), 5.06 (s, 1H), 4.86 (s, 1H), 2.51-2.59 (m, 2H), 1.36 (s, 18H), 1.16 (t, *J* = 7.5, 3H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.0 (s), 151.6 (s), 145.4 (s), 137.1 (s), 135.4 (s), 134.6 (s), 130.4 (s), 129.6 (s), 129.4 (s), 128.1 (s), 127.8 (s), 126.0 (s), 125.9 (s), 114.8 (s), 56.2 (s), 34.4 (s), 30.4 (s), 23.1 (s), 14.2 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₇O₂ [M+H]⁺: 417.2789, found: 417.2783.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3-methoxyphenyl)(phenyl)methyl)phenol (**6f**)



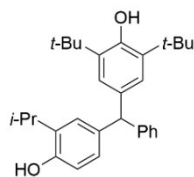
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6f** (73.7 mg, 0.18 mmol, 88%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.24-7.28 (m, 2H), 7.16-7.20 (m, 1H), 7.10-7.11 (m, 2H), 6.91 (s, 2H), 6.81-6.83 (m, 1H), 6.64-6.65 (m, 1H), 6.56-6.59 (m, 1H), 5.50 (s, 1H), 5.36 (s, 1H), 5.08 (s, 1H), 3.77 (s, 3H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 146.2 (s), 145.2 (s), 143.8 (s), 136.9 (s), 135.4 (s), 134.3 (s), 129.3 (s), 128.1 (s), 126.0 (s), 125.9 (s), 122.2 (s), 113.9 (s), 112.1 (s), 56.4 (s), 55.9 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₅O₃ [M+H]⁺: 419.2581, found: 419.2575.

5-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-[1,1'-biphenyl]-2-ol (**6g**)



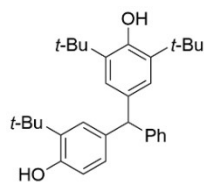
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6g** (76.3 mg, 0.16 mmol, 82%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.40-7.49 (m, 4H), 7.34-7.37 (m, 1H), 7.25-7.28 (m, 2H), 7.13-7.19 (m, 3H), 6.97-7.03 (m, 2H), 6.93 (m, 2H), 6.87-6.89 (m, 1H), 5.40 (s, 1H), 5.09 (s, 1H), 5.06 (s, 1H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 150.6 (s), 145.1 (s), 137.3 (s), 135.5 (s), 134.3 (s), 131.1 (s), 130.1 (s), 129.3 (s), 129.2 (s), 129.1 (s), 128.4 (s), 128.2 (s), 127.8 (s), 127.7 (s), 126.0 (s), 125.9 (s), 115.5 (s), 56.1 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₃H₃₇O₂ [M+H]⁺: 465.2789, found: 465.2782.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3-isopropylphenyl)(phenyl)methyl)phenol (6h)



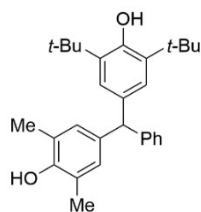
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6h** (75.9 mg, 0.18 mmol, 88%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.24-7.27 (m, 2H), 7.15-7.19 (m, 1H), 7.09-7.11 (m, 2H), 6.95-6.96 (m, 1H), 6.90 (s, 2H), 6.75-6.78 (m, 1H), 6.62-6.64 (m, 1H), 5.36 (s, 1H), 5.06 (s, 1H), 4.71 (s, 1H), 3.10-3.21 (m, 1H), 1.35 (s, 18H), 1.17 (d, *J* = 6.9, 6H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.0 (s), 150.9 (s), 145.5 (s), 137.1 (s), 135.3 (s), 134.6 (s), 133.9 (s), 129.3 (s), 128.1 (s), 127.6 (s), 127.5 (s), 126.0 (s), 125.9 (s), 114.9 (s), 56.2 (s), 34.4 (s), 30.4 (s), 27.1 (s), 22.6 (s). HRMS (ESI) *m/z*: calcd. for C₃₀H₃₉O₂ [M+H]⁺: 431.2945, found: 431.2940.

2,6-Di-*tert*-butyl-4-((3-(*tert*-butyl)-4-hydroxyphenyl)(phenyl)methyl)phenol (6i)



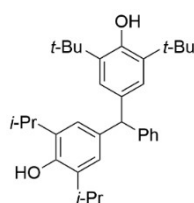
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6i** (77.4 mg, 0.17 mmol, 87%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.27 (m, 2H), 7.15-7.18 (m, 1H), 7.10-7.11 (m, 2H), 7.02-7.03 (m, 1H), 6.91 (s, 2H), 6.76-6.79 (m, 1H), 6.54-6.56 (m, 1H), 5.34 (s, 1H), 5.06 (s, 1H), 4.74 (s, 1H), 1.36 (s, 18H), 1.33 (s, 9H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.4 (s), 152.0 (s), 145.6 (s), 136.5 (s), 135.5 (s), 135.3 (s), 134.6 (s), 129.3 (s), 128.3 (s), 128.1 (s), 127.7 (s), 126.0 (s), 125.8 (s), 116.2 (s), 56.3 (s), 34.6 (s), 34.4 (s), 30.4 (s), 29.6 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₄₁O₂ [M+H]⁺: 445.3102, found: 445.3097.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3,5-dimethylphenyl)(phenyl)methyl)phenol (6j)



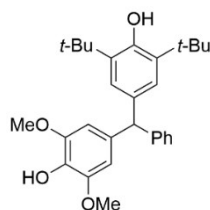
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6j** (68.4 mg, 0.16 mmol, 82%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.29 (m, 2H), 7.14-7.18 (m, 1H), 7.09-7.11 (m, 2H), 6.89-6.91 (m, 2H), 6.73 (s, 2H), 5.30 (s, 1H), 5.05 (s, 1H), 4.48 (s, 1H), 2.17 (s, 6H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.0 (s), 150.4 (s), 145.4 (s), 136.4 (s), 135.4 (s), 134.6 (s), 129.5 (s), 129.3 (s), 128.1 (s), 126.0 (s), 125.8 (s), 122.6 (s), 56.2 (s), 34.4 (s), 30.4 (s), 16.0 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₇O₂ [M+H]⁺: 417.2789, found: 417.2784.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3,5-diisopropylphenyl)(phenyl)methyl)phenol (6k)



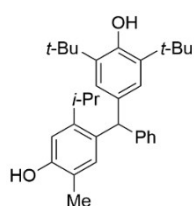
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6k** (74.7 mg, 0.16 mmol, 79%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.26 (m, 2H), 7.10-7.17 (m, 3H), 6.92 (s, 2H), 6.80 (s, 2H), 5.34 (s, 1H), 5.05 (s, 1H), 4.65 (s, 1H), 3.05-3.16 (m, 2H), 1.36 (s, 18H), 1.17 (d, *J* = 6.9 Hz, 12H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 151.9 (s), 148.1 (s), 136.4 (s), 135.3 (s), 134.8 (s), 133.1 (s), 129.3 (s), 128.0 (s), 126.0 (s), 125.8 (s), 124.6 (s), 56.6 (s), 34.4 (s), 30.4 (s), 27.3 (s), 22.8 (s). HRMS (ESI) *m/z*: calcd. for C₃₃H₄₅O₂ [M+H]⁺: 473.3415, found: 473.3410.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3,5-dimethoxyphenyl)(phenyl)methyl)phenol (**6l**)



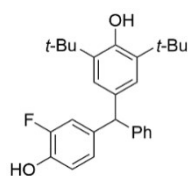
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6l** (46.7 mg, 0.41 mmol, 52%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.25-7.29 (m, 2H), 7.17-7.20 (m, 1H), 7.09-7.11 (m, 2H), 6.92 (s, 2H), 6.34 (s, 2H), 5.39 (s, 1H), 5.35 (s, 1H), 5.08 (s, 1H), 3.76 (s, 6H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 152.1 (s), 146.8 (s), 145.1 (s), 135.9 (s), 135.5 (s), 134.1 (s), 132.9 (s), 129.3 (s), 128.1 (s), 126.0 (s), 125.9 (s), 106.4 (s), 56.8 (s), 56.2 (s), 34.4 (s), 30.4 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{29}\text{H}_{37}\text{O}_4$ $[\text{M}+\text{H}]^+$: 449.2687, found: 449.2681.

4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-5-isopropyl-2-methylphenol (**6m**)



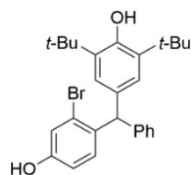
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6m** (73.9 mg, 0.17 mmol, 83%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.22-7.26 (m, 2H), 7.14-7.18 (m, 1H), 7.06-7.07 (m, 2H), 6.82 (s, 2H), 6.79 (s, 1H), 6.55 (s, 1H), 5.62 (s, 1H), 5.03 (s, 1H), 4.58 (s, 1H), 3.09-3.19 (m, 1H), 2.10 (s, 3H), 1.34 (s, 18H), 1.03 (dd, J = 8.7, 6.8 Hz, 6H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 152.2 (s), 151.9 (s), 146.3 (s), 145.3 (s), 135.3 (s), 134.7 (s), 133.8 (s), 132.3 (s), 129.6 (s), 128.0 (s), 126.2 (s), 125.8 (s), 120.1 (s), 112.1 (s), 51.8 (s), 34.3 (s), 30.4 (s), 28.4 (s), 24.0 (s), 15.5 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{31}\text{H}_{41}\text{O}_2$ $[\text{M}+\text{H}]^+$: 445.3102, found: 445.3098.

2,6-Di-*tert*-butyl-4-((3-fluoro-4-hydroxyphenyl)(phenyl)methyl)phenol (**6n**)



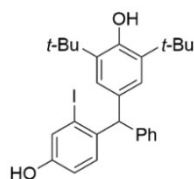
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6n** (70.0 mg, 0.17 mmol, 86%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.26-7.29 (m, 2H), 7.18-7.21 (m, 1H), 7.08-7.10 (m, 2H), 6.87-6.92 (m, 3H), 6.76-6.83 (m, 2H), 5.36 (s, 1H), 5.18 (s, 1H), 5.09 (s, 1H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 152.2 (s), 150.8 (d, $J(\text{C},\text{F})$ = 235.6 Hz), 144.4 (s), 141.6 (d, $J(\text{C},\text{F})$ = 8.1 Hz), 138.1 (d, $J(\text{C},\text{F})$ = 5.1 Hz), 135.6 (s), 133.8 (s), 129.3 (s), 128.2 (s), 126.2 (s), 125.9 (s), 125.6 (d, $J(\text{C},\text{F})$ = 3.3 Hz), 116.7 (d, $J(\text{C},\text{F})$ = 2.0 Hz), 116.4 (d, $J(\text{C},\text{F})$ = 18.5 Hz), 55.9 (s), 34.4 (s), 30.3 (s); ^{19}F NMR (376 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = -141.0 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{32}\text{FO}_2$ $[\text{M}+\text{H}]^+$: 407.2381, found: 407.2373.

4-((2-Bromo-4-hydroxyphenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (**6o**)



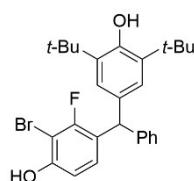
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6o** (82.2 mg, 0.18 mmol, 88%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.25-7.28 (m, 2H), 7.17-7.21 (m, 1H), 7.05-7.10 (m, 3H), 6.80-6.83 (m, 3H), 6.68-6.71 (m, 1H), 5.75 (s, 1H), 5.08 (s, 1H), 4.87 (s, 1H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 154.3 (s), 152.14 (s), 136.5 (s), 135.5 (s), 133.3 (s), 131.8 (s), 130.8 (s), 129.5 (s), 128.1 (s), 126.2 (s), 126.1 (s), 125.5 (s), 124.0 (s), 119.7 (s), 118.8 (s), 114.4 (s), 114.2 (s), 55.0 (s), 34.3 (s), 30.3 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{32}\text{BrO}_2$ $[\text{M}+\text{H}]^+$: 467.1581, found: 467.1576.

2,6-Di-*tert*-butyl-4-((4-hydroxy-2-iodophenyl)(phenyl)methyl)phenol (**6p**)



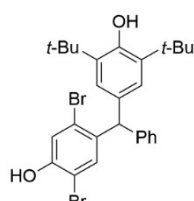
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6p** (89.6 mg, 0.17 mmol, 87%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.36-7.37 (m, 1H), 7.18-7.28 (m, 3H), 7.05-7.07 (m, 2H), 6.73-6.83 (m, 4H), 5.63 (s, 1H), 5.08 (s, 1H), 4.73 (s, 1H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 154.1 (s), 152.1 (s), 143.6 (s), 139.7 (s), 135.5 (s), 133.5 (s), 130.9 (s), 129.6 (s), 128.1 (s), 126.3 (s), 126.1 (s), 115.2 (s), 114.9 (s), 102.2 (s), 59.6 (s), 34.3 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₂IO₂ [M+H]⁺: 515.1442, found: 515.1437.

3-Bromo-4-((3,5-di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-2-fluorophenol (**6q**)



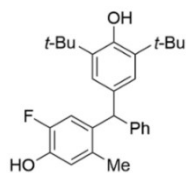
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6q** (72.8 mg, 0.15 mmol, 75%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.24-7.29 (m, 2H), 7.16-7.22 (m, 1H), 7.07-7.09 (m, 2H), 6.87 (s, 2H), 6.79-6.83 (m, 1H), 6.71-6.76 (m, 1H), 5.66 (s, 1H), 5.53 (s, 1H), 5.11 (s, 1H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 156.8 (d, *J*(C,F) = 246.0 Hz), 152.3 (s), 152.0 (s), 143.1 (s), 135.6 (s), 132.4 (s), 130.0 (d, *J*(C,F) = 5.5 Hz), 129.1 (s), 128.3 (s), 126.4 (s), 125.9 (s), 125.1 (d, *J*(C,F) = 15.6 Hz), 110.7 (d, *J*(C,F) = 3.4 Hz), 98.3 (d, *J*(C,F) = 24.1 Hz), 49.0 (s), 34.4 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -108.1 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₁BrFO₂ [M+H]⁺: 485.1486, found: 485.1480.

2,5-Dibromo-4-((3,5-di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)phenol (**6r**)



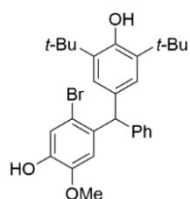
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6r** (86.4 mg, 0.16 mmol, 79%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.17-7.31 (m, 4H), 7.03-7.07 (m, 3H), 6.82 (s, 2H), 5.71 (s, 1H), 5.44 (s, 1H), 5.12 (s, 1H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.4 (s), 151.1 (s), 142.9 (s), 138.0 (s), 135.6 (s), 133.5 (s), 132.6 (s), 129.4 (s), 128.3 (s), 126.4 (s), 126.1 (s), 125.0 (s), 120.2 (s), 109.1 (s), 55.1 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₁Br₂O₂ [M+H]⁺: 547.0665, found: 547.0661.

4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-2-fluoro-5-methylphenol (**6s**)



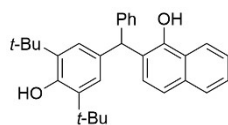
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6s** (69.2 mg, 0.16 mmol, 82%) as a yellow solid. mp: 137.8-138.2 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.27 (m, 2H), 7.16-7.20 (m, 1H), 7.03-7.04 (m, 2H), 6.77-6.80 (m, 3H), 6.51-6.54 (m, 1H), 5.44 (s, 1H), 5.08 (s, 1H), 5.01 (s, 1H), 2.11 (s, 3H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.2 (s), 149.1 (d, *J*(C,F) = 232.5 Hz), 143.8 (s), 141.1 (d, *J*(C,F) = 14.4 Hz), 136.3 (d, *J*(C,F) = 4.8 Hz), 135.6 (s), 133.5 (s), 133.1 (d, *J*(C,F) = 3.5 Hz), 129.4 (s), 128.2 (s), 126.2 (s), 126.1 (s), 118.8 (d, *J*(C,F) = 1.5 Hz), 116.5 (d, *J*(C,F) = 19.0 Hz), 52.9 (s), 34.4 (s), 30.4 (s), 19.4 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -144.8 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₄FO₂ [M+H]⁺: 421.2538, found: 421.2531.

5-Bromo-4-((3,5-di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-2-methoxyphenol (**6t**)



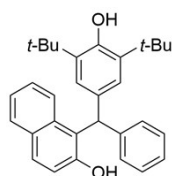
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **6t** (82.5 mg, 0.17 mmol, 83%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.25-7.28 (m, 2H), 7.17-7.21 (m, 1H), 7.13 (s, 1H), 7.06-7.08 (m, 2H), 6.89 (s, 2H), 6.45 (s, 1H), 5.74 (s, 1H), 5.51 (s, 1H), 5.10 (s, 1H), 3.64 (s, 3H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.2 (s), 145.7 (s), 144.5 (s), 143.7 (s), 135.6 (s), 133.5 (s), 133.3 (s), 129.3 (s), 128.2 (s), 126.2 (s), 126.1 (s), 118.6 (s), 115.9 (s), 113.3 (s), 55.9 (s), 55.4 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₄BrO₃ [M+H]⁺: 497.1686, found: 497.1681.

2-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)naphthalen-1-ol (**6u**)



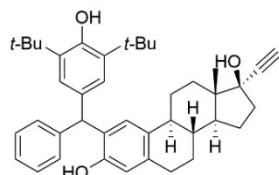
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6u** (44.8 mg, 0.10 mmol, 51%) as a colorless oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 8.20-8.22 (m, 1H), 7.97-7.99 (m, 1H), 7.38-7.45 (m, 2H), 7.22-7.25 (m, 2H), 7.10-7.18 (m, 3H), 6.88 (m, 2H), 6.74-6.76 (m, 1H), 6.68-6.70 (m, 1H), 6.08 (s, 1H), 5.26 (s, 1H), 5.06 (s, 1H), 1.33 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.1 (s), 150.2 (s), 144.8 (s), 135.5 (s), 134.3 (s), 133.5 (s), 133.1 (s), 129.5 (s), 128.2 (s), 127.4 (s), 126.5 (s), 126.4 (s), 126.0 (s), 124.8 (s), 124.7 (s), 124.5 (s), 122.1 (s), 107.9 (s), 52.7 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₃₅O₂ [M+H]⁺: 439.2632, found: 439.2628.

1-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)naphthalen-2-ol (**6v**)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6v** (56.2 mg, 0.13 mmol, 64%) as a colorless oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 8.01-8.04 (m, 1H), 7.77-7.79 (m, 1H), 7.71-7.73 (m, 1H), 7.40-7.44 (m, 1H), 7.29-7.33 (m, 3H), 7.22-7.27 (m, 3H), 7.06-7.08 (m, 1H), 7.01 (s, 2H), 6.29 (s, 1H), 5.40 (s, 1H), 5.20 (s, 1H), 1.32 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.2 (s), 153.1 (s), 142.1 (s), 136.8 (s), 133.5 (s), 131.8 (s), 129.6 (s), 129.5 (s), 129.0 (s), 128.9 (s), 128.7 (s), 126.9 (s), 126.7 (s), 125.7 (s), 123.1 (s), 122.9 (s), 120.3 (s), 120.0 (s), 48.6 (s), 34.5 (s), 30.2 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₃₅O₂ [M+H]⁺: 439.2632, found: 439.2625.

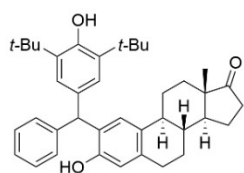
(8*R*,9*S*,13*S*,14*S*,17*R*)-2-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-17-ethynyl-13-methyl-7,8,9,11,12,13,14,15,16,17-decahydro-6*H*-cyclopenta[*a*]phenanthrene-3,17-diol (**6w**)



According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 2:1) gave product **6w** (77.9 mg, 0.13 mmol, 66%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.28-7.31 (m, 2H), 7.15-7.24 (m, 3H), 6.93-6.97 (m, 2H), 6.74-6.75 (m, 1H), 6.53 (s, 1H), 5.46-5.53 (s, 1H), 5.10-5.11 (s, 1H), 4.52-4.54 (s, 1H), 2.76-2.82 (m, 2H), 2.56-2.57 (m, 1H), 2.28-2.35 (m, 1H), 2.10-2.19 (m, 1H), 1.97-2.04 (m, 2H), 1.73-1.90 (m, 3H), 1.60-1.70 (m, 2H), 1.43 (s, 1H), 1.36-1.37 (s, 18H), 0.85-0.90 (m, 5H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.5 (s), 152.4 (s), 151.5 (s), 151.4 (s), 143.4 (s), 143.3 (s), 136.3 (s), 136.2 (s), 135.9 (s), 135.8 (s), 132.7 (s), 132.6 (s), 132.2 (s), 132.1 (s), 129.2 (s), 129.1 (s),

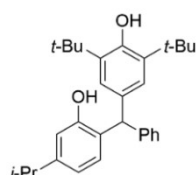
128.5 (s), 128.1 (s), 127.8 (s), 127.7 (s), 127.5 (s), 126.5 (s), 126.0 (s), 116.2 (s), 116.1 (s), 87.6 (s), 77.9 (s), 74.0 (s), 51.8 (s), 51.3 (s), 49.5 (s), 47.2 (s), 47.1 (s), 43.6 (s), 43.5 (s), 39.4 (s), 39.0 (s), 34.4 (s), 32.7 (s), 30.3 (s), 30.2 (s), 29.2 (s), 27.3 (s), 27.2 (s), 26.9 (s), 26.4 (s), 26.3 (s), 22.8 (s), 12.8 (s), 12.7 (s). HRMS (ESI) m/z : calcd. for $C_{41}H_{51}O_3$ $[M+H]^+$: 591.3833, found: 591.3828.

(8*R*,9*S*,13*S*,14*S*)-2-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-3-hydroxy-13-methyl-6,7,8,9,11,12,13,14,15,16-decahydro-17*H*-cyclopenta[*a*]phenanthren-17-one (6x)



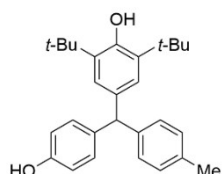
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 2:1) gave product **6x** (99.4 mg, 0.18 mmol, 88%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.14-7.31 (m, 5H), 6.93-6.97 (m, 2H), 6.75-6.76 (m, 1H), 6.54 (s, 1H), 5.48-5.54 (s, 1H), 5.11-5.12 (s, 1H), 4.76-4.80 (s, 1H), 2.82-2.86 (m, 2H), 2.44-2.51 (m, 1H), 1.95-2.18 (m, 4H), 1.82-1.86 (m, 1H), 1.52-1.63 (m, 2H), 1.35-1.36 (s, 18H), 0.78-0.97 (m, 6H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 152.5 (s), 152.4 (s), 151.7 (s), 151.6 (s), 143.4 (s), 143.3 (s), 136.0 (s), 135.9 (s), 135.8 (s), 135.7 (s), 132.7 (s), 132.5 (s), 131.7 (s), 131.5 (s), 129.2 (s), 129.0 (s), 128.5 (s), 128.3 (s), 128.0 (s), 127.6 (s), 127.5 (s), 126.5 (s), 126.0 (s), 125.9 (s), 116.3 (s), 116.2 (s), 51.7 (s), 51.2 (s), 50.4 (s), 48.1 (s), 48.0 (s), 44.1 (s), 44.0 (s), 38.4 (s), 38.3 (s), 35.9 (s), 34.4 (s), 31.5 (s), 30.4 (s), 30.3 (s), 29.1 (s), 29.0 (s), 27.0 (s), 26.6 (s), 26.5 (s), 25.9 (s), 25.8 (s), 21.6 (s), 13.9 (s), 13.8 (s). HRMS (ESI) m/z : calcd. for $C_{39}H_{49}O_3$ $[M+H]^+$: 565.3677, found: 565.3671.

2,6-Di-*tert*-butyl-4-((2-hydroxy-4-isopropylphenyl)(phenyl)methyl)phenol (6ae)



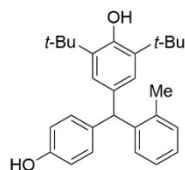
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **6ae** (69.8 mg, 0.16 mmol, 81%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.28-7.31 (m, 2H), 7.19-7.25 (m, 1H), 7.13-7.17 (m, 2H), 6.93 (m, 2H), 6.67-6.73 (m, 3H), 5.50 (s, 1H), 5.12 (s, 1H), 4.63 (s, 1H), 2.78-2.88 (m, 3H), 1.35 (s, 18H), 1.22 (d, J = 6.9, 6H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 153.4 (s), 152.5 (s), 149.0 (s), 143.2 (s), 135.9 (s), 132.4 (s), 130.1 (s), 129.3 (s), 128.5 (s), 128.1 (s), 126.6 (s), 126.0 (s), 118.7 (s), 114.3 (s), 51.2 (s), 34.4 (s), 33.6 (s), 30.4 (s), 24.0 (s). HRMS (ESI) m/z : calcd. for $C_{30}H_{39}O_2$ $[M+H]^+$: 431.2945, found: 431.2940.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(*p*-tolyl)methyl)phenol (7a)



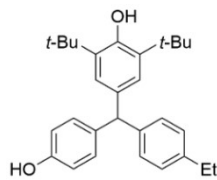
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **7a** (73.3 mg, 0.18 mmol, 91%) as a yellow oil. 1H NMR (400 MHz, $CDCl_3$, 25 °C, TMS): δ = 7.05-7.07 (m, 2H), 6.95-6.99 (m, 4H), 6.89 (s, 2H), 6.70-6.73 (m, 2H), 5.33 (s, 1H), 5.05 (s, 1H), 4.79 (s, 1H), 2.31 (s, 3H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, $CDCl_3$, 25 °C, TMS): δ = 153.7 (s), 152.0 (s), 142.1 (s), 137.5 (s), 135.4 (s), 135.3 (s), 134.6 (s), 130.5 (s), 129.2 (s), 128.8 (s), 125.9 (s), 114.9 (s), 55.7 (s), 34.4 (s), 30.4 (s), 21.0 (s). HRMS (ESI) m/z : calcd. for $C_{28}H_{35}O_2$ $[M+H]^+$: 403.2632, found: 403.2625.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(*o*-tolyl)methyl)phenol (7b)



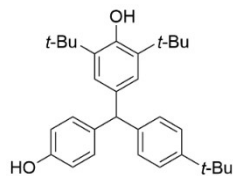
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **7b** (71.7 mg, 0.18 mmol, 89%) as a yellow solid. mp: 156.5-156.7 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.06-7.14 (m, 3H), 6.90-6.92 (m, 2H), 6.80-6.82 (m, 3H), 6.70-6.74 (m, 2H), 5.49 (s, 1H), 5.05 (s, 1H), 4.76 (s, 1H), 2.21 (s, 3H), 1.34 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.6 (s), 152.0 (s), 143.5 (s), 136.5 (s), 136.4 (s), 135.4 (s), 134.0 (s), 130.6 (s), 130.2 (s), 129.2 (s), 126.2 (s), 126.0 (s), 125.6 (s), 115.0 (s), 52.6 (s), 34.3 (s), 30.4 (s), 20.0 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₅O₂ [M+H]⁺: 403.2632, found: 403.2627.

2,6-Di-*tert*-butyl-4-((4-ethylphenyl)(4-hydroxyphenyl)methyl)phenol (7c)



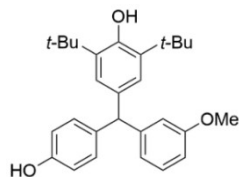
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **7c** (73.4 mg, 0.18 mmol, 88%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.08-7.10 (m, 2H), 6.92-7.02 (m, 4H), 6.90 (s, 2H), 6.71-6.73 (m, 2H), 5.33 (s, 1H), 5.05 (s, 1H), 4.86 (s, 1H), 2.58-2.64 (m, 2H), 1.35 (s, 18H), 1.20 (d, *J* = 7.6, 3H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.6 (s), 152.0 (s), 142.3 (s), 141.8 (s), 137.6 (s), 135.4 (s), 134.6 (s), 130.5 (s), 129.2 (s), 127.6 (s), 126.0 (s), 114.9 (s), 55.7 (s), 34.4 (s), 30.4 (s), 28.4 (s), 15.6 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₇O₂ [M+H]⁺: 417.2789, found: 417.2783.

2,6-Di-*tert*-butyl-4-((4-(*tert*-butyl)phenyl)(4-hydroxyphenyl)methyl)phenol (7d)



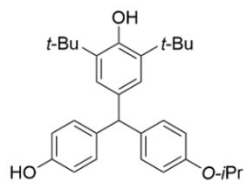
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **7d** (80.1 mg, 0.18 mmol, 90%) as a yellow solid. mp: 160.8-161.1 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.26-7.28 (m, 2H), 6.97-7.03 (m, 4H), 6.90 (s, 2H), 6.71-6.75 (m, 2H), 5.33 (s, 1H), 5.06 (s, 1H), 4.72 (s, 1H), 1.36 (s, 18H), 1.30 (s, 9H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.7 (s), 152.0 (s), 148.7 (s), 142.0 (s), 137.5 (s), 135.3 (s), 134.6 (s), 130.5 (s), 128.8 (s), 126.0 (s), 125.0 (s), 114.9 (s), 55.6 (s), 34.4 (s), 34.3 (s), 31.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₄₁O₂ [M+H]⁺: 445.3102, found: 445.3096.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(3-methoxyphenyl)methyl)phenol (7e)



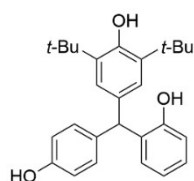
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 10:1) gave product **7e** (73.7 mg, 0.18 mmol, 88%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.15-7.19 (m, 1H), 6.95-9.97 (m, 2H), 6.89-6.90 (m, 2H), 6.67-6.74 (m, 5H), 5.33 (s, 1H), 5.12 (s, 1H), 5.07 (s, 1H), 3.72 (s, 3H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 159.4 (s), 153.8 (s), 152.1 (s), 146.9 (s), 137.0 (s), 135.5 (s), 134.3 (s), 130.5 (s), 129.1 (s), 126.0 (s), 122.1 (s), 115.4 (s), 115.0 (s), 111.2 (s), 56.0 (s), 55.2 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₅O₃ [M+H]⁺: 419.2581, found: 419.2575.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(4-isopropoxyphenyl)methyl)phenol (7f)



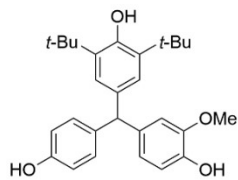
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7f** (76.9 mg, 0.17 mmol, 86%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.95-6.99 (m, 4H), 6.88 (s, 2H), 7.78-7.80 (m, 2H), 6.72-6.74 (m, 2H), 5.31 (s, 1H), 5.05 (s, 1H), 4.76 (s, 1H), 4.46-4.55 (m, 1H), 1.35 (s, 18H), 1.31 (d, *J* = 6.0, 6H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 171.3 (s), 156.0 (s), 153.9 (s), 151.9 (s), 137.2 (s), 135.4 (s), 134.9 (s), 130.4 (s), 130.2 (s), 125.9 (s), 115.6 (s), 114.9 (s), 60.5 (s), 34.3 (s), 30.3 (s), 22.1 (s), 21.1 (s). HRMS (ESI) *m/z*: calcd. for C₃₀H₃₉O₃ [M+H]⁺: 447.2894, found: 447.2890.

2,6-Di-*tert*-butyl-4-((2-hydroxyphenyl)(4-hydroxyphenyl)methyl)phenol (7h)



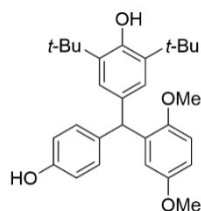
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7h** (64.8 mg, 0.16 mmol, 80%) as a yellow oil. ¹H NMR (400 MHz, DMSO-*d*₆, 25 °C, TMS): δ = 9.16 (s, 2H), 6.96-7.00 (m, 2H), 6.72-6.82 (m, 4H), 6.67-6.69 (m, 4H), 5.95 (s, 1H), 1.29 (s, 18H); ¹³C NMR (100 MHz, DMSO-*d*₆, 25 °C, TMS): δ = 159.8 (s), 159.7 (s), 156.8 (s), 143.7 (s), 143.6 (s), 139.7 (s), 136.3 (s), 134.6 (s), 134.5 (s), 131.8 (s), 130.4 (s), 123.5 (s), 123.4 (s), 120.1 (s), 47.5 (s), 39.6 (s), 35.7 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₃O₃ [M+H]⁺: 405.2425, found: 405.2420.

2,6-Di-*tert*-butyl-4-((4-hydroxy-3-methoxyphenyl)(4-hydroxyphenyl)methyl)phenol (7i)



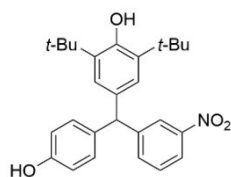
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7i** (68.73 mg, 0.16 mmol, 79%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.95-6.97 (m, 2H), 6.89 (s, 2H), 6.80-6.82 (m, 2H), 6.71-6.74 (m, 2H), 6.62-6.63 (m, 1H), 6.55-6.58 (m, 1H), 5.49 (s, 1H), 5.29 (s, 1H), 5.06 (s, 1H), 4.76 (s, 1H), 3.76 (s, 3H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.7 (s), 152.0 (s), 146.2 (s), 143.7 (s), 137.6 (s), 137.2 (s), 135.4 (s), 134.6 (s), 130.4 (s), 125.9 (s), 122.1 (s), 114.9 (s), 113.9 (s), 112.1 (s), 55.9 (s), 55.6 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₅O₄ [M+H]⁺: 435.2530, found: 435.2523.

2,6-Di-*tert*-butyl-4-((2,5-dimethoxyphenyl)(4-hydroxyphenyl)methyl)phenol (7j)



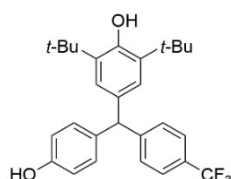
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7j** (71.82 mg, 0.16 mmol, 80%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.94-6.96 (m, 2H), 6.88-6.91 (m, 2H), 6.78-6.84 (m, 2H), 6.78-6.72 (m, 2H), 6.47-6.48 (m, 1H), 5.71 (s, 1H), 5.03 (s, 1H), 4.74 (s, 1H), 3.67 (s, 3H), 3.64 (s, 3H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.5 (s), 153.4 (s), 151.9 (s), 151.6 (s), 136.8 (s), 135.4 (s), 135.3 (s), 134.1 (s), 130.4 (s), 126.1 (s), 117.0 (s), 114.8 (s), 112.0 (s), 111.0 (s), 56.5 (s), 55.6 (s), 48.8 (s), 34.3 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₉H₃₇O₄ [M+H]⁺: 449.2687, found: 449.2681.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(3-nitrophenyl)methyl)phenol (**7k**)



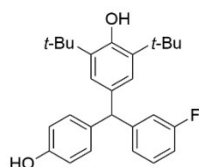
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7k** (71.2 mg, 0.16 mmol, 82%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 8.00-8.07 (m, 2H), 7.41-7.46 (m, 2H), 6.94-6.98 (m, 2H), 6.56 (s, 2H), 6.75-6.79 (m, 2H), 5.46 (s, 1H), 5.13 (s, 1H), 4.85 (s, 1H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 154.2 (s), 152.5 (s), 148.3 (s), 147.5 (s), 135.9 (s), 135.6 (s), 135.4 (s), 133.0 (s), 130.4 (s), 129.0 (s), 125.8 (s), 124.1 (s), 121.3 (s), 115.4 (s), 55.7 (s), 34.4 (s), 30.3 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{32}\text{NO}_4$ $[\text{M}+\text{H}]^+$: 434.2326, found: 434.2321.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(4-(trifluoromethyl)phenyl)methyl)phenol (**7l**)



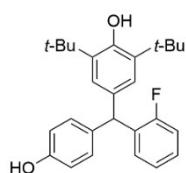
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7l** (76.8 mg, 0.17 mmol, 84%) as a yellow solid. mp: 112.3-112.5 $^\circ\text{C}$. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.50-7.53 (m, 2H), 7.20-7.26 (m, 2H), 6.93-6.96 (m, 2H), 6.86 (s, 2H), 6.73-6.77 (m, 2H), 5.42 (s, 1H), 5.11 (s, 1H), 4.76 (s, 1H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 154.0 (s), 152.3 (s), 149.3 (s), 136.3 (s), 135.8 (s), 132.0 (d, $J(\text{C},\text{F})$ = 297.7 Hz), 130.5 (s), 129.7 (s), 129.6 (s), 125.9 (s), 125.1 (q, $J(\text{C},\text{F})$ = 3.7 Hz), 120.8 (s), 115.2 (s), 55.8 (s), 34.4 (s), 30.3 (s). ^{19}F NMR (376 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = -62.3 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{28}\text{H}_{32}\text{F}_3\text{O}_2$ $[\text{M}+\text{H}]^+$: 457.2349, found: 457.2345.

2,6-Di-*tert*-butyl-4-((3-fluorophenyl)(4-hydroxyphenyl)methyl)phenol (**7m**)



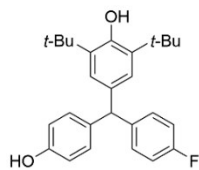
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7m** (70.3 mg, 0.17 mmol, 86%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.19-7.25 (m, 1H), 6.94-6.98 (m, 2H), 6.84-6.91 (m, 4H), 6.73-6.82 (m, 3H), 5.36 (s, 1H), 5.10 (s, 1H), 4.82 (s, 1H), 1.36 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 162.9 (d, $J(\text{C},\text{F})$ = 243.7 Hz), 153.9 (s), 152.2 (s), 147.9 (s), 136.5 (s), 135.6 (s), 133.8 (s), 130.5 (s), 129.45 (d, $J(\text{C},\text{F})$ = 8.2 Hz), 125.9 (s), 125.0 (s), 116.2 (d, $J(\text{C},\text{F})$ = 21.4 Hz), 115.1 (s), 112.9 (d, $J(\text{C},\text{F})$ = 21.0 Hz), 55.7 (s), 34.4 (s), 30.3 (s); ^{19}F NMR (376 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = -113.7 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{32}\text{FO}_2$ $[\text{M}+\text{H}]^+$: 407.2381, found: 407.2373.

2,6-Di-*tert*-butyl-4-((2-fluorophenyl)(4-hydroxyphenyl)methyl)phenol (**7n**)



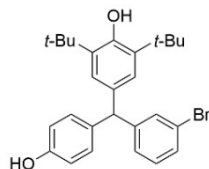
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7n** (65.9 mg, 0.16 mmol, 81%) as a yellow oil. ^1H NMR (400 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 7.15-7.21 (m, 1H), 6.92-7.05 (m, 5H), 6.86-6.88 (m, 2H), 6.73-6.75 (m, 2H), 5.67 (s, 1H), 5.08 (s, 1H), 4.74 (s, 1H), 1.35 (s, 18H); ^{13}C NMR (100 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = 160.7 (d, $J(\text{C},\text{F})$ = 244.8 Hz), 153.8 (s), 152.2 (s), 135.9 (s), 135.5 (s), 133.1 (s), 132.2 (d, $J(\text{C},\text{F})$ = 14.3 Hz), 130.7 (d, $J(\text{C},\text{F})$ = 4.0 Hz), 130.3 (s), 127.8 (d, $J(\text{C},\text{F})$ = 8.3 Hz), 125.9 (s), 123.7 (d, $J(\text{C},\text{F})$ = 3.5 Hz), 115.3 (s), 115.1 (d, $J(\text{C},\text{F})$ = 9.2 Hz), 48.4 (s), 34.4 (s), 30.3 (s); ^{19}F NMR (376 MHz, CDCl_3 , 25 $^\circ\text{C}$, TMS): δ = -116.7 (s). HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{32}\text{FO}_2$ $[\text{M}+\text{H}]^+$: 407.2381, found: 407.2376.

2,6-Di-*tert*-butyl-4-((4-fluorophenyl)(4-hydroxyphenyl)methyl)phenol (**7o**)



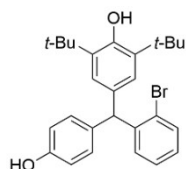
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7o** (69.2 mg, 0.17 mmol, 85%) as a yellow solid. mp: 162.3-162.7 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.15-7.21 (m, 1H), 6.92-7.05 (m, 5H), 6.86-6.88 (m, 2H), 6.73-6.75 (m, 2H), 5.67 (s, 1H), 5.08 (s, 1H), 4.74 (s, 1H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 161.3 (d, *J*(C,F) = 242.8 Hz), 153.8 (s), 152.1 (s), 140.9 (s), 137.0 (s), 135.5 (s), 134.3 (s), 130.71 (d, *J*(C,F) = 7.8 Hz), 130.5 (s), 125.6 (s), 115.0 (d, *J*(C,F) = 8.0 Hz), 114.8 (s), 55.2 (s), 34.4 (s), 30.3 (s); ¹⁹F NMR (376 MHz, CDCl₃, 25 °C, TMS): δ = -117.5 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₂FO₂ [M+H]⁺: 407.2381, found: 407.2375.

4-((3-Bromophenyl)(4-hydroxyphenyl)methyl)-2,6-di-*tert*-butylphenol (**7p**)



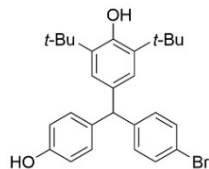
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7p** (80.2 mg, 0.17 mmol, 86%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.31-7.33 (m, 1H), 7.26-7.27 (m, 1H), 7.11-7.15 (m, 1H), 7.01-7.03 (m, 1H), 6.94-6.96 (m, 2H), 6.86 (s, 2H), 6.73-6.77 (m, 2H), 5.33 (s, 1H), 5.10 (s, 1H), 4.71 (s, 1H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.9 (s), 152.3 (s), 136.4 (s), 135.6 (s), 133.6 (s), 132.4 (s), 130.7 (s), 129.7 (s), 129.2 (s), 128.0 (s), 125.9 (s), 122.4 (s), 115.1 (s), 100.0 (s), 55.7 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₂BrO₂ [M+H]⁺: 467.1581, found: 467.1577.

4-((2-Bromophenyl)(4-hydroxyphenyl)methyl)-2,6-di-*tert*-butylphenol (**7q**)



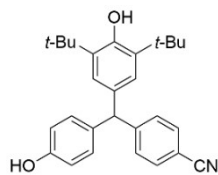
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7q** (75.7 mg, 0.16 mmol, 81%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.53-7.55 (m, 1H), 7.18-7.22 (m, 1H), 7.03-7.07 (m, 1H), 6.92-6.97 (m, 3H), 6.84 (s, 2H), 6.72-6.75 (m, 2H), 5.76 (s, 1H), 5.08 (s, 1H), 4.84 (s, 1H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.8 (s), 152.2 (s), 144.4 (s), 135.6 (s), 135.5 (s), 133.3 (s), 133.0 (s), 131.1 (s), 130.7 (s), 127.7 (s), 127.1 (s), 126.2 (s), 125.5 (s), 115.0 (s), 55.0 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₂BrO₂ [M+H]⁺: 467.1581, found: 467.1573.

4-((4-Bromophenyl)(4-hydroxyphenyl)methyl)-2,6-di-*tert*-butylphenol (**7r**)



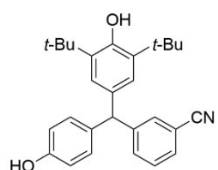
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7r** (80.2 mg, 0.17 mmol, 86%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.36-7.39 (m, 2H), 6.93-6.98 (m, 4H), 6.85 (s, 2H), 6.72-6.76 (m, 2H), 5.32 (s, 1H), 5.08 (s, 1H), 4.83 (s, 1H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.9 (s), 152.2 (s), 144.3 (s), 136.6 (s), 135.6 (s), 133.9 (s), 131.2 (s), 131.1 (s), 130.5 (s), 129.7 (s), 125.8 (s), 120.8 (s), 119.9 (s), 115.1 (s), 55.4 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₂BrO₂ [M+H]⁺: 467.1581, found: 467.1575.

4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(4-hydroxyphenyl)methyl)benzonitrile (7s)



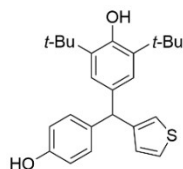
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7s** (63.8 mg, 0.15 mmol, 77%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.55-7.57 (m, 2H), 7.20-7.22 (m, 2H), 6.92-6.94 (m, 2H), 6.83 (s, 2H), 6.75-6.78 (m, 2H), 5.41 (s, 1H), 5.13 (s, 1H), 4.95 (s, 1H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 154.2 (s), 152.4 (s), 150.9 (s), 135.8 (s), 135.5 (s), 133.0 (s), 132.0 (s), 130.5 (s), 130.1 (s), 125.8 (s), 119.1 (s), 115.3 (s), 109.8 (s), 56.0 (s), 34.4 (s), 30.3 (s); HRMS (ESI) *m/z*: calcd. for C₂₈H₃₂NO₂ [M+H]⁺: 414.2428, found: 414.2424.

3-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(4-hydroxyphenyl)methyl)benzonitrile (7t)



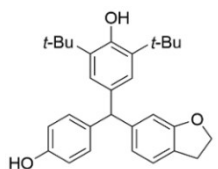
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7t** (65.4 mg, 0.16 mmol, 79%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.48-7.50 (m, 1H), 7.36-7.39 (m, 3H), 6.92-6.94 (m, 2H), 6.76-6.82 (m, 2H), 5.39 (s, 1H), 5.13 (s, 1H), 4.83 (s, 1H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 154.2 (s), 152.5 (s), 146.8 (s), 135.9 (s), 135.6 (s), 133.9 (s), 133.1 (s), 132.8 (s), 130.5 (s), 129.9 (s), 128.9 (s), 125.8 (s), 119.1 (s), 115.3 (s), 112.1 (s), 55.5 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₂NO₂ [M+H]⁺: 414.2428, found: 414.2421.

2,6-Di-*tert*-butyl-4-((4-hydroxyphenyl)(thiophen-3-yl)methyl)phenol (7u)



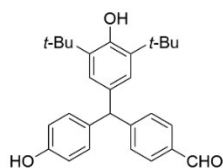
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7u** (64.8 mg, 0.16 mmol, 82%) as a yellow solid. mp: 121.8-122.1 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.23-7.25 (m, 1H), 7.00-7.02 (m, 2H), 6.92-6.94 (m, 2H), 6.85-6.86 (m, 1H), 6.70-6.75 (m, 3H), 5.33 (s, 1H), 5.07 (s, 1H), 4.70 (s, 1H), 1.37 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.8 (s), 152.2 (s), 146.2 (s), 137.1 (s), 135.5 (s), 134.5 (s), 130.1 (s), 128.9 (s), 125.5 (s), 125.2 (s), 122.3 (s), 115.0 (s), 51.8 (s), 34.4 (s), 30.4 (s). HRMS (ESI) *m/z*: calcd. for C₂₅H₃₁O₂S [M+H]⁺: 395.2040, found: 395.2035.

2,6-Di-*tert*-butyl-4-((2,3-dihydrobenzofuran-6-yl)(4-hydroxyphenyl)methyl)phenol (7x)



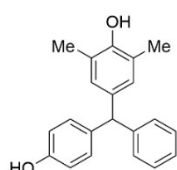
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7x** (73.3 mg, 0.17 mmol, 85%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 6.93-6.97 (m, 3H), 6.89 (s, 2H), 6.80-6.83 (m, 1H), 6.71-6.75 (m, 2H), 6.67-6.69 (m, 1H), 5.43 (s, 1H), 5.06 (s, 1H), 4.79 (s, 1H), 4.54 (t, *J* = 8.7 Hz, 2H), 3.14 (t, *J* = 8.7 Hz, 2H), 1.36 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 158.3 (s), 153.6 (s), 152.0 (s), 137.7 (s), 137.4 (s), 135.4 (s), 134.9 (s), 130.5 (s), 128.8 (s), 126.8 (s), 125.9 (s), 125.8 (s), 114.9 (s), 108.7 (s), 71.2 (s), 55.4 (s), 34.4 (s), 30.4 (s), 29.8 (s); HRMS (ESI) *m/z*: calcd. for C₂₉H₃₅O₃ [M+H]⁺: 431.2581, found: 431.2573.

4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(4-hydroxyphenyl)methyl)benzaldehyde (**7z**)



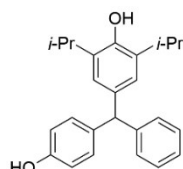
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7z** (55.2 mg, 0.13 mmol, 66%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 9.98 (s, 1H), 7.78-7.80 (m, 2H), 7.26-7.29 (m, 2H), 6.94-6.96 (m, 2H), 6.86 (s, 2H), 6.75-6.78 (m, 2H), 5.45 (s, 1H), 5.12 (s, 1H), 5.02 (s, 1H), 1.35 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 192.2 (s), 154.1 (s), 152.6 (s), 152.4 (s), 136.0 (s), 135.7 (s), 134.5 (s), 133.4 (s), 130.5 (s), 130.0 (s), 129.8 (s), 125.9 (s), 115.2 (s), 34.4 (s), 30.3 (s). HRMS (ESI) *m/z*: calcd. for C₂₈H₃₃O₃ [M+H]⁺: 417.2425, found: 417.2420.

4-((4-Hydroxyphenyl)(phenyl)methyl)-2,6-dimethylphenol (**7ab**)



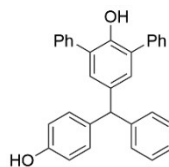
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7ab** (51.9 mg, 0.17 mmol, 85%) as a yellow solid. mp: 127.5-127.9 °C. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.25-7.29 (m, 2H), 7.17-7.21 (m, 1H), 7.09-7.10 (m, 2H), 6.94-6.97 (m, 2H), 6.70-6.75 (m, 4H), 5.35 (s, 1H), 4.80 (s, 1H), 4.53 (s, 1H), 2.17 (s, 6H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.8 (s), 150.6 (s), 144.7 (s), 136.8 (s), 135.8 (s), 130.5 (s), 129.5 (s), 129.3 (s), 128.2 (s), 126.1 (s), 122.8 (s), 115.1 (s), 55.3 (s), 16.1 (s); HRMS (ESI) *m/z*: calcd. for C₂₁H₂₁O₂ [M+H]⁺: 305.1537, found: 305.1531.

4-((4-Hydroxyphenyl)(phenyl)methyl)-2,6-diisopropylphenol (**7ac**)



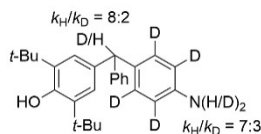
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7ac** (60.1 mg, 0.17 mmol, 83%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.24-7.28 (m, 2H), 7.15-7.20 (m, 1H), 7.09-7.10 (m, 2H), 6.95-6.98 (m, 2H), 6.77 (s, 2H), 6.70-6.75 (m, 2H), 5.40 (s, 1H), 4.77 (s, 1H), 4.68 (s, 1H), 3.00-3.14 (m, 2H), 1.17 (d, *J* = 6.9 Hz, 12H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.7 (s), 148.3 (s), 145.0 (s), 137.1 (s), 135.9 (s), 133.3 (s), 130.5 (s), 129.3 (s), 128.2 (s), 126.0 (s), 124.5 (s), 115.0 (s), 55.8 (s), 27.3 (s), 22.7 (s); HRMS (ESI) *m/z*: calcd. for C₂₅H₂₉O₂ [M+H]⁺: 361.2163, found: 361.2159.

5'-((4-Hydroxyphenyl)(phenyl)methyl)-[1,1':3',1''-terphenyl]-2'-ol (**7ad**)



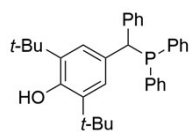
According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 5:1) gave product **7ad** (62.6 mg, 0.15 mmol, 73%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.48-7.50 (m, 4H), 7.40-7.44 (m, 4H), 7.24-7.35 (m, 4H), 7.16-7.21 (m, 3H), 7.02-7.04 (m, 4H), 6.72-6.74 (m, 2H), 5.48 (s, 1H), 5.35 (s, 1H), 4.87 (s, 1H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.9 (s), 147.7 (s), 144.4 (s), 137.7 (s), 136.5 (s), 136.4 (s), 130.80 (s), 130.54 (s), 129.4 (s), 129.3 (s), 128.83 (s), 128.5 (s), 128.4 (s), 127.6 (s), 126.23 (s), 115.2 (s), 55.45 (s). HRMS (ESI) *m/z*: calcd. for C₃₁H₂₅O₂ [M+H]⁺: 429.1850, found: 429.1846.

4-((4-Aminophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3a')



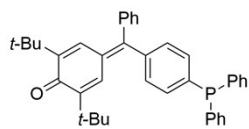
According to the general procedure, work-up and flash column chromatography (Hexane/EtOAc: 10:1) gave product **3a'** (68.9 mg, 0.16 mmol, 82%) as a yellow oil. ¹H NMR (400 MHz, Acetone-*d*₆, 25 °C, TMS): δ = 7.11-7.17 (m, 2H), 6.99-7.07 (m, 3H), 6.85 (s, 2H), 5.76-5.81 (s, 1H), 5.35 (s, 0.8H), 4.35 (s, 1.4H), 1.22 (s, 18H); ¹³C NMR (100 MHz, Acetone-*d*₆, 25 °C, TMS): δ = 152.2 (s), 145.3 (s), 137.1 (s), 135.1 (s), 129.2 (s), 129.1 (s), 128.1 (s), 127.9 (s), 125.9 (s), 125.7 (s), 125.6 (s), 119.0 (s), 56.0 (s), 34.3 (s), 29.8 (s). HRMS (ESI) *m/z*: calcd. for C₂₇H₃₄N₂O [M+H]⁺: 388.2635, found: 388.2629.

2,6-Di-*tert*-butyl-4-((diphenylphosphaneyl)(phenyl)methyl)phenol (9a)



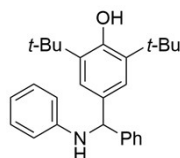
According to the general procedure, work-up and flash column chromatography (Hexane/EtOAc: 20:1) gave product **9a** (85.6 mg, 0.18 mmol, 89%) as a colorless oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.62-7.68 (m, 2H), 7.14-7.51 (m, 13H), 6.94-6.95 (m, 2H), 5.06 (s, 1H), 4.89 (d, *J* = 17.5 Hz, 1H), 1.29 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 152.9 (s), 138.1 (s), 135.3 (d, *J*(C,P) = 1.6 Hz), 133.2 (d, *J*(C,P) = 17.1 Hz), 133.0 (d, *J*(C,P) = 9.3 Hz), 131.7 (d, *J*(C,P) = 2.4 Hz), 130.9 (d, *J*(C,P) = 2.4 Hz), 130.8 (d, *J*(C,P) = 2.4 Hz), 130.0 (d, *J*(C,P) = 6.3 Hz), 129.8 (d, *J*(C,P) = 81.5 Hz), 129.3 (d, *J*(C,P) = 78.8 Hz), 129.1 (d, *J*(C,P) = 10.4 Hz), 128.4 (d, *J*(C,P) = 9.9 Hz), 128.1 (d, *J*(C,P) = 9.6 Hz), 127.0 (d, *J*(C,P) = 5.1 Hz), 126.9 (d, *J*(C,P) = 1.8 Hz), 51.2 (d, *J*(C,P) = 29.0 Hz), 34.3 (s), 30.2 (s). ³¹P NMR (160 MHz, CDCl₃, 25 °C, TMS): δ = 22.5. HRMS (ESI) *m/z*: calcd. for C₃₃H₃₈OP[M+H]⁺: 481.2655, found: 481.2651.

2,6-Di-*tert*-butyl-4-((4-(diphenylphosphaneyl)phenyl)(phenyl)methylene)cyclohexa-2,5-dien-1-one (9b)



According to the general procedure, work-up and flash column chromatography (Hexane/EtOAc: 40:1) gave product **9b** (94.4 mg, 0.17 mmol, 85%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.53-7.60 (m, 2H), 7.37-7.48 (m, 7H), 7.29-7.30 (m, 10H), 7.06 (s, 1H), 7.02 (s, 1H), 1.32 (d, *J* = 14.0 Hz, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 186.6 (s), 149.4 (s), 147.8 (s), 142.6 (s), 137.3 (d, *J*(C,P) = 10.9 Hz), 136.0 (s), 135.2 (s), 133.8 (d, *J*(C,P) = 19.5 Hz), 133.3 (d, *J*(C,P) = 9.6 Hz), 132.0 (s), 131.3 (d, *J*(C,P) = 2.3 Hz), 130.4 (s), 129.3 (d, *J*(C,P) = 57.7 Hz), 129.0 (d, *J*(C,P) = 25.3 Hz), 128.9 (d, *J*(C,P) = 4.3 Hz), 128.6 (d, *J*(C,P) = 6.9 Hz), 127.9 (s), 35.5 (s), 35.1 (s), 29.6 (s), 29.5 (s). ³¹P NMR (160 MHz, CDCl₃, 25 °C, TMS): δ = -5.37. HRMS (ESI) *m/z*: calcd. for C₃₉H₄₀OP[M+H]⁺: 555.2812, found: 555.2807.

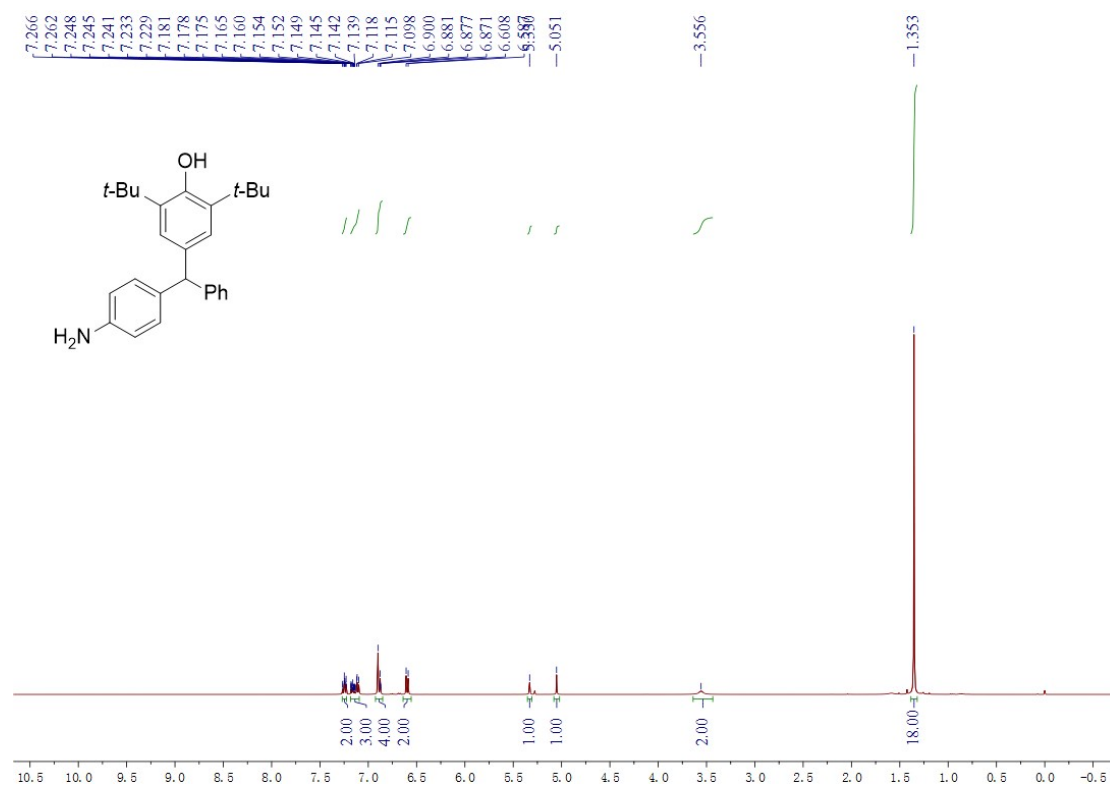
2,6-Di-*tert*-butyl-4-(phenyl(phenylamino)methyl)phenol (3a'')

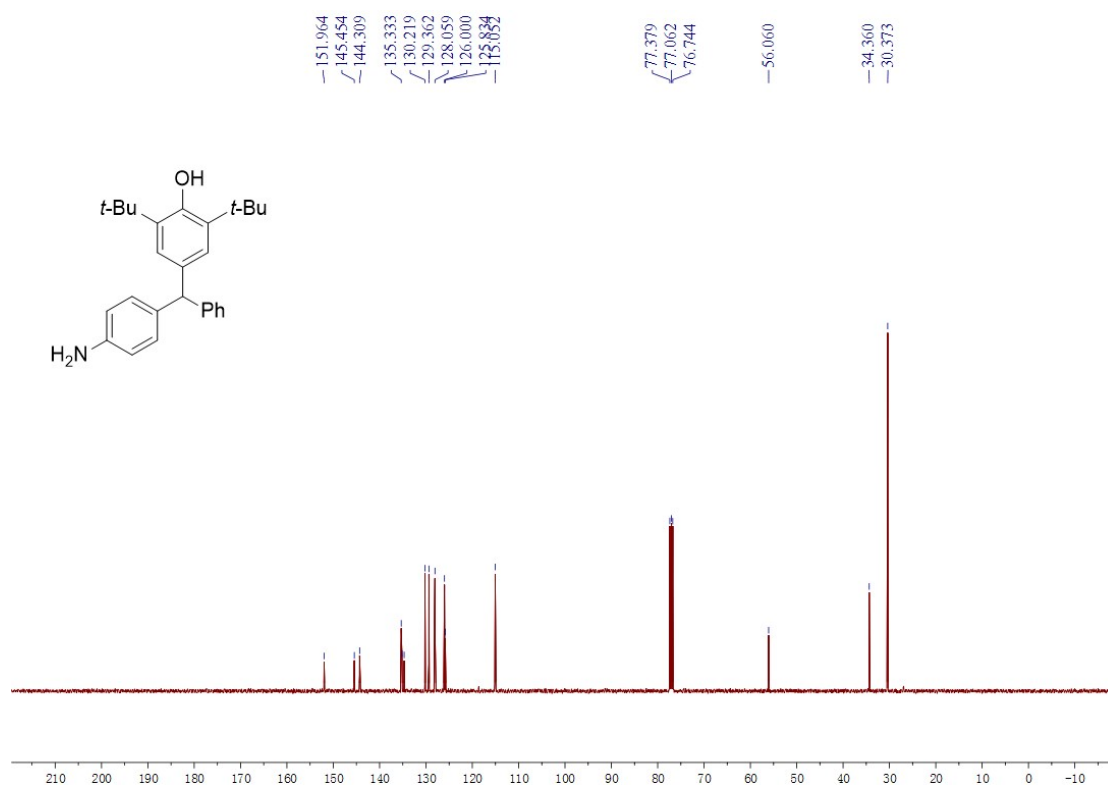


According to the general procedure, work-up and flash column chromatography (hexane/EtOAc: 40:1) gave product **3a''** (58.3 mg, 0.15 mmol, 75%) as a yellow oil. ¹H NMR (400 MHz, CDCl₃, 25 °C, TMS): δ = 7.39-7.41 (m, 2H), 7.30-7.34 (m, 2H), 7.21-7.25 (m, 1H), 7.07-7.12 (m, 4H), 6.65-6.68 (m, 1H), 6.53-6.55 (m, 2H), 5.40 (s, 1H), 5.16 (s, 1H), 4.24 (s, 1H), 1.38 (s, 18H); ¹³C NMR (100 MHz, CDCl₃, 25 °C, TMS): δ = 153.1 (s), 147.7 (s), 143.1 (s), 136.0 (s), 134.1 (s), 129.1 (s), 128.6 (s), 127.1 (s), 126.9 (s), 124.6 (s), 117.3 (s), 113.4 (s), 63.2 (s), 34.4 (s), 30.3 (s).

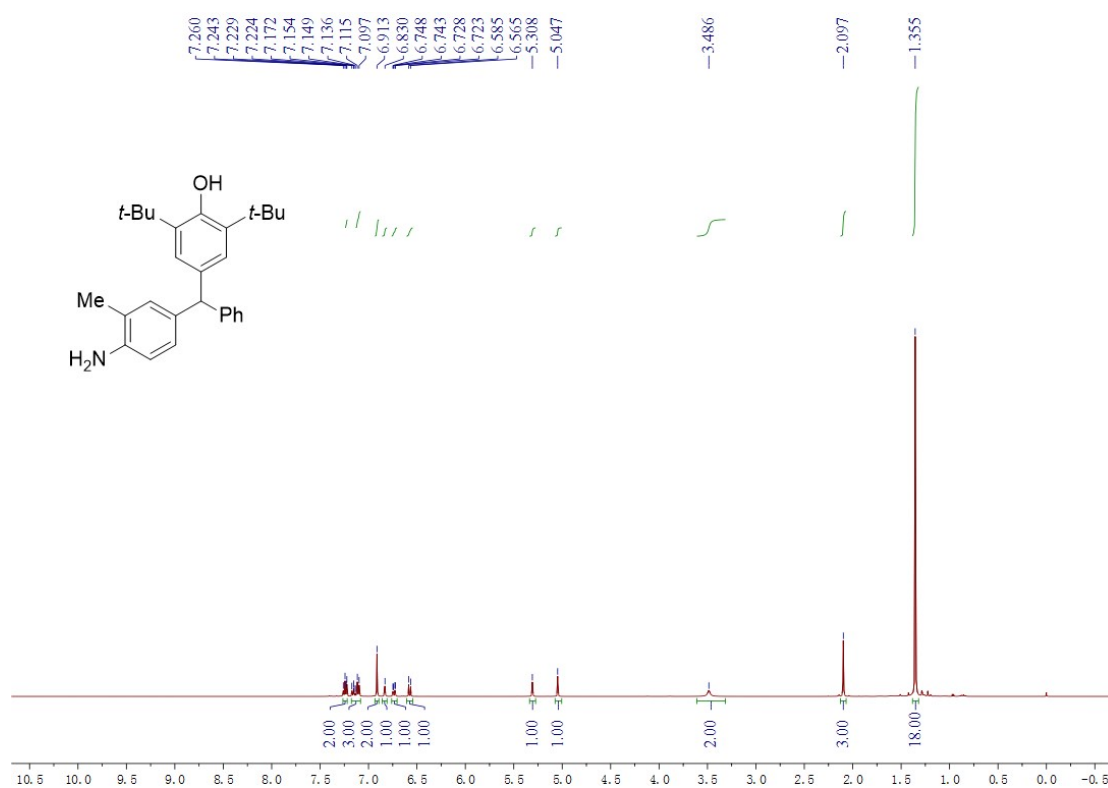
^1H , ^{13}C , ^{31}P and ^{19}F NMR spectra

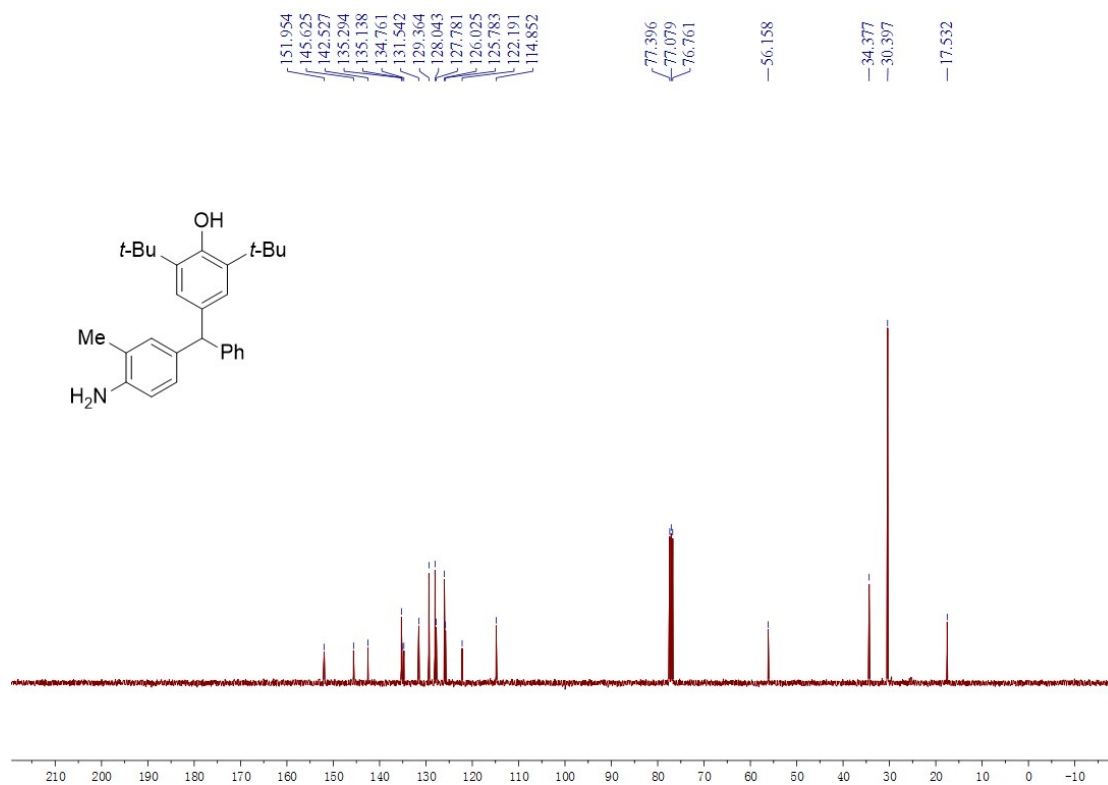
4-((4-Aminophenyl)(phenyl)methyl)-2,6-di-*tert*-butylphenol (3a)



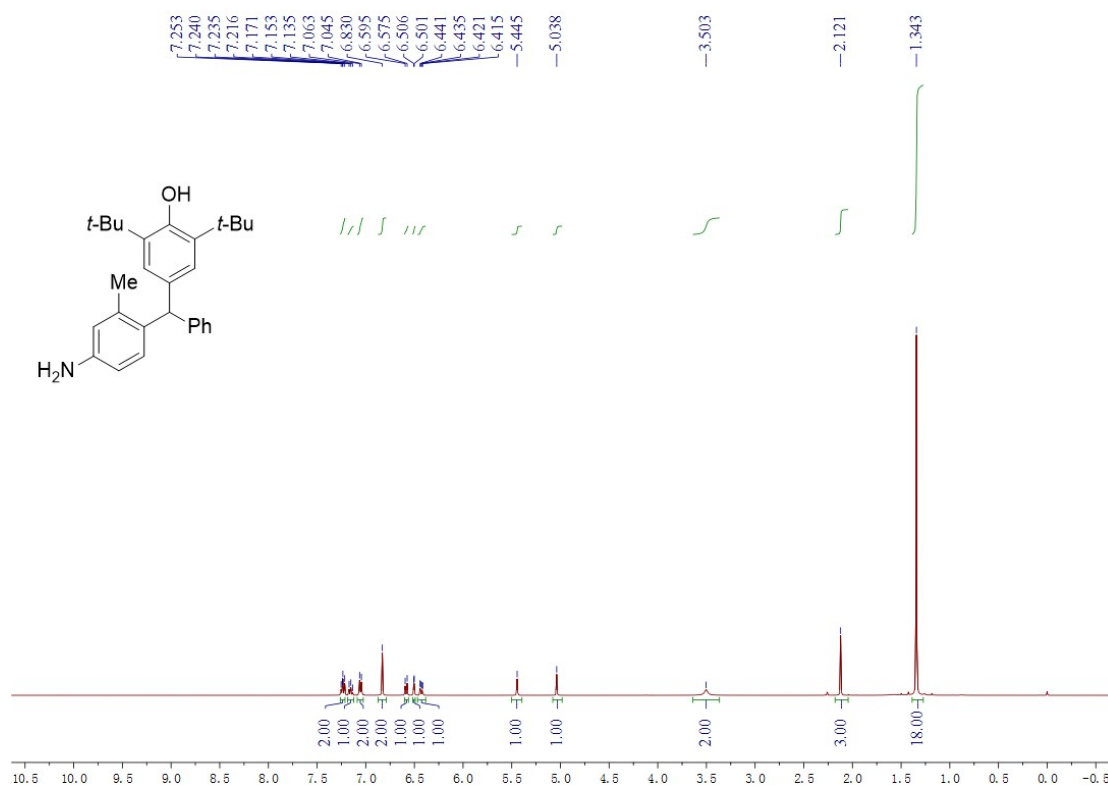


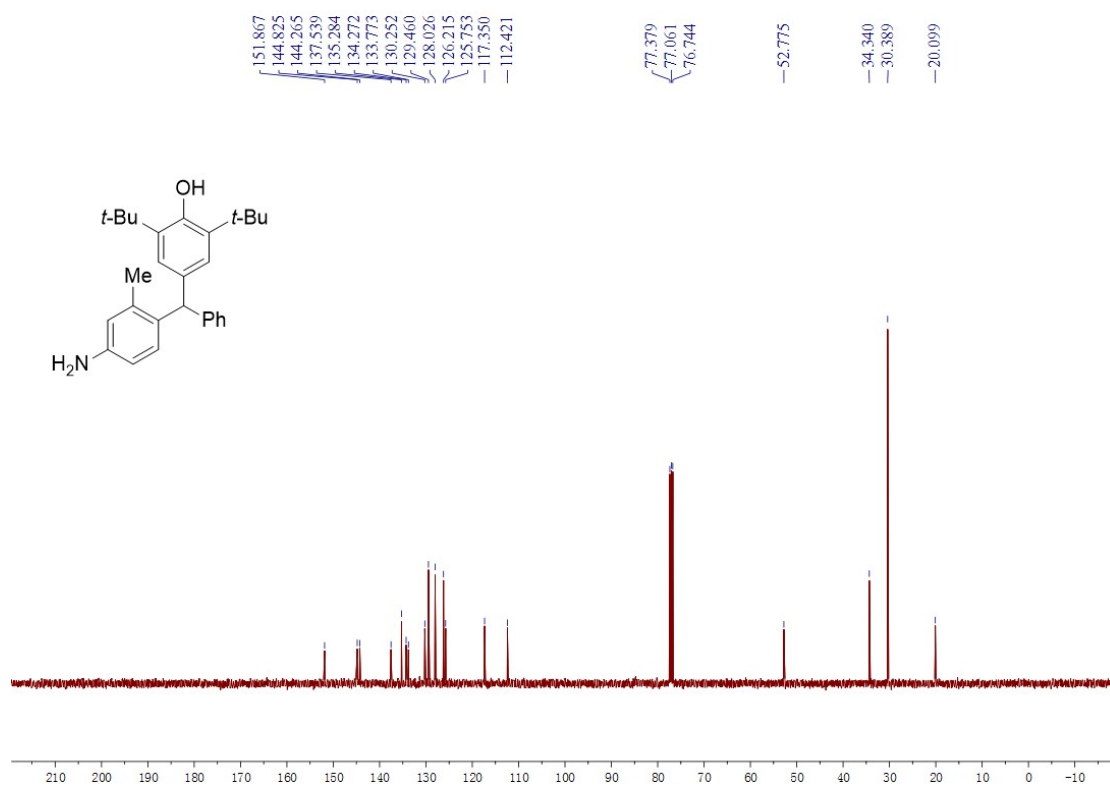
4-((4-Amino-3-methylphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3b)



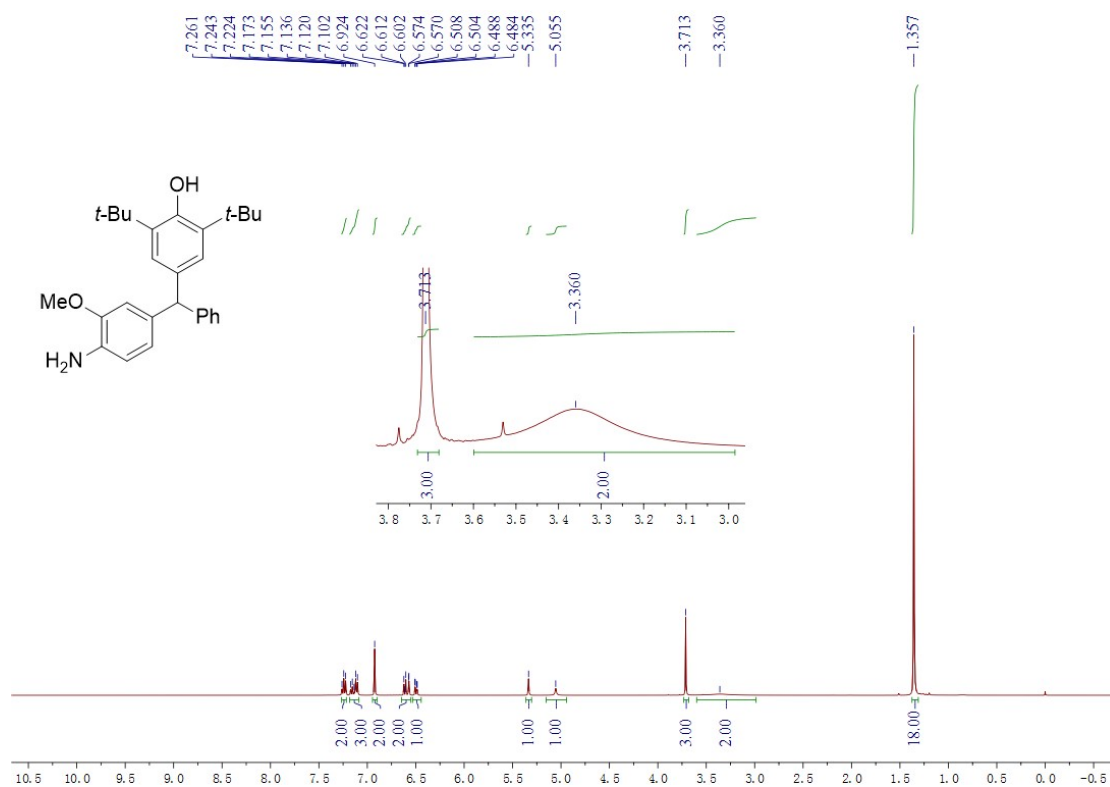


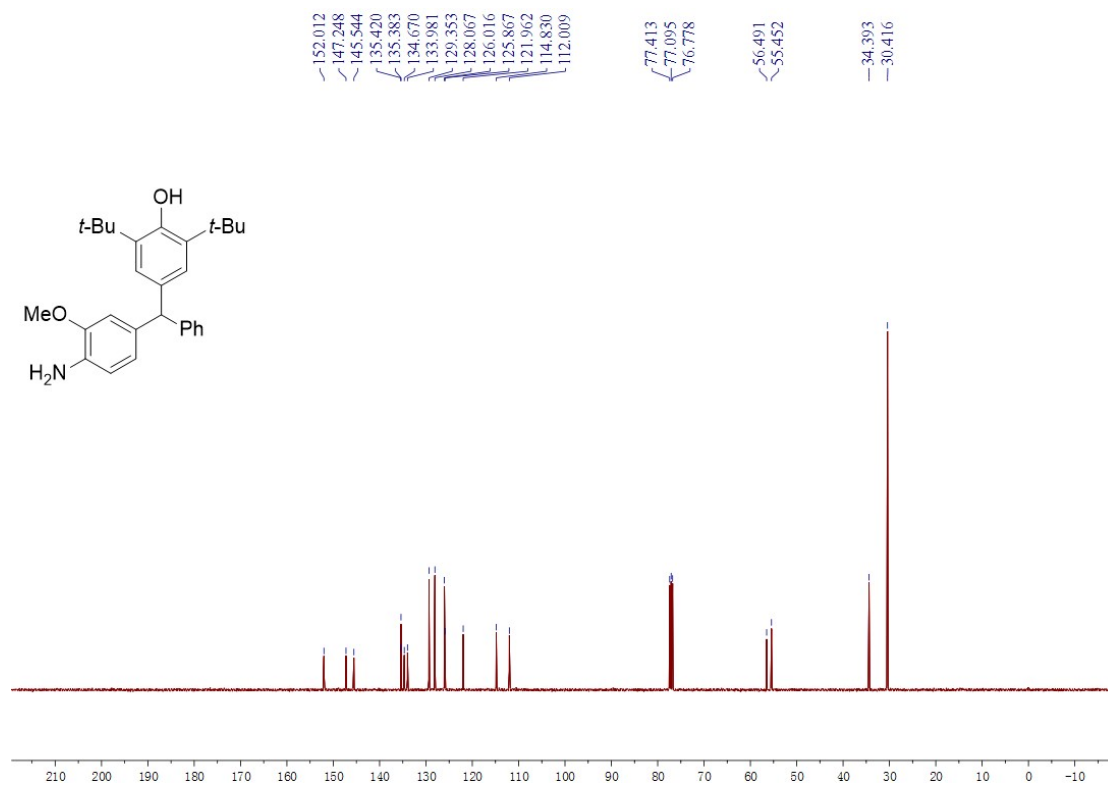
4-((4-Amino-2-methylphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3c)



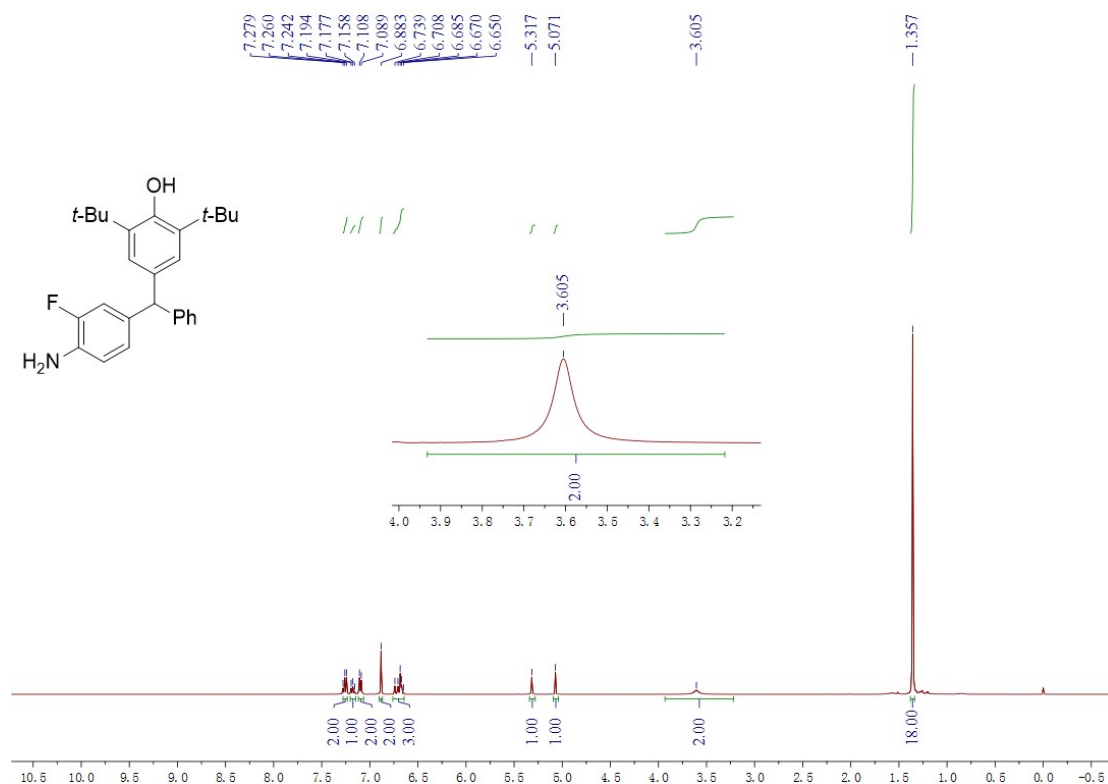


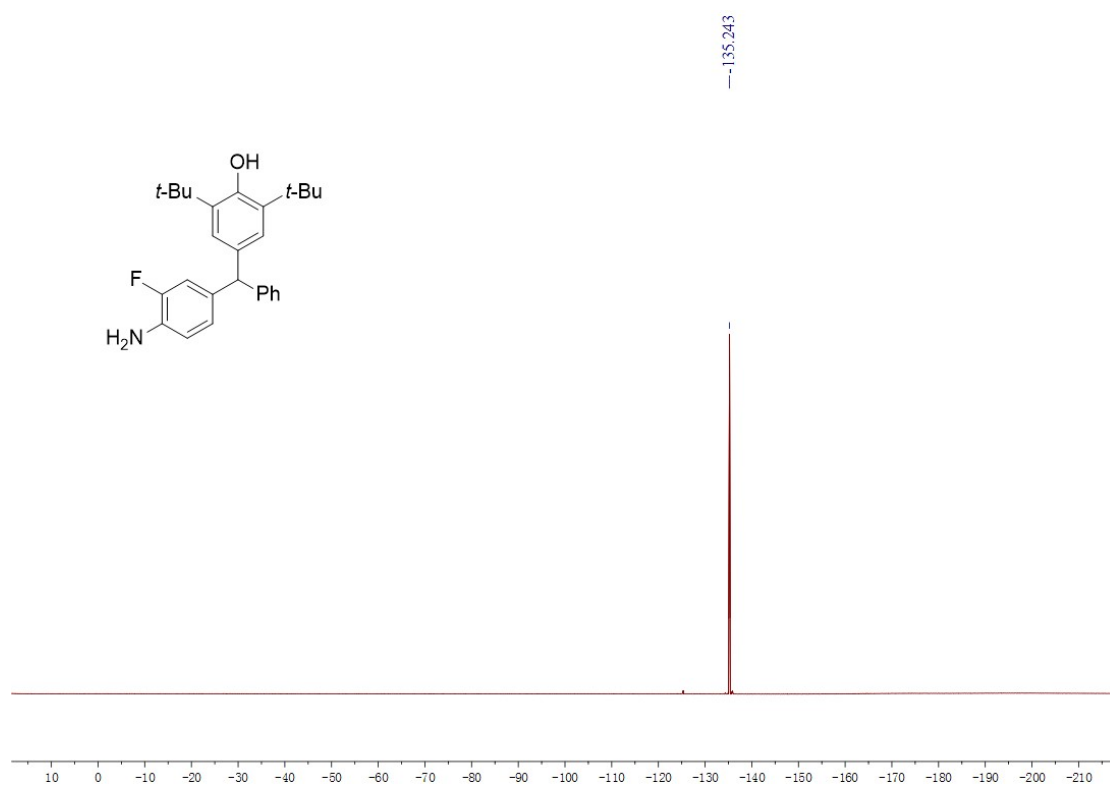
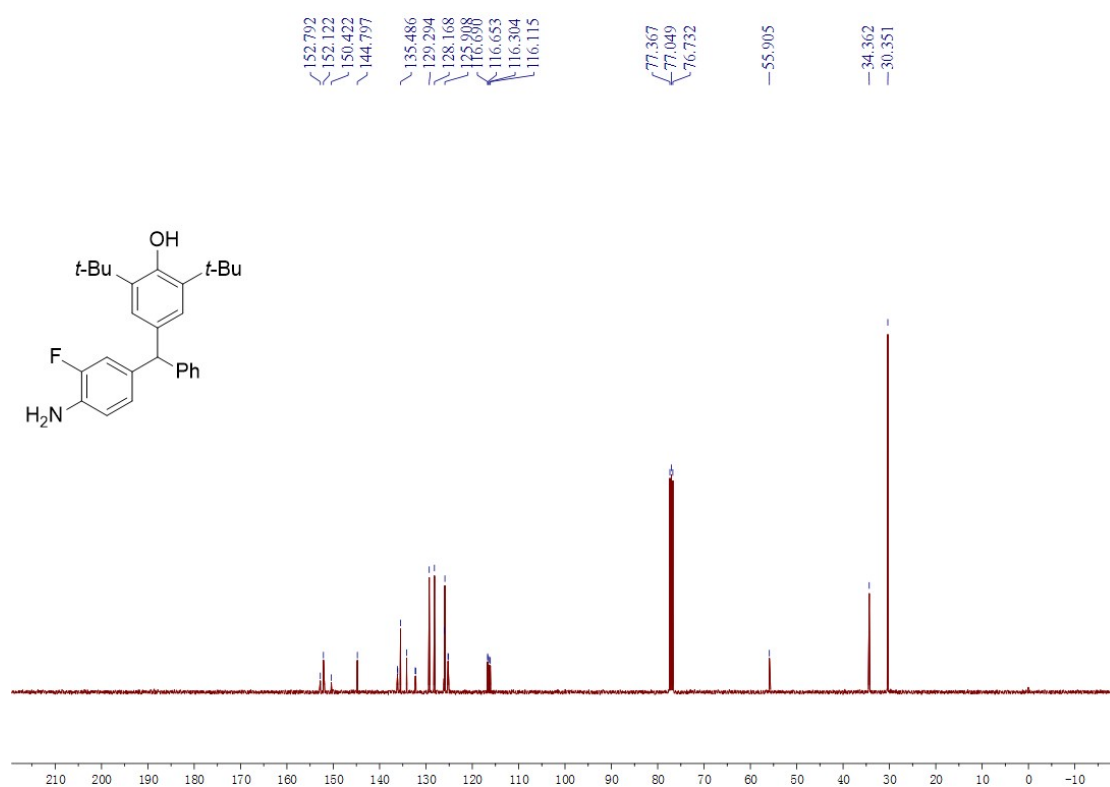
4-((4-Amino-3-methoxyphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3d)



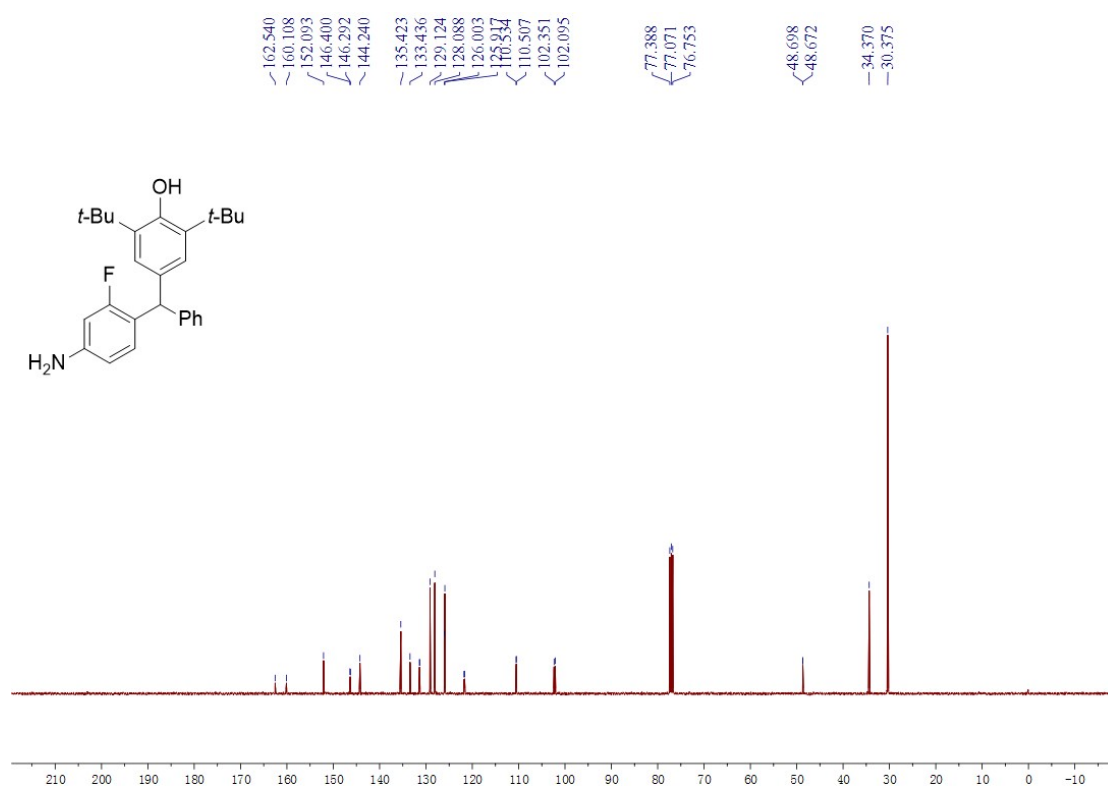
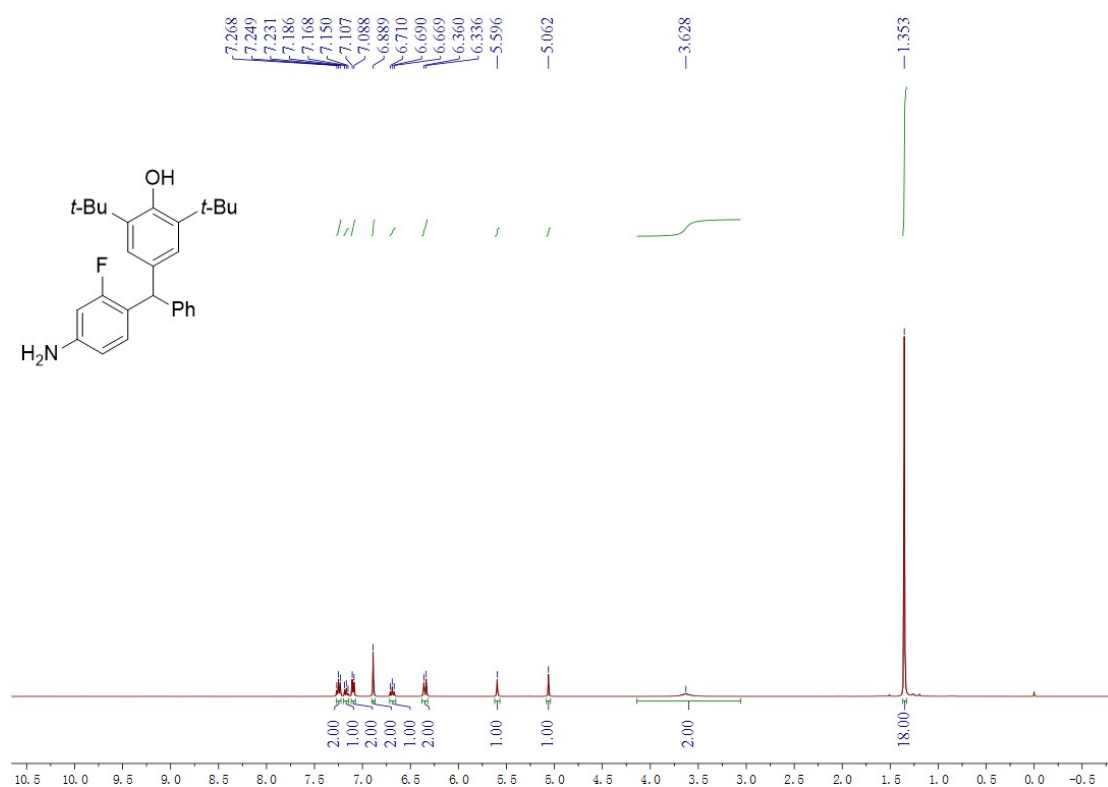


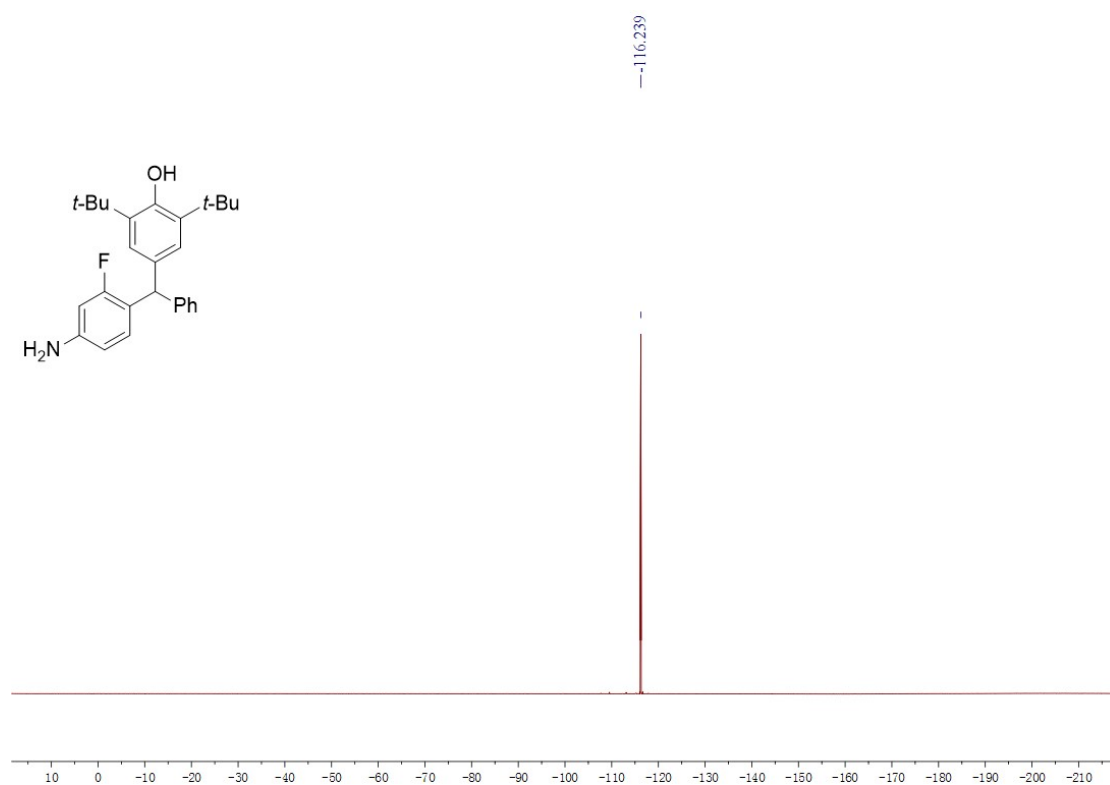
4-((4-Amino-3-fluorophenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3e)



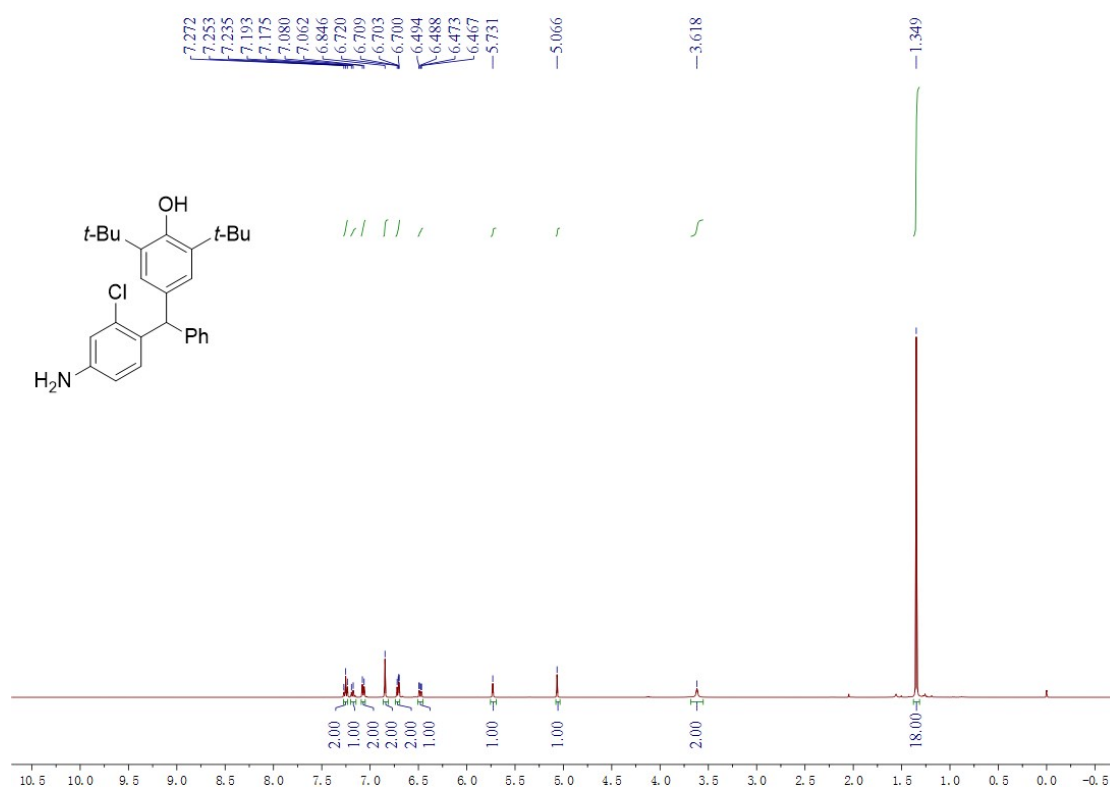


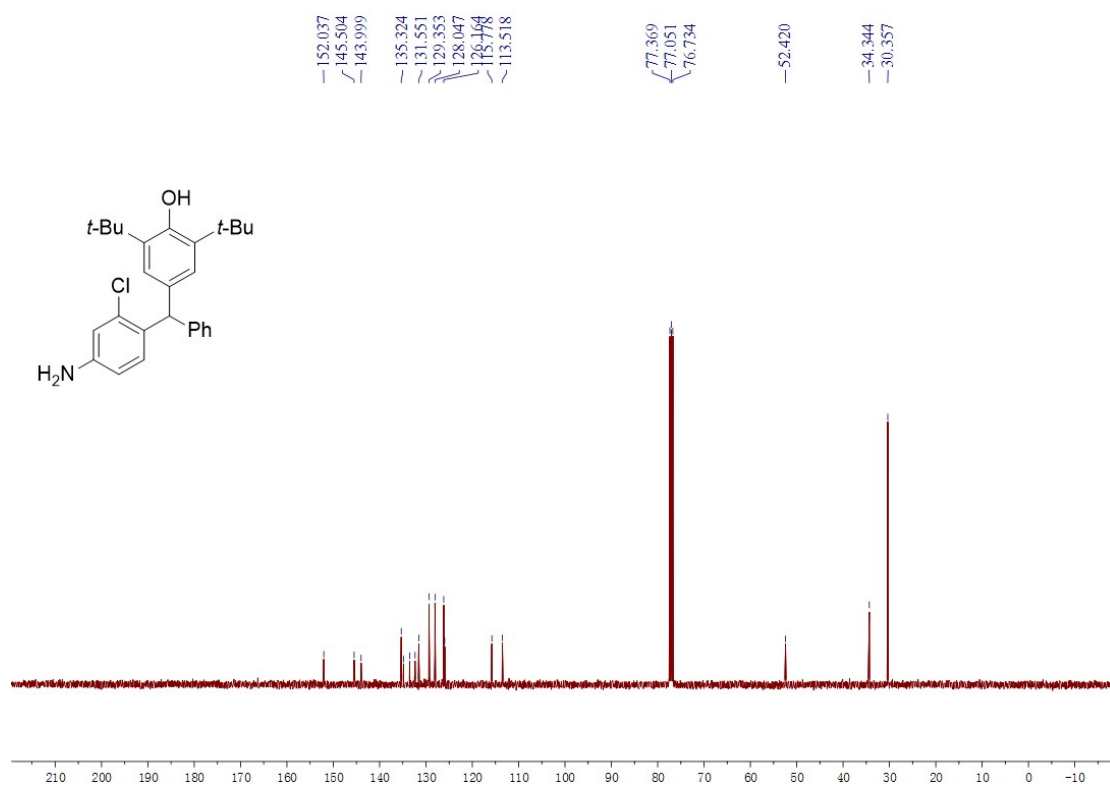
4-((4-Amino-2-fluorophenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3f)



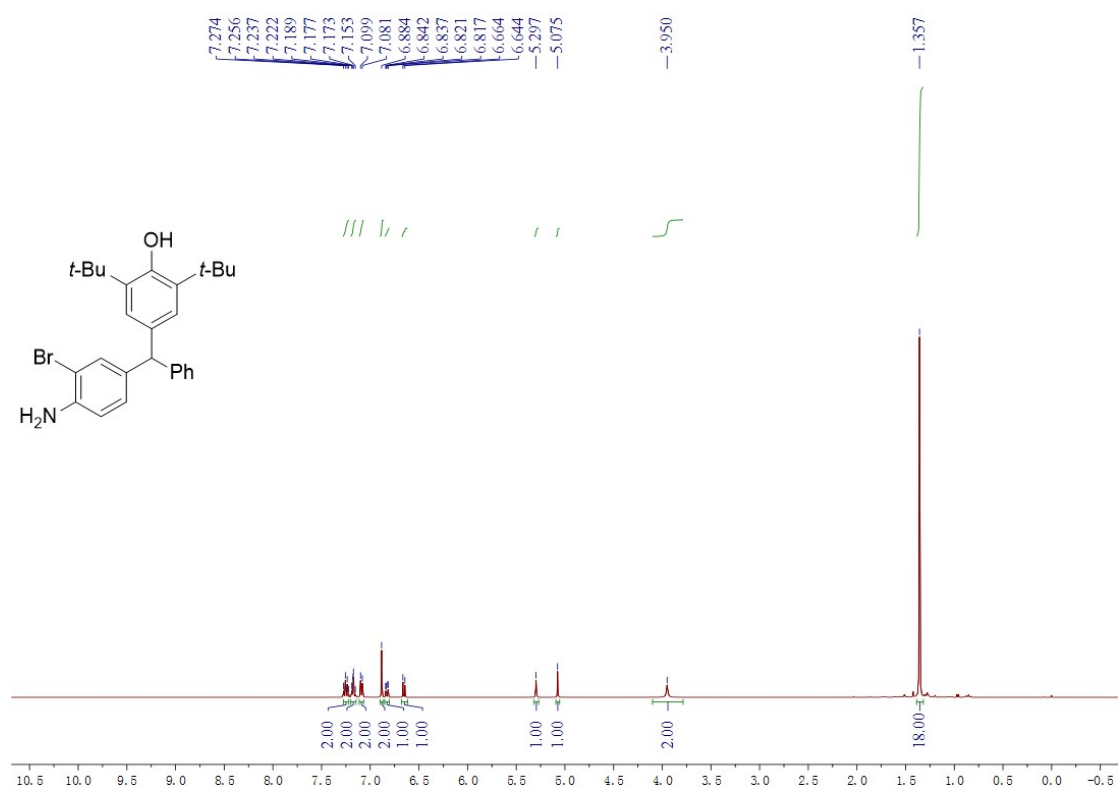


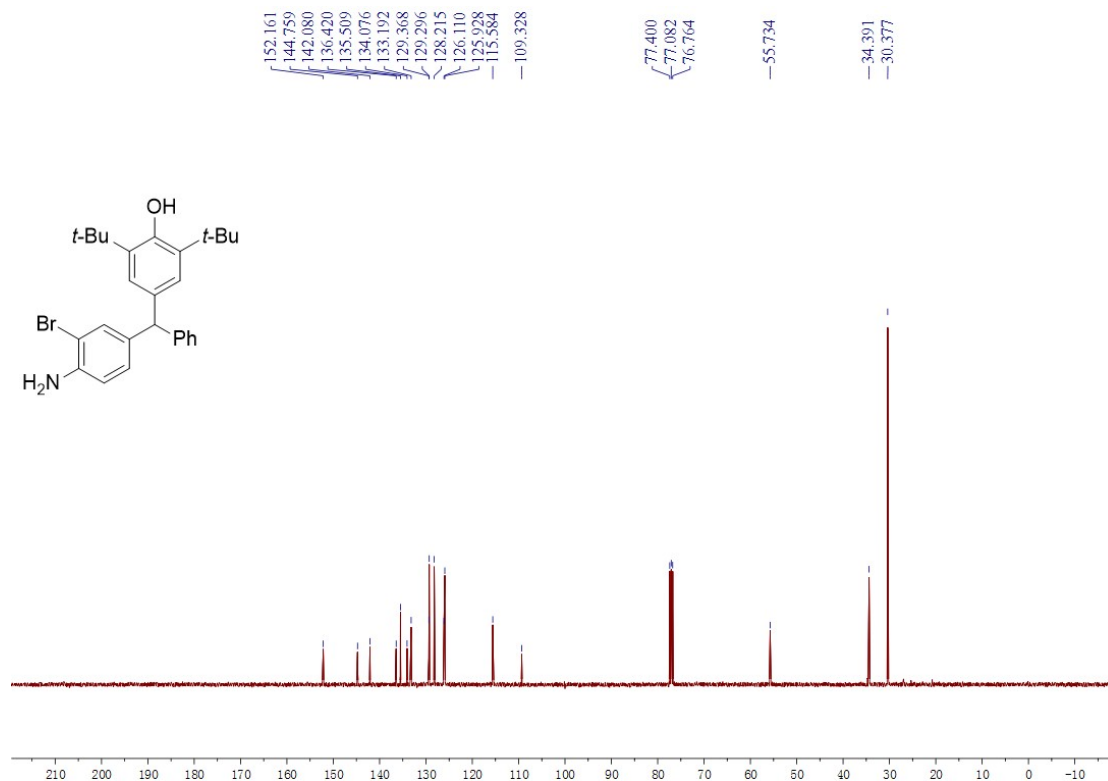
4-((4-Amino-2-chlorophenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3g)



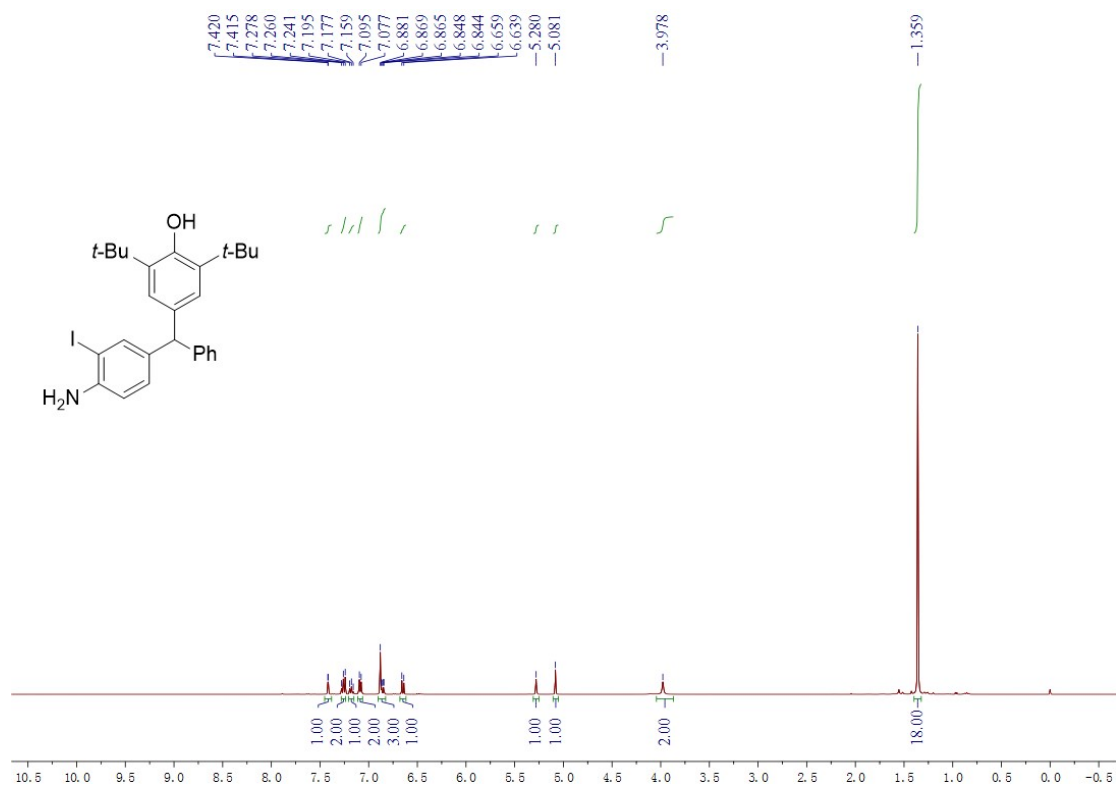


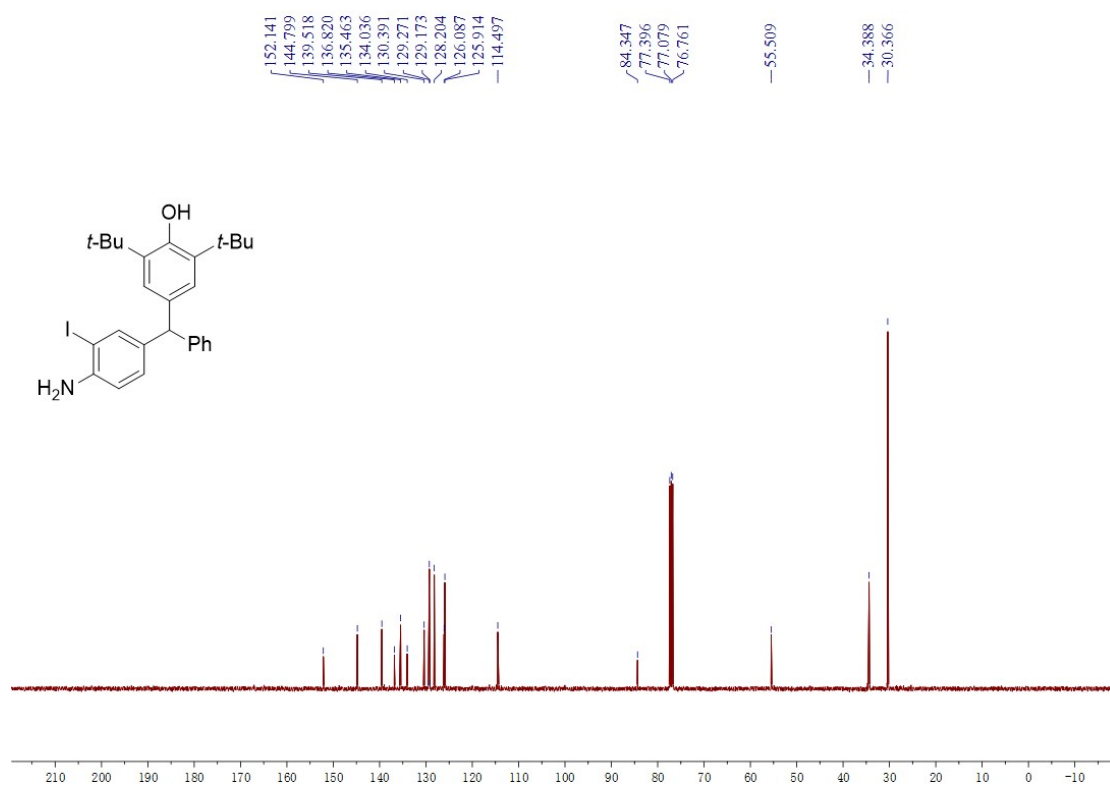
4-((4-Amino-3-bromophenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3h)



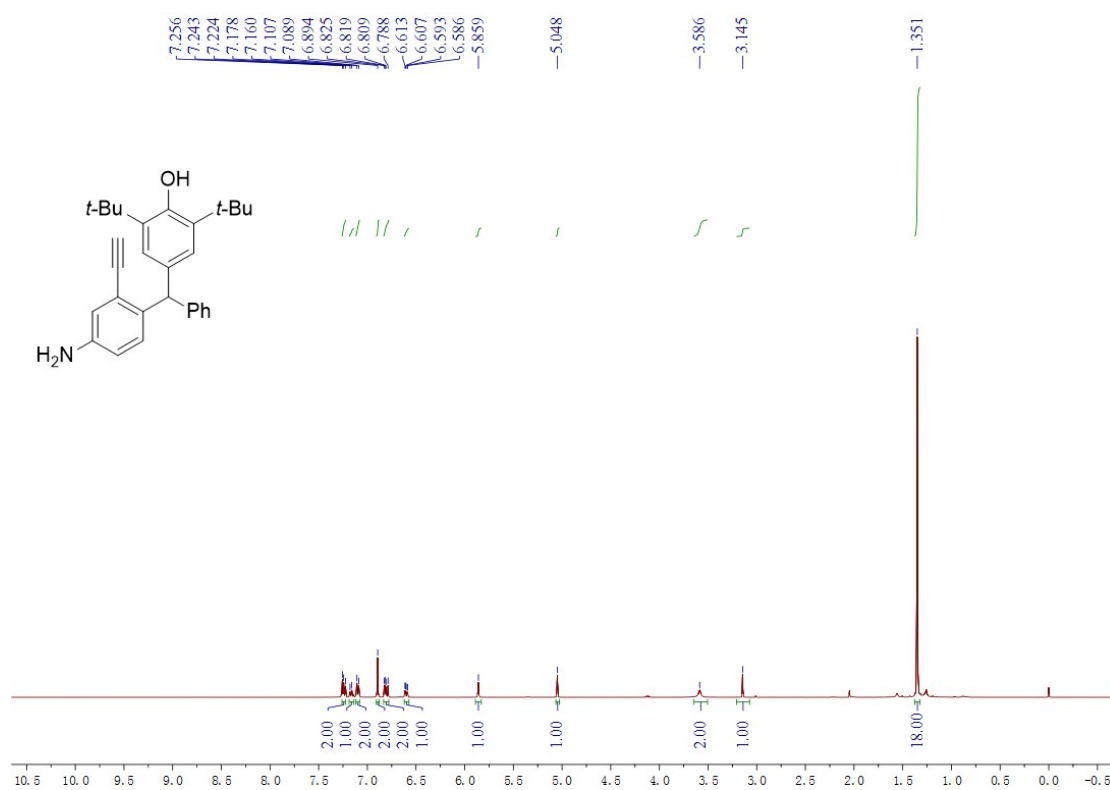


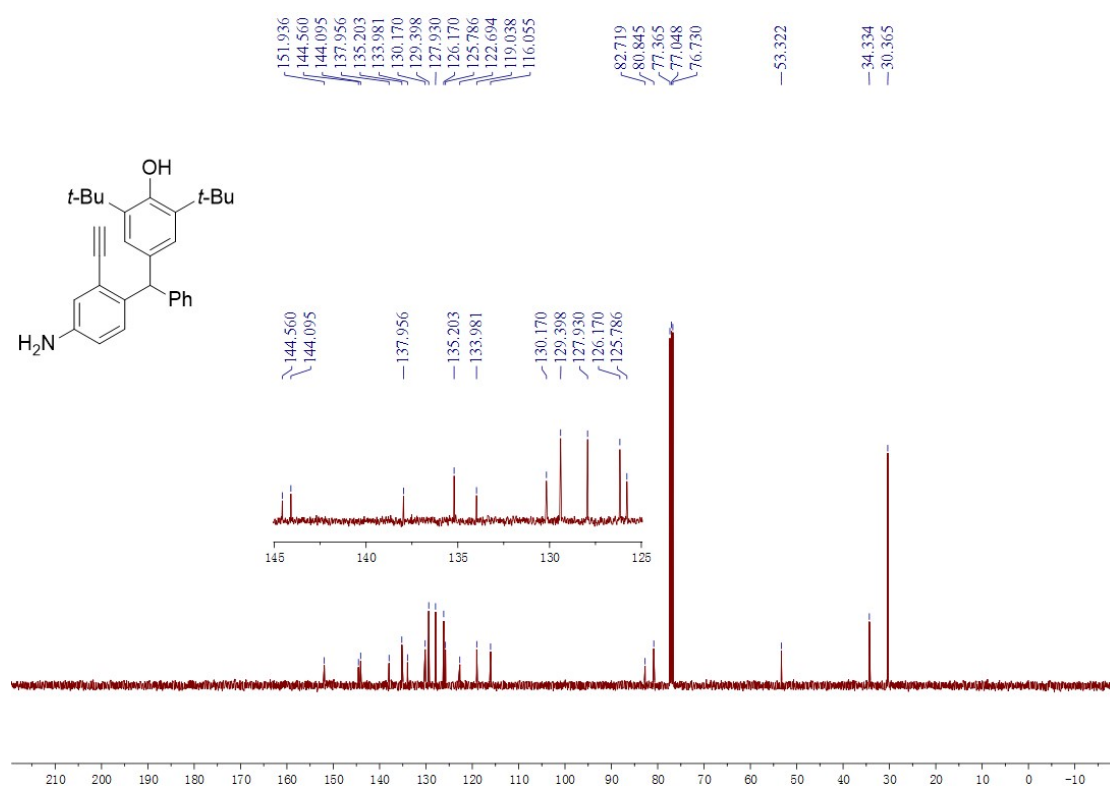
4-((4-Amino-3-iodophenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3i)



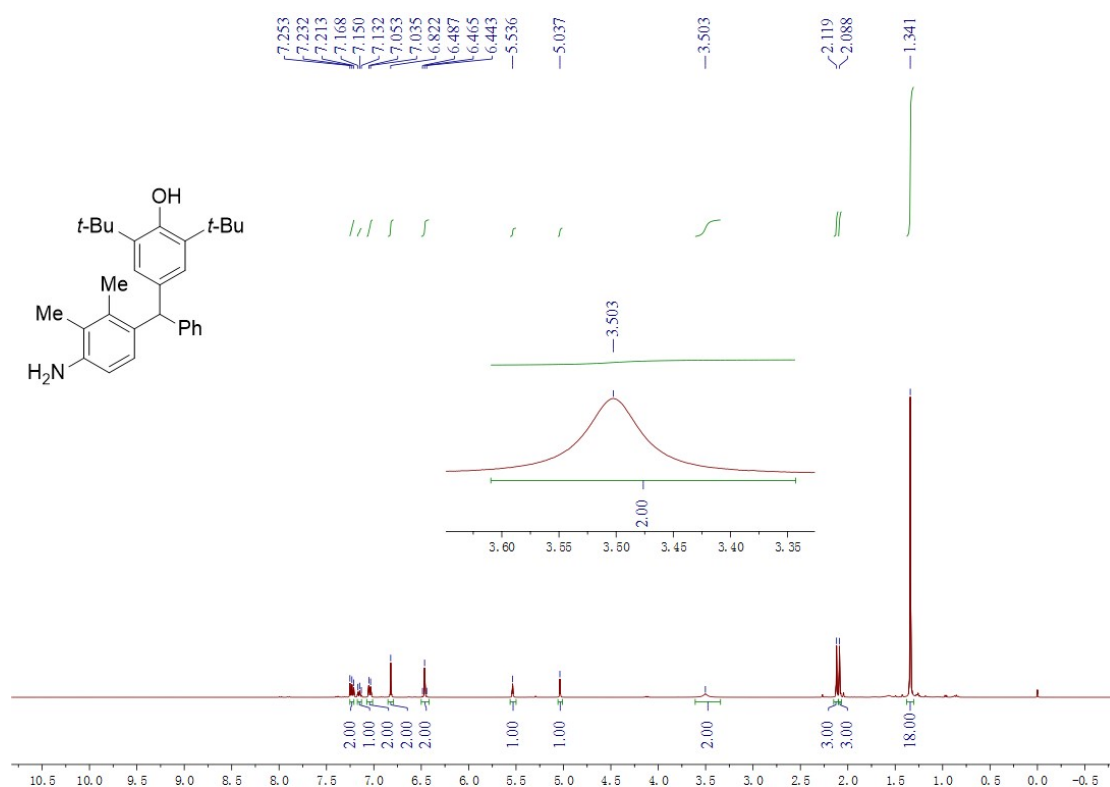


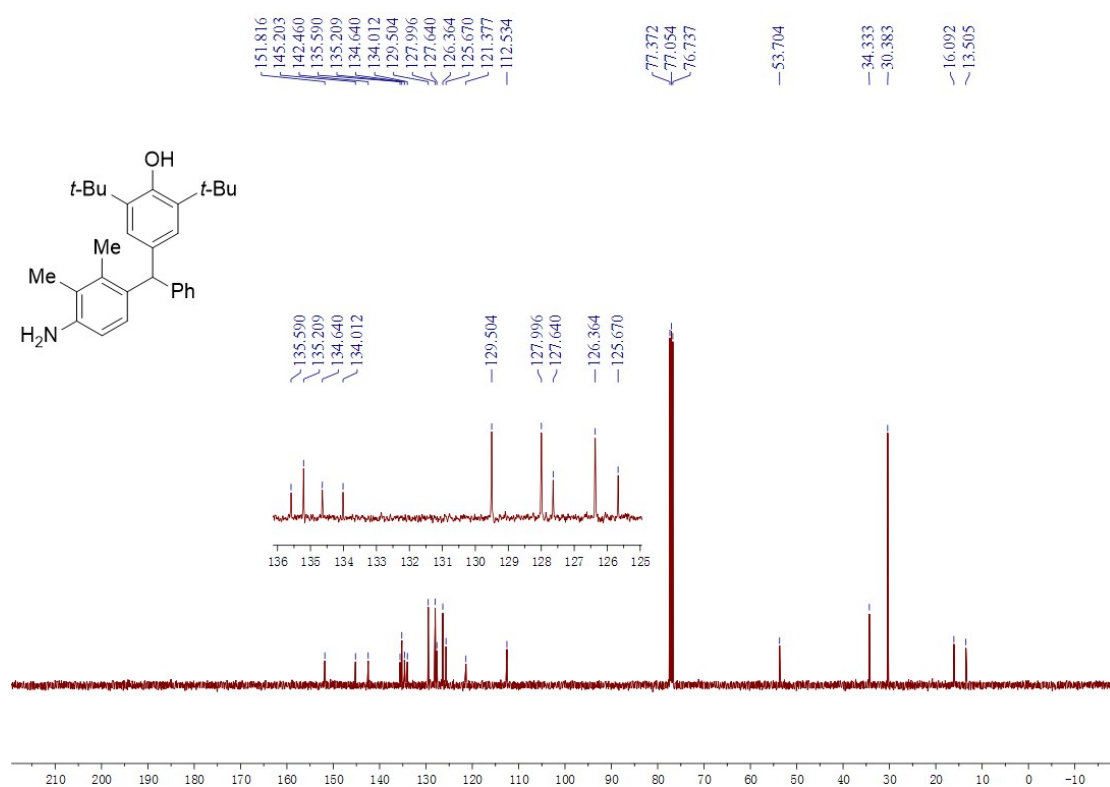
4-((4-Amino-2-ethynylphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3j)



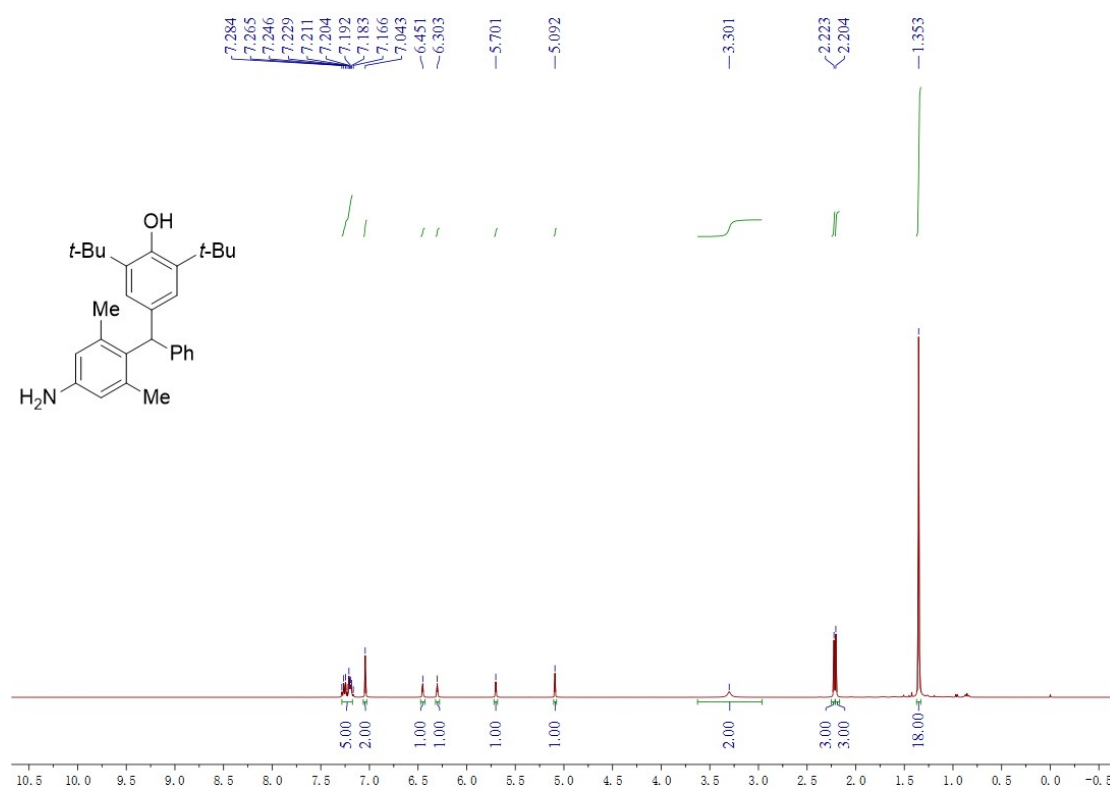


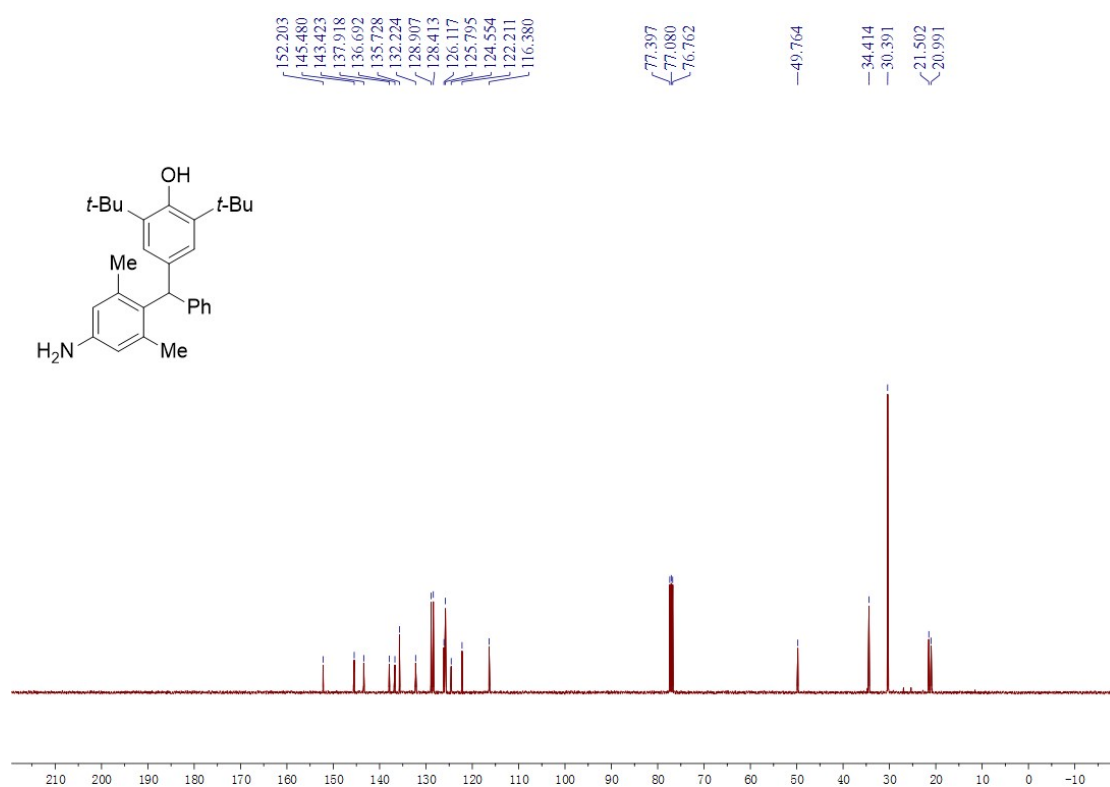
4-((4-Amino-2,3-dimethylphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3k)



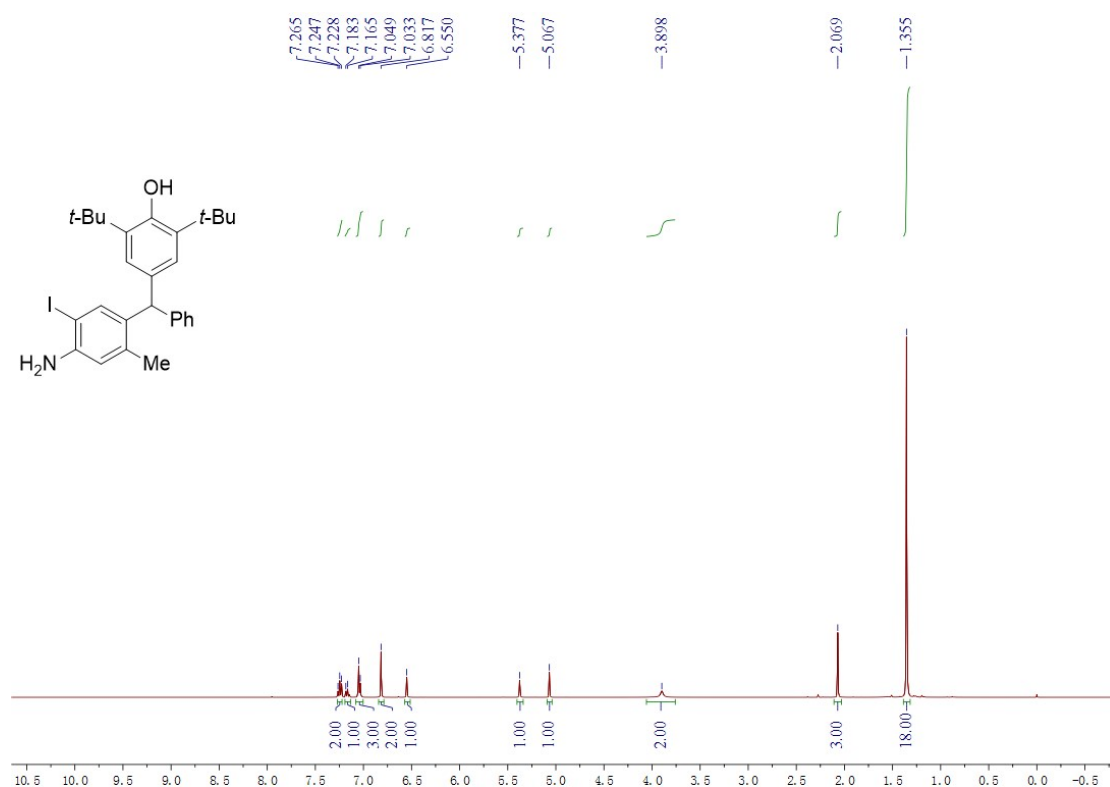


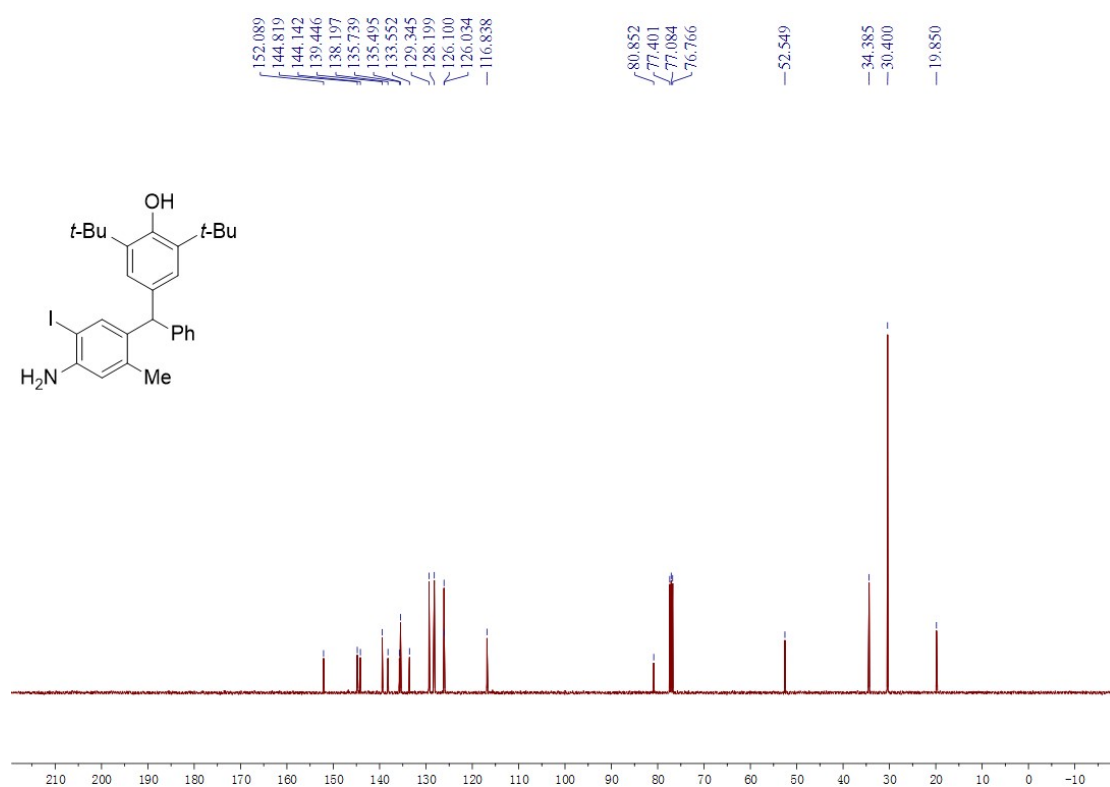
4-((4-Amino-2,6-dimethylphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3l)



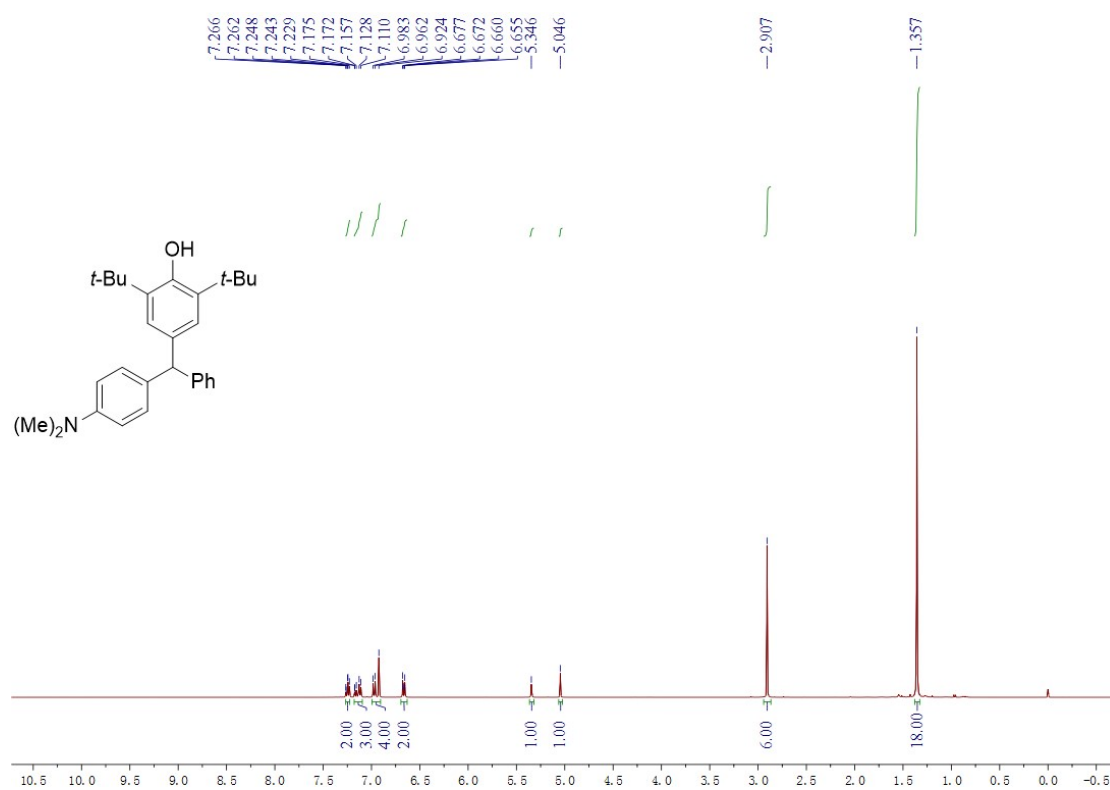


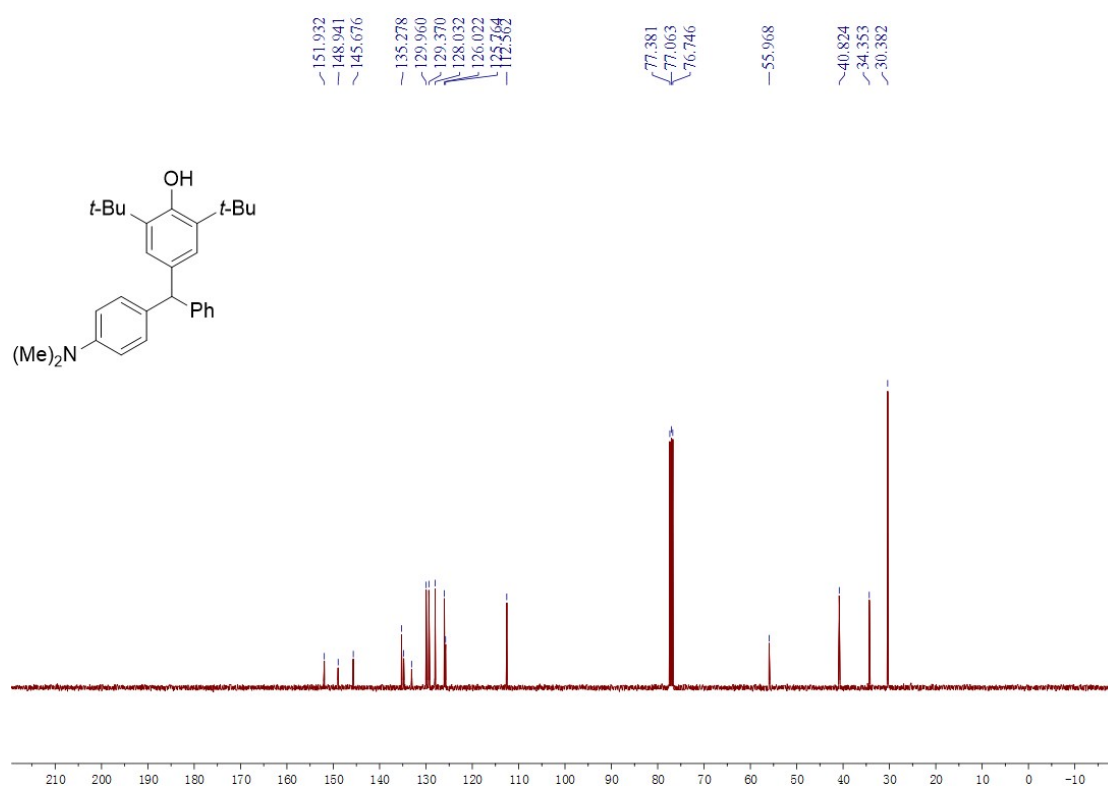
4-((4-Amino-5-iodo-2-methylphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3m)



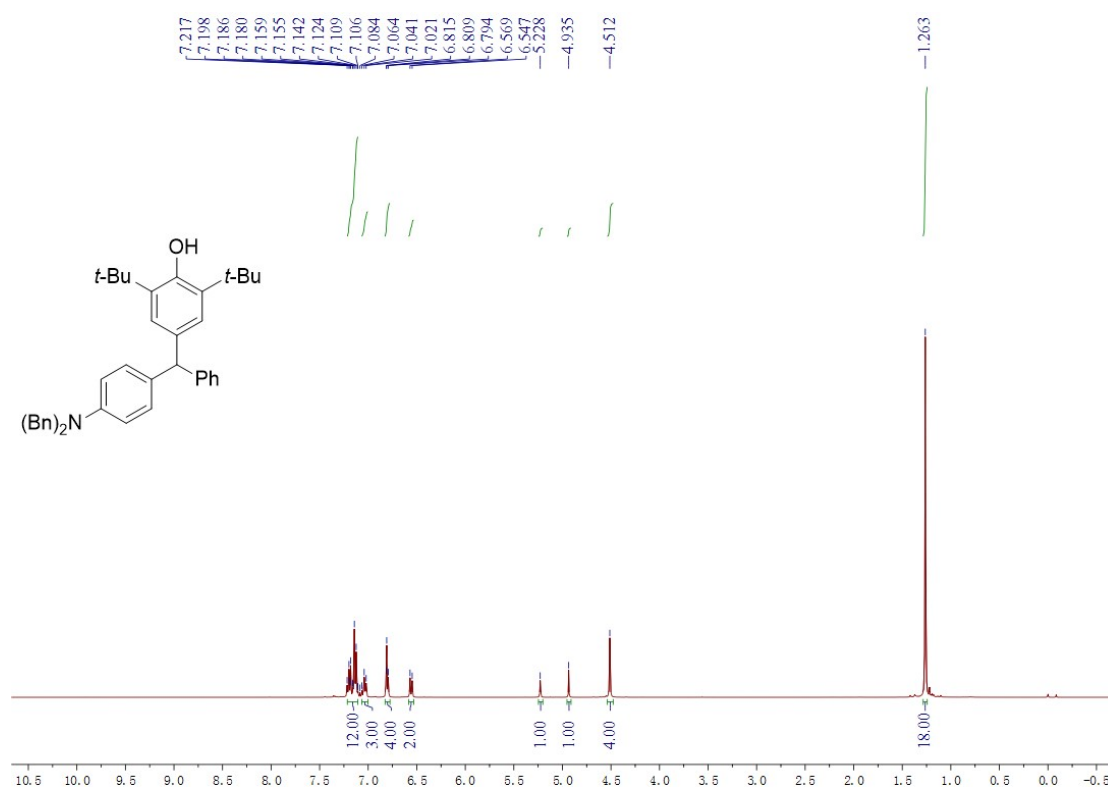


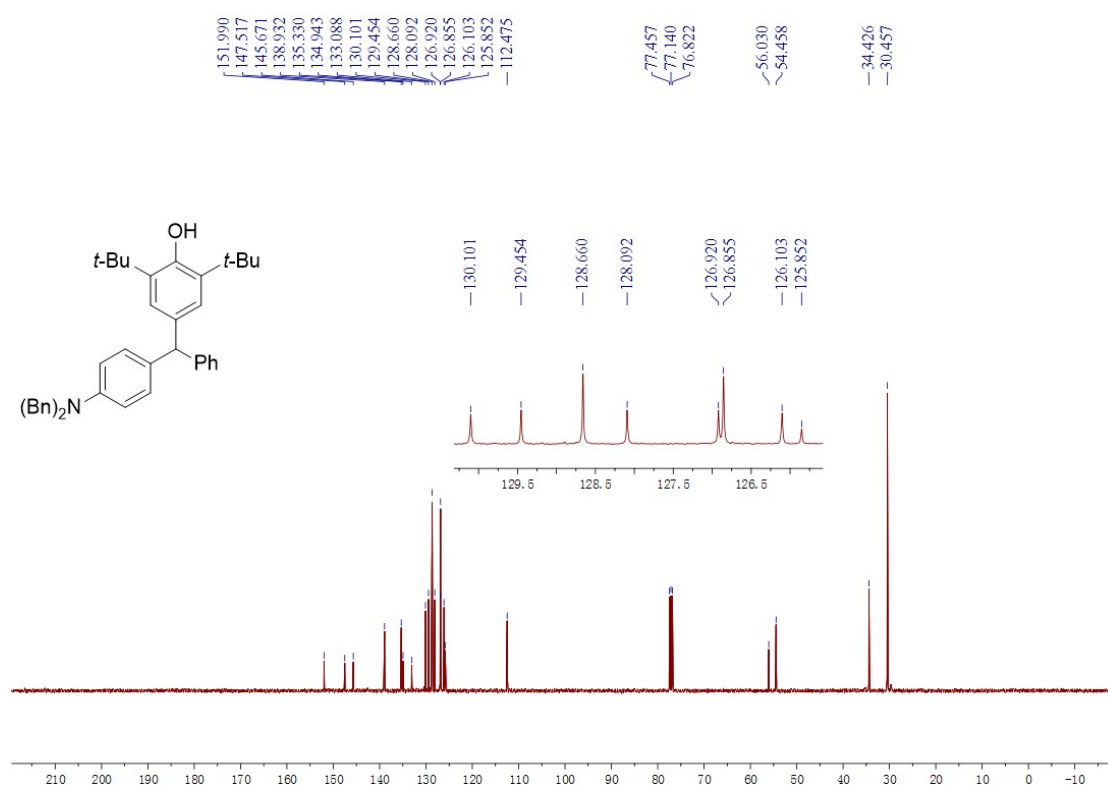
2,6-Di-tert-butyl-4-((4-(dimethylamino)phenyl)(phenyl)methyl)phenol (3n)



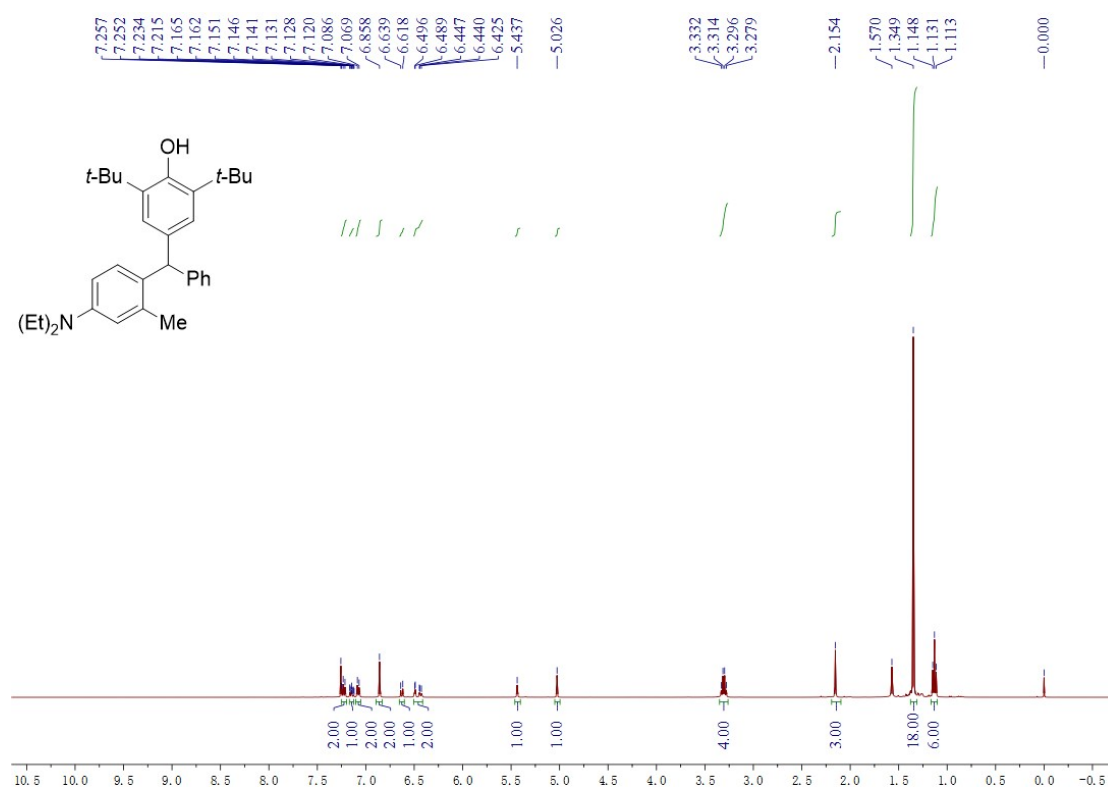


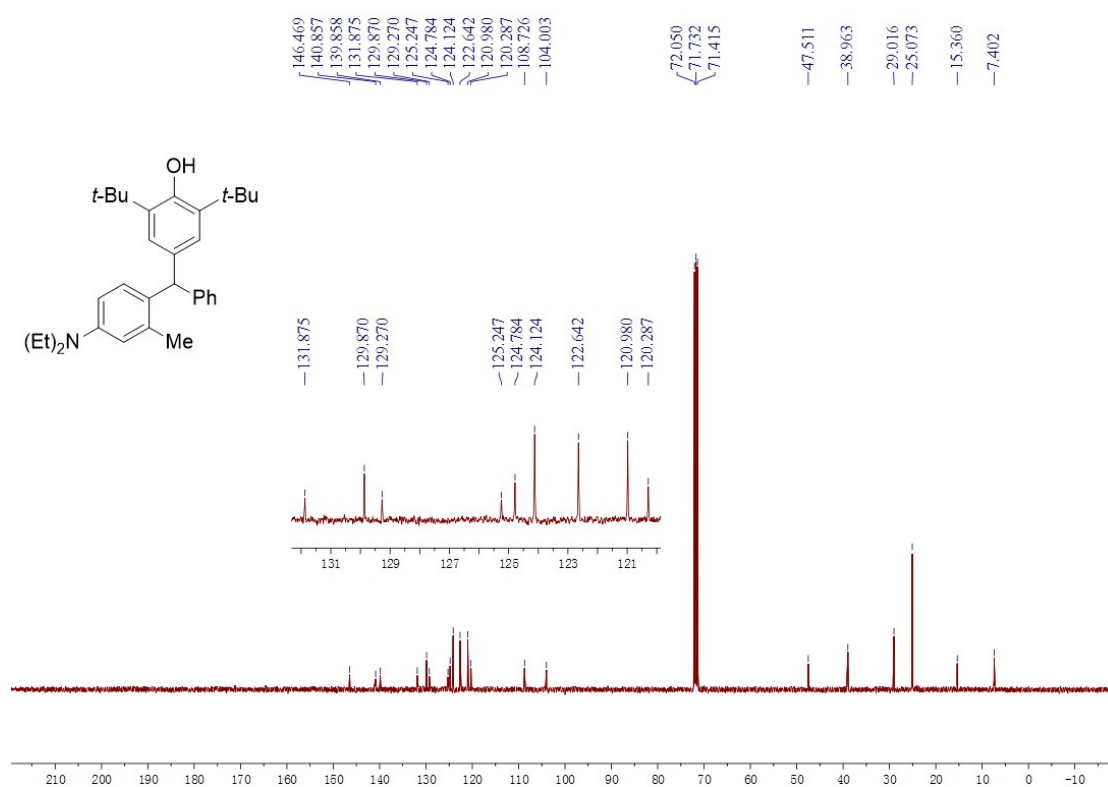
2,6-Di-tert-butyl-4-((4-(dibenzylamino)phenyl)(phenyl)methyl)phenol (3o)



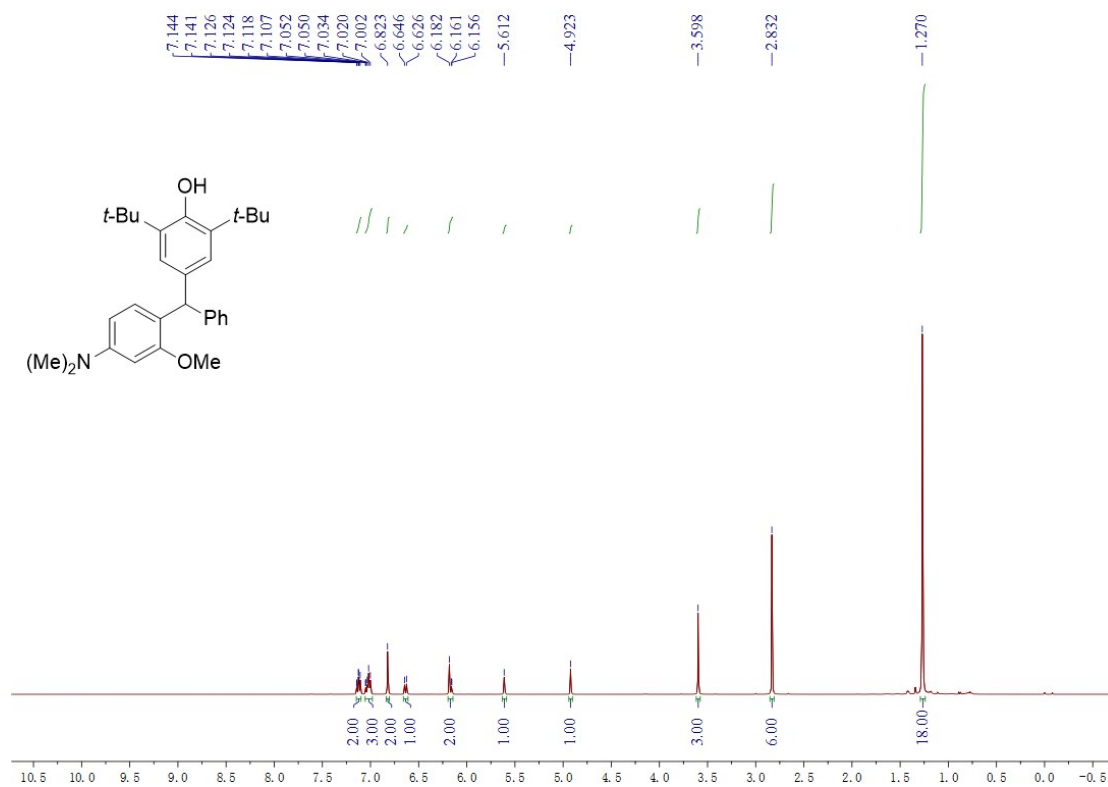


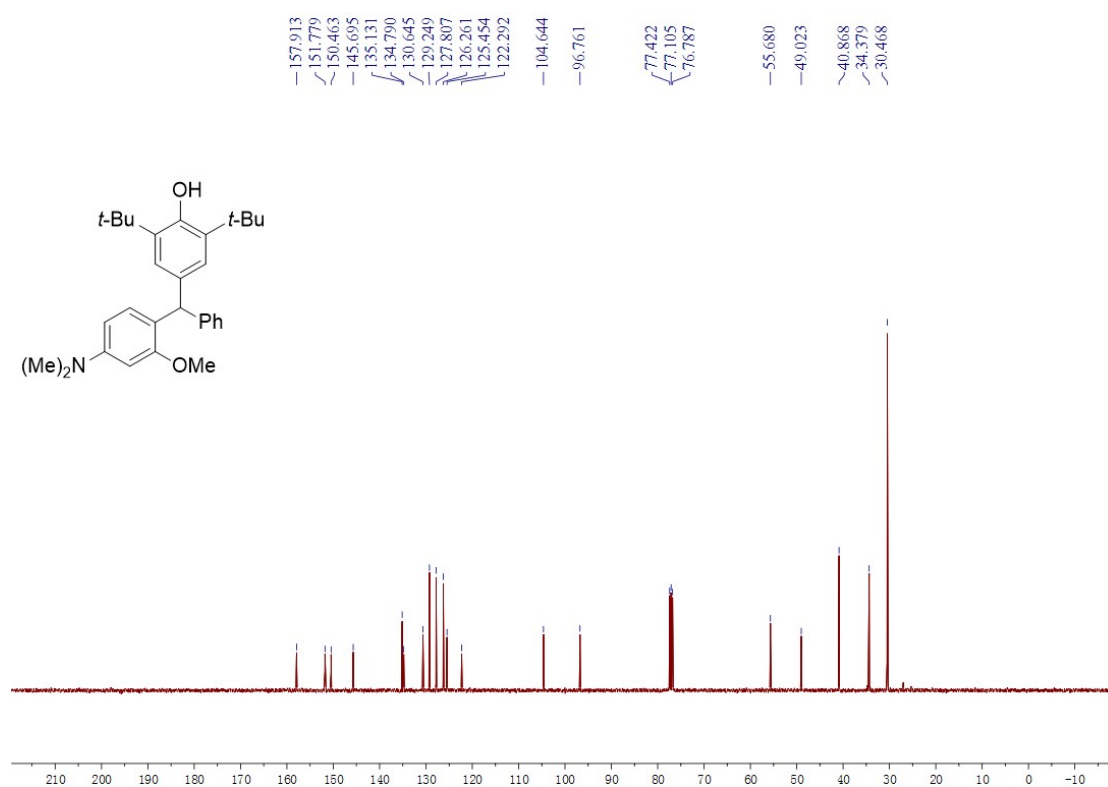
2,6-Di-tert-butyl-4-((4-(diethylamino)-2-methylphenyl)(phenyl)methyl)phenol (3p)



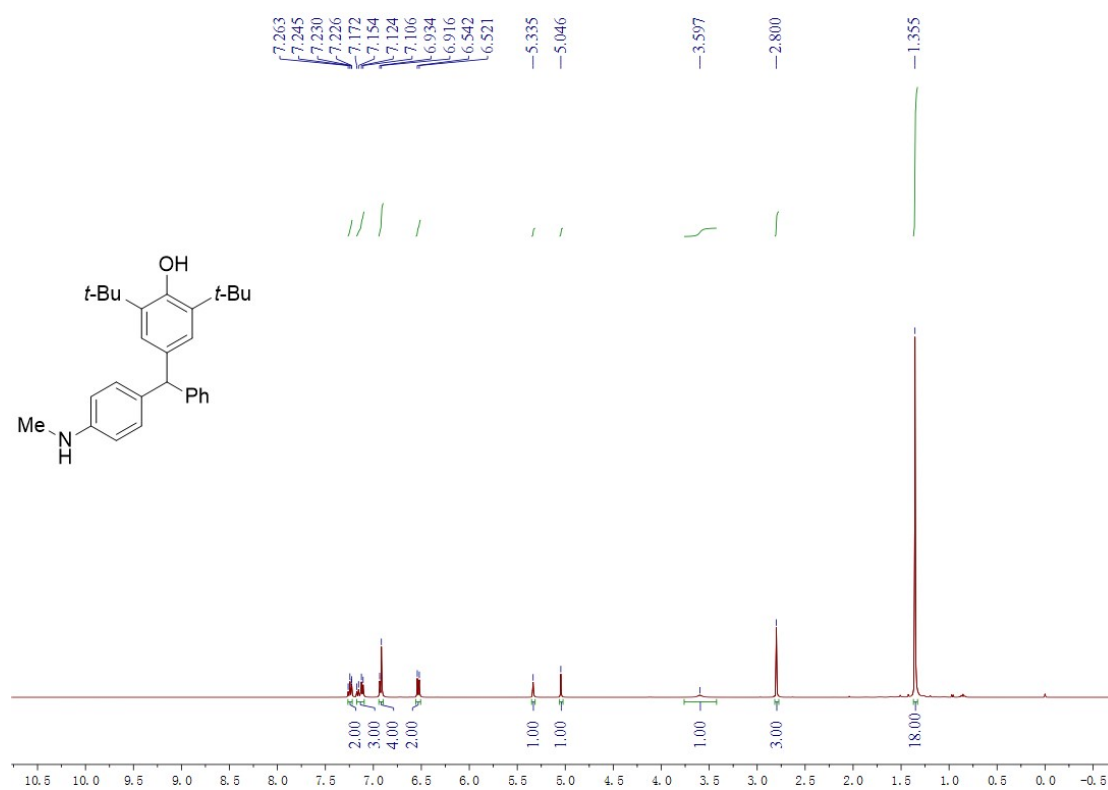


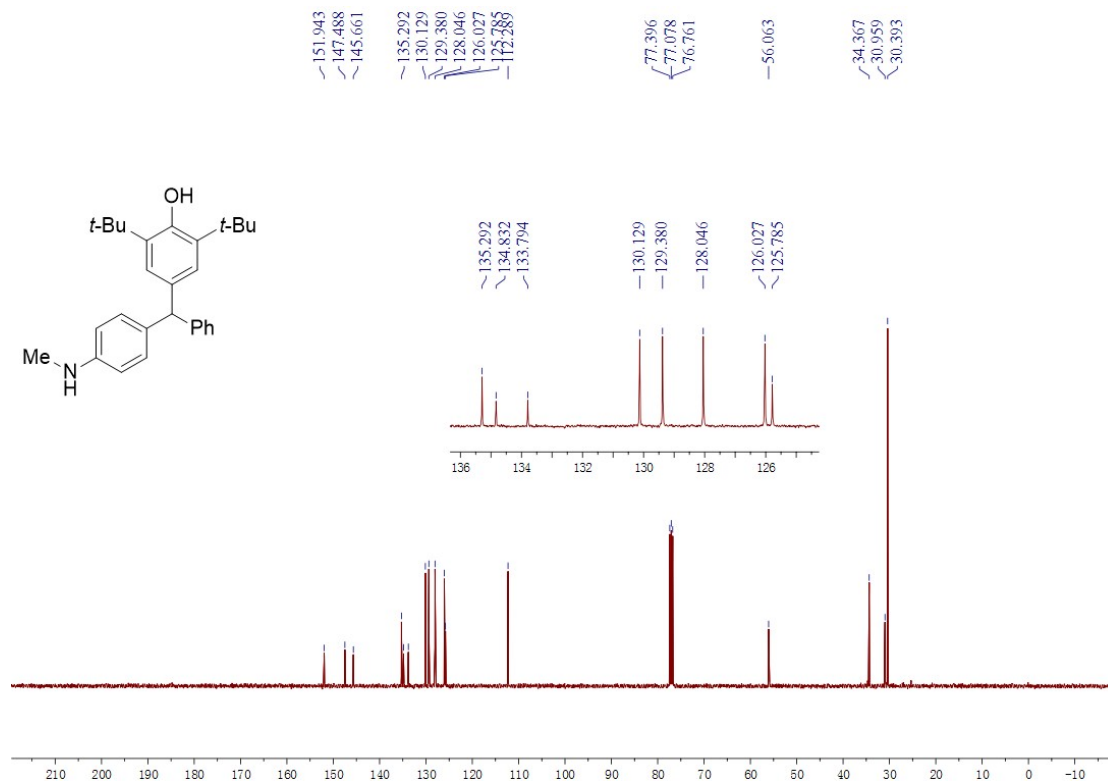
2,6-Di-tert-butyl-4-((4-(dimethylamino)-2-methoxyphenyl)(phenyl)methyl)phenol (3q)



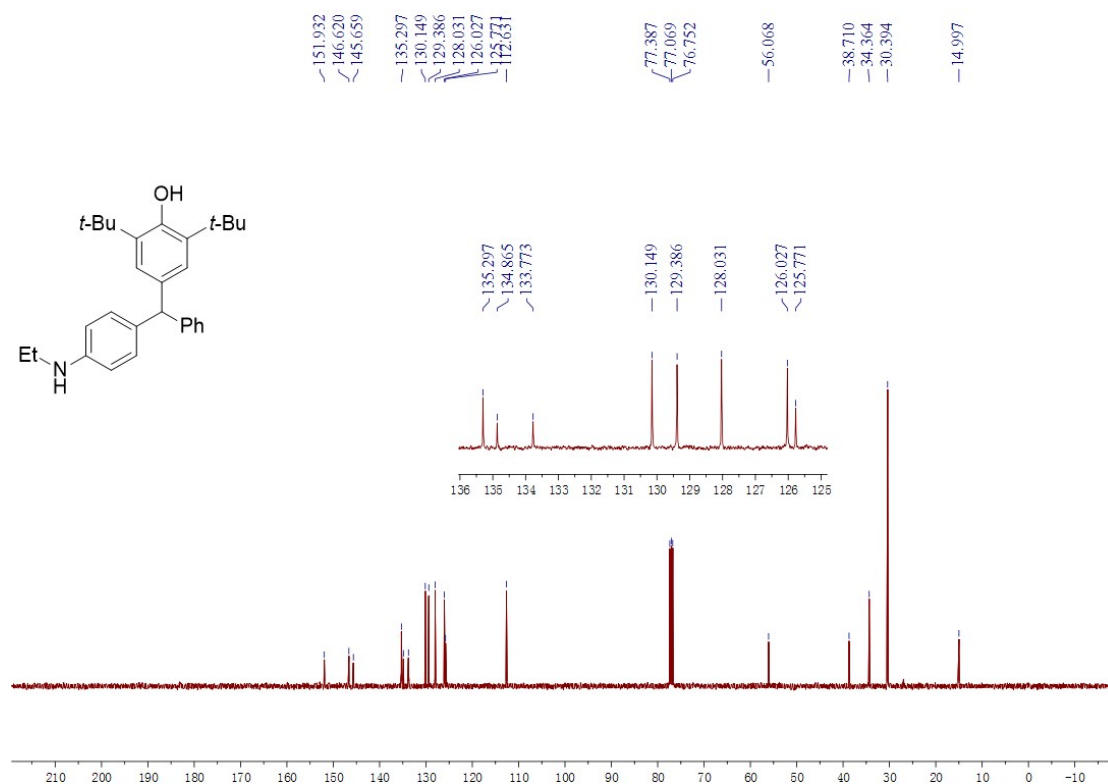


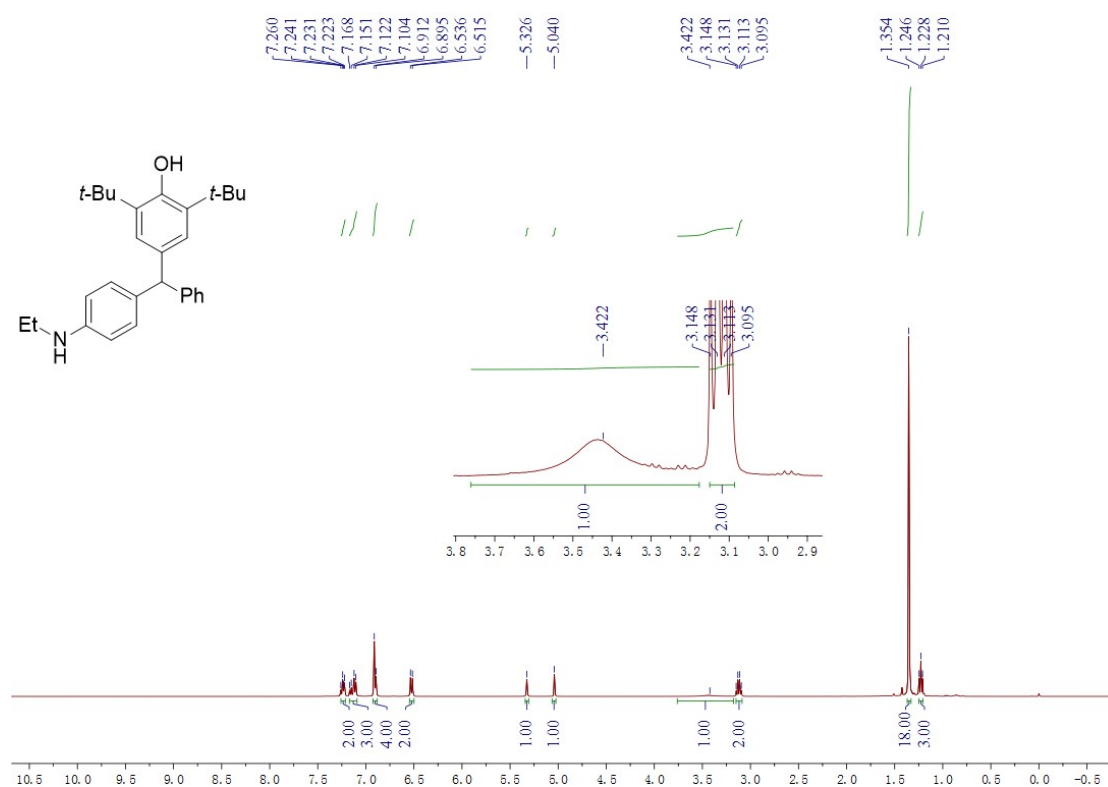
2,6-Di-tert-butyl-4-((4-(methylamino)phenyl)(phenyl)methyl)phenol (3r)



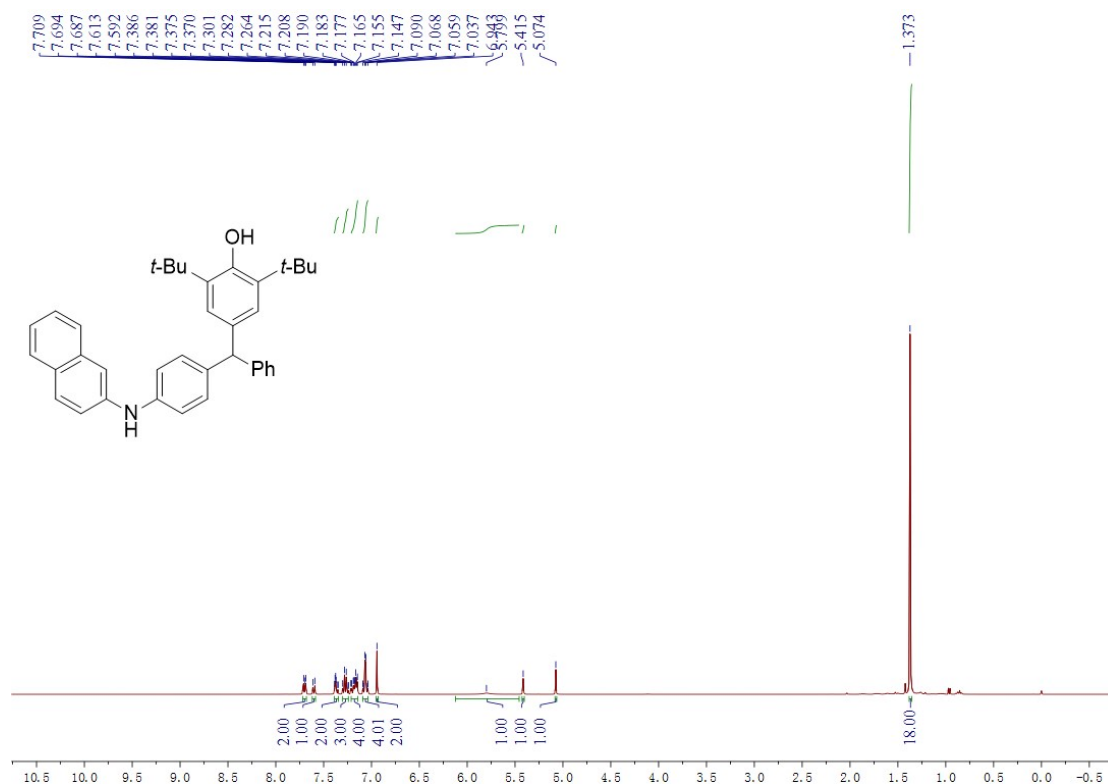


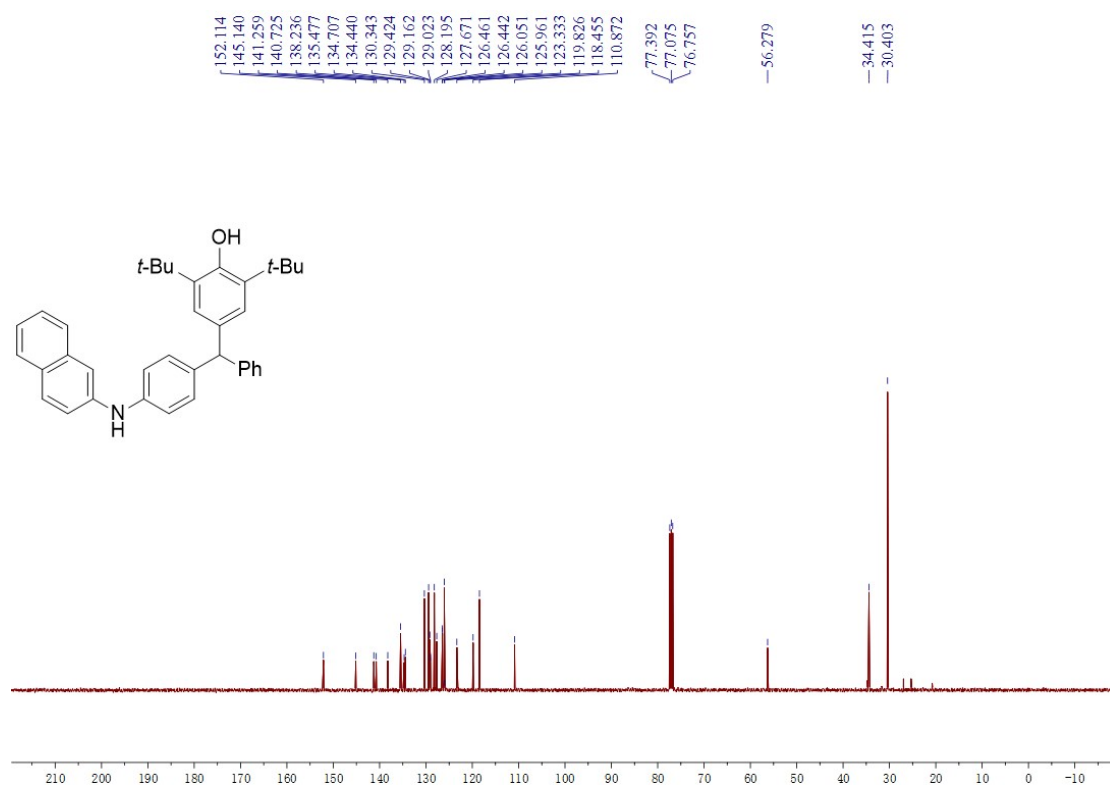
2,6-Di-*tert*-butyl-4-((4-(ethylamino)phenyl)(phenyl)methyl)phenol (3s)



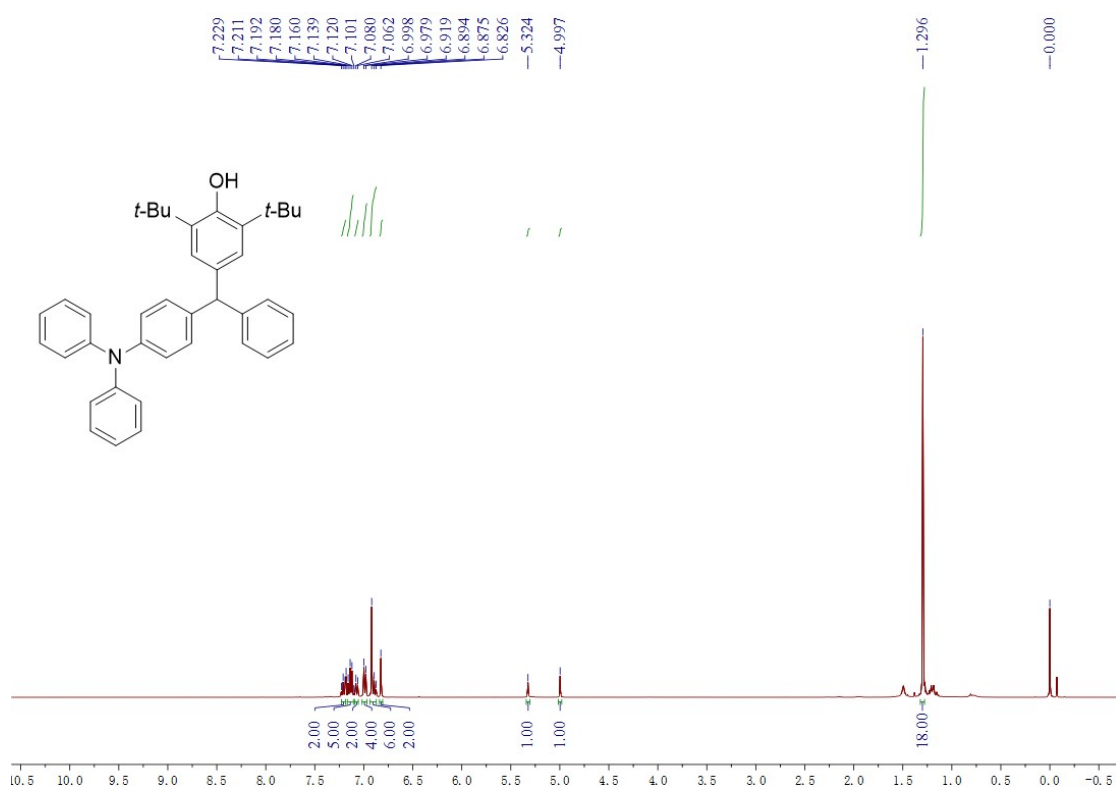


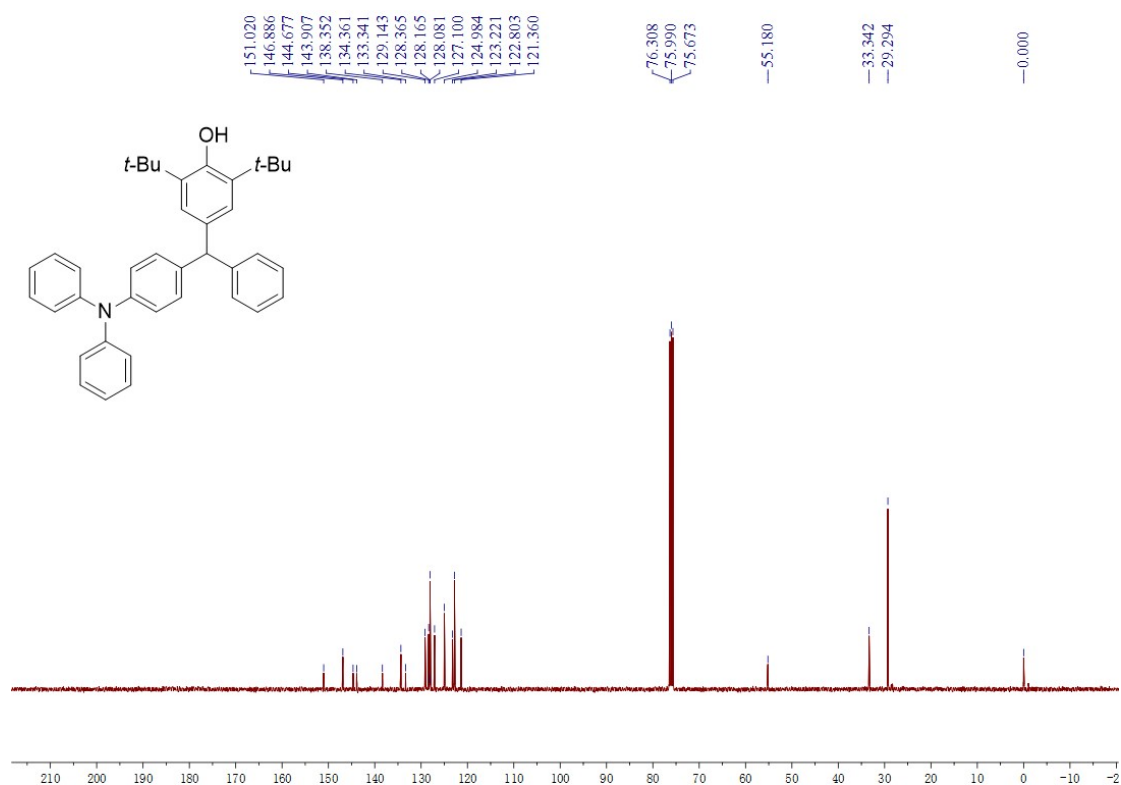
2,6-Di-tert-butyl-4-((4-(naphthalen-2-ylamino)phenyl)(phenyl)methyl)phenol (3t)



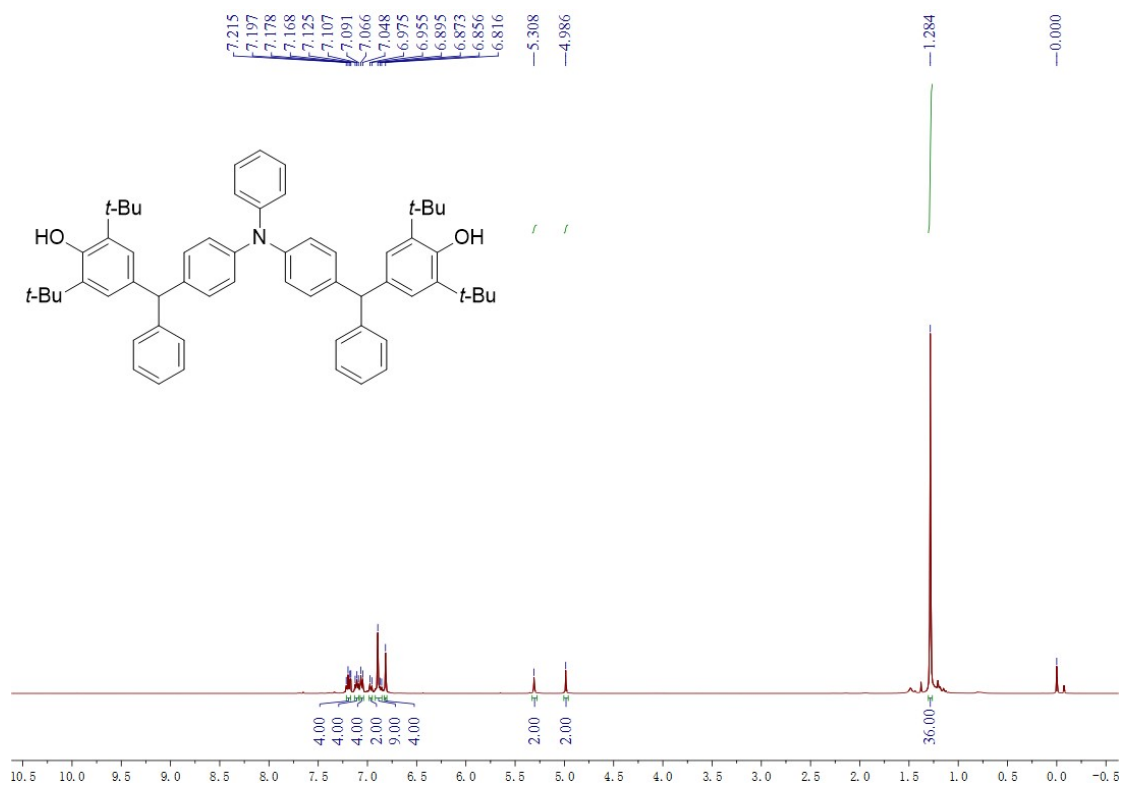


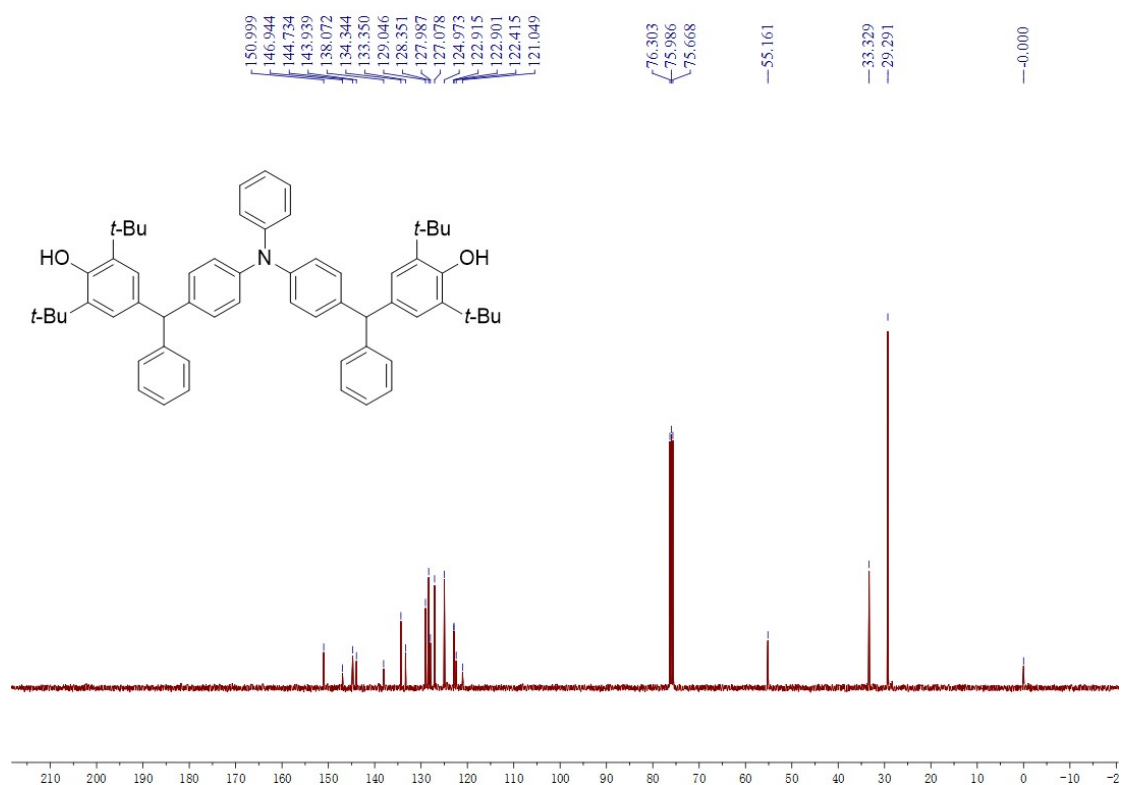
2,6-Di-tert-butyl-4-((4-(diphenylamino)phenyl)(phenyl)methyl)phenol (3u)



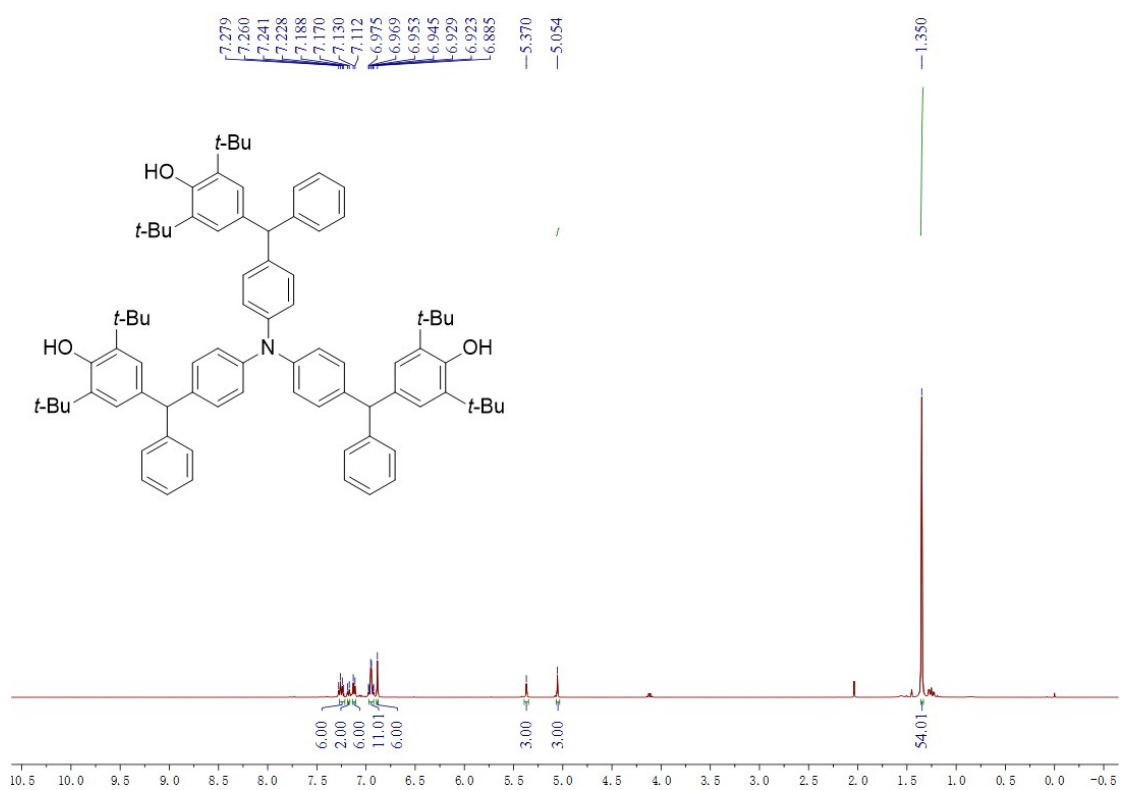


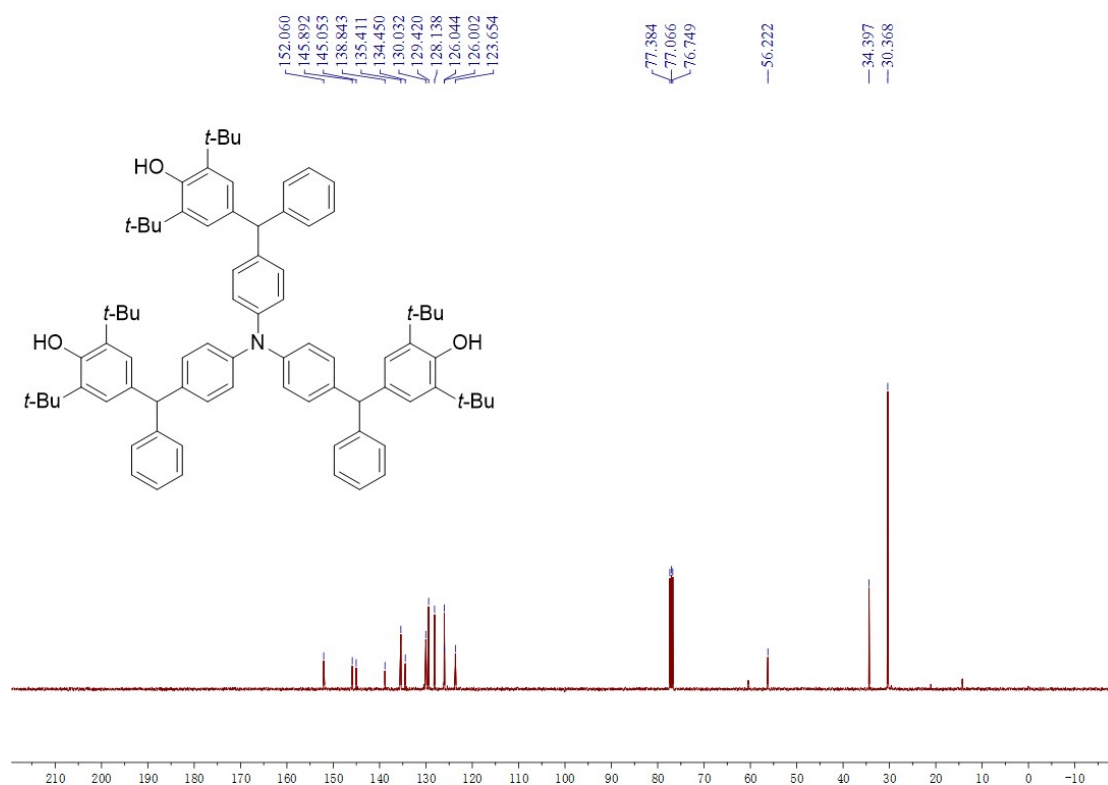
4,4'-(((Phenylazanediy)bis(4,1-phenylene))bis(phenylmethylene))bis(2,6-di-tert-butylphenol) (3u')



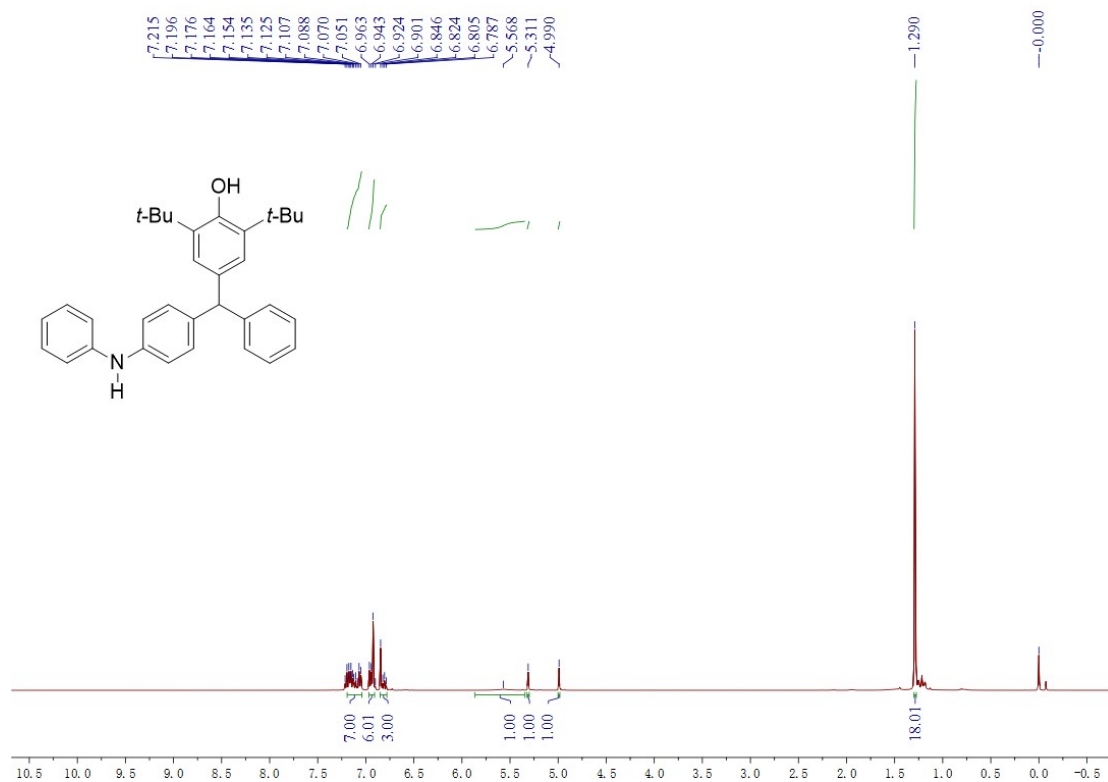


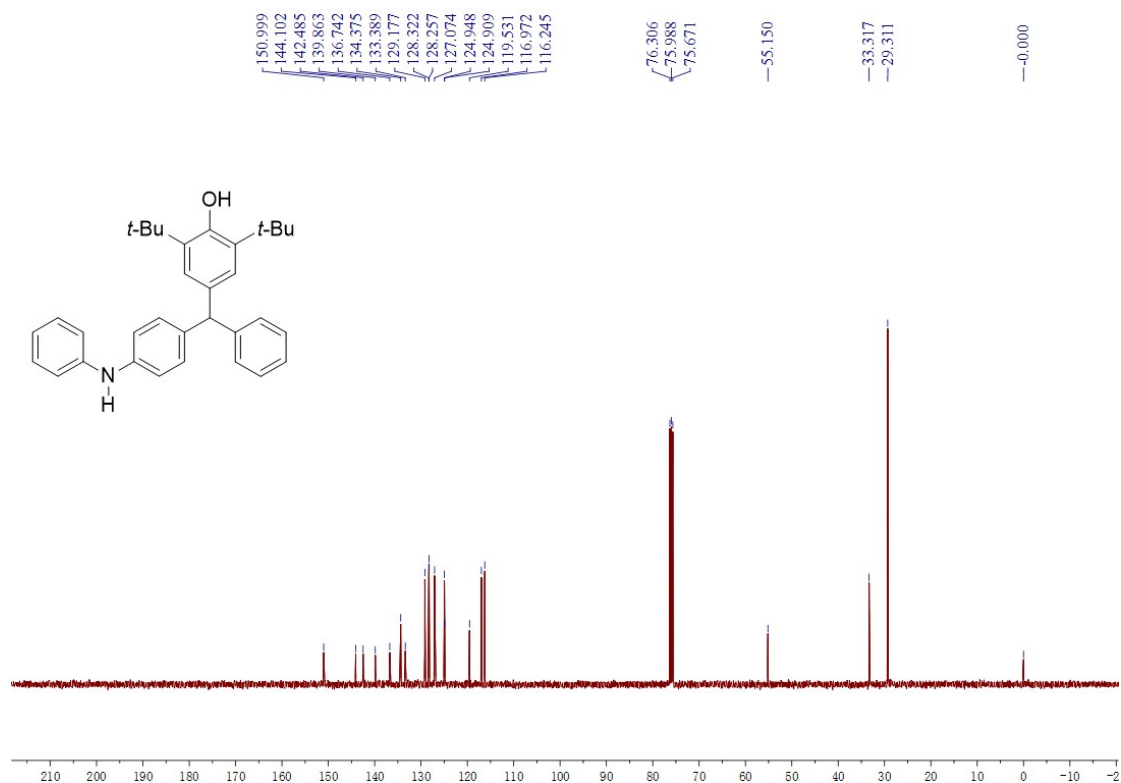
4,4',4''-((Nitrilotris(benzene-4,1-diyl))tris(phenylmethylene))tris(2,6-di-tert-butylphenol) (3u'')



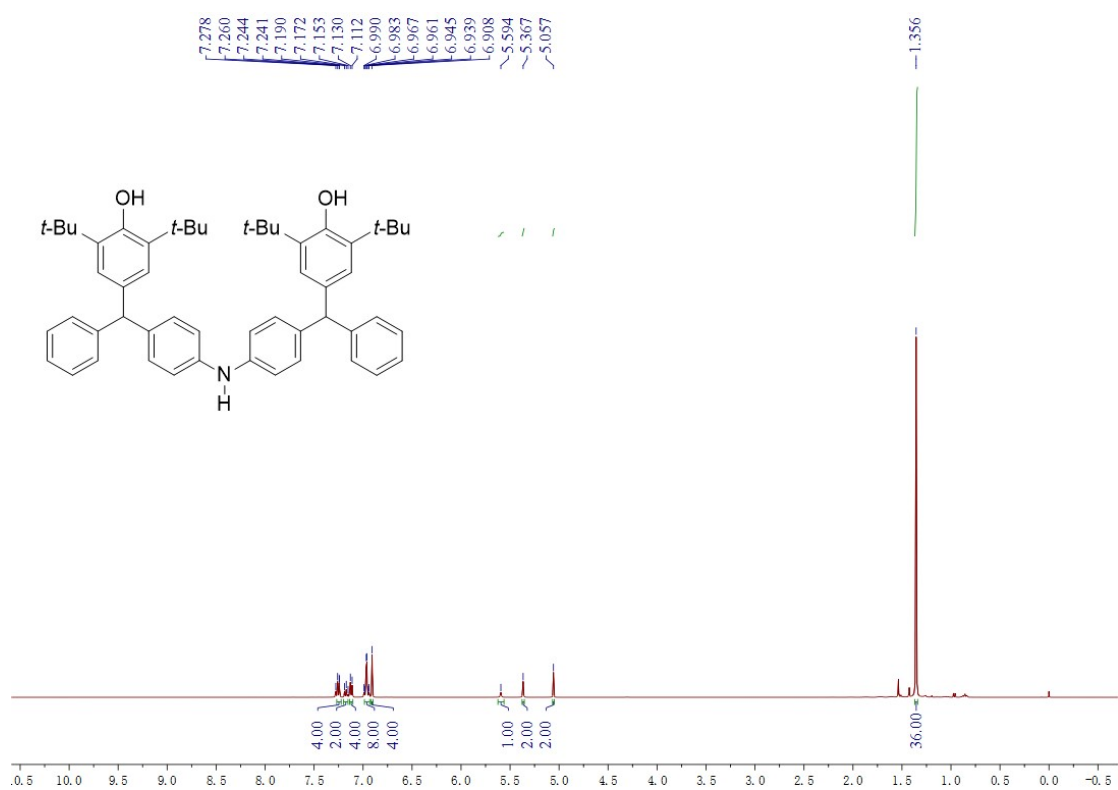


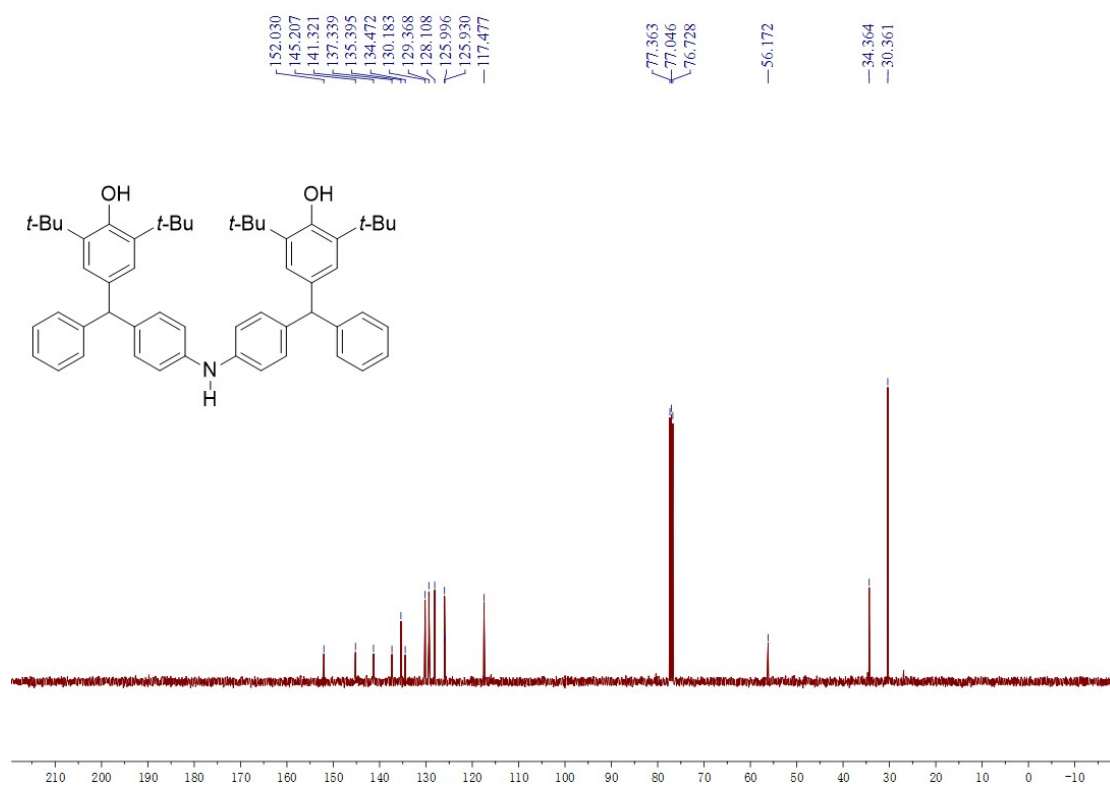
2,6-Di-tert-butyl-4-(phenyl(4-(phenylamino)phenyl)methyl)phenol (3v)



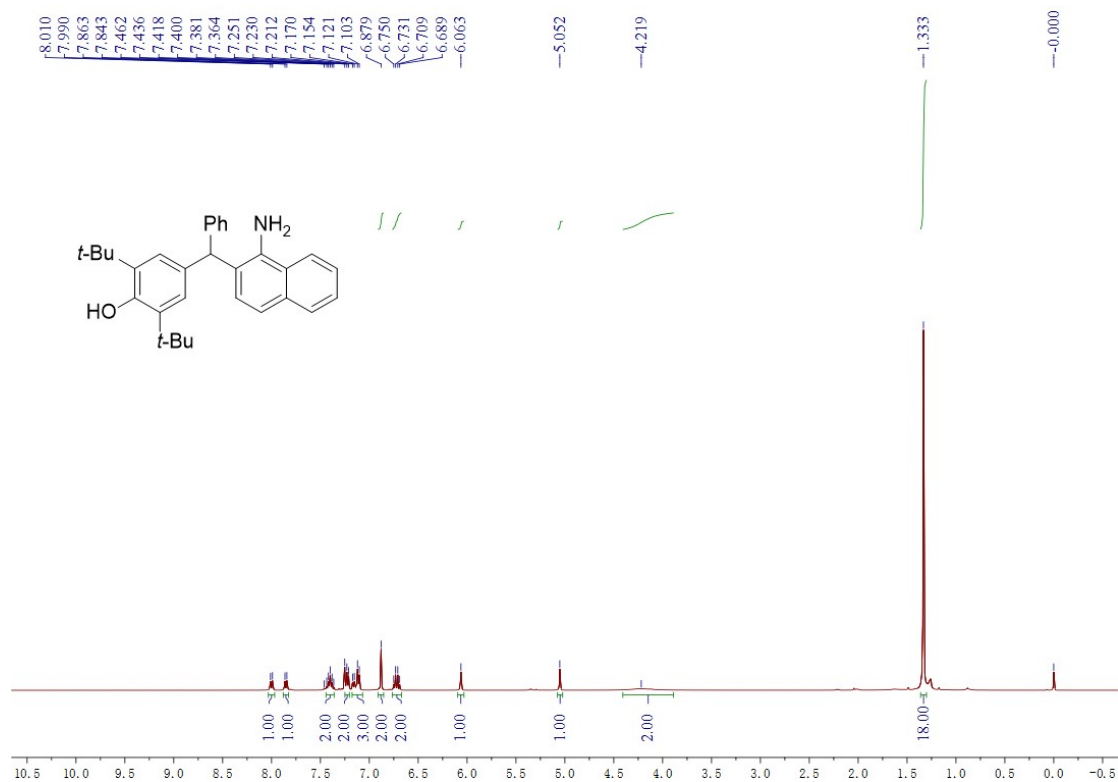


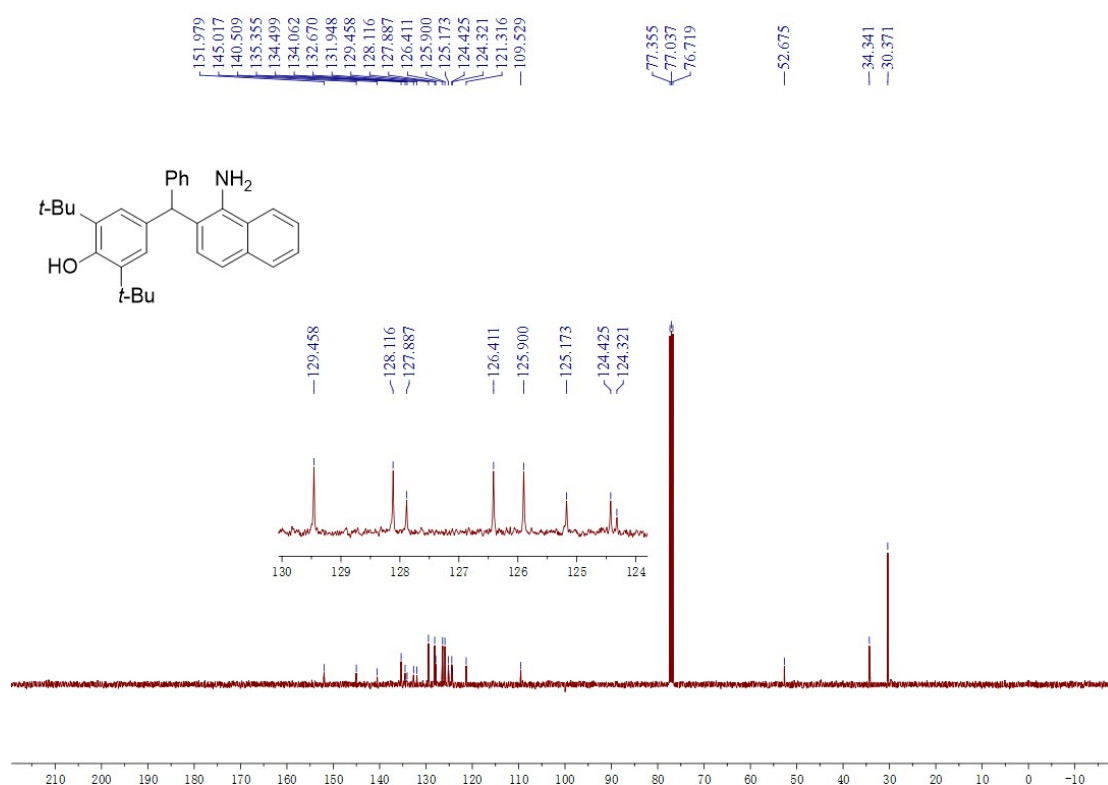
4,4'-((Azanediylbis(4,1-phenylene))bis(phenylmethylene))bis(2,6-di-tert-butylphenol) (3v')



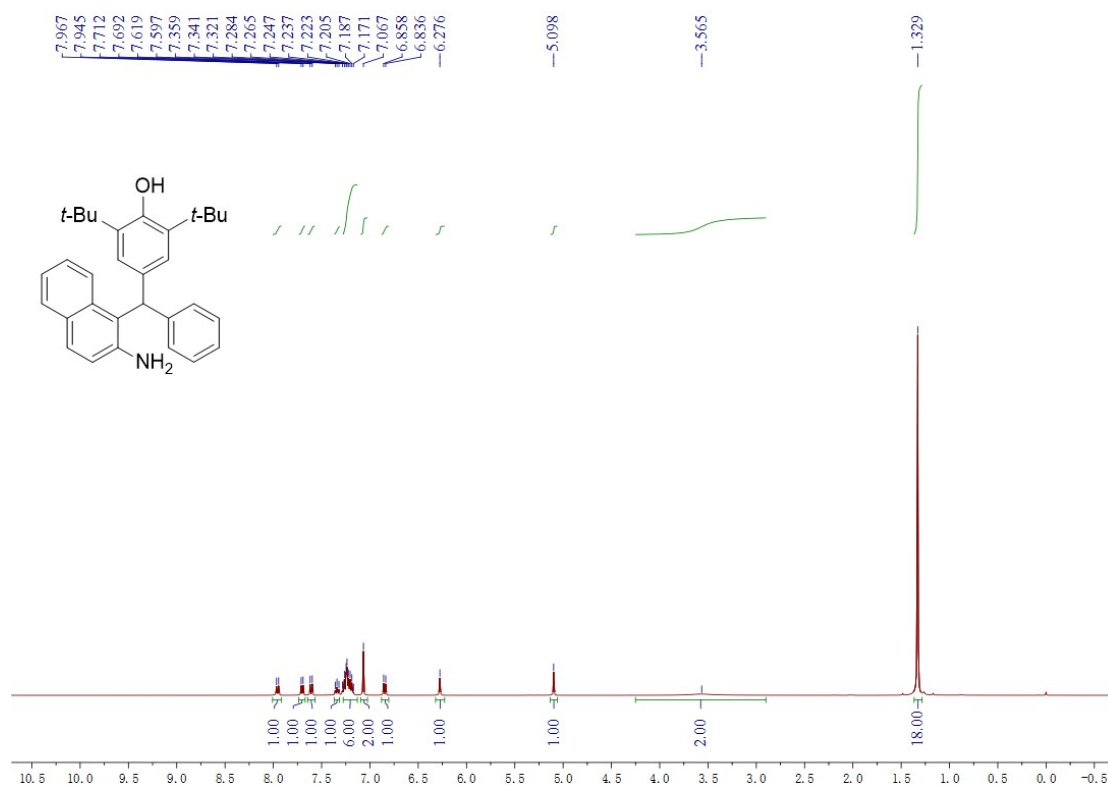


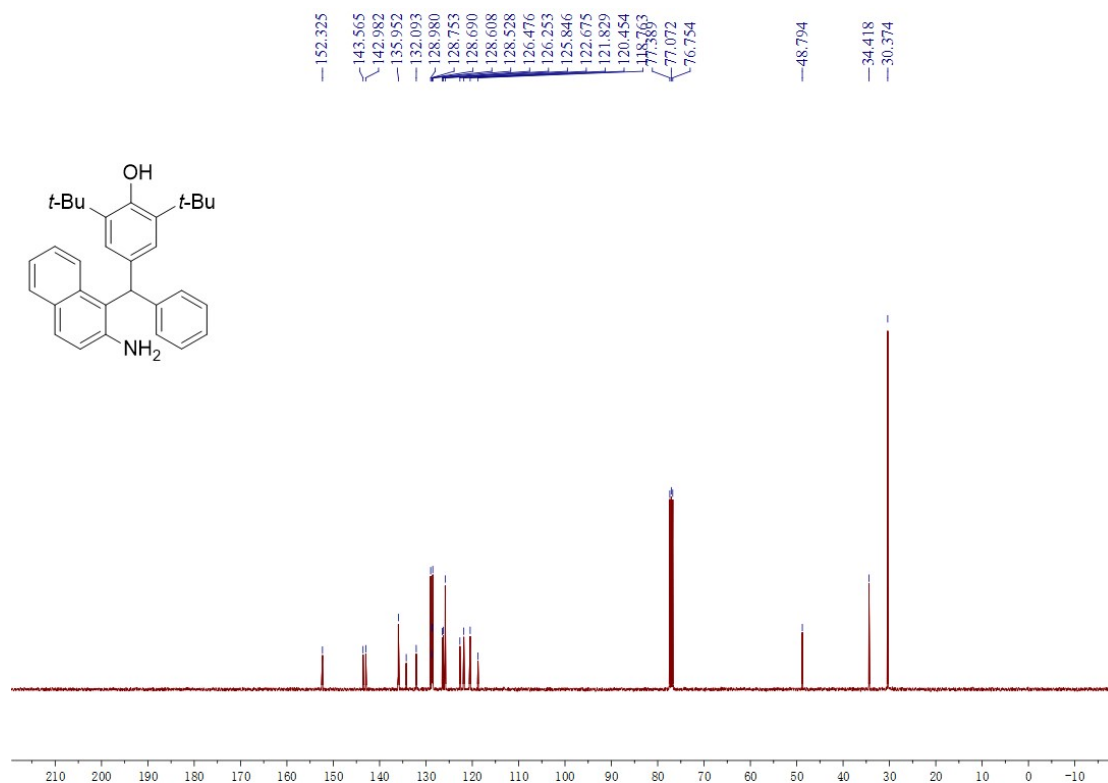
4-((1-Aminonaphthalen-2-yl)(phenyl)methyl)-2,6-di-tert-butylphenol (3w)



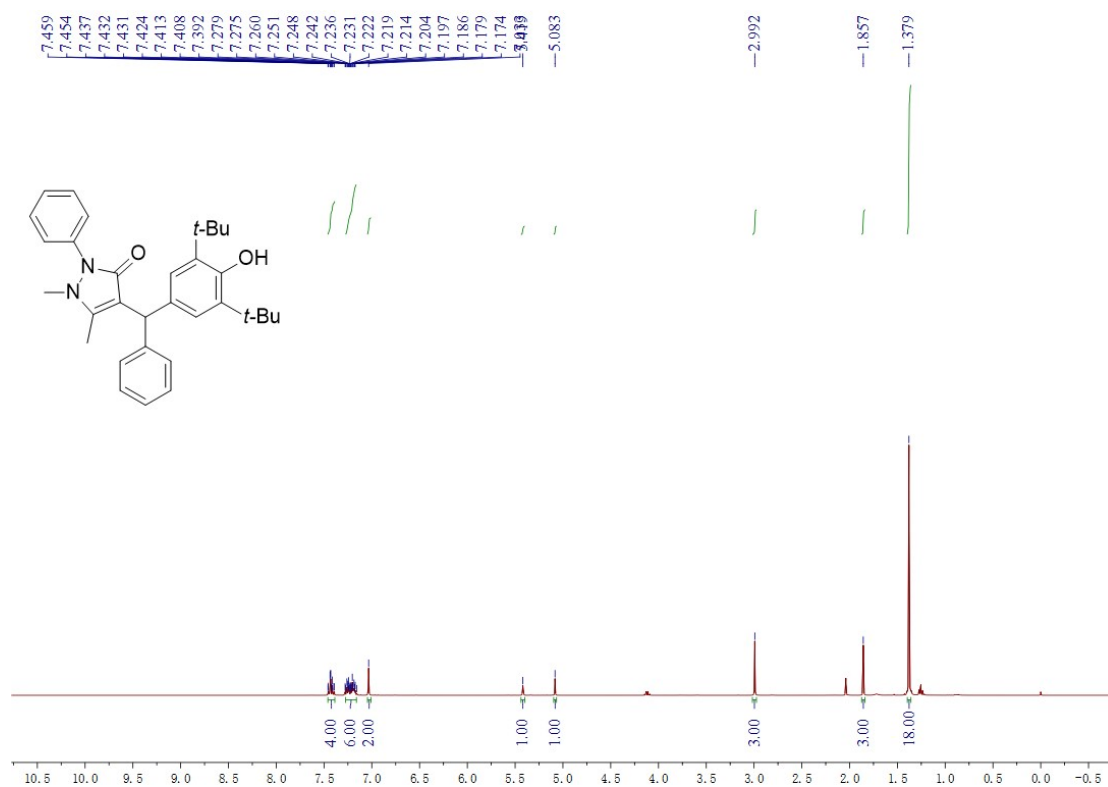


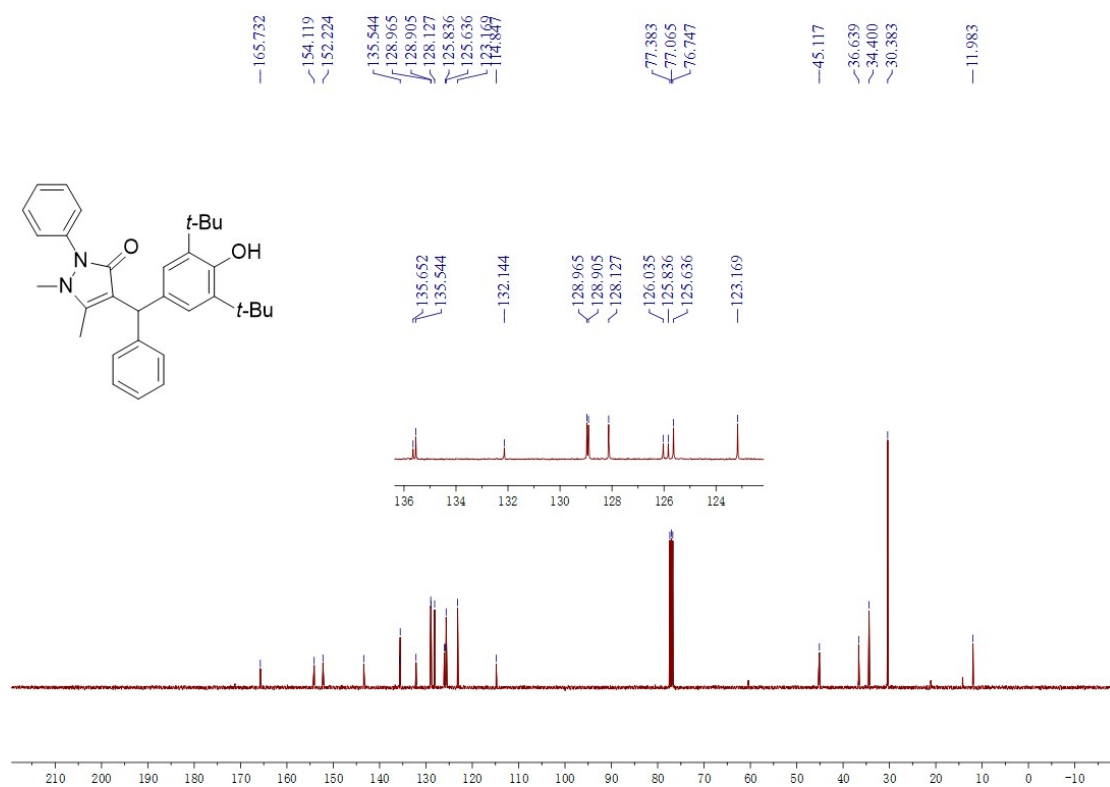
4-((2-Aminonaphthalen-1-yl)(phenyl)methyl)-2,6-di-tert-butylphenol (3x)



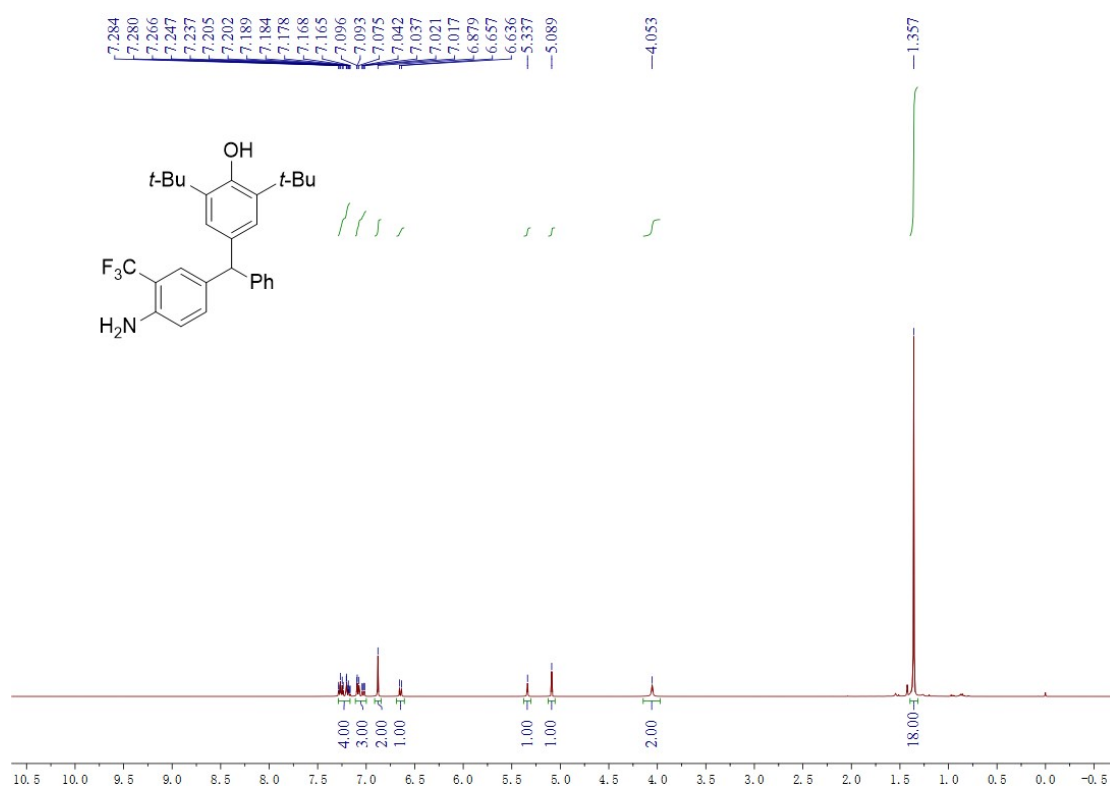


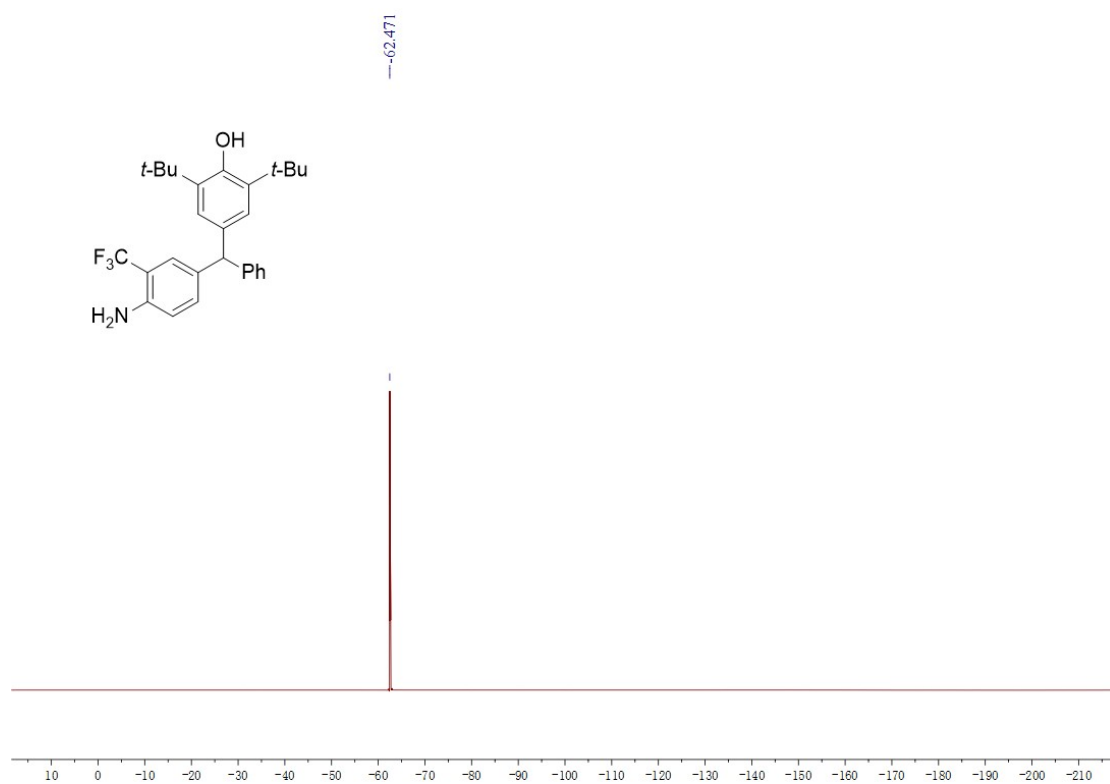
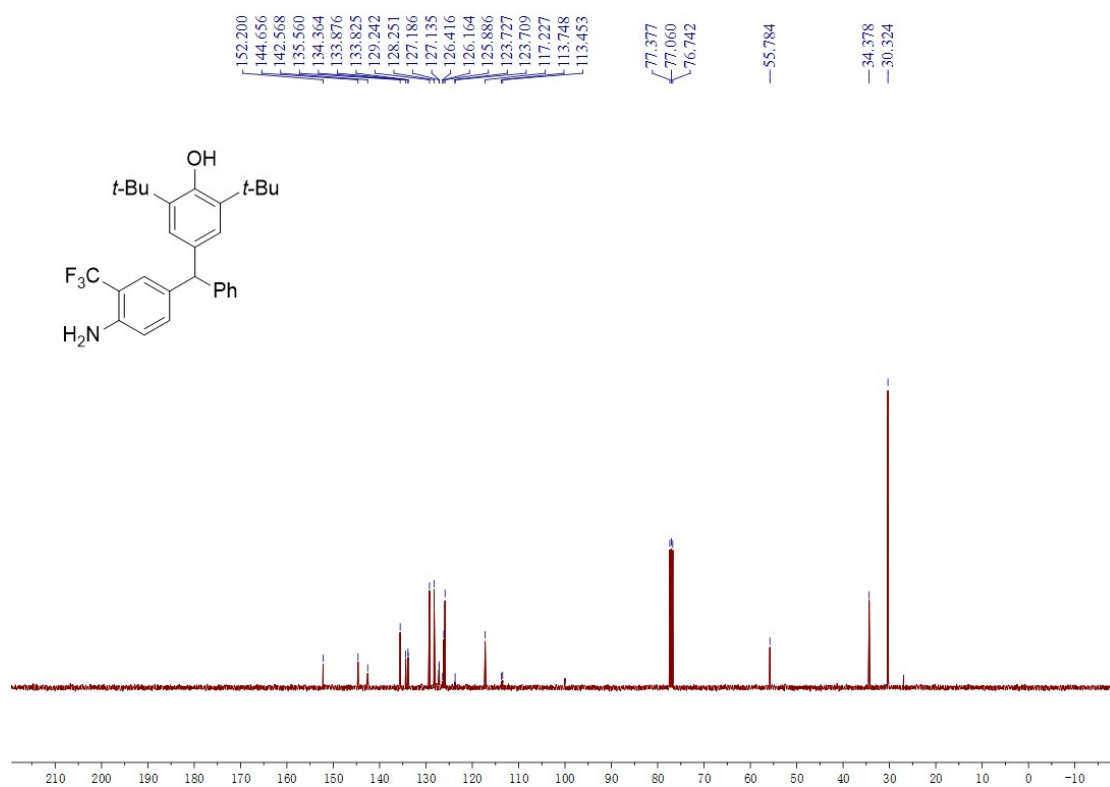
4-((3,5-Di-tert-butyl-4-hydroxyphenyl)(phenyl)methyl)-1,5-dimethyl-2-phenyl-1,2-dihydro-3H-pyrazol-3-one (3y)



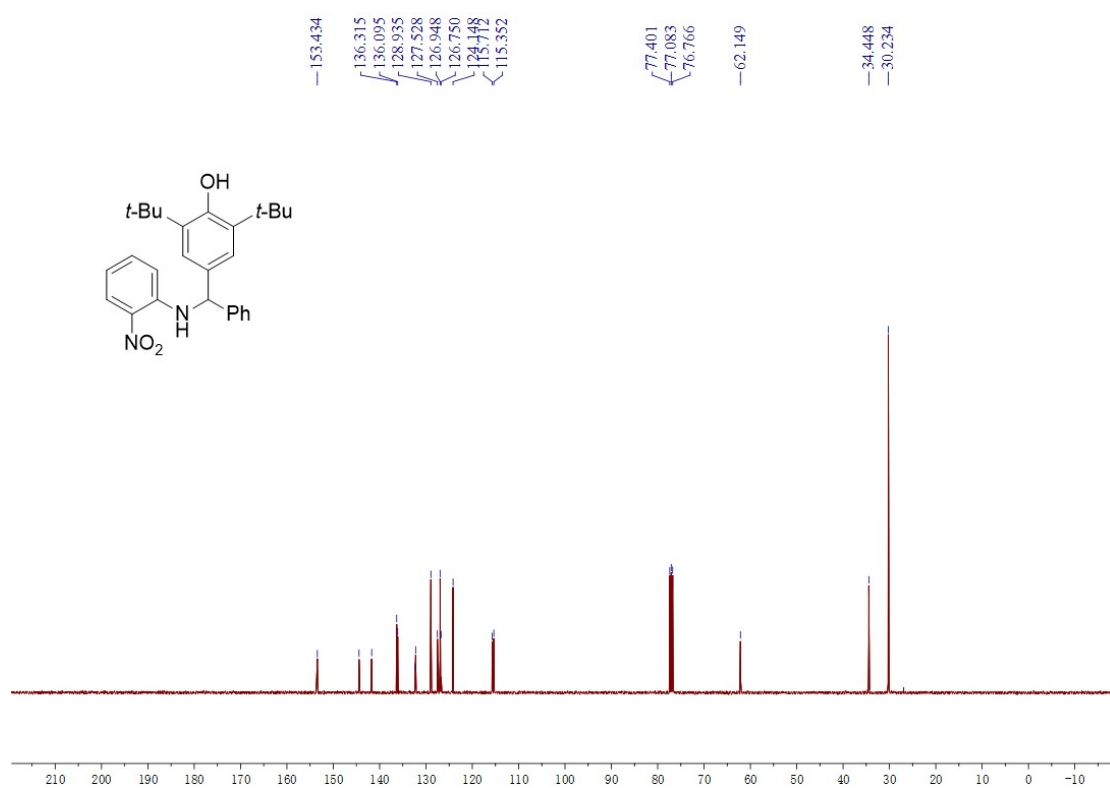
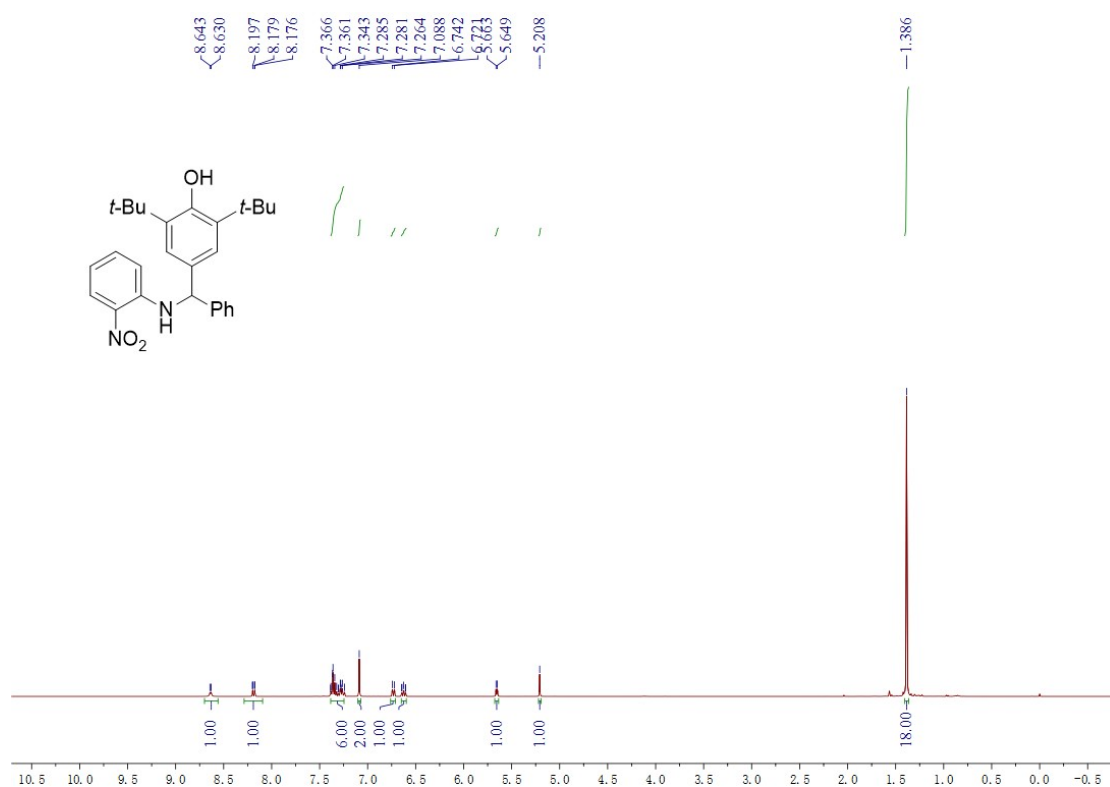


4-((4-Amino-3-(trifluoromethyl)phenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3z)

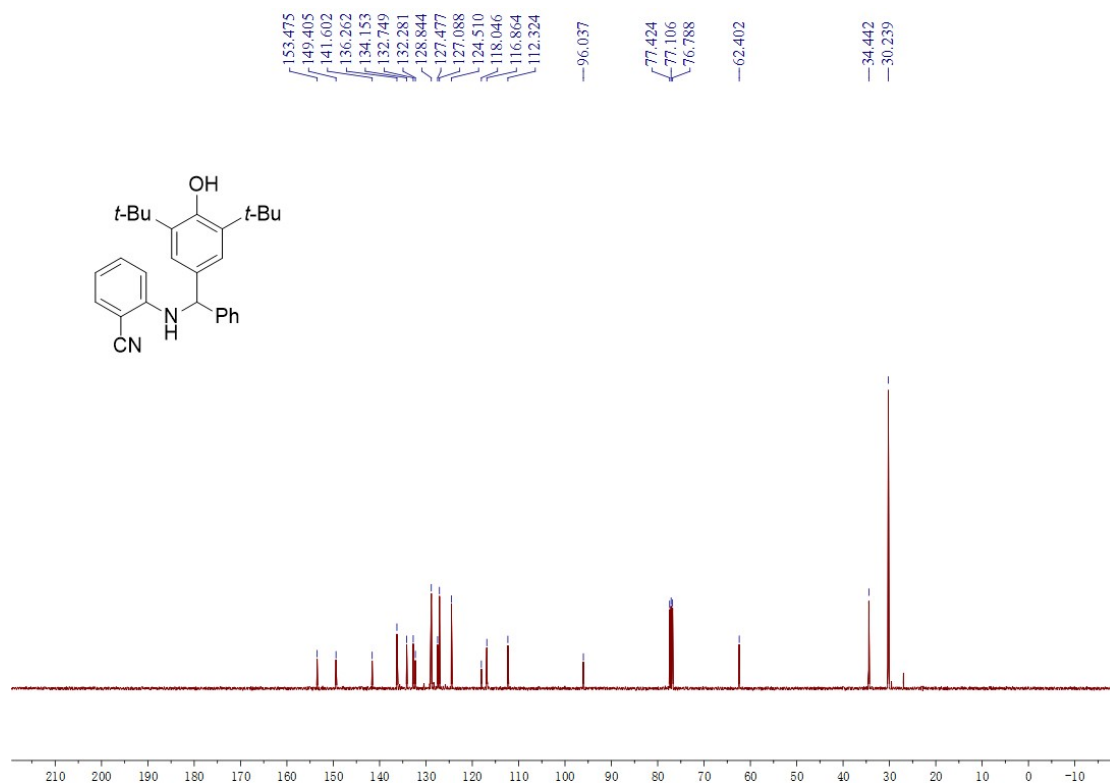
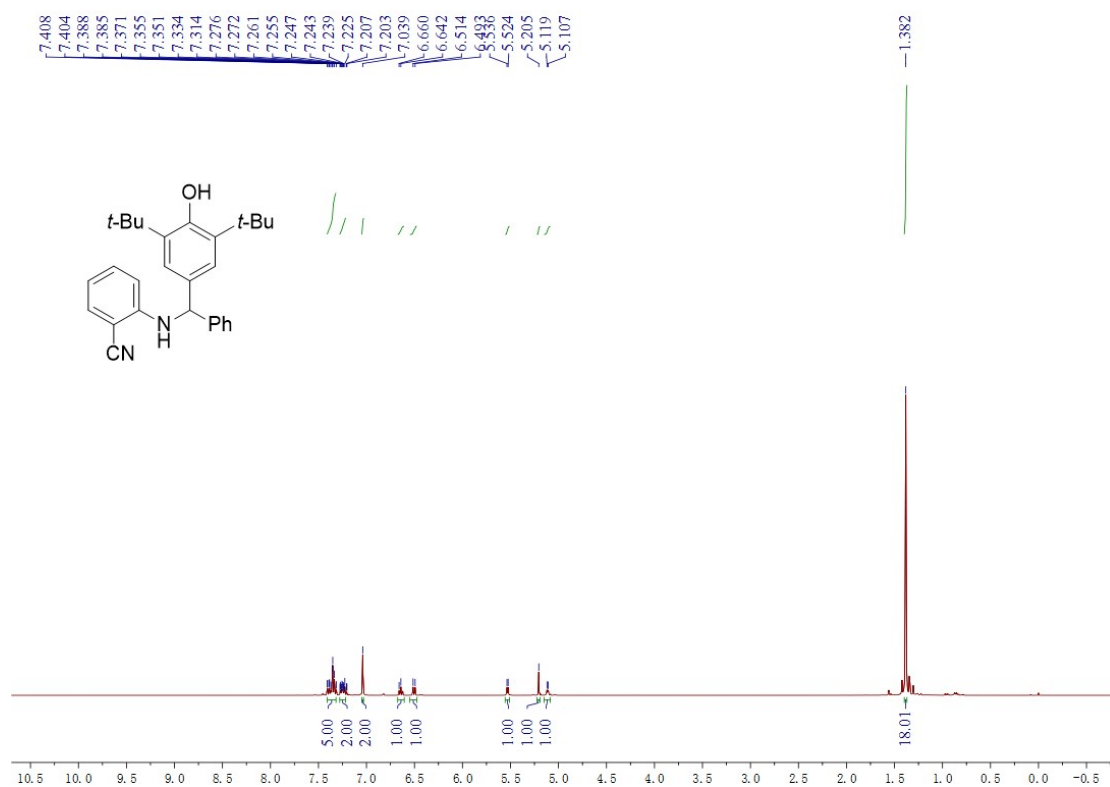




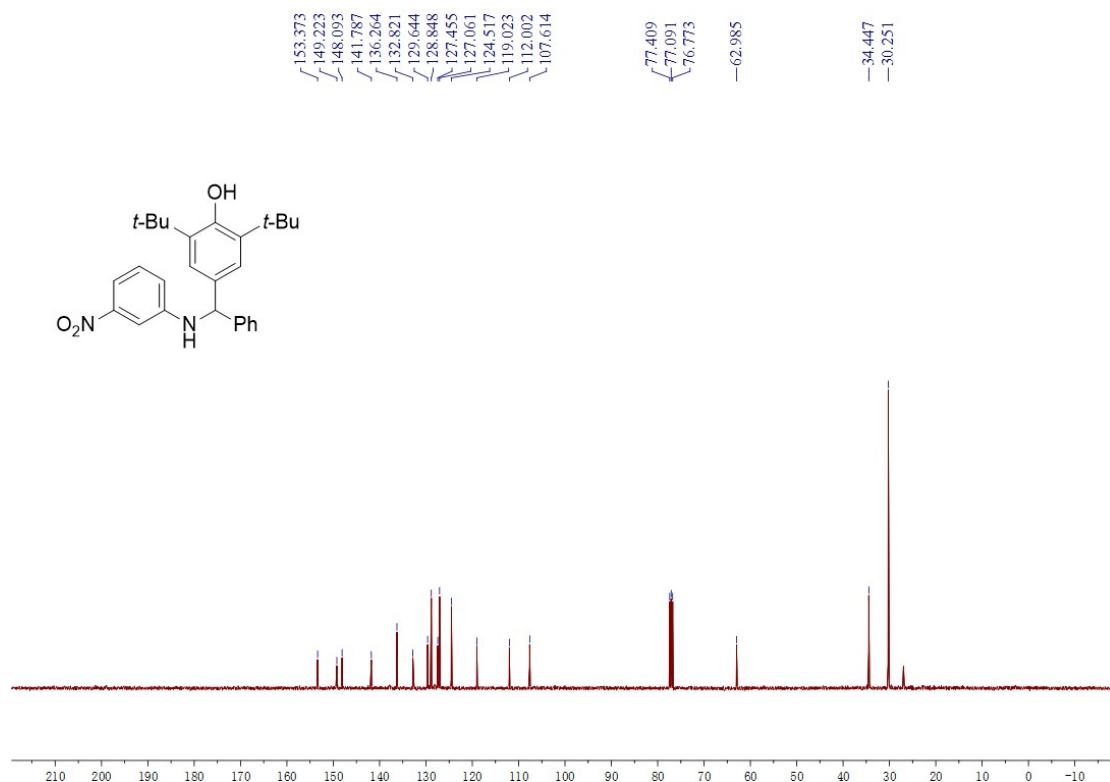
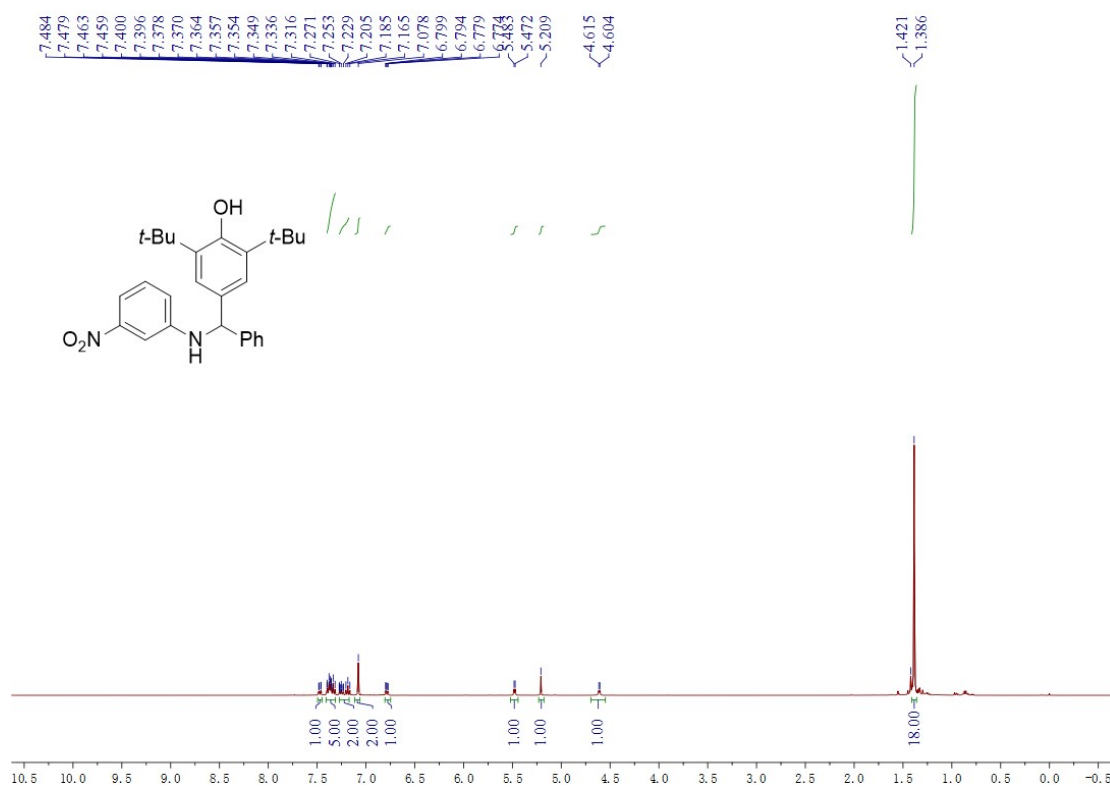
2,6-Di-*tert*-butyl-4-(((2-nitrophenyl)amino)(phenyl)methyl)phenol (3aa')



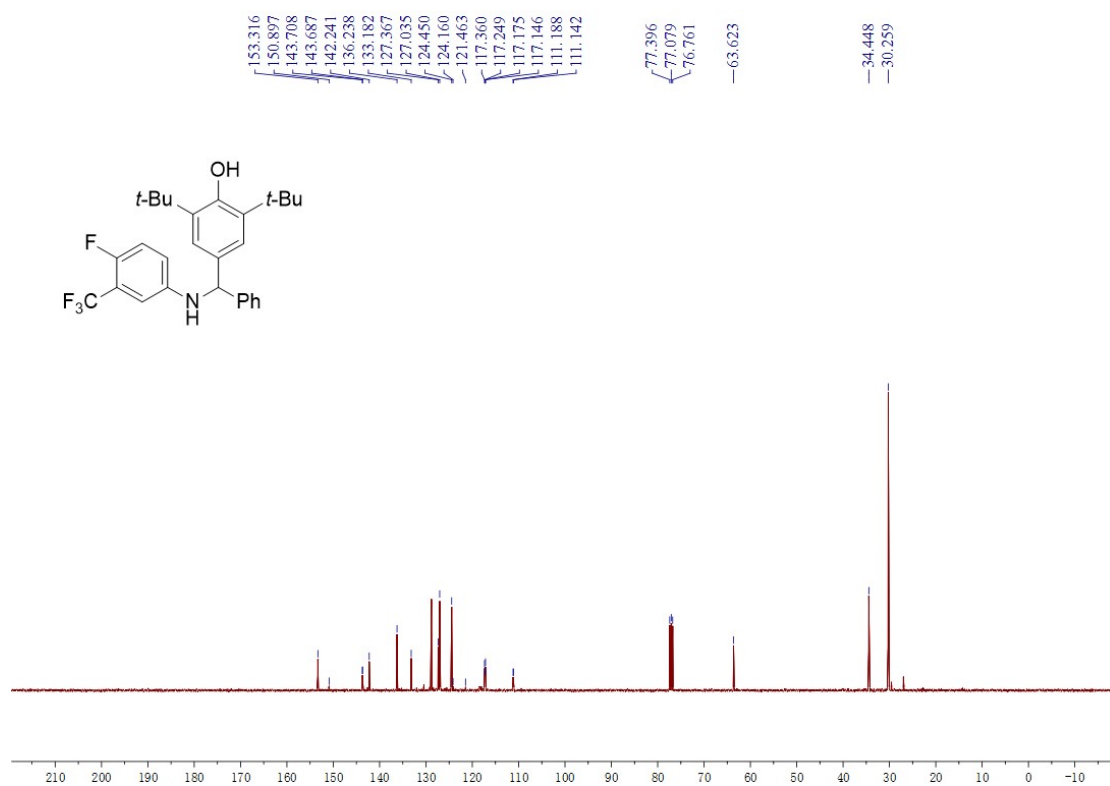
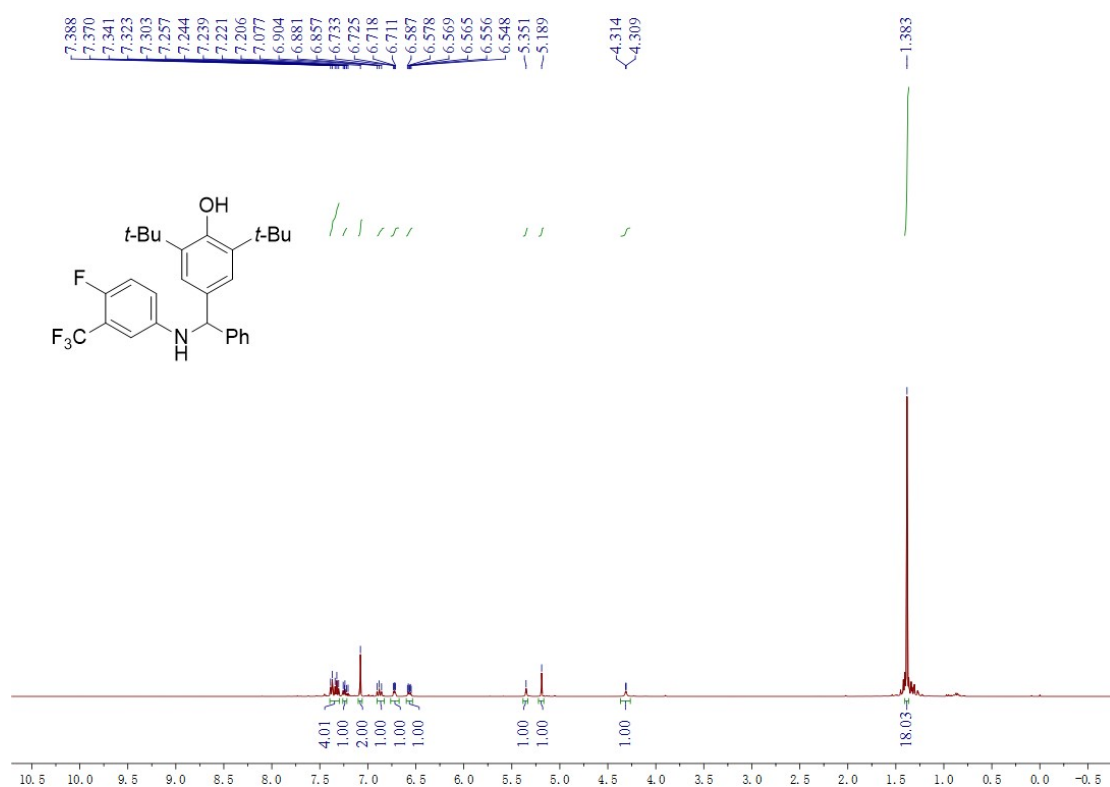
2-(((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)amino)benzonitrile (3ab')

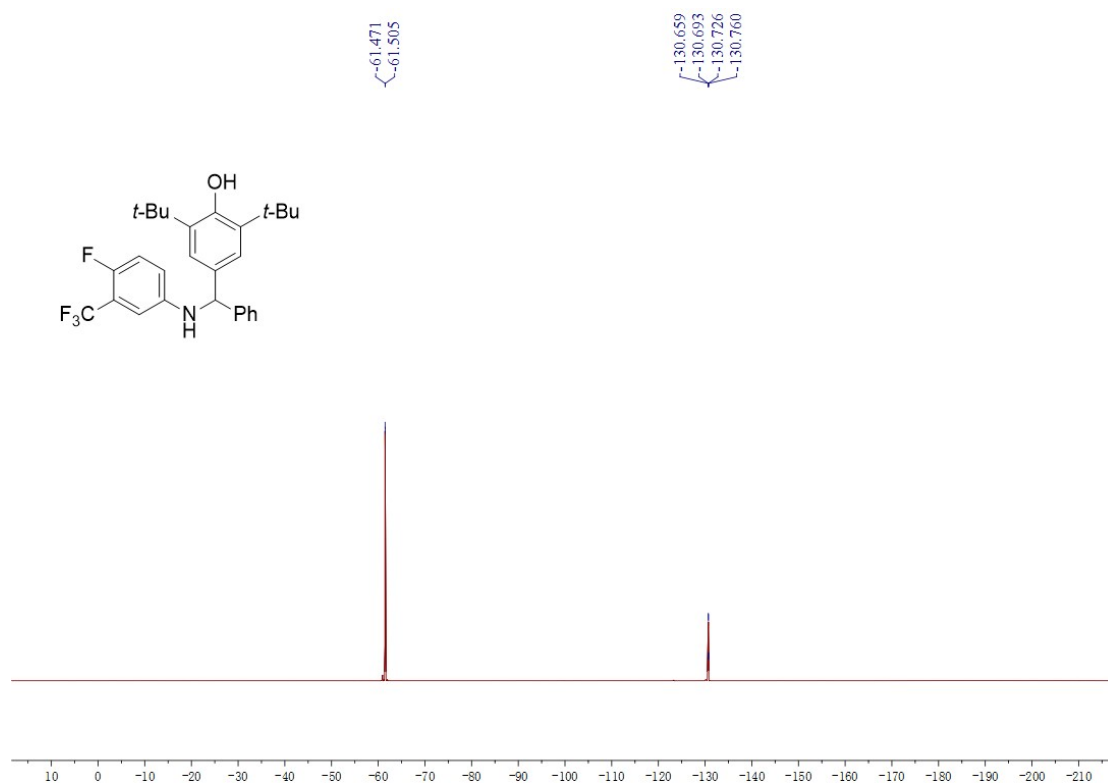


2,6-Di-tert-butyl-4-(((3-nitrophenyl)amino)(phenyl)methyl)phenol (3ac')

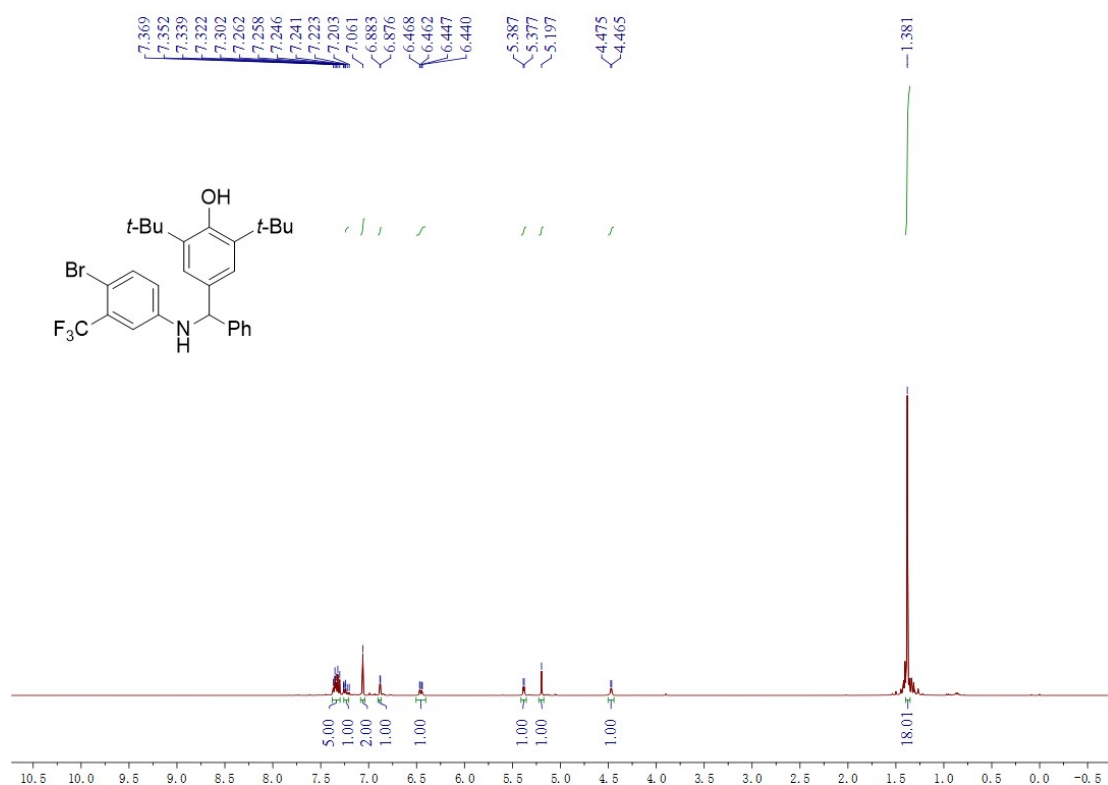


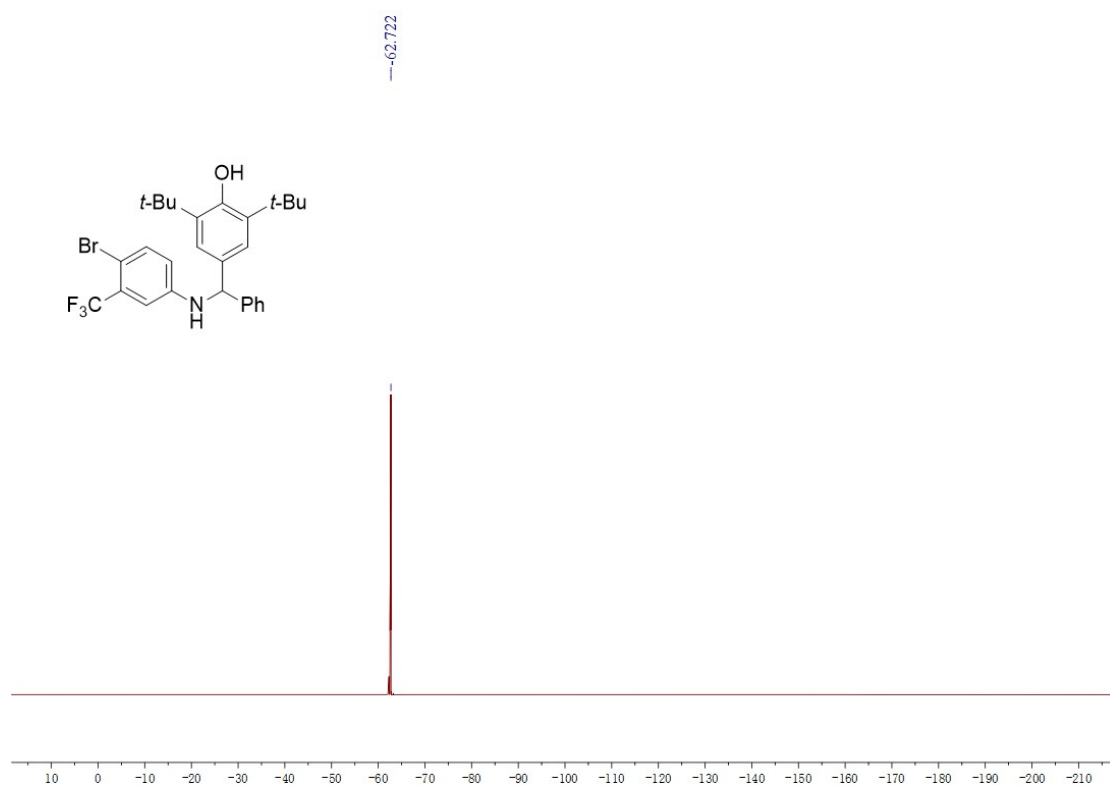
2,6-Di-tert-butyl-4-(((4-fluoro-3-(trifluoromethyl)phenyl)amino)(phenyl)methyl)phenol (3ad')





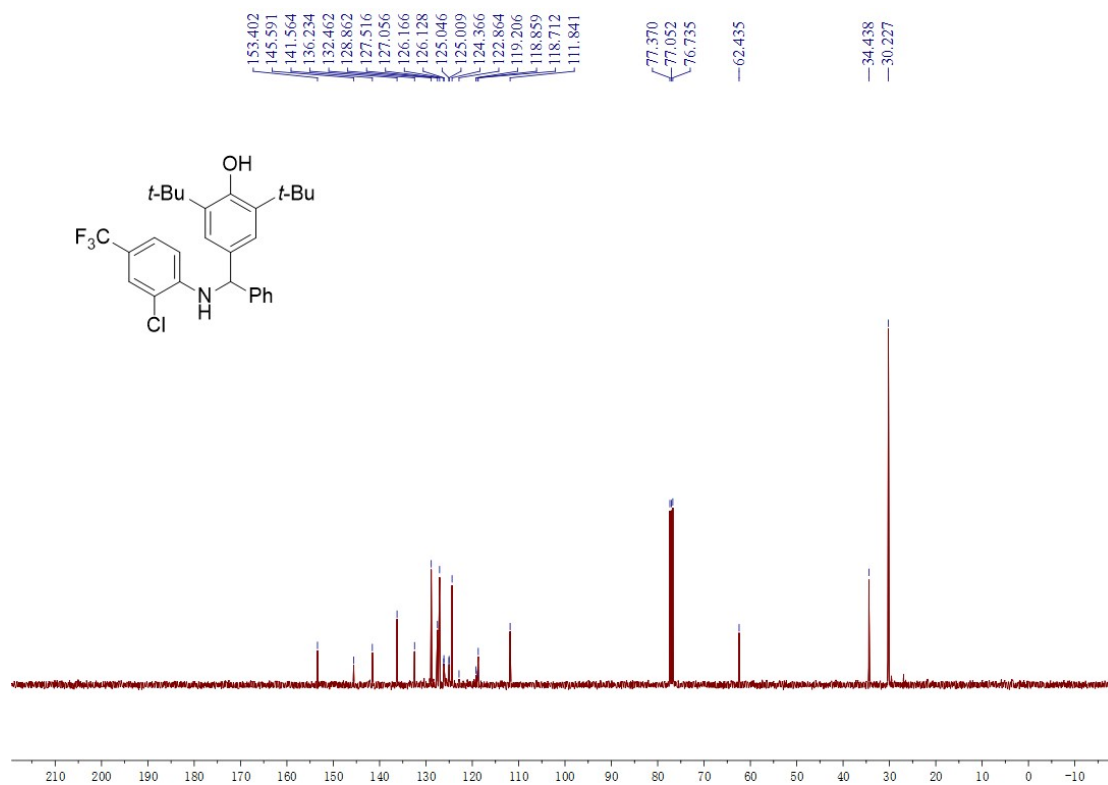
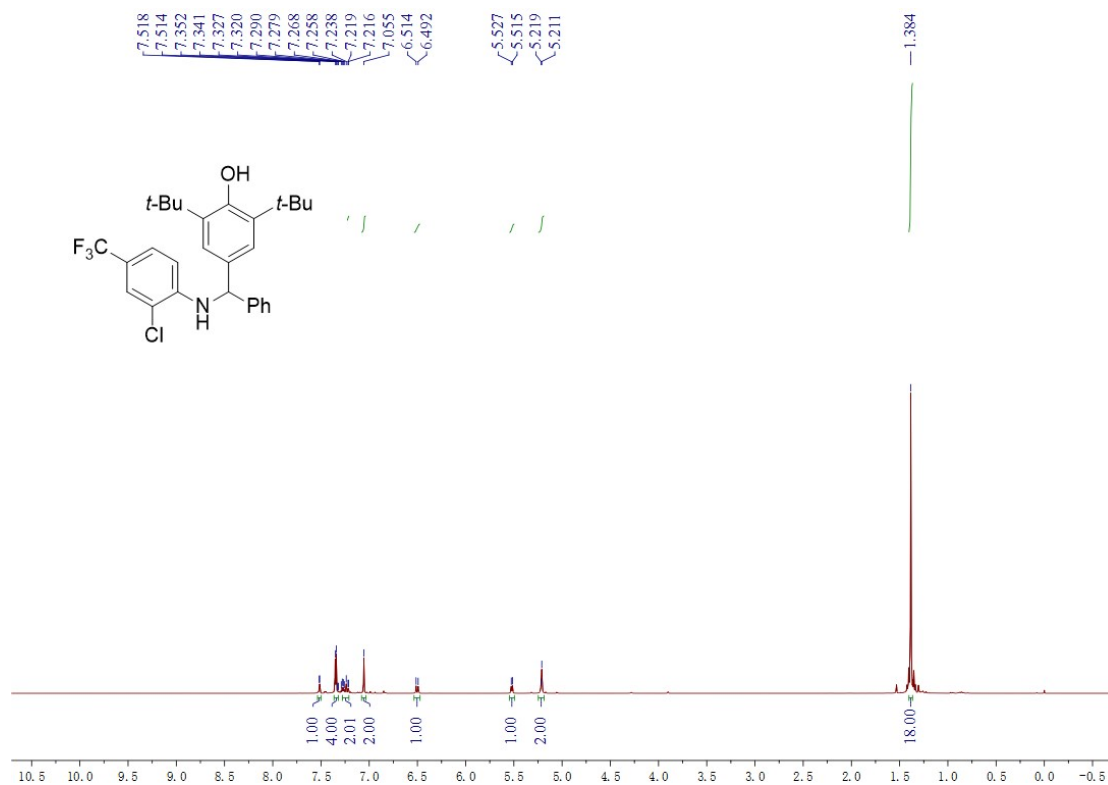
4-(((4-Bromo-3-(trifluoromethyl)phenyl)amino)(phenyl)methyl)-2,6-di-tert-butylphenol (3ae')

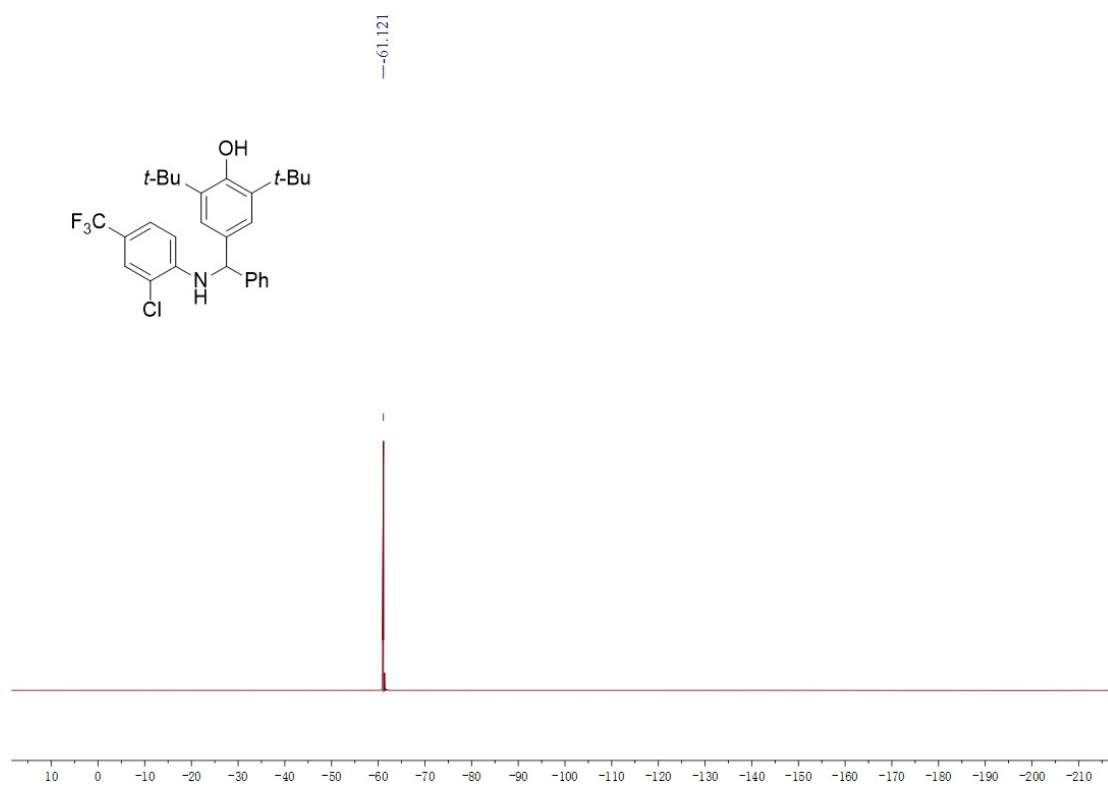




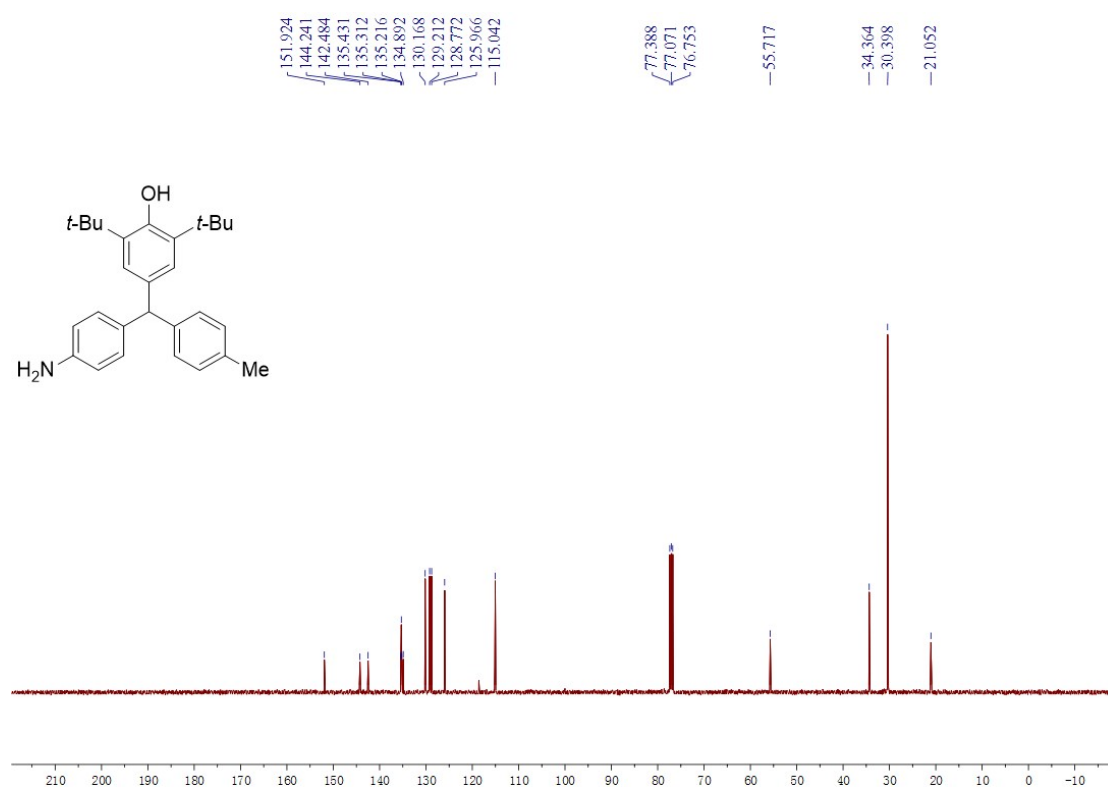
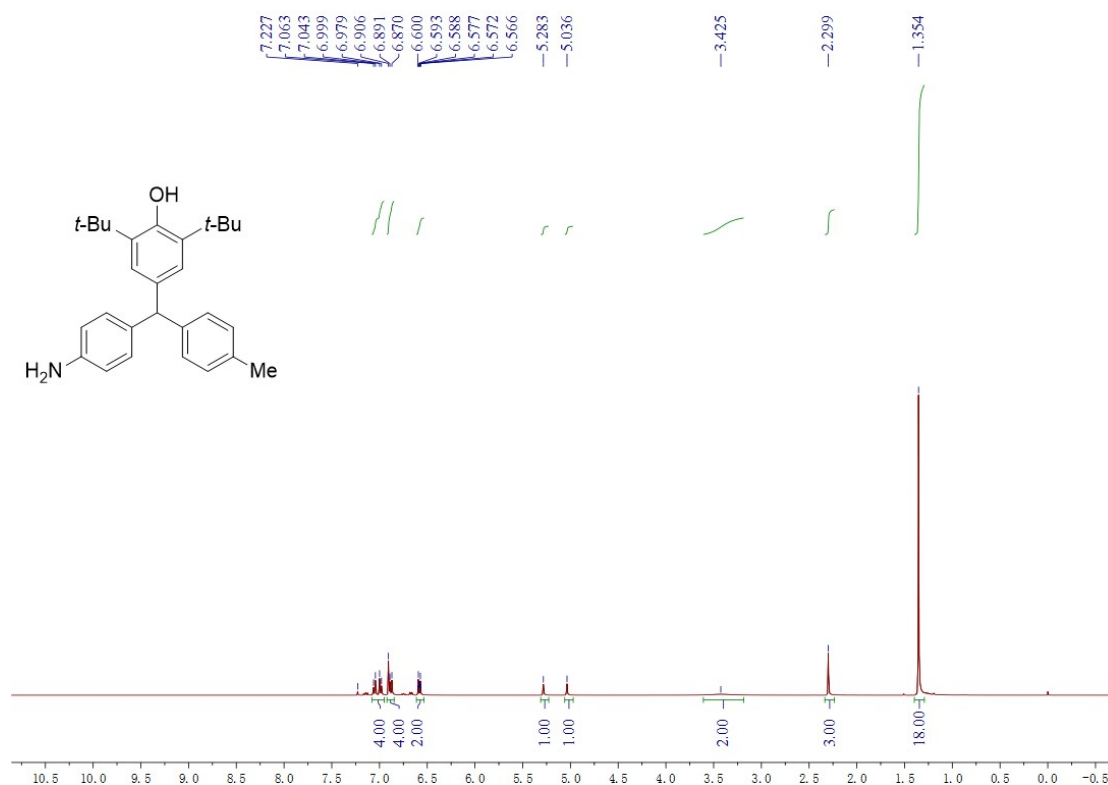
2,6-Di-tert-butyl-4-(((2-chloro-4-(trifluoromethyl)phenyl)amino)(phenyl)methyl)ph-

enol (3af')

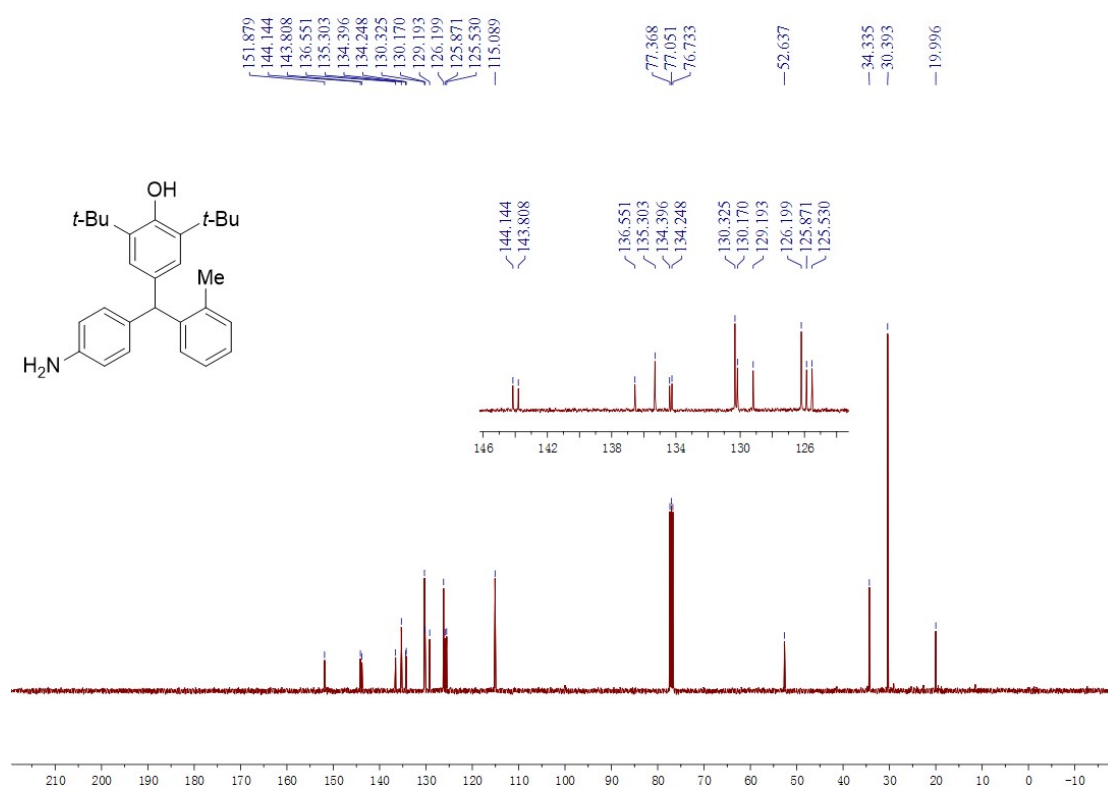
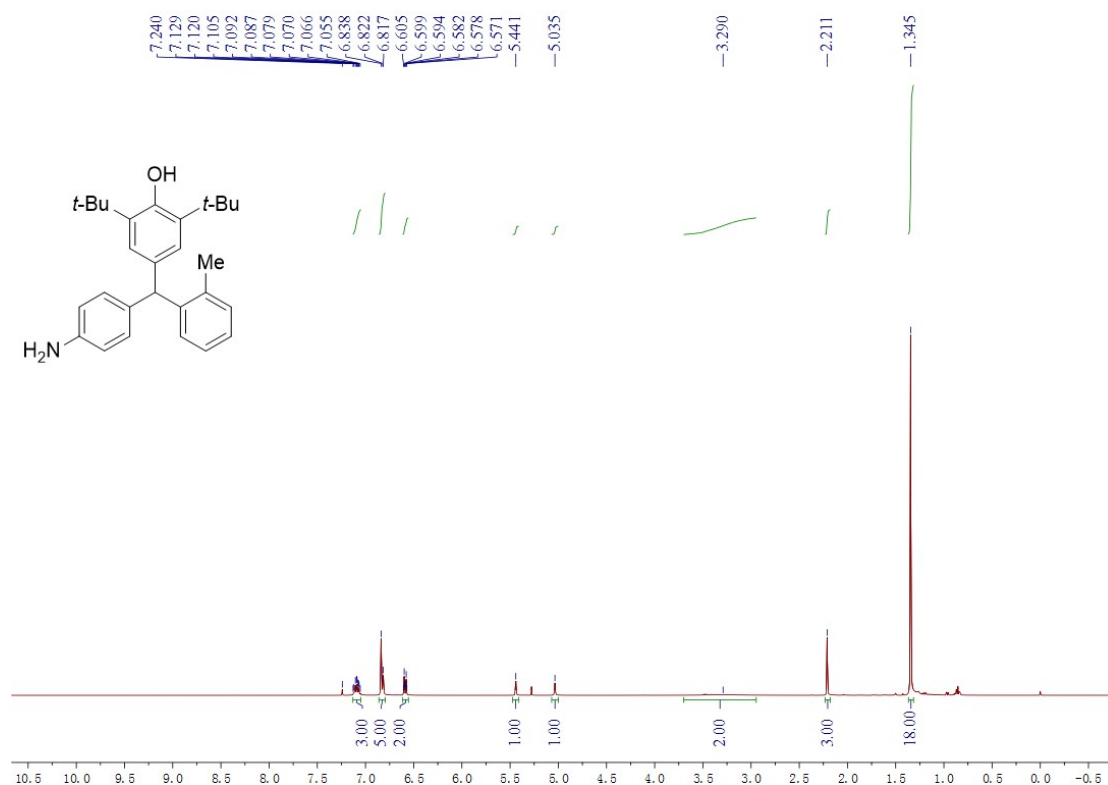




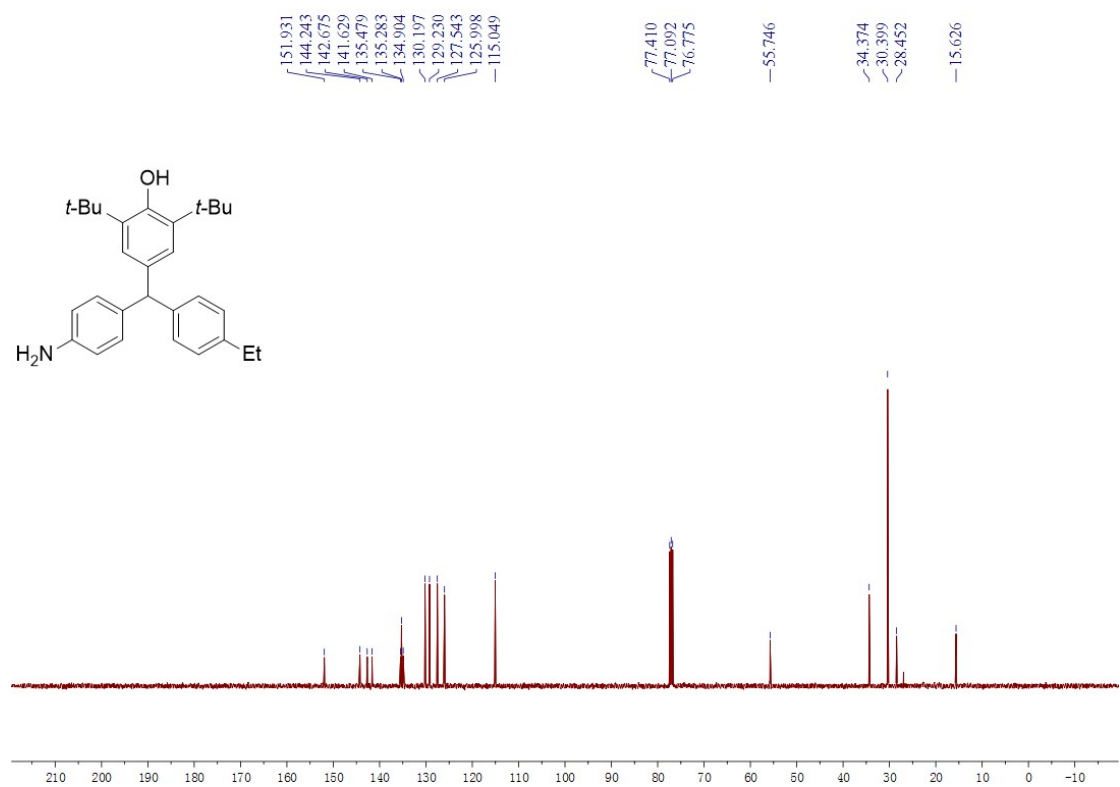
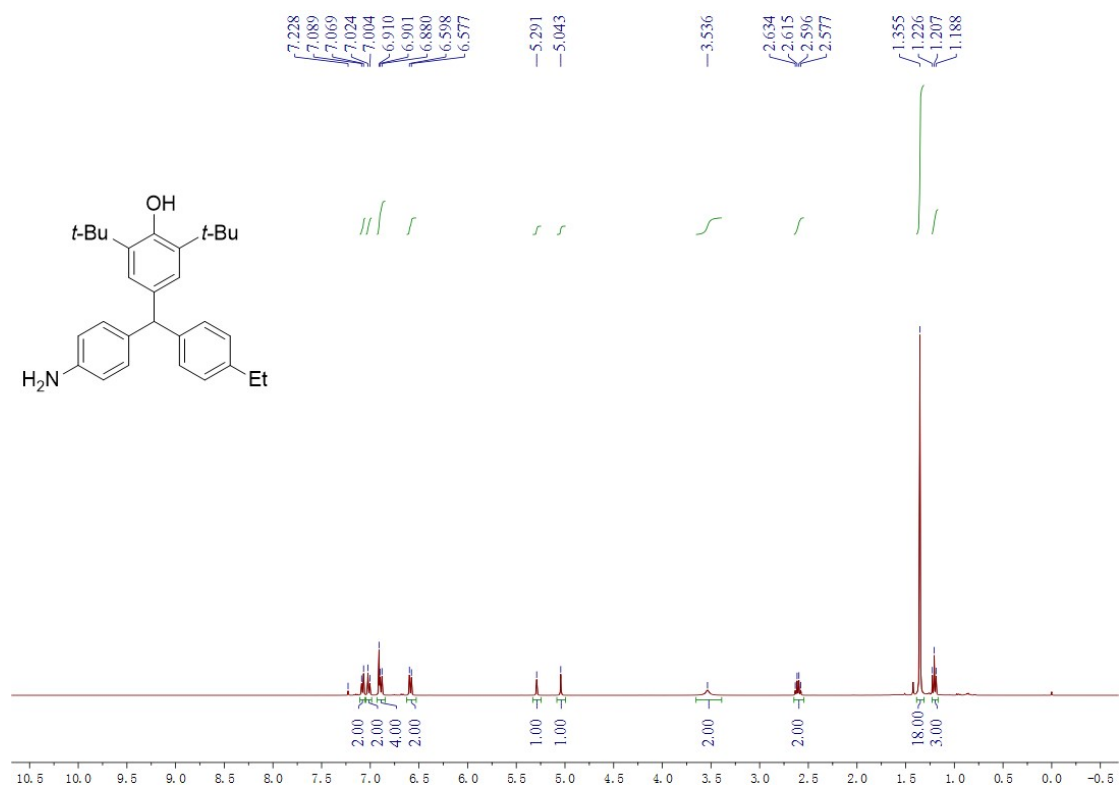
4-((4-Aminophenyl)(p-tolyl)methyl)-2,6-di-tert-butylphenol (4a)



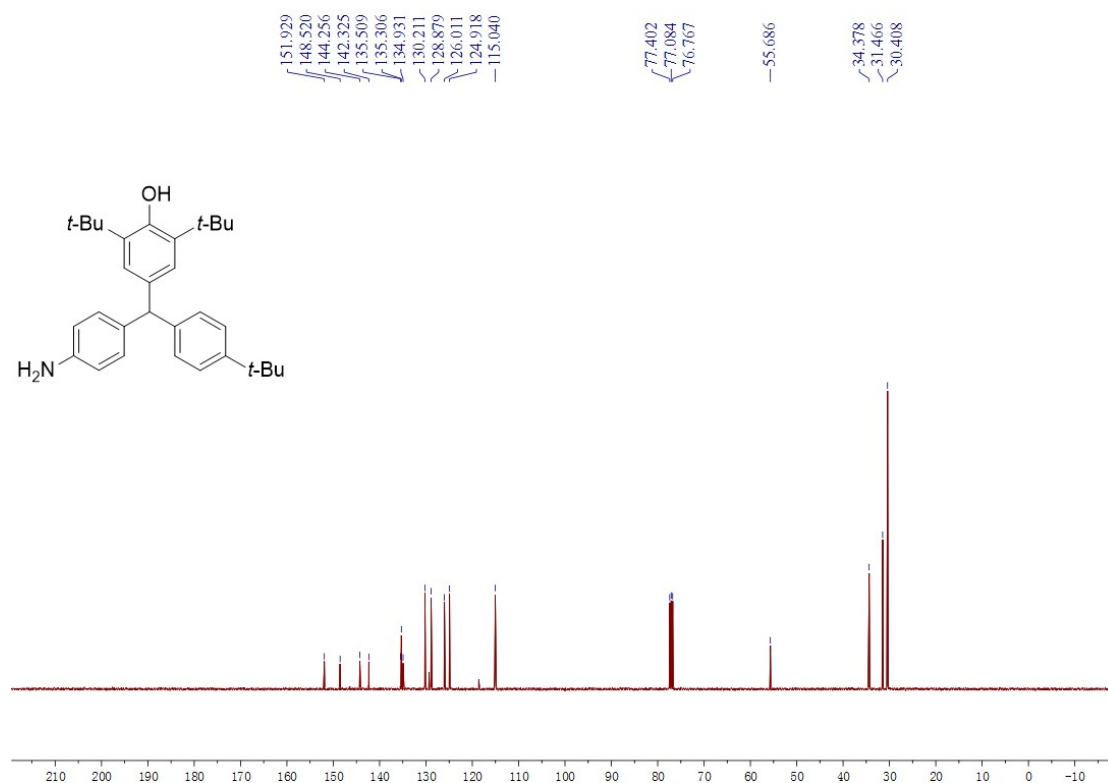
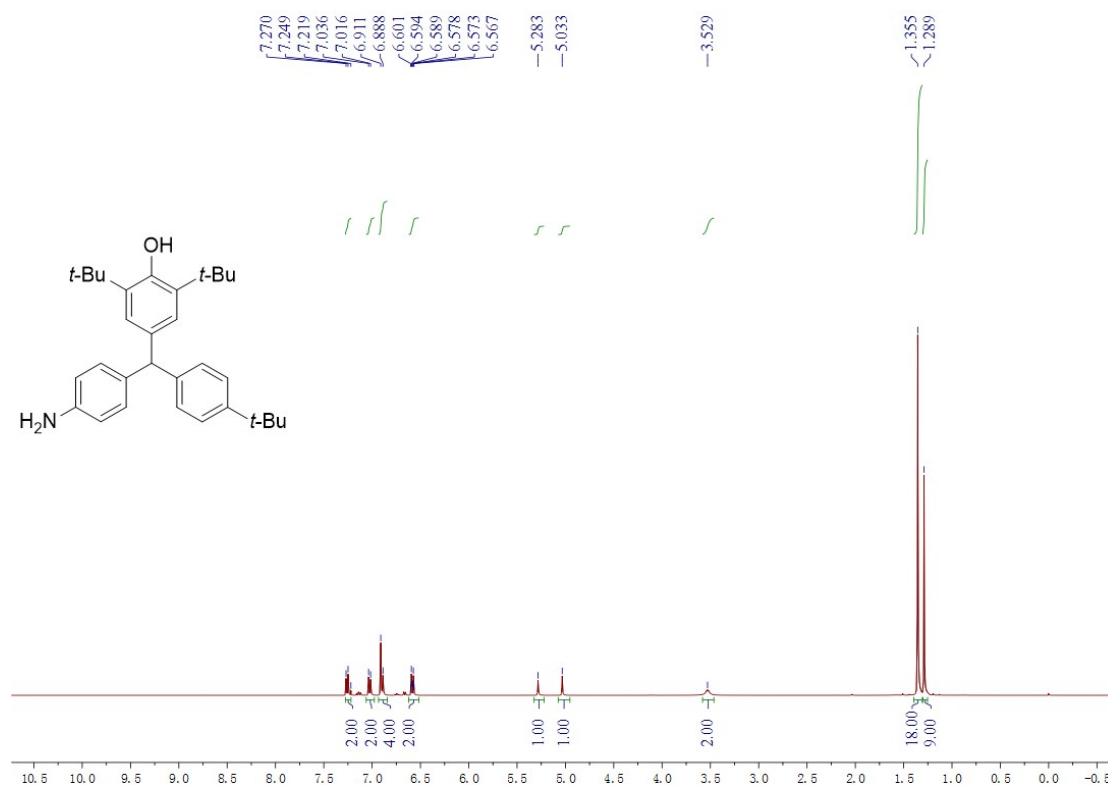
4-((4-Aminophenyl)(o-tolyl)methyl)-2,6-di-tert-butylphenol (4b)



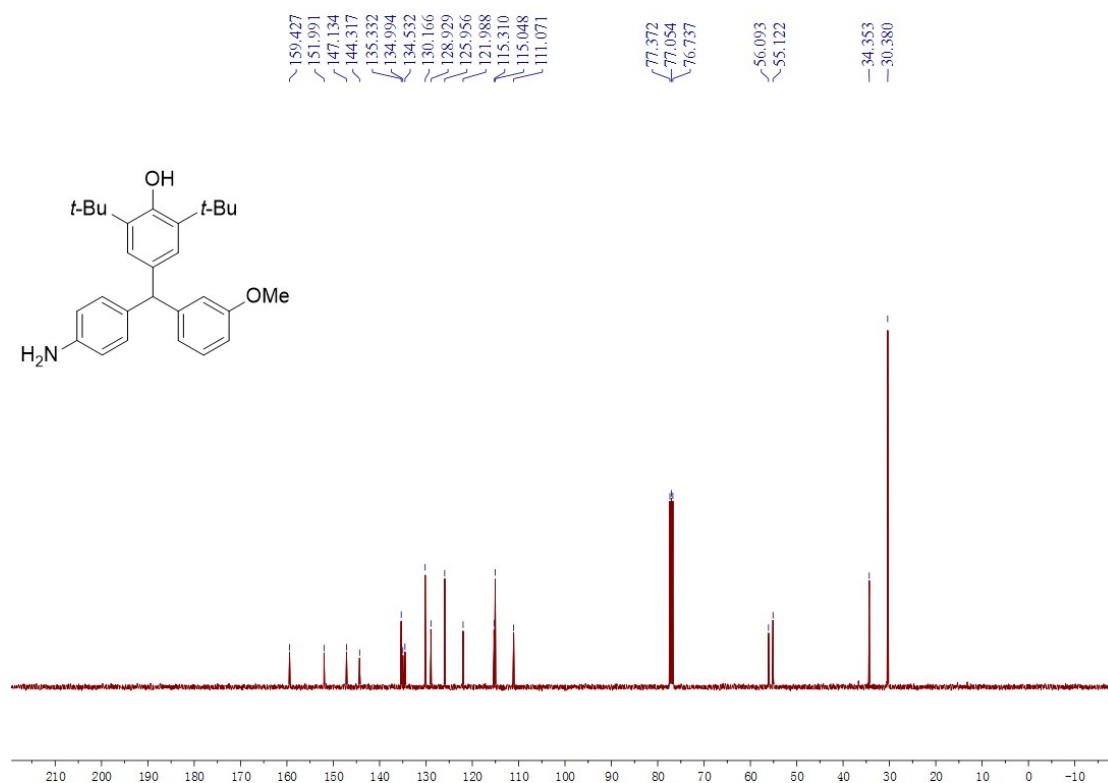
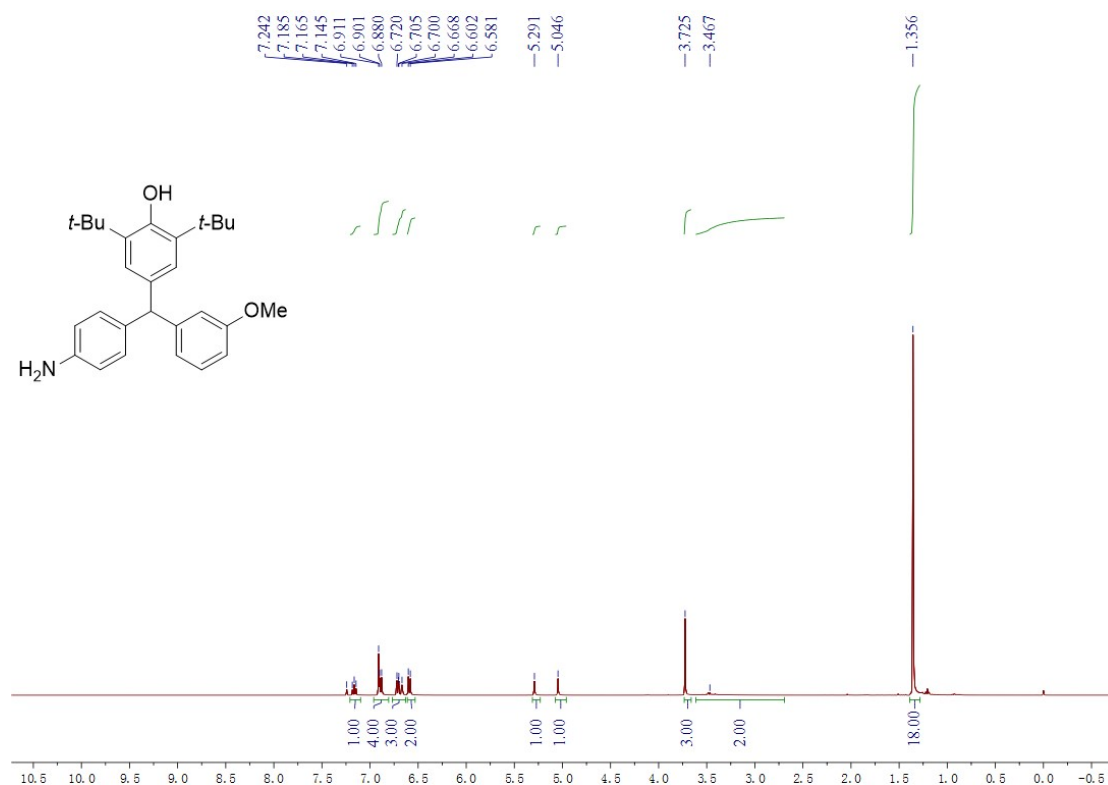
4-((4-Aminophenyl)(4-ethylphenyl)methyl)-2,6-di-tert-butylphenol (4c)



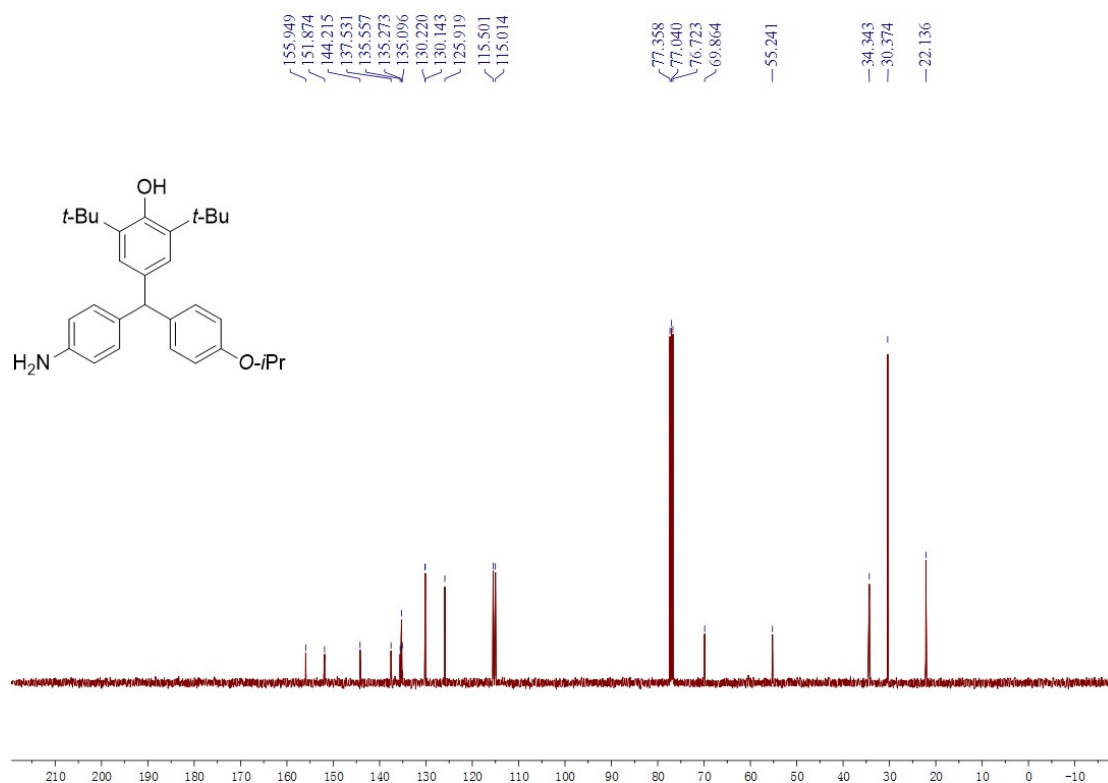
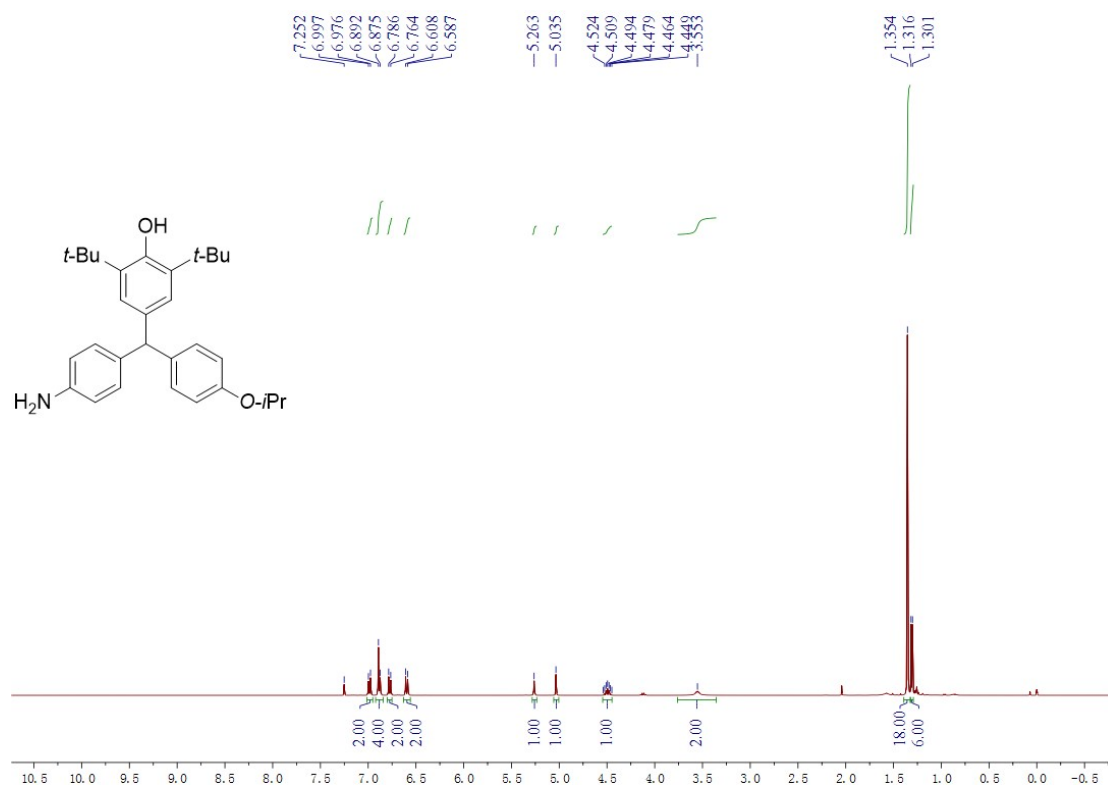
4-((4-Aminophenyl)(4-(tert-butyl)phenyl)methyl)-2,6-di-tert-butylphenol (4d)



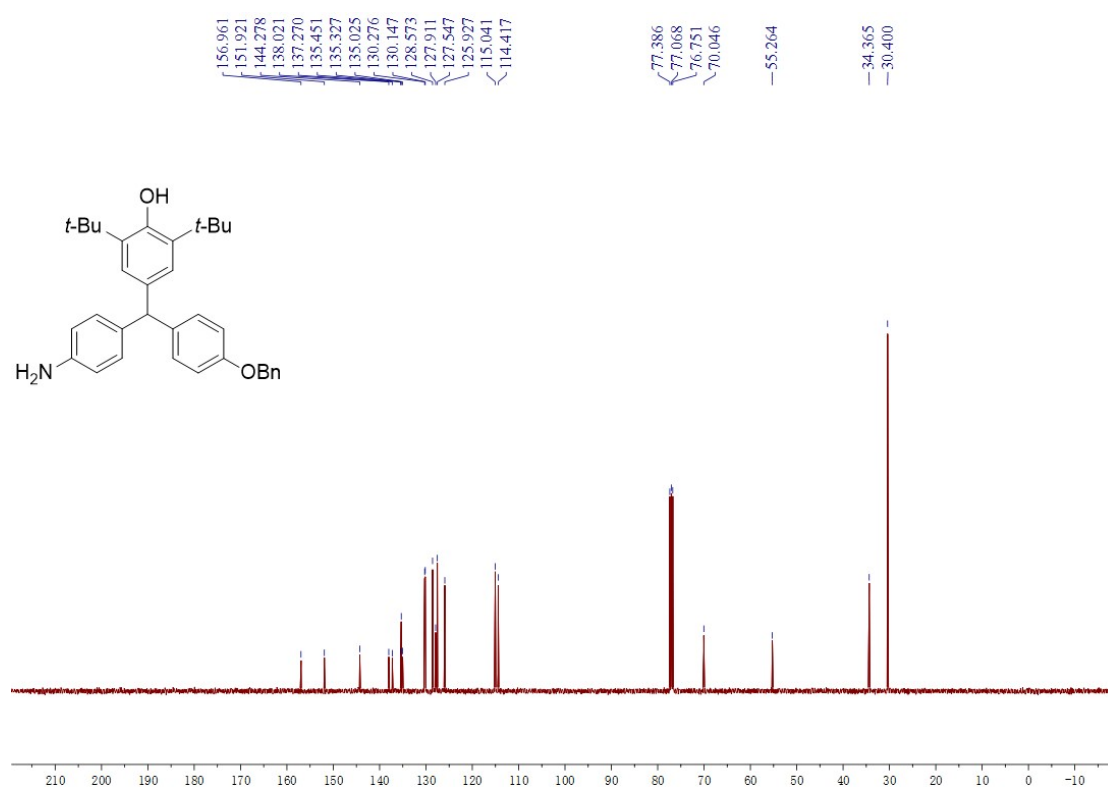
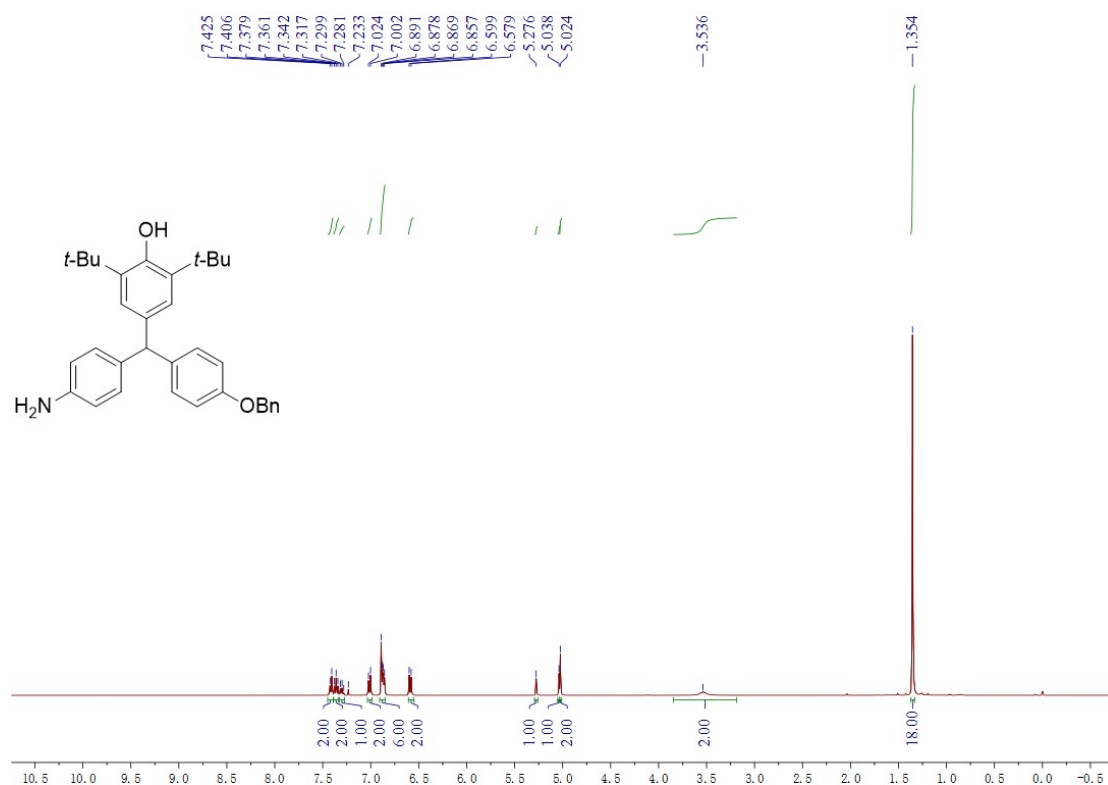
4-((4-Aminophenyl)(3-methoxyphenyl)methyl)-2,6-di-tert-butylphenol (4e)



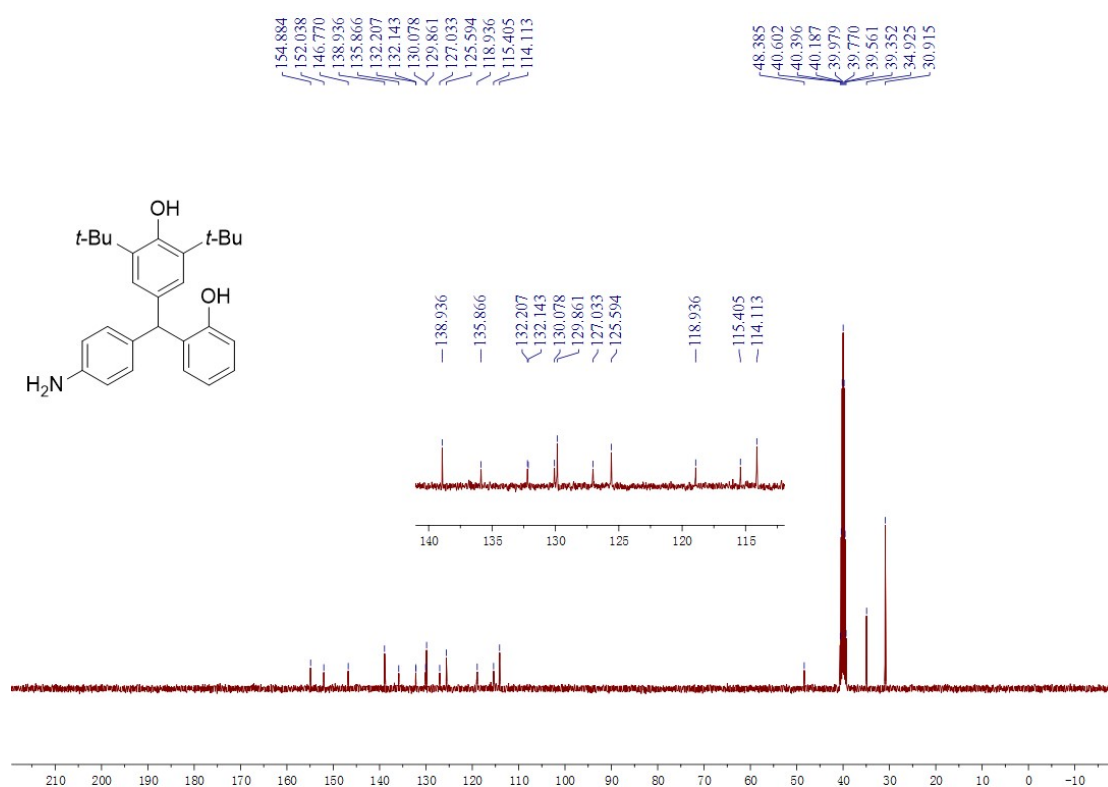
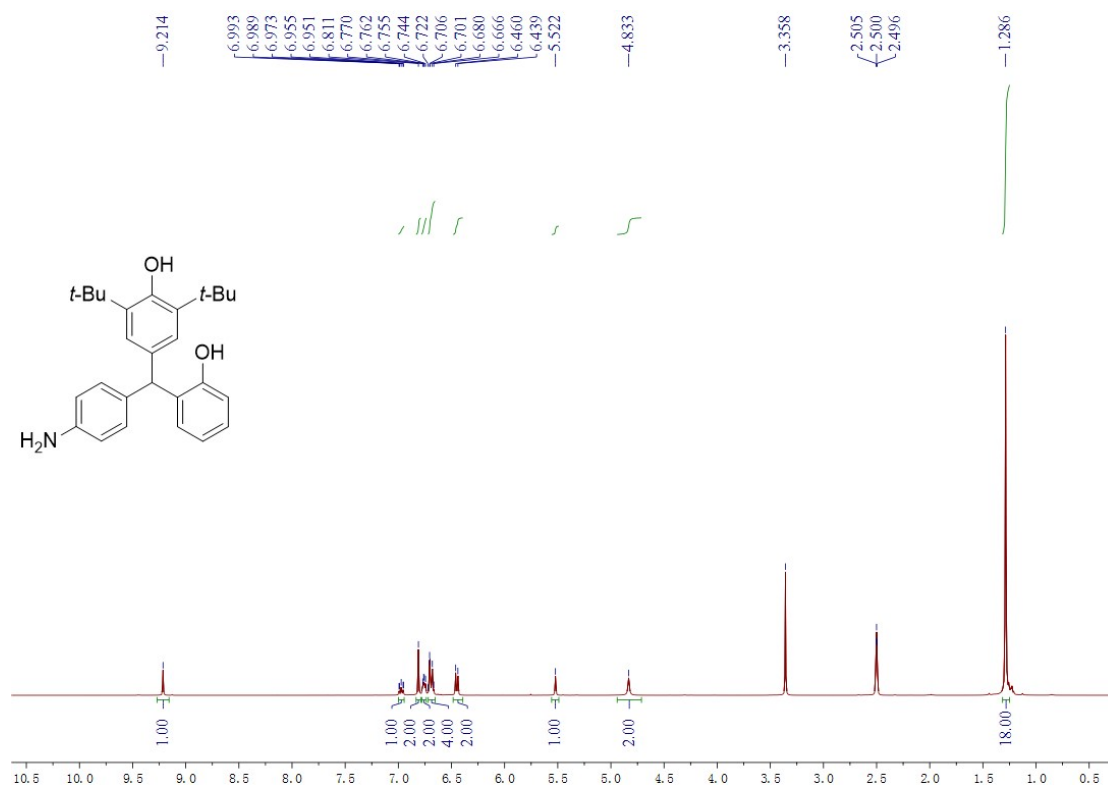
4-((4-Aminophenyl)(4-isopropoxyphenyl)methyl)-2,6-di-tert-butylphenol (4f)



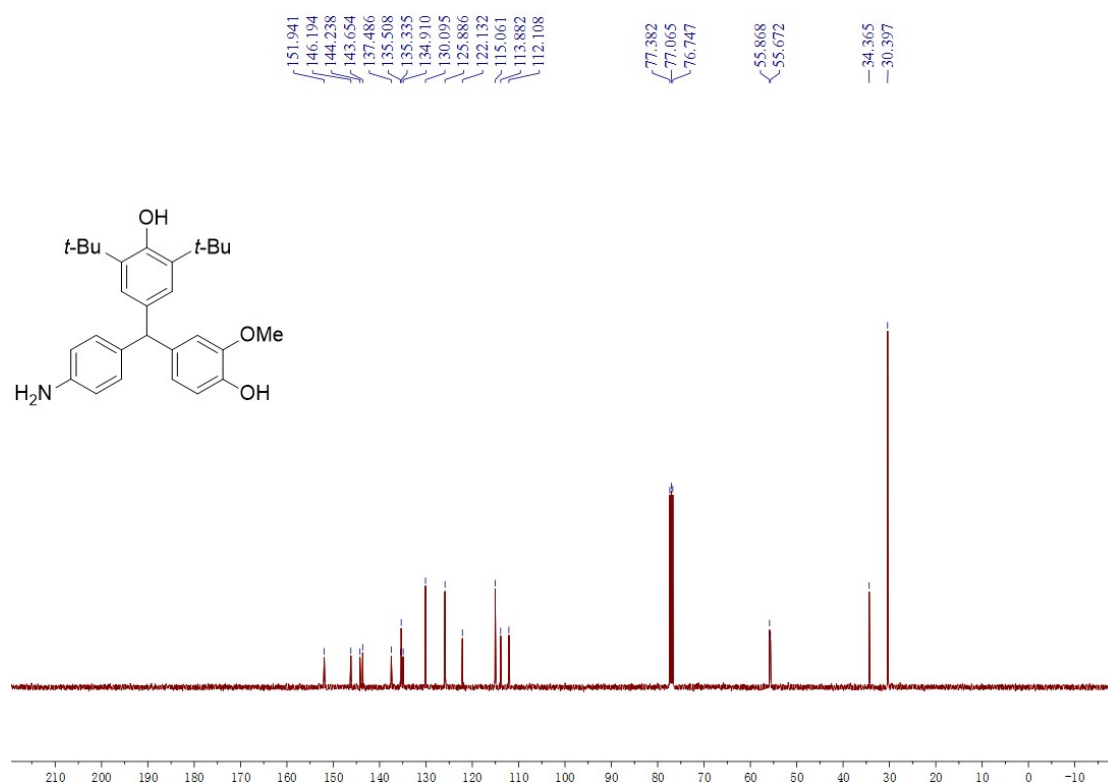
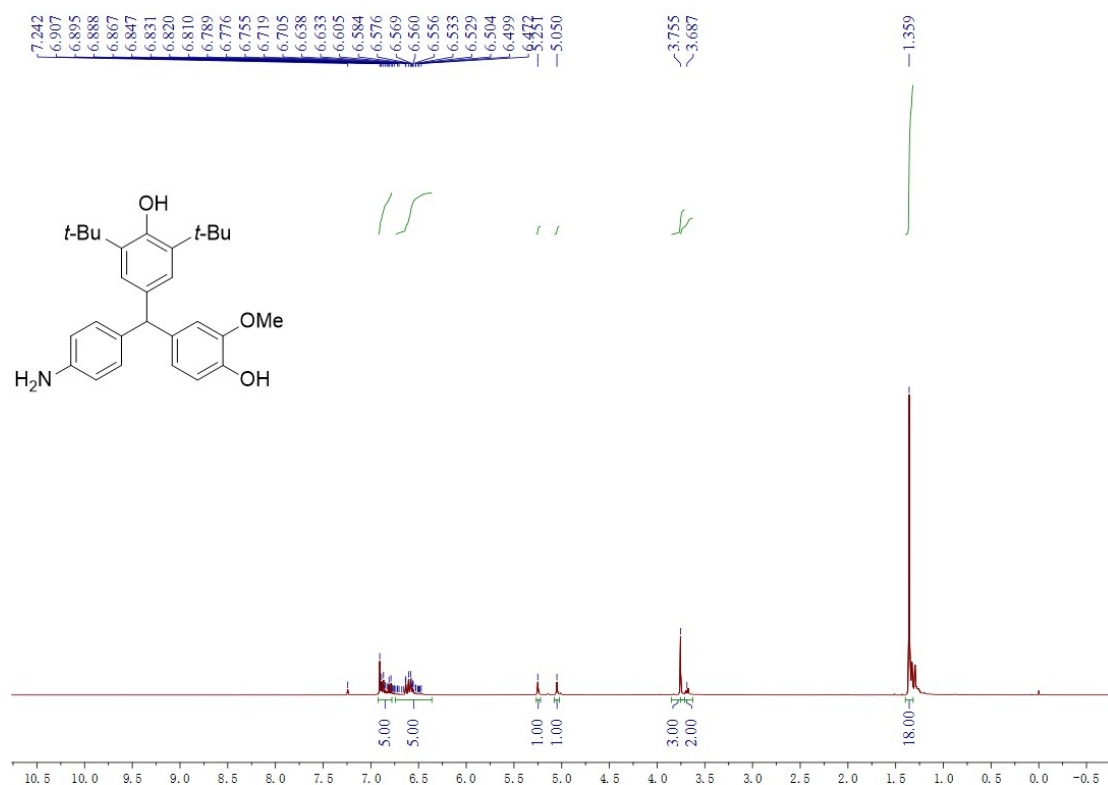
4-((4-Aminophenyl)(4-(benzyloxy)phenyl)methyl)-2,6-di-tert-butylphenol (4g)



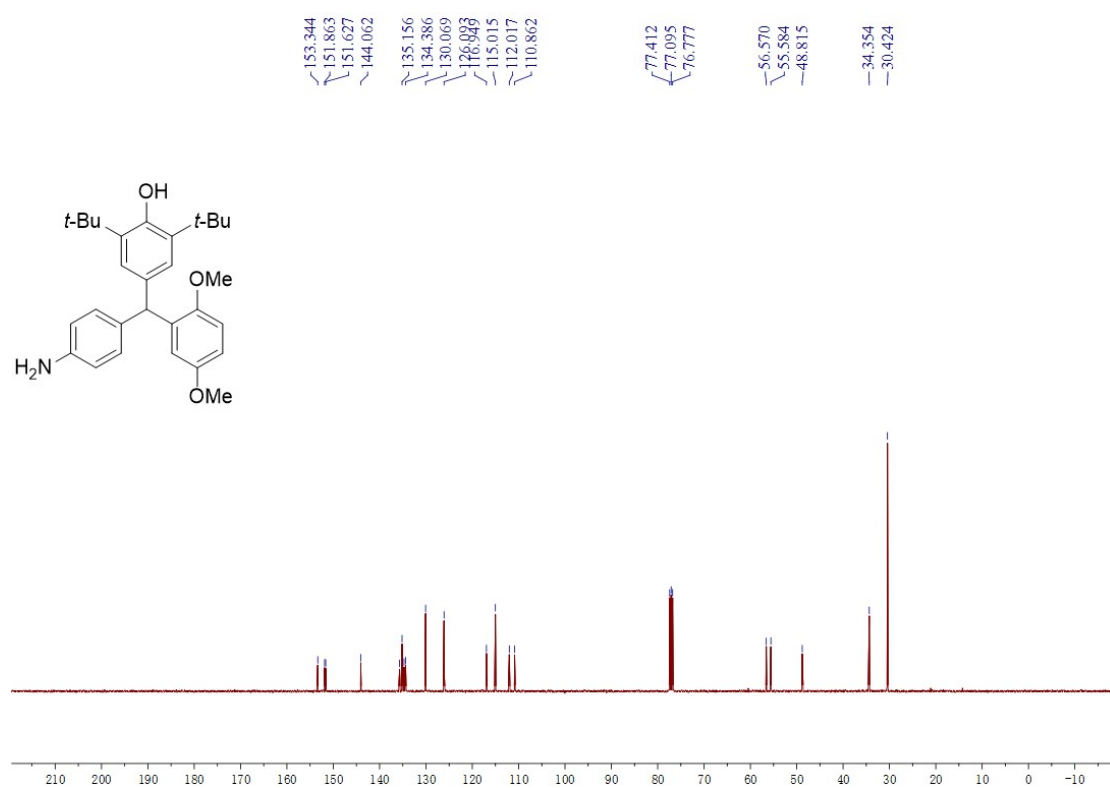
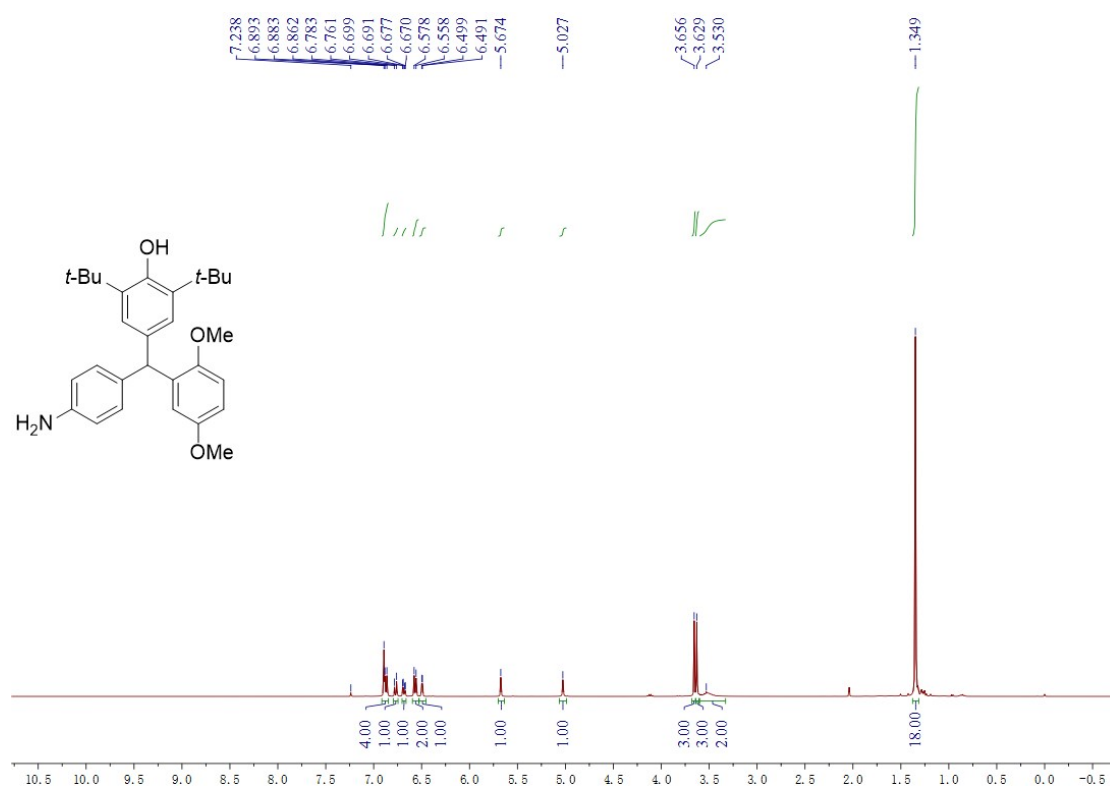
4-((4-Aminophenyl)(2-hydroxyphenyl)methyl)-2,6-di-tert-butylphenol (4h)



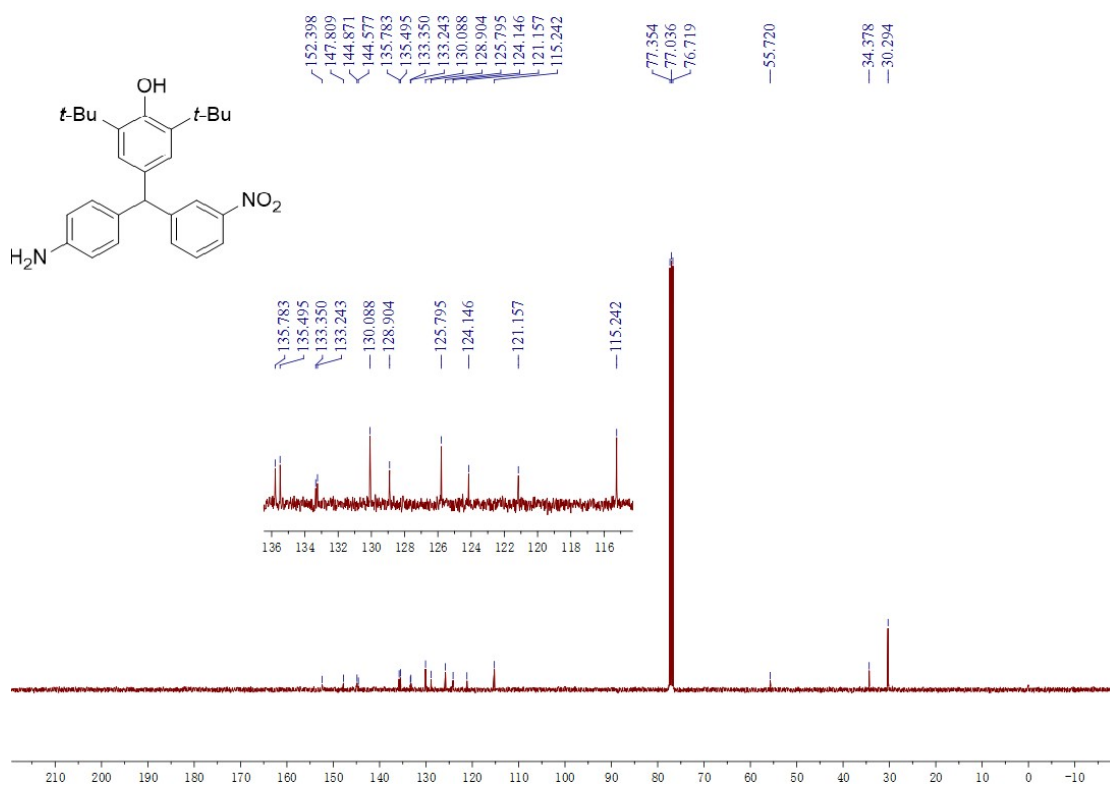
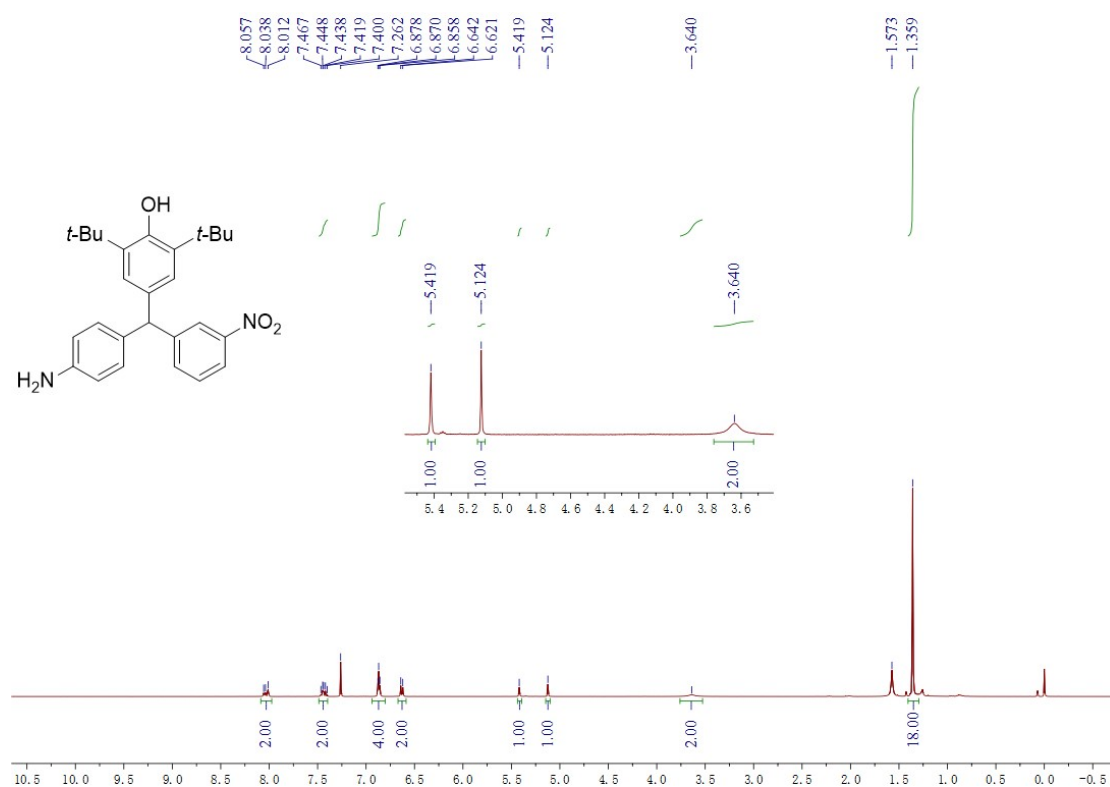
4-((4-Aminophenyl)(4-hydroxy-3-methoxyphenyl)methyl)-2,6-di-tert-butylphenol (4i)



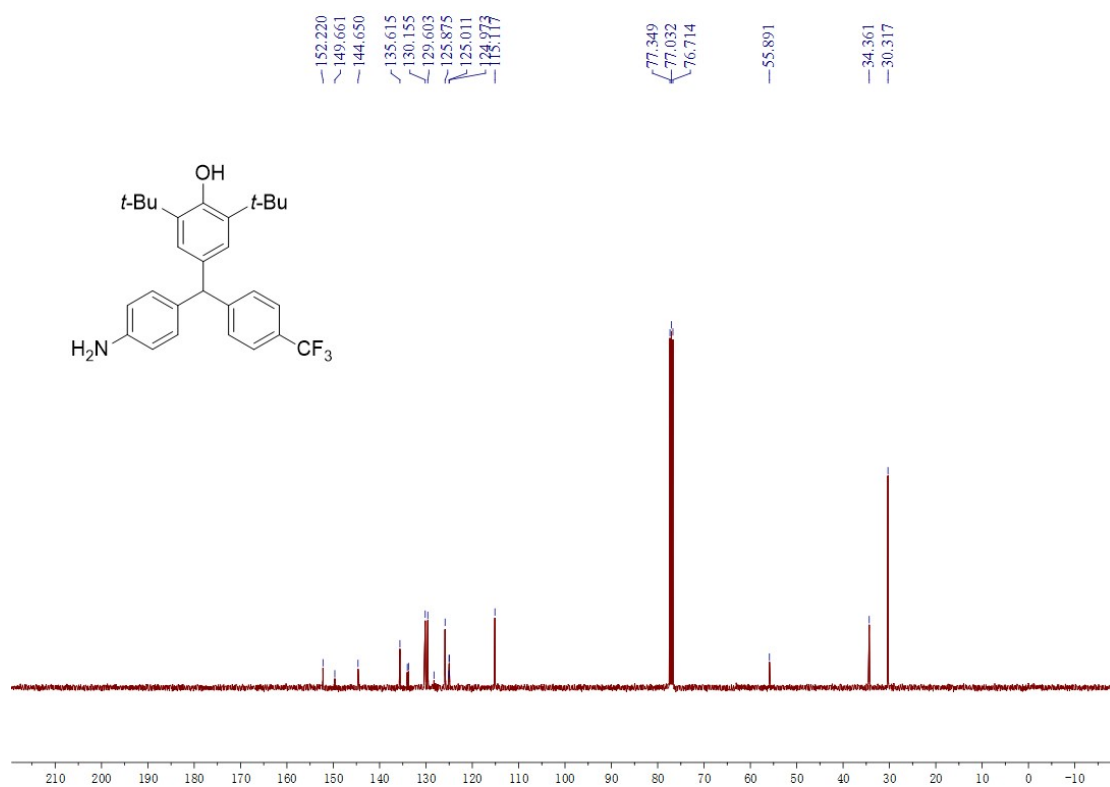
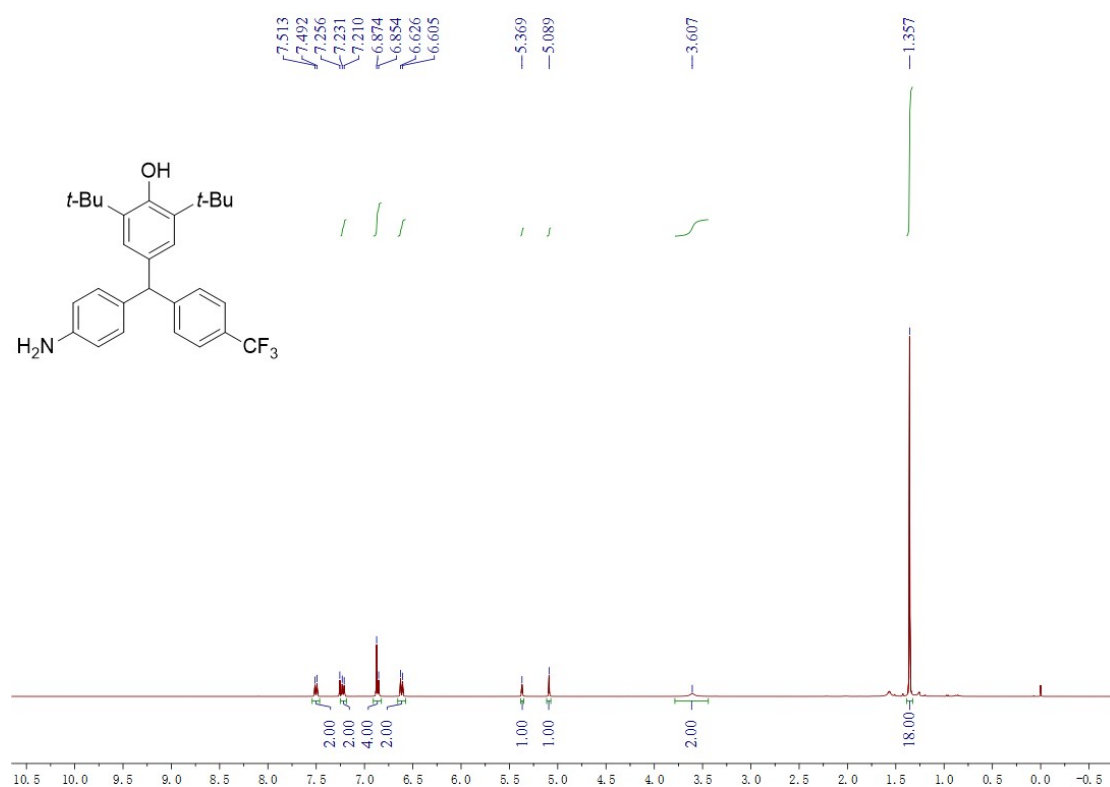
4-((4-Aminophenyl)(2,5-dimethoxyphenyl)methyl)-2,6-di-tert-butylphenol (4j)

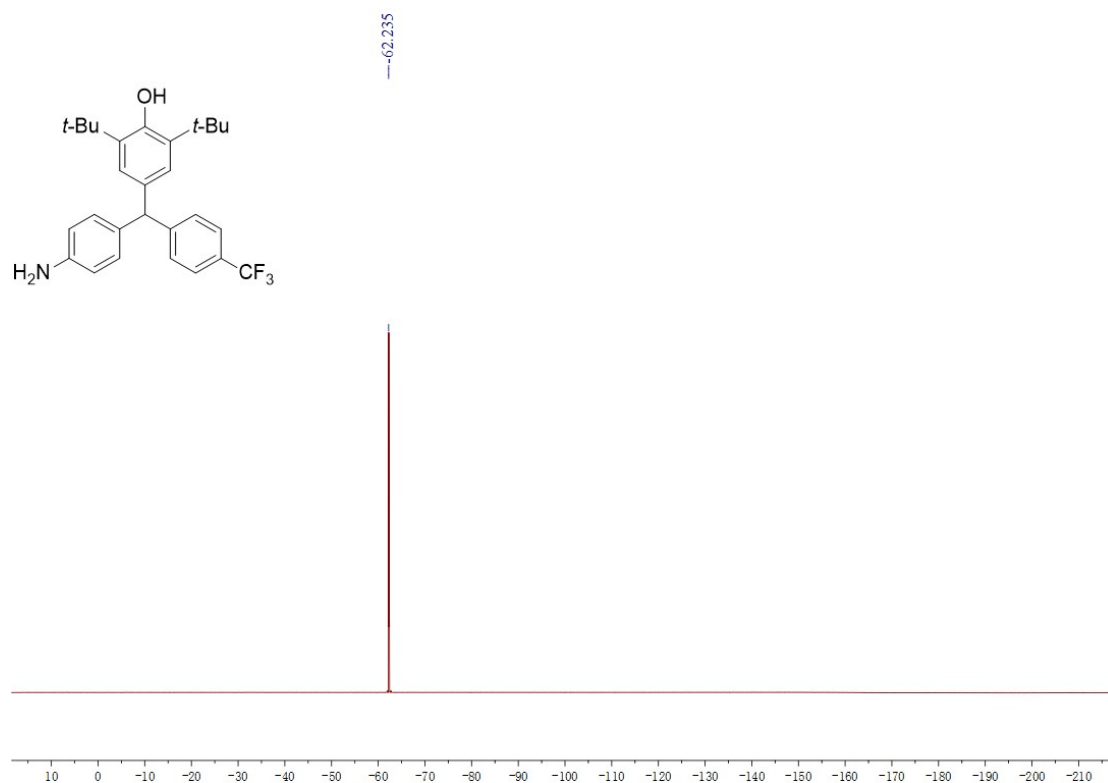


4-((4-Aminophenyl)(3-nitrophenyl)methyl)-2,6-di-tert-butylphenol (4k)

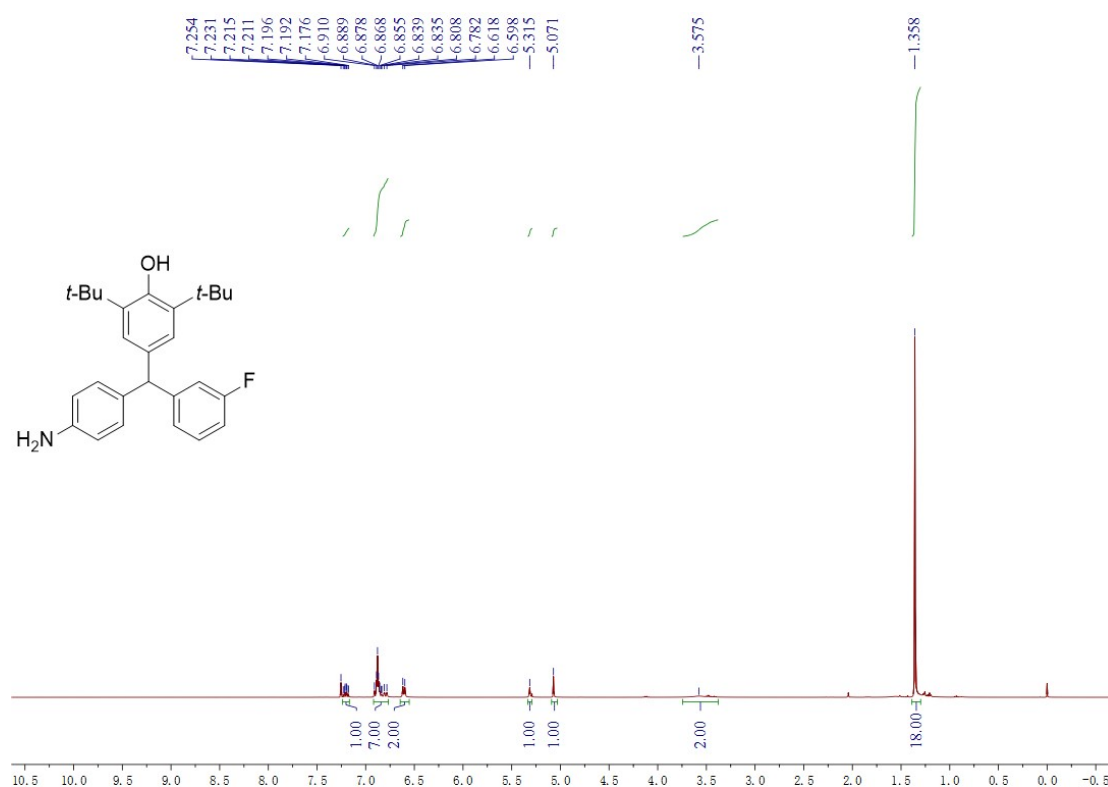


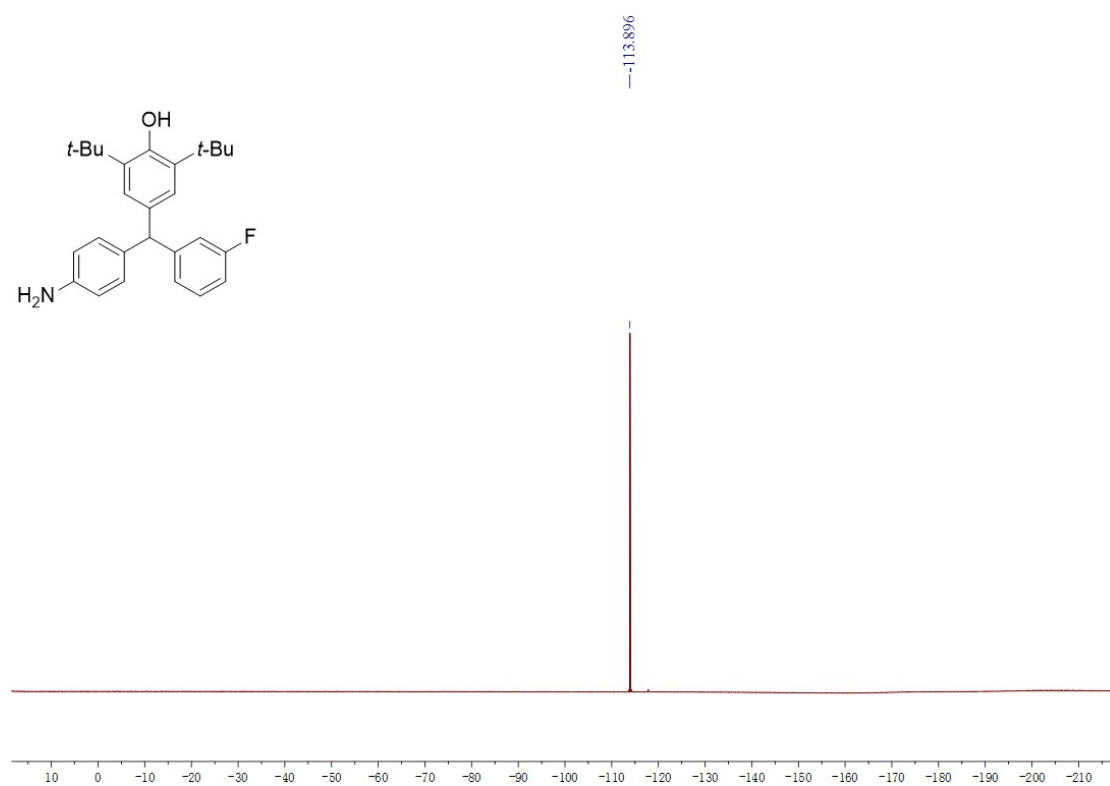
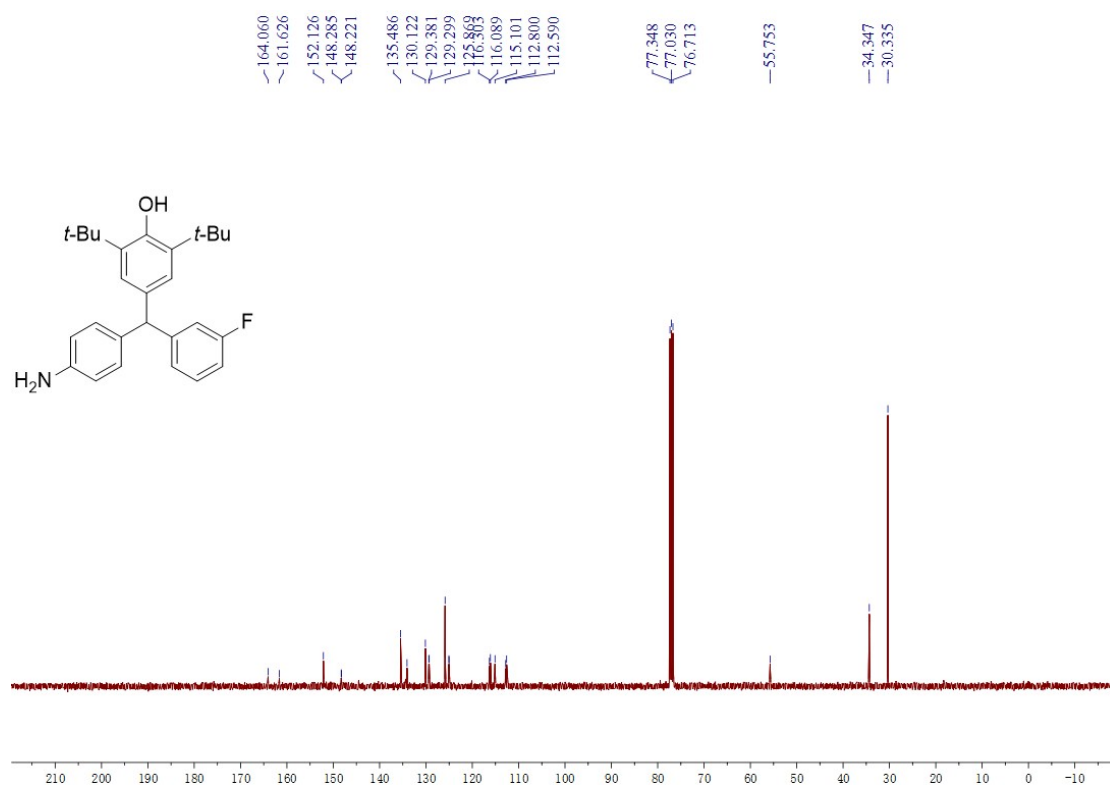
4-((4-Aminophenyl)(4-(trifluoromethyl)phenyl)methyl)-2,6-di-tert-butylphenol (4l)



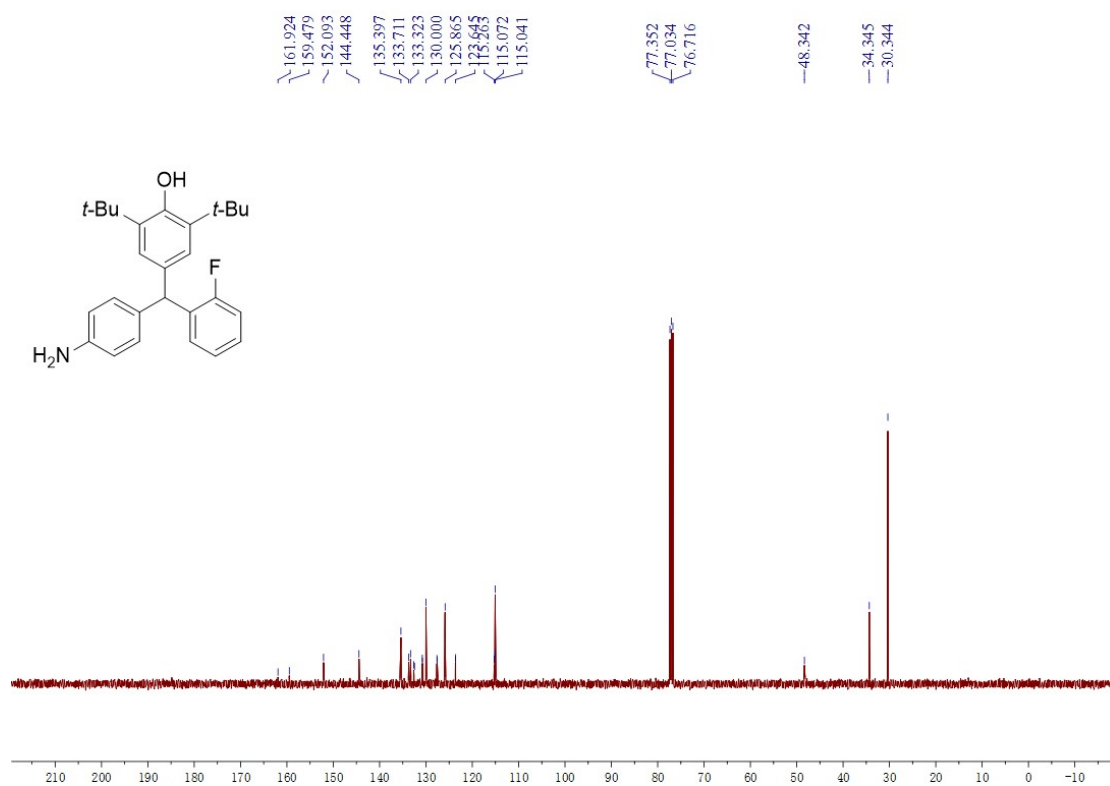
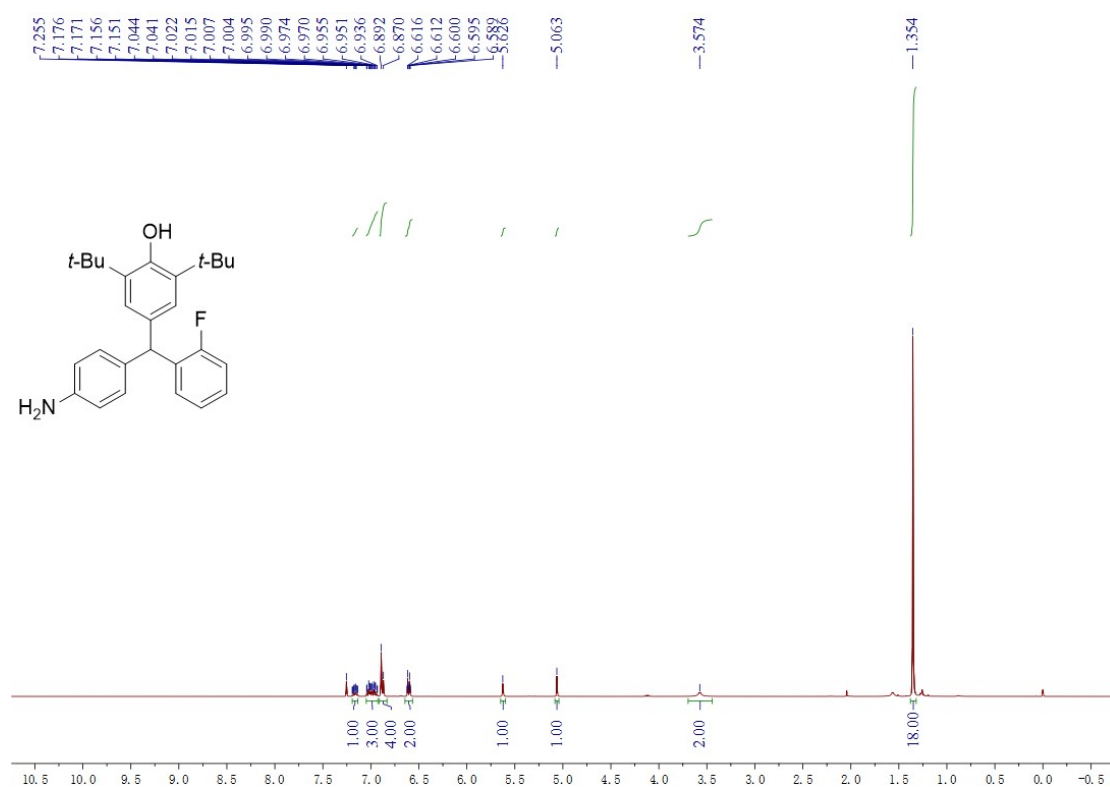


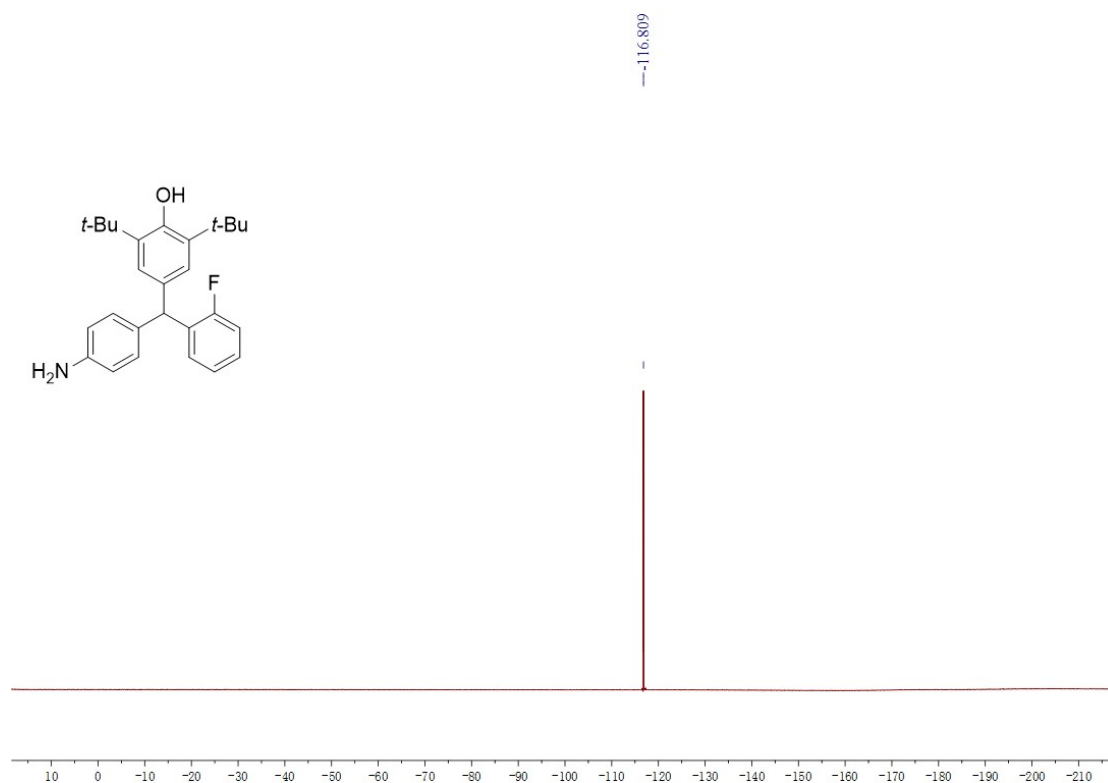
4-((4-Aminophenyl)(3-fluorophenyl)methyl)-2,6-di-tert-butylphenol (4m)



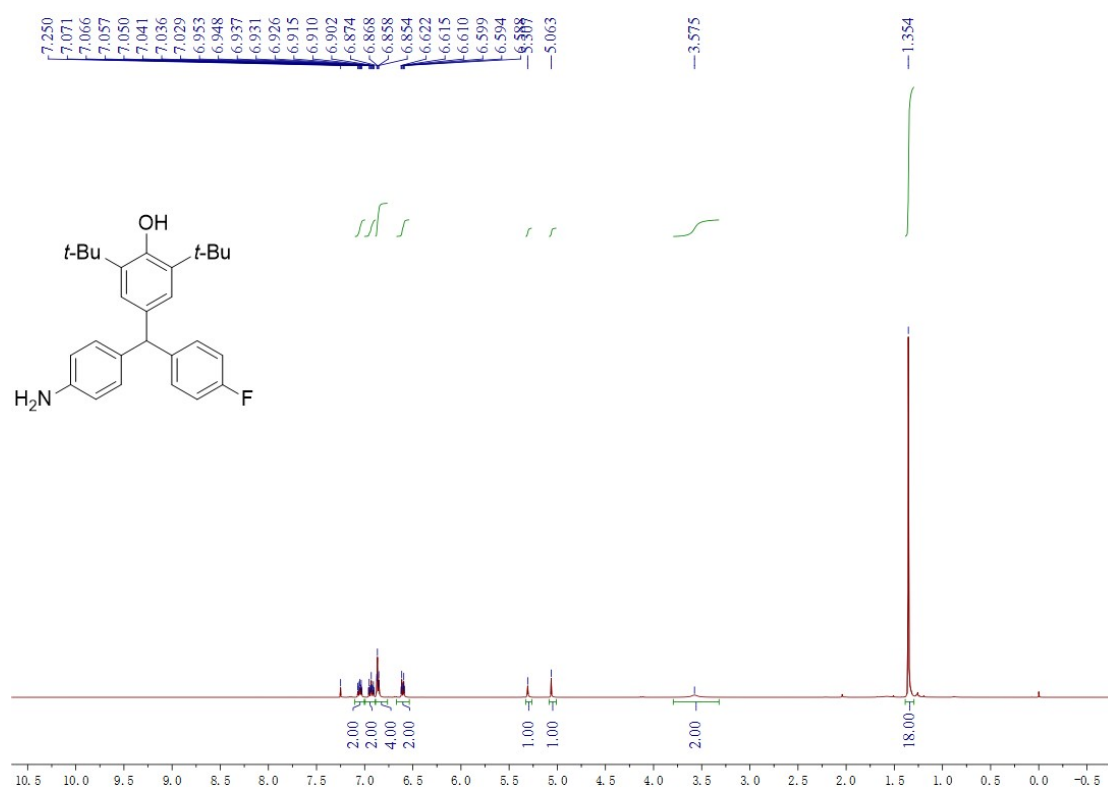


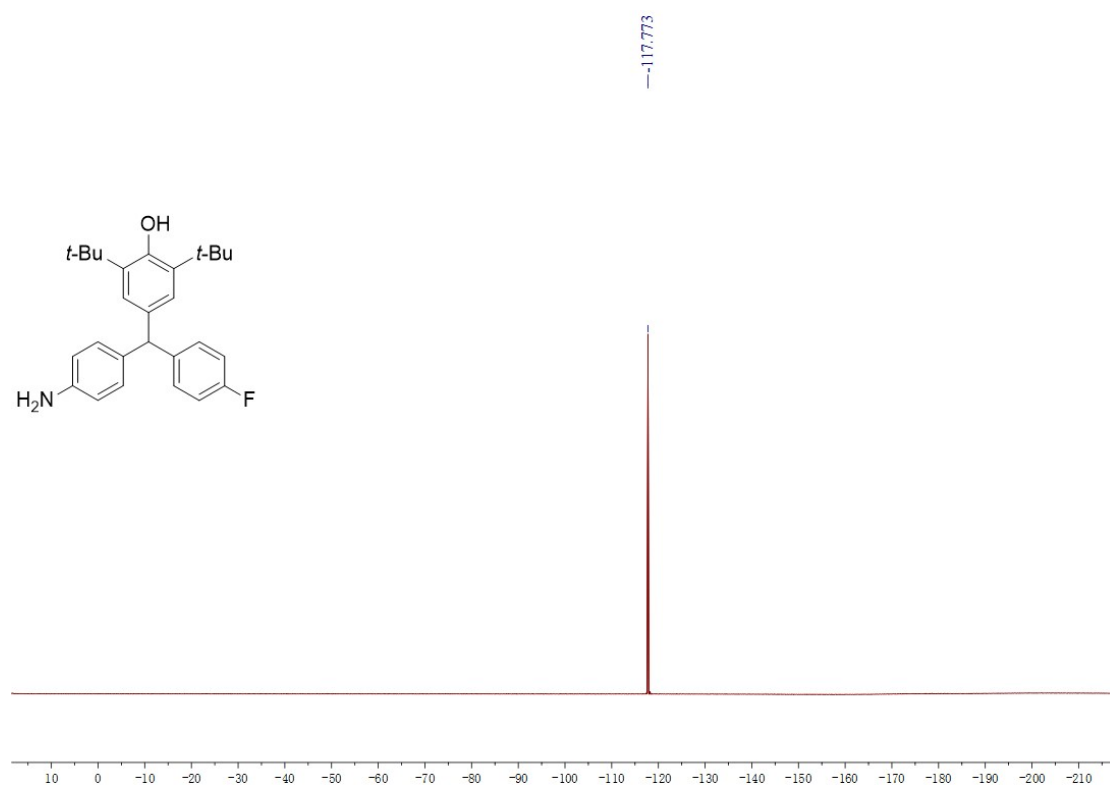
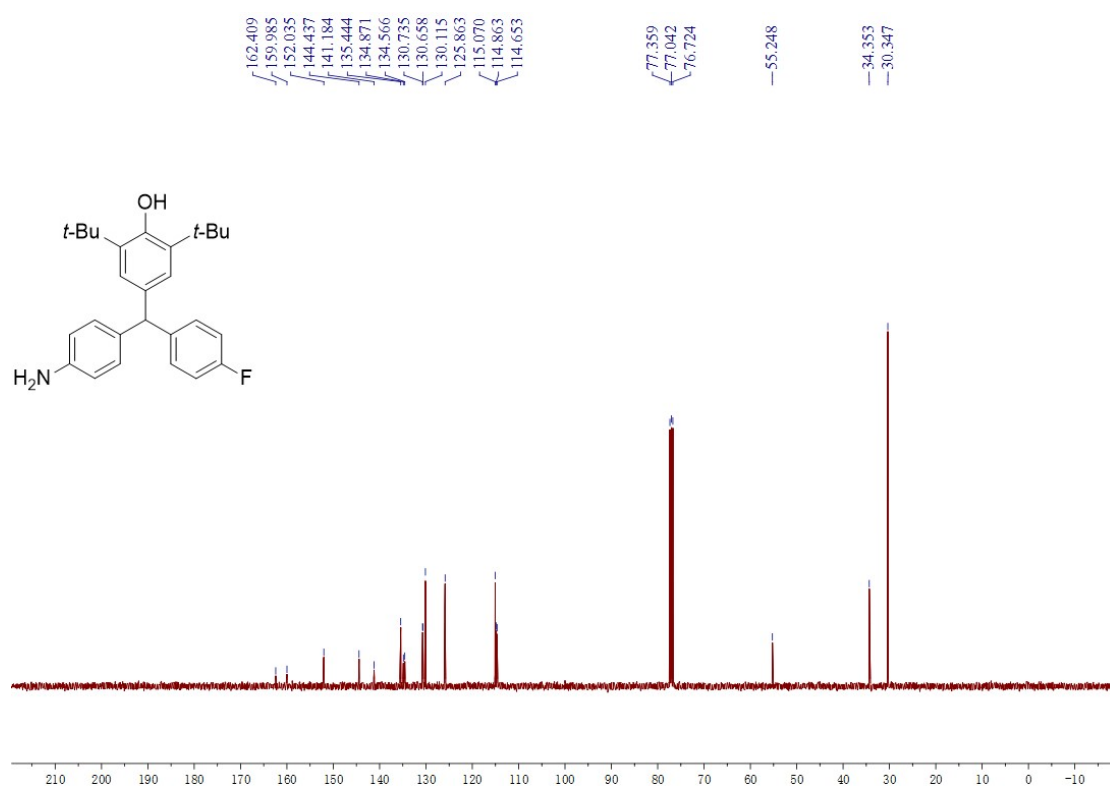
4-((4-Aminophenyl)(2-fluorophenyl)methyl)-2,6-di-tert-butylphenol (4n)



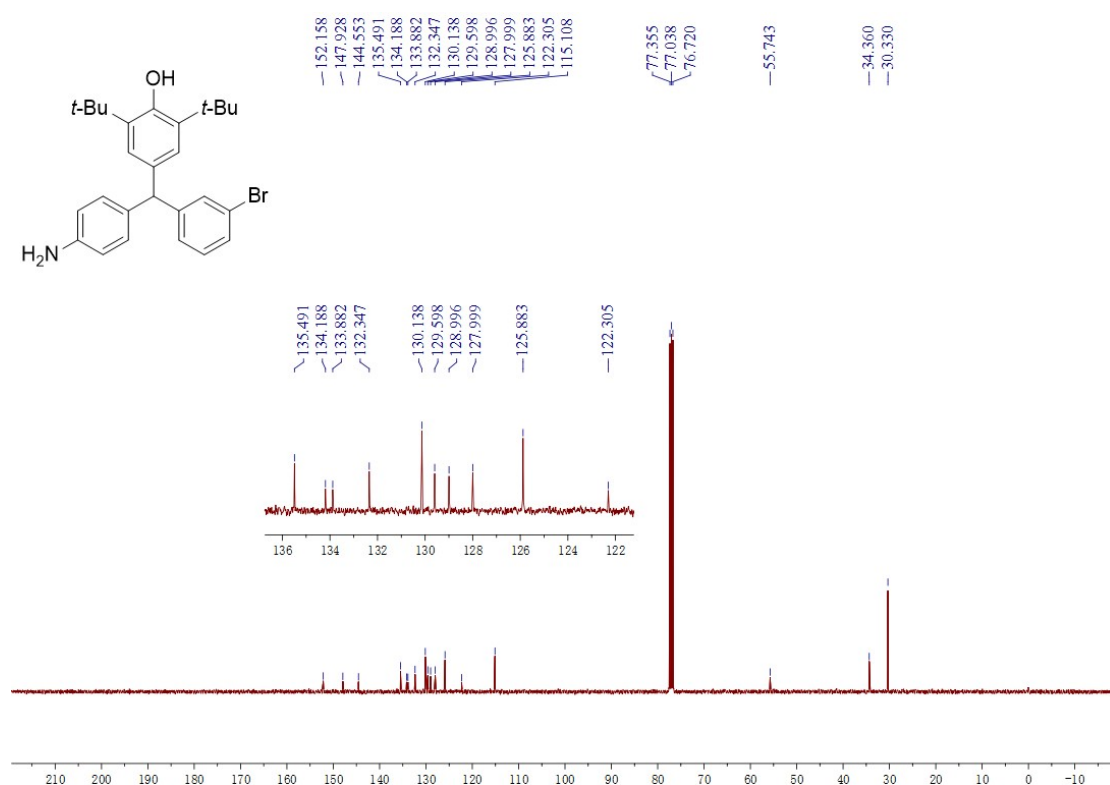
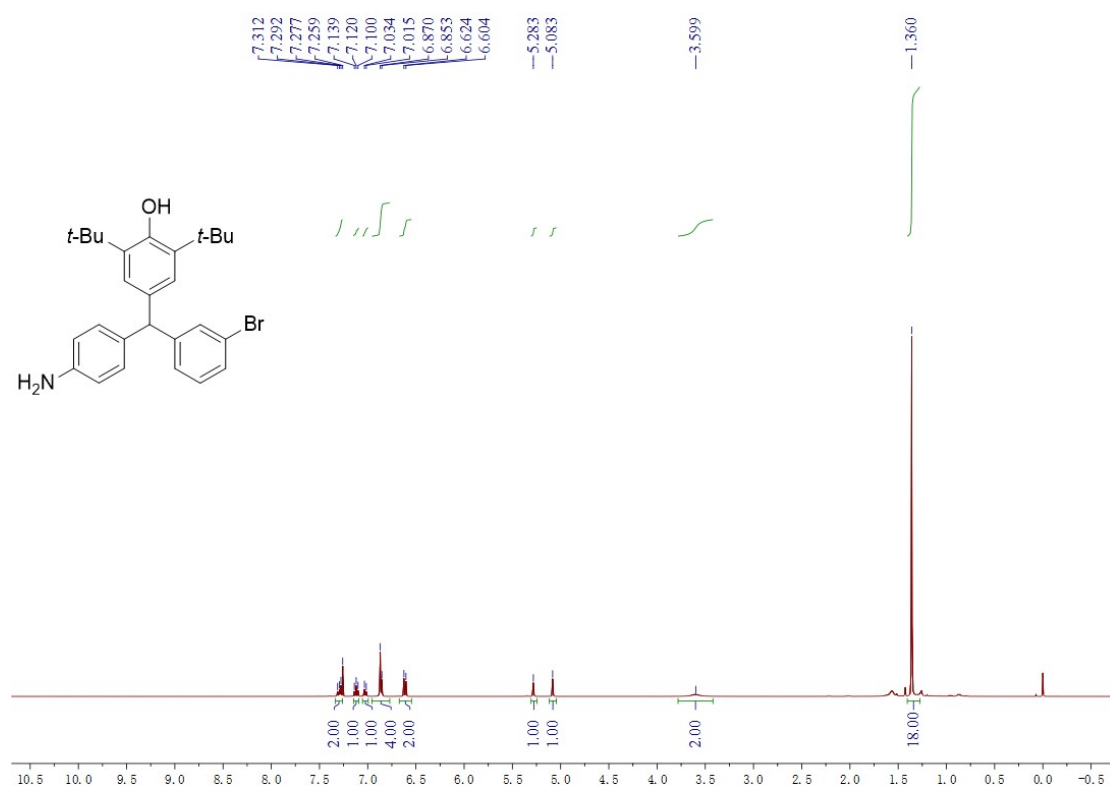


4-((4-Aminophenyl)(4-fluorophenyl)methyl)-2,6-di-tert-butylphenol (4o)

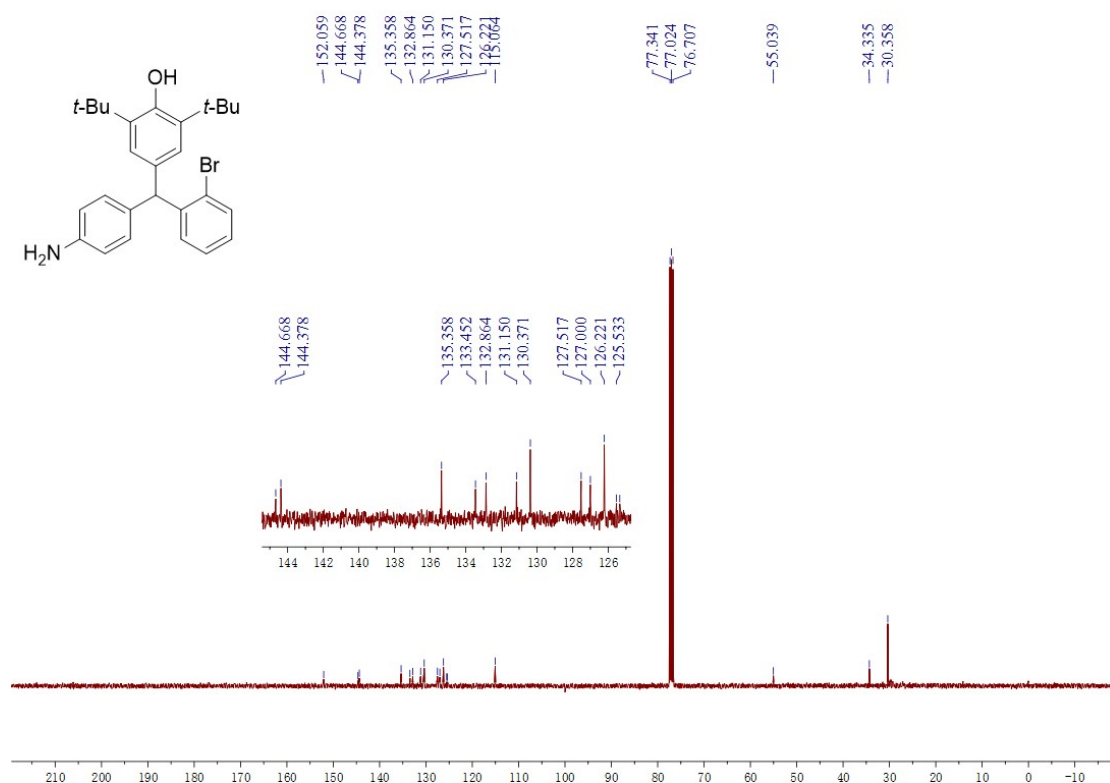
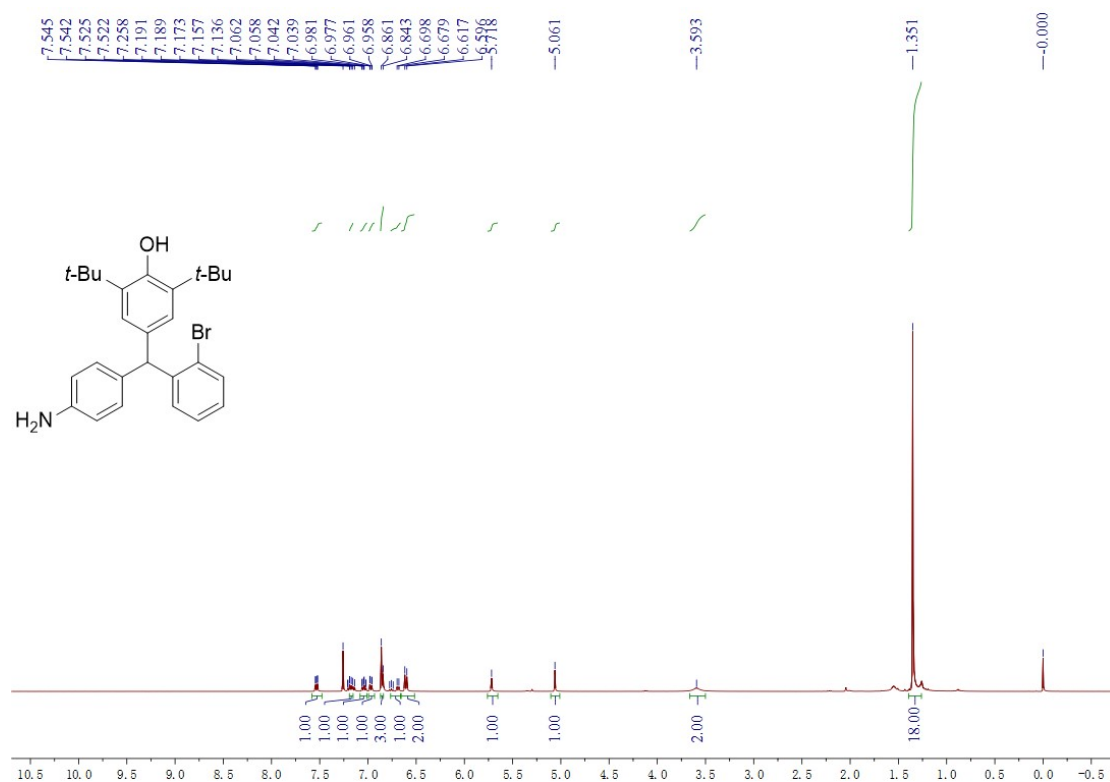




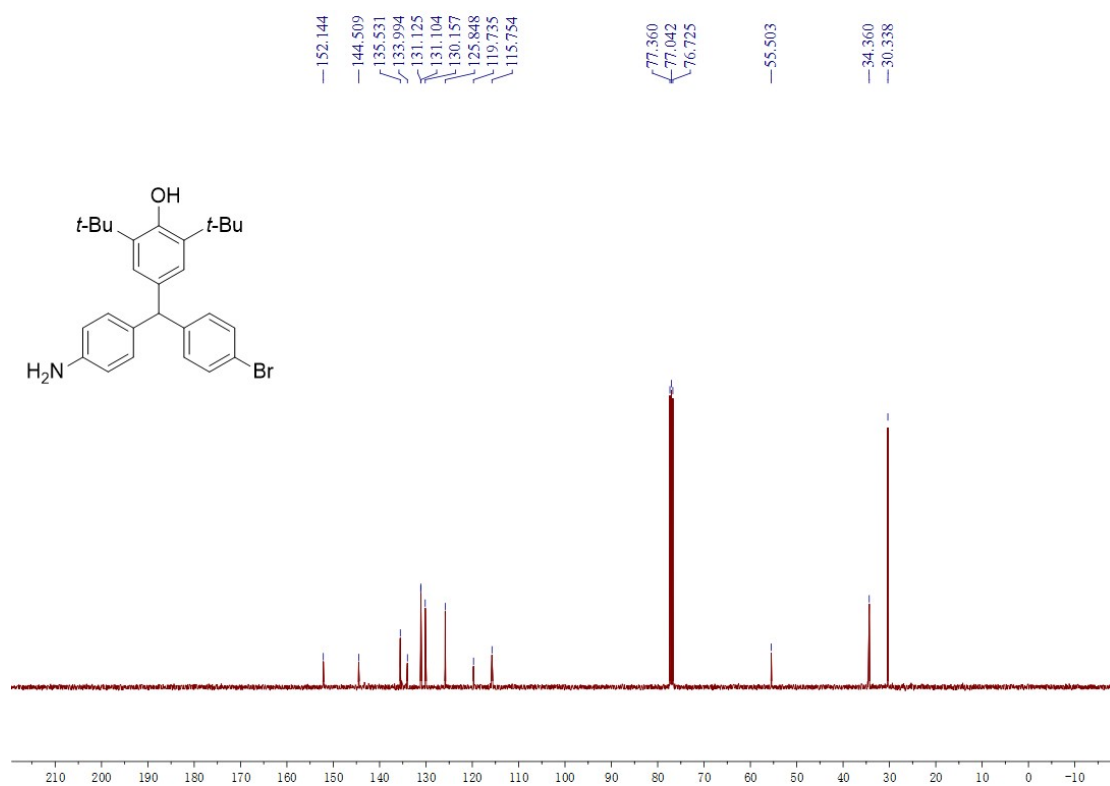
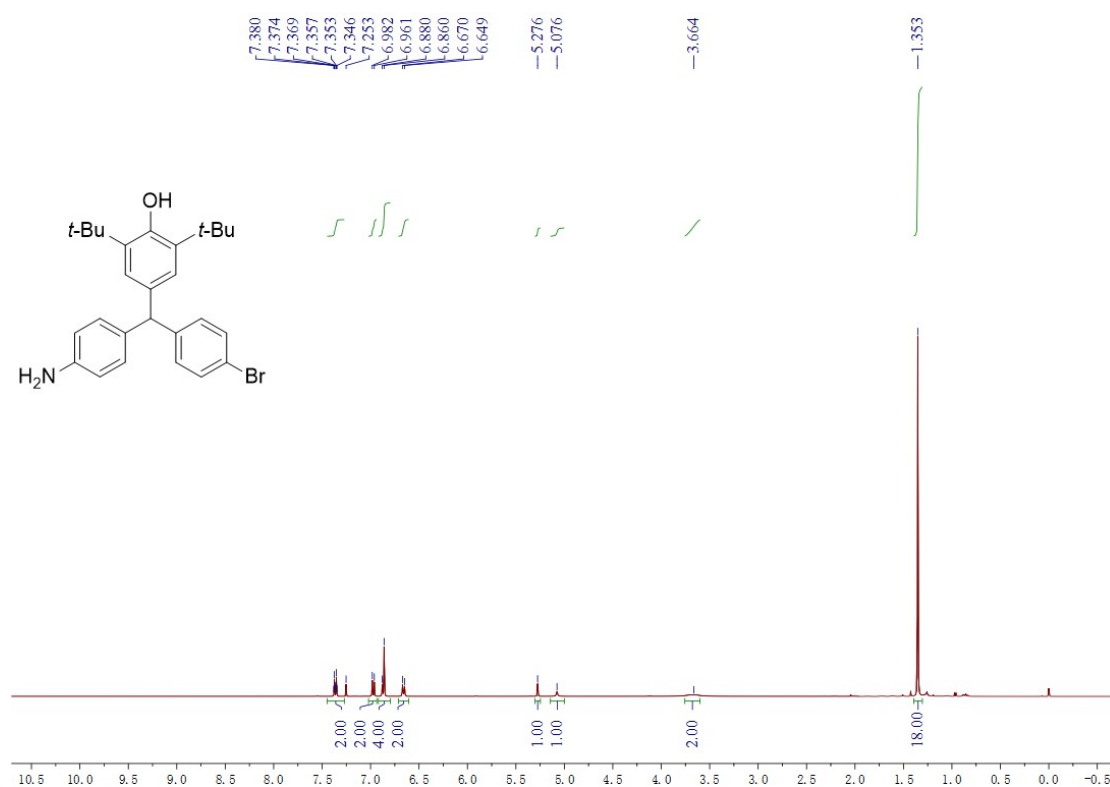
4-((4-Aminophenyl)(3-bromophenyl)methyl)-2,6-di-tert-butylphenol (4p)



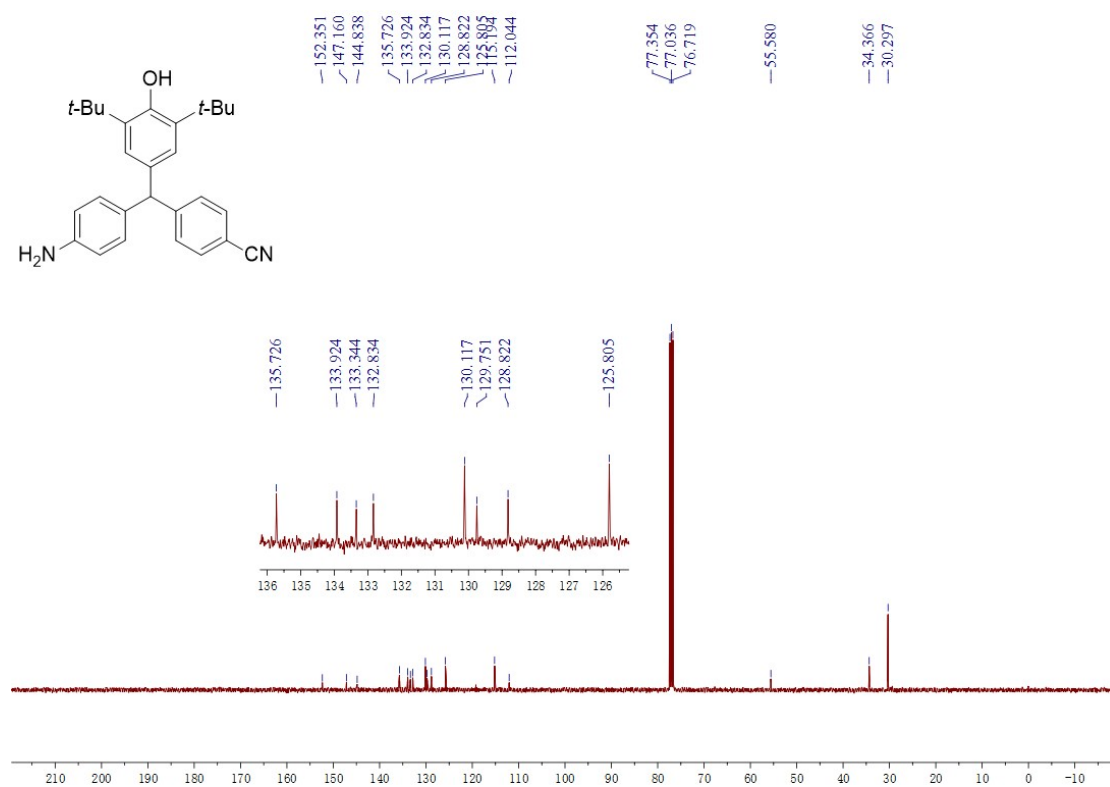
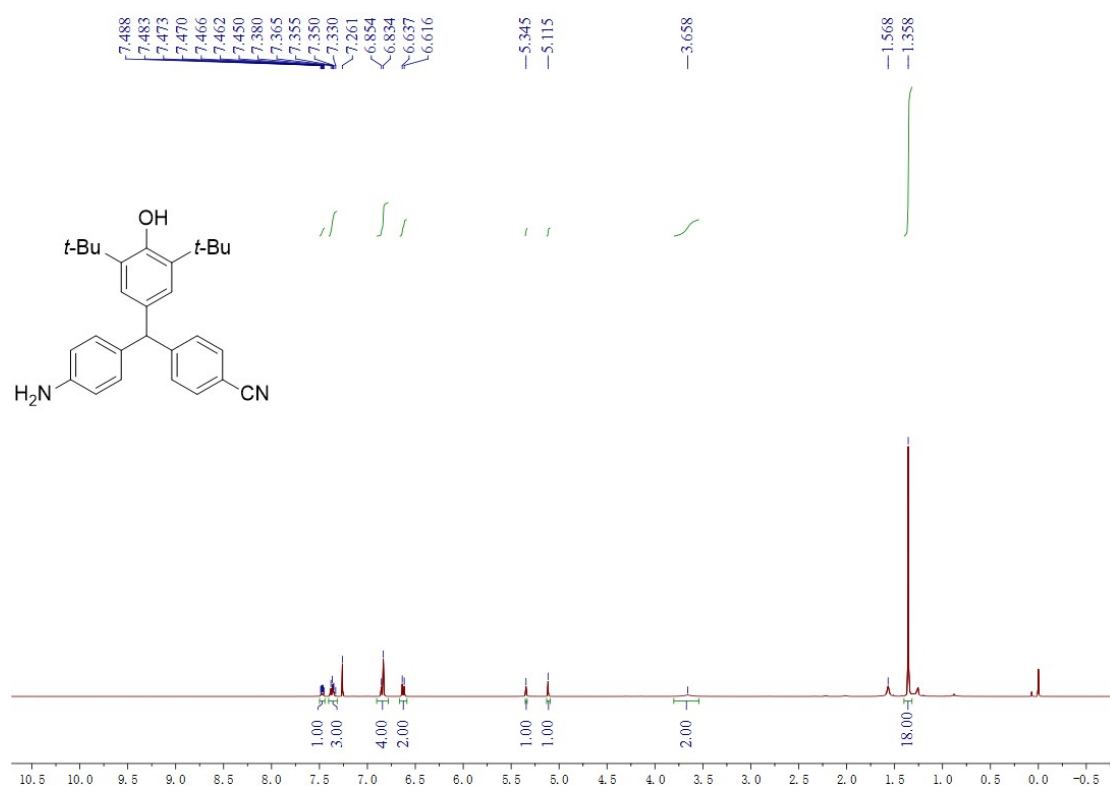
4-((4-Aminophenyl)(2-bromophenyl)methyl)-2,6-di-tert-butylphenol (4q)



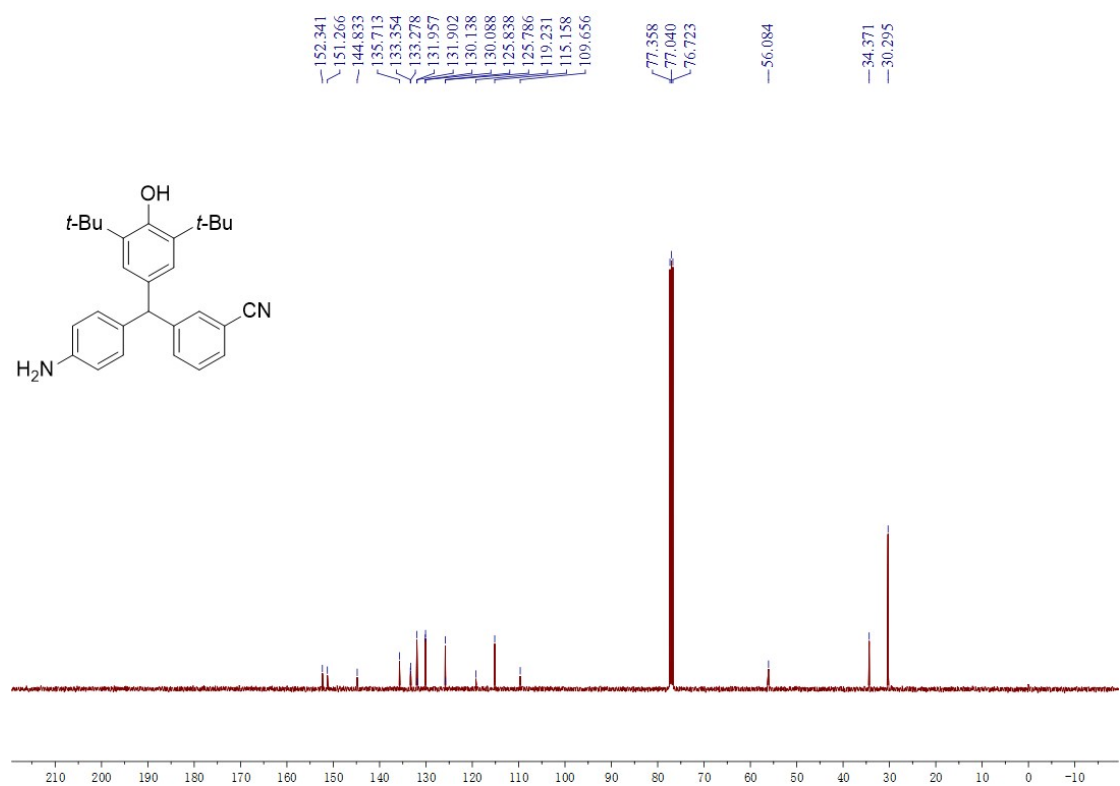
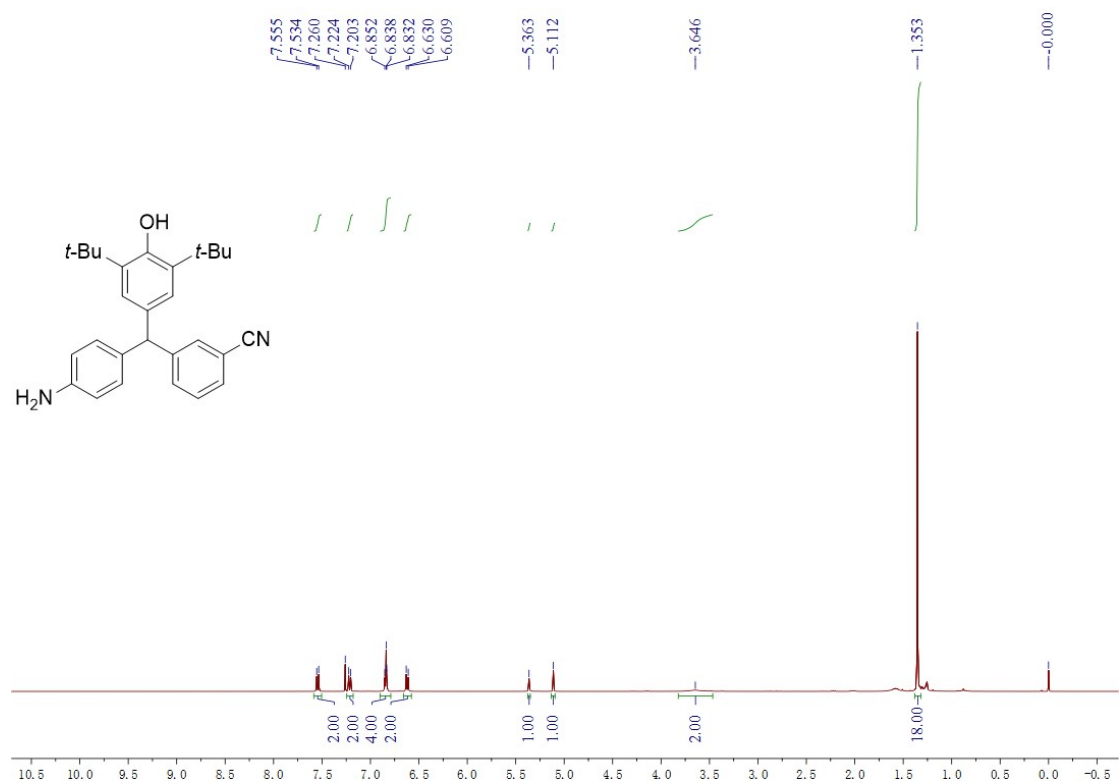
4-((4-Aminophenyl)(4-bromophenyl)methyl)-2,6-di-tert-butylphenol (4r)



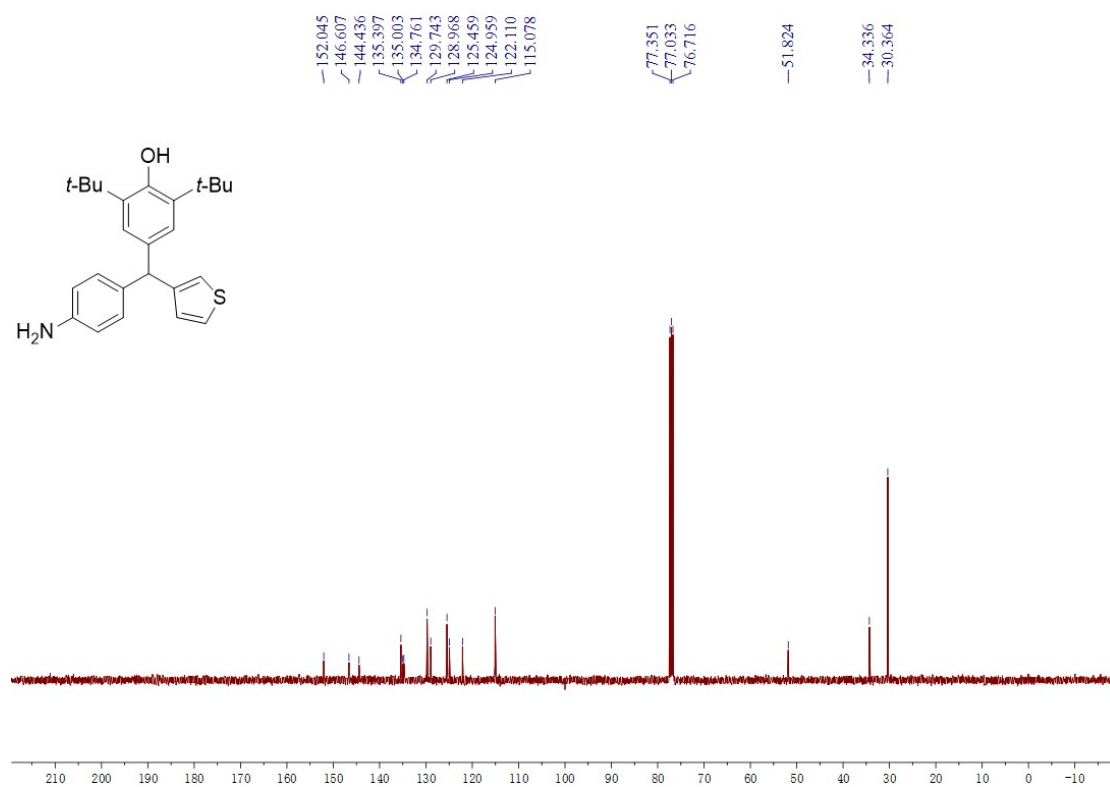
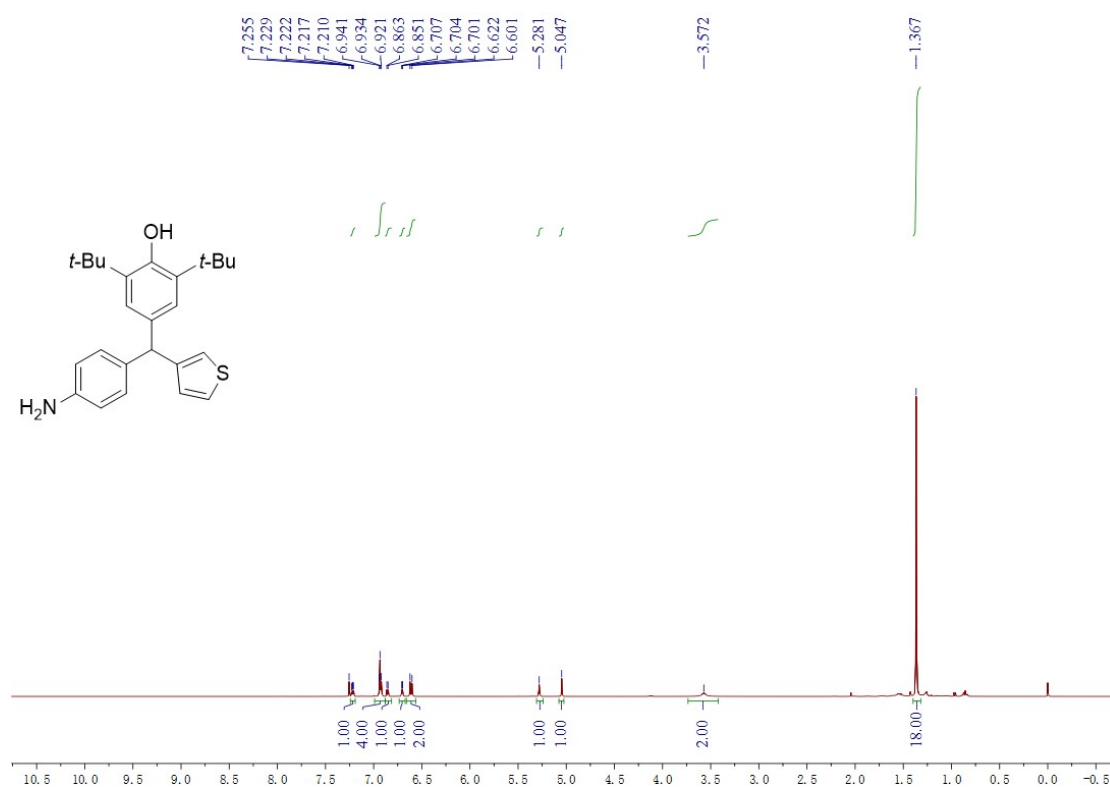
4-((4-Aminophenyl)(3,5-di-tert-butyl-4-hydroxyphenyl)methyl)benzonitrile (4s)



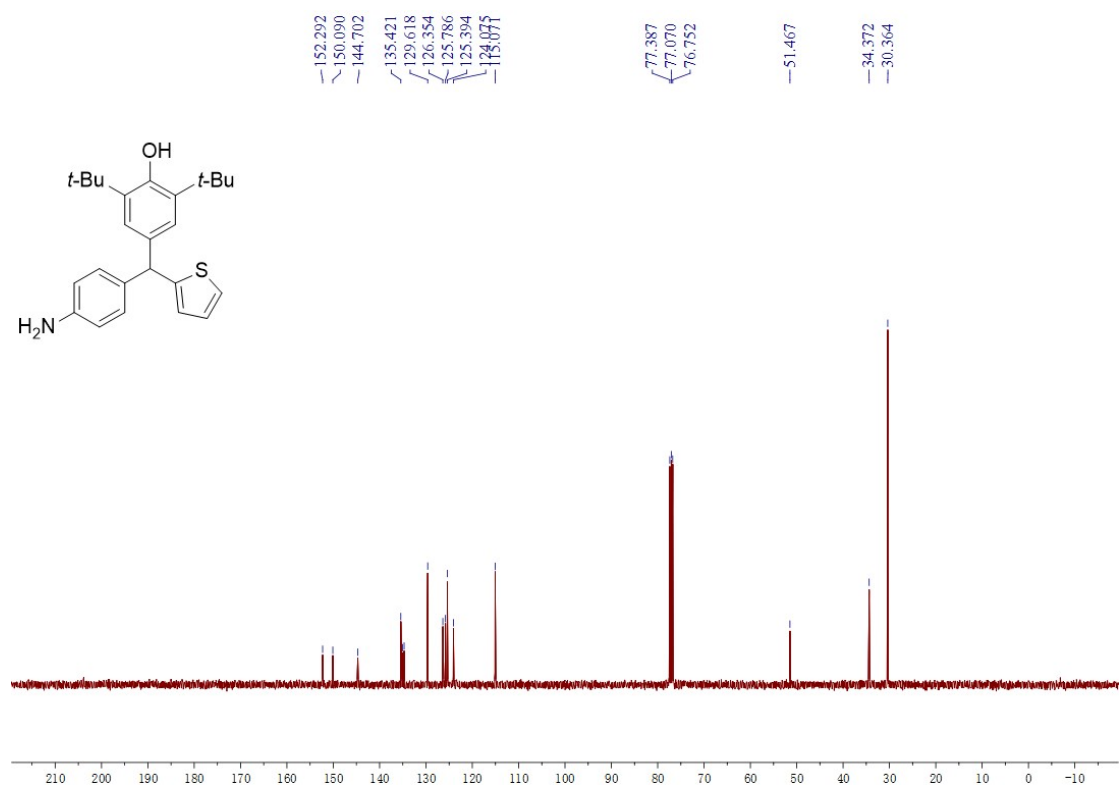
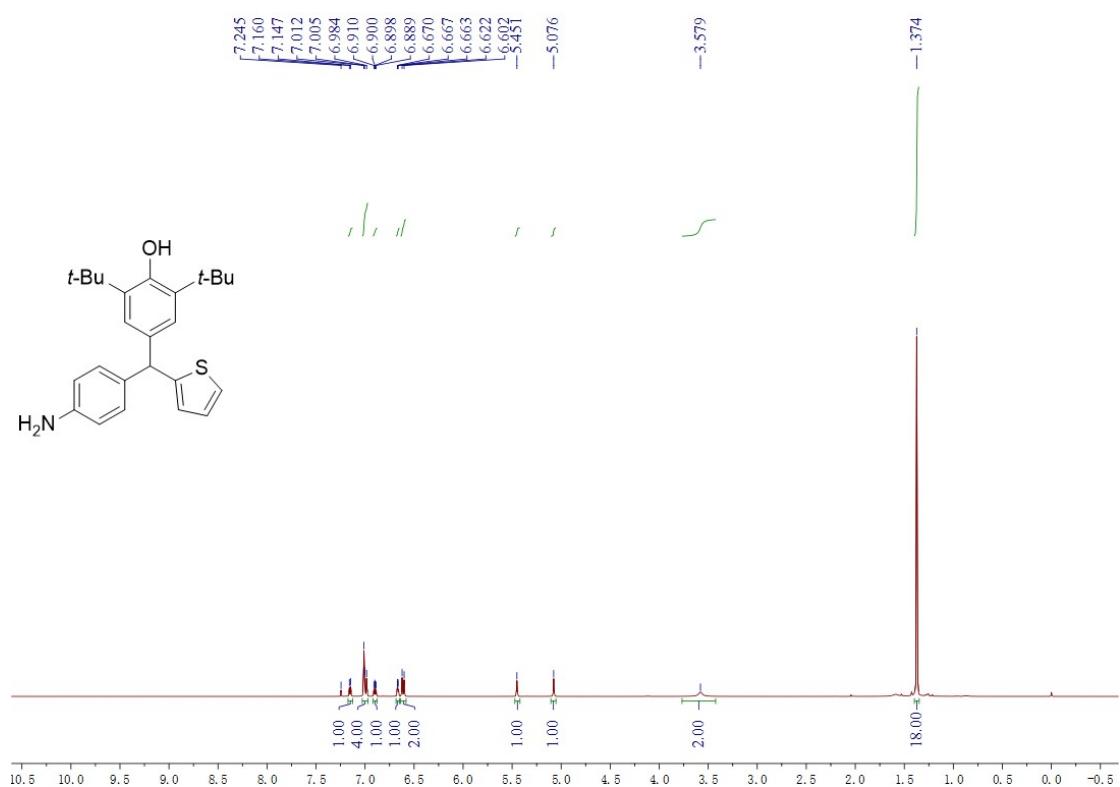
3-((4-Aminophenyl)(3,5-di-tert-butyl-4-hydroxyphenyl)methyl)benzonitrile (4t)



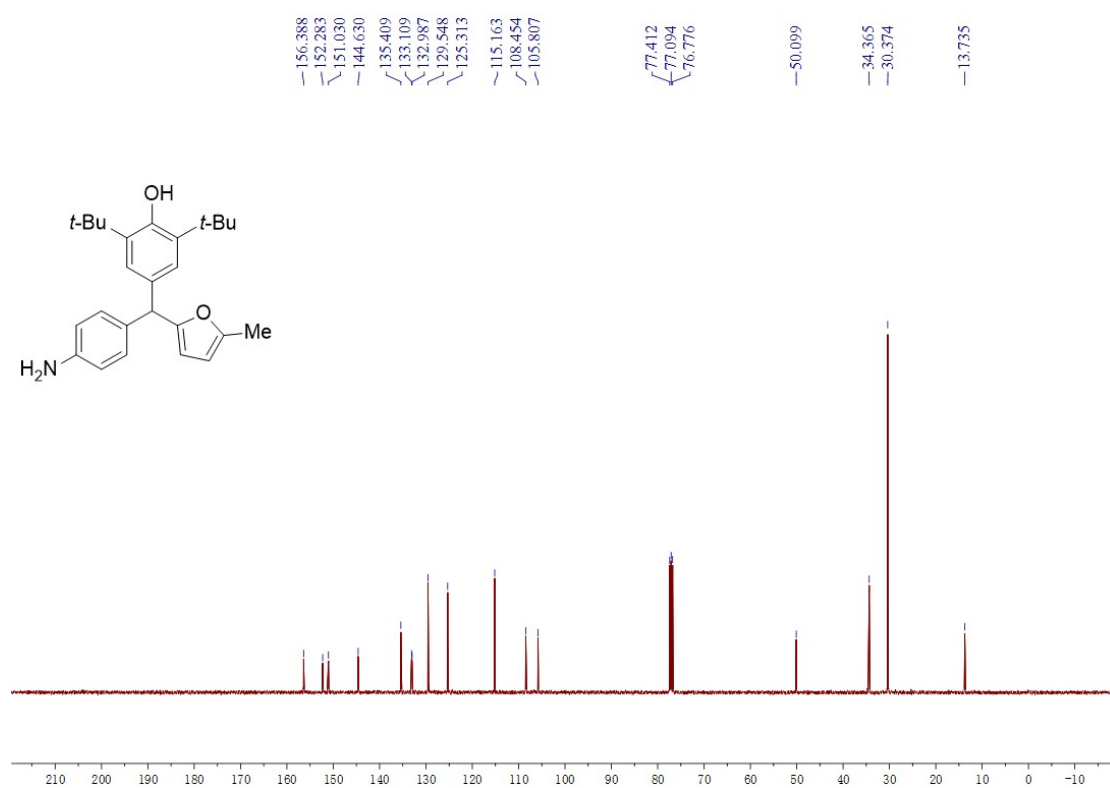
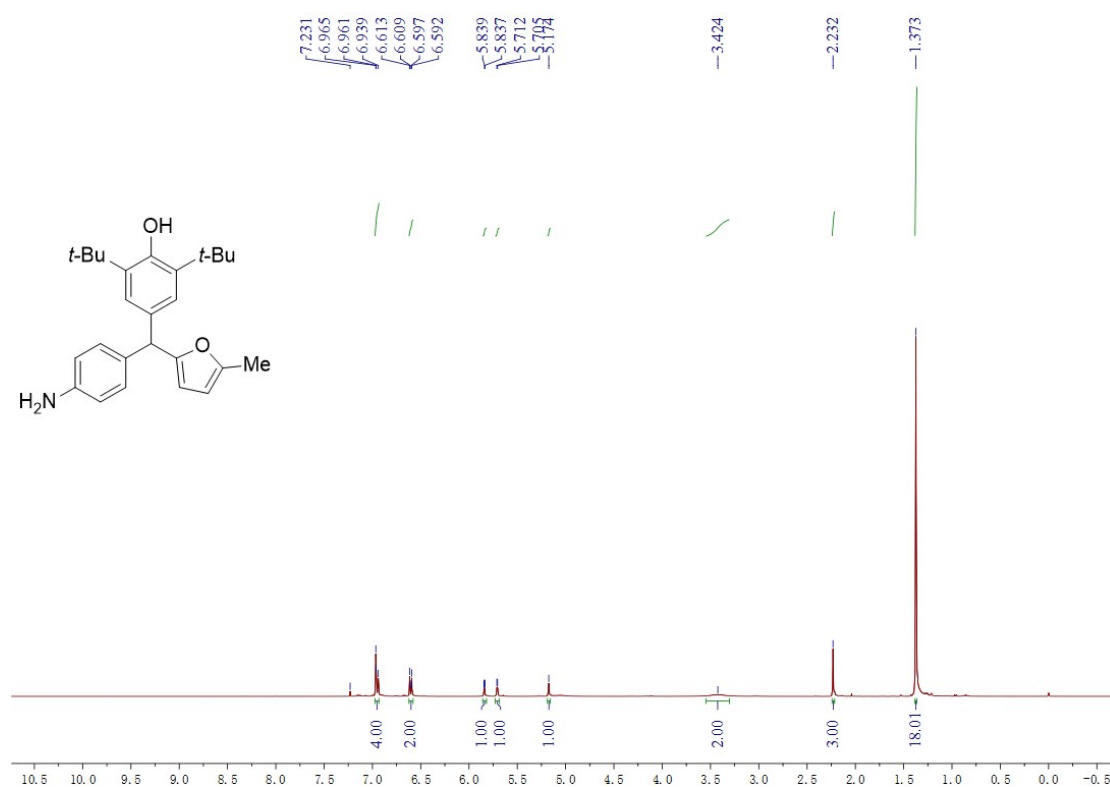
4-((4-Aminophenyl)(thiophen-3-yl)methyl)-2,6-di-tert-butylphenol (4u)



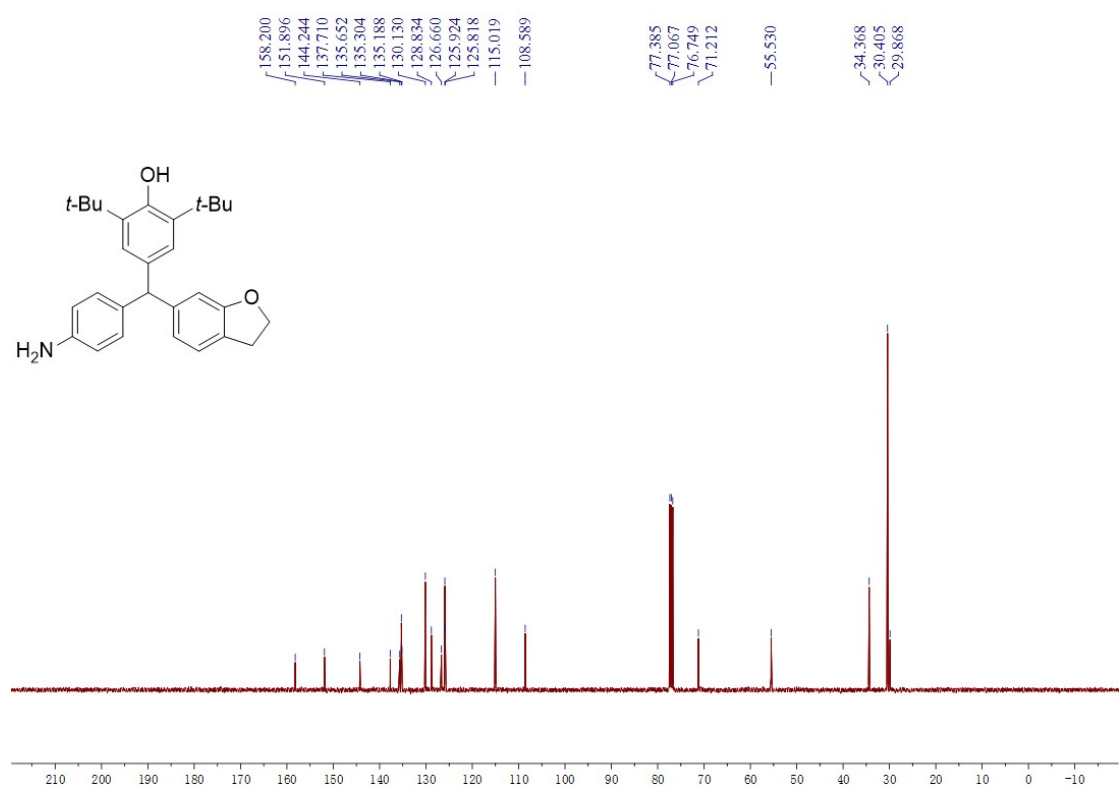
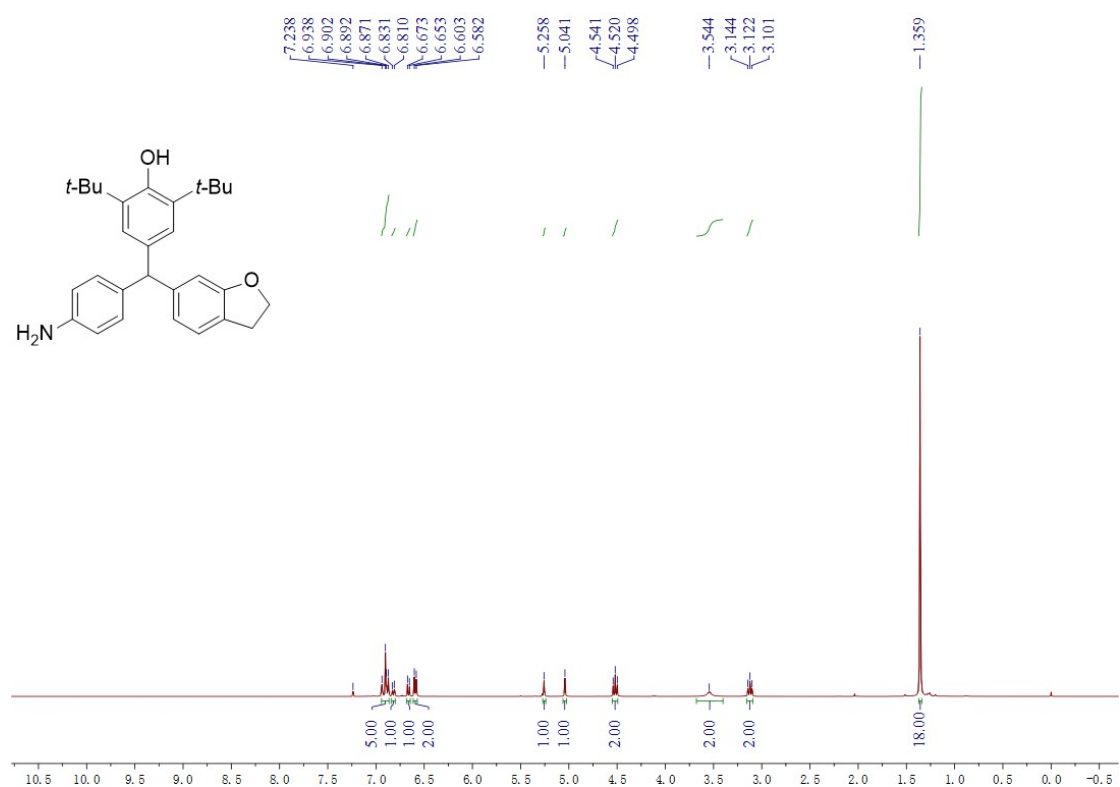
4-((4-Aminophenyl)(thiophen-2-yl)methyl)-2,6-di-tert-butylphenol (4v)



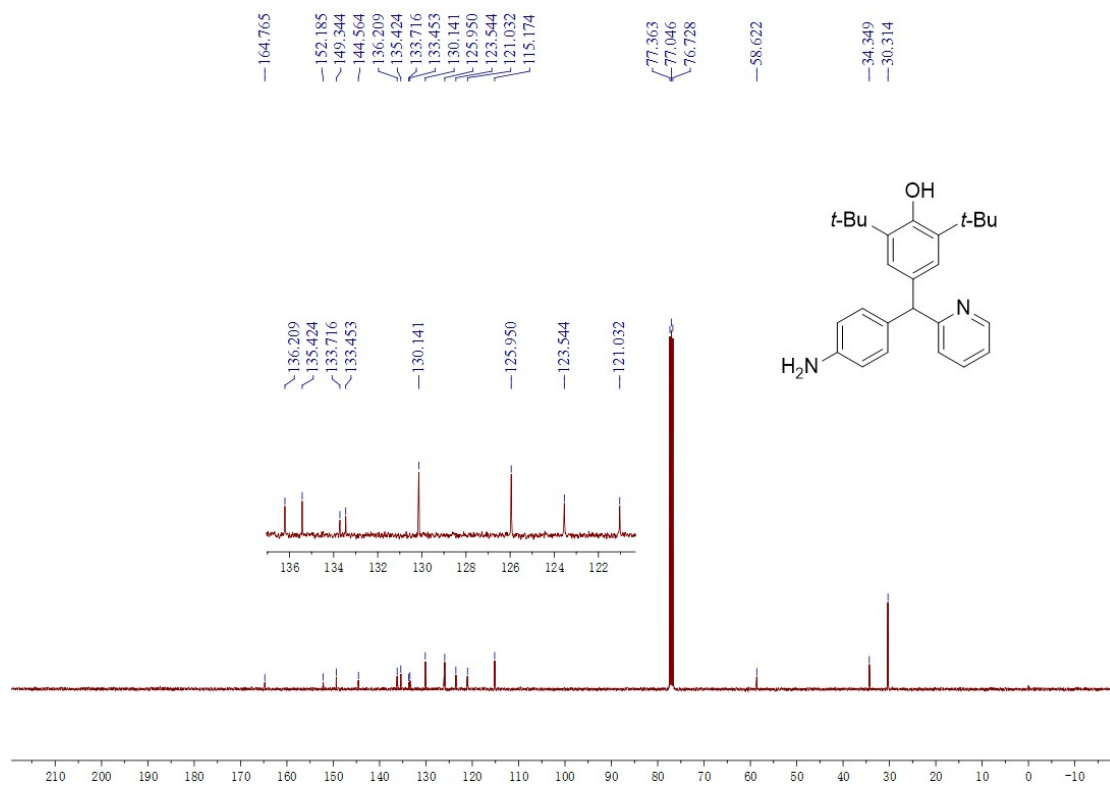
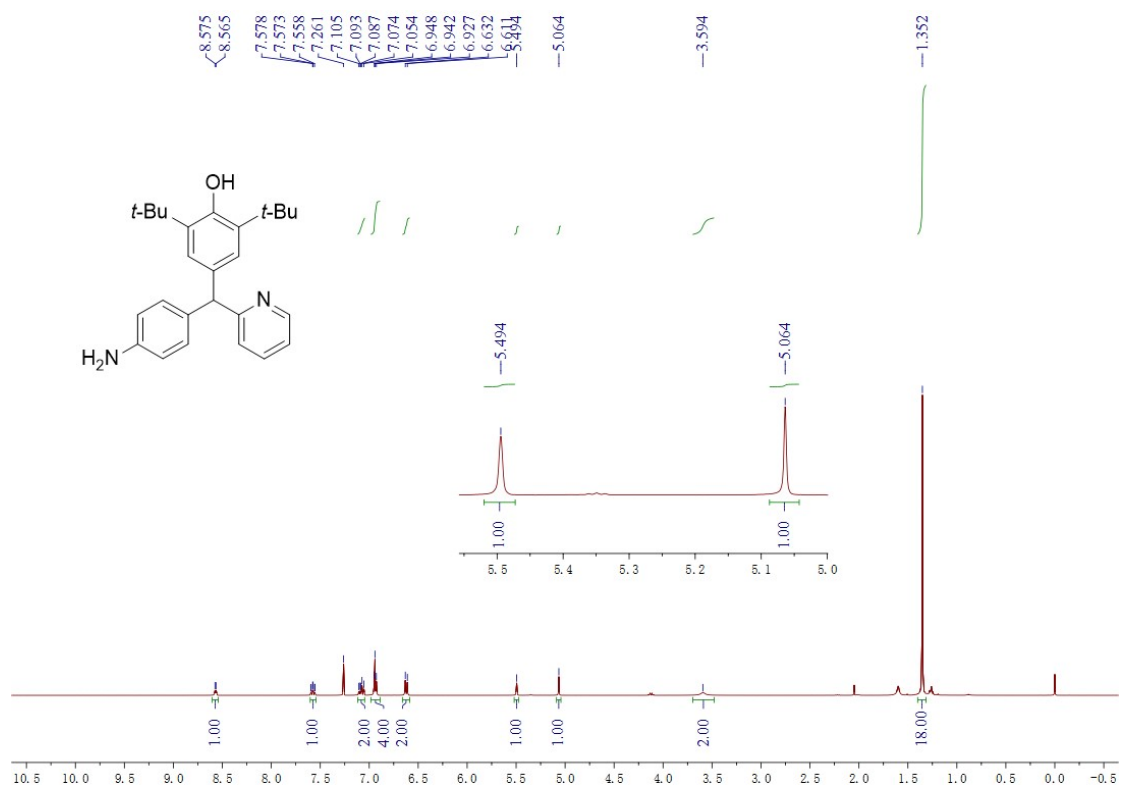
4-((4-Aminophenyl)(5-methylfuran-2-yl)methyl)-2,6-di-tert-butylphenol (4w)



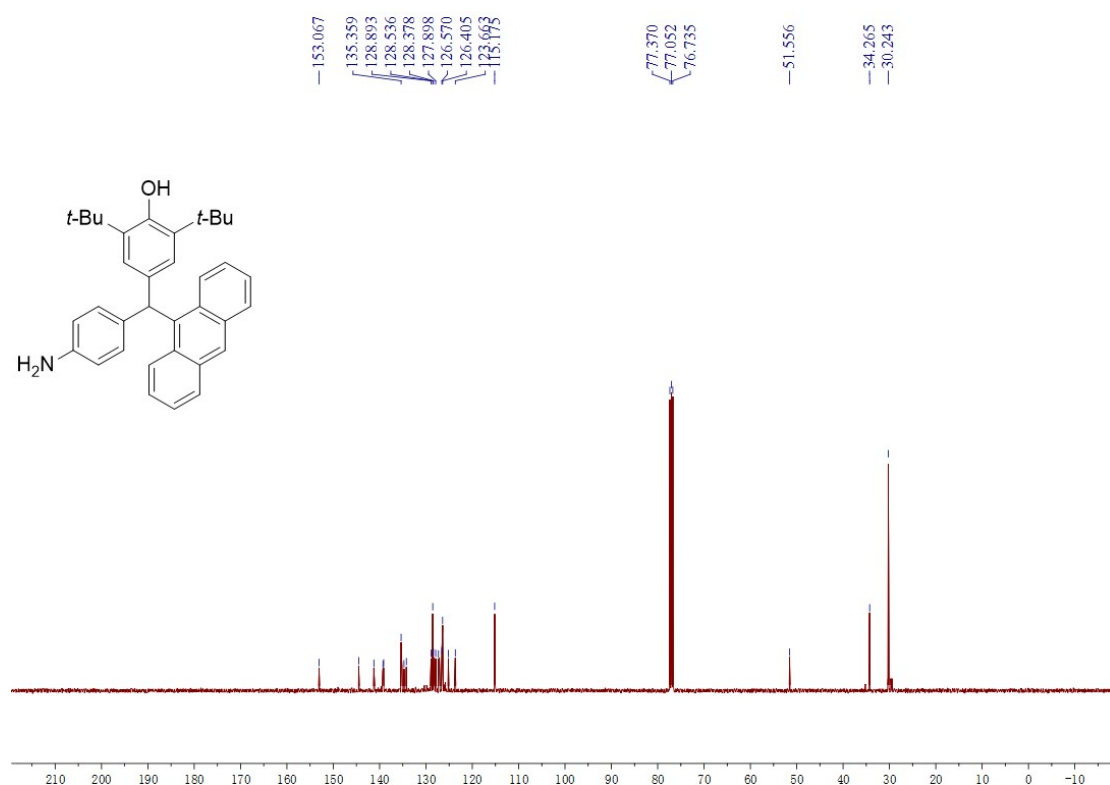
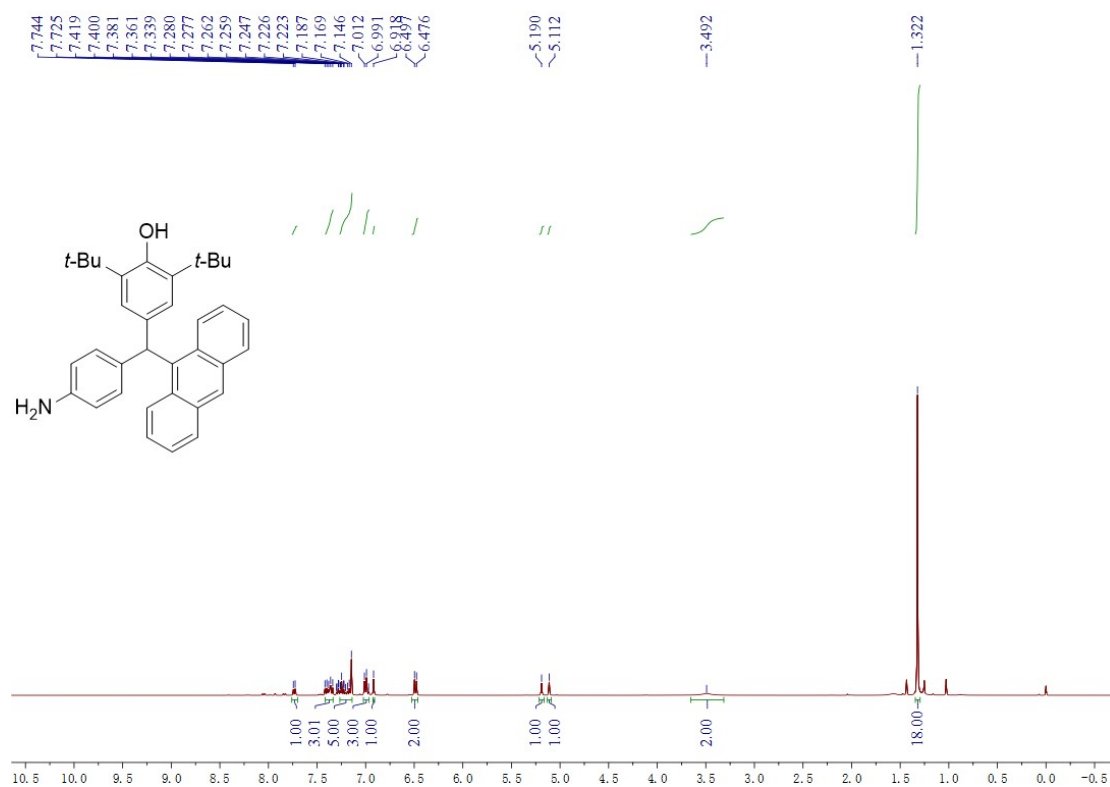
4-((4-Aminophenyl)(2,3-dihydrobenzofuran-6-yl)methyl)-2,6-di-tert-butylphenol (4x)



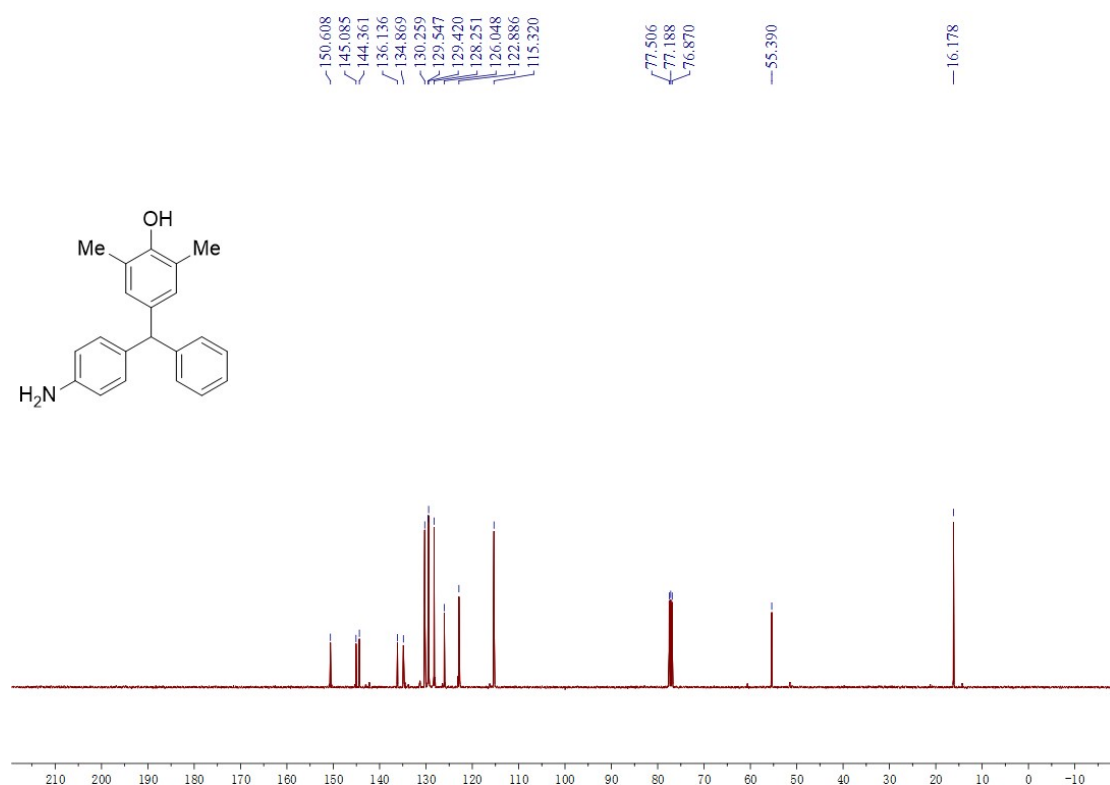
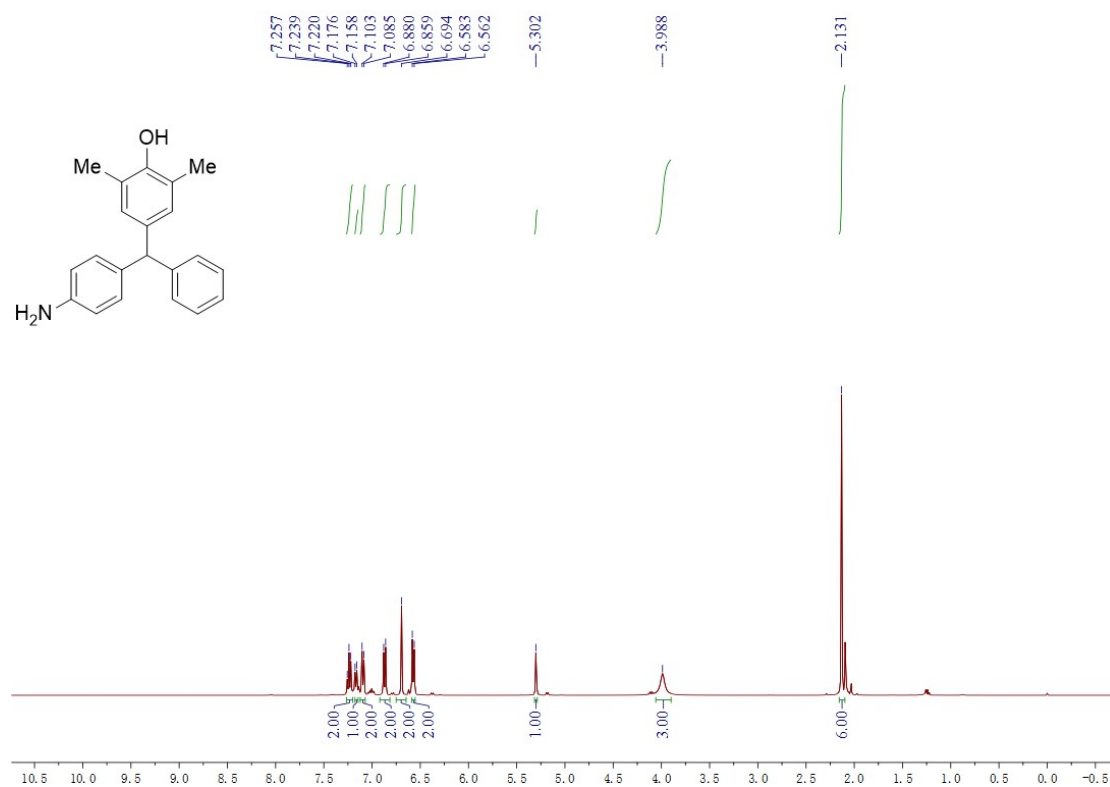
4-((4-Aminophenyl)(pyridin-2-yl)methyl)-2,6-di-tert-butylphenol (4y)



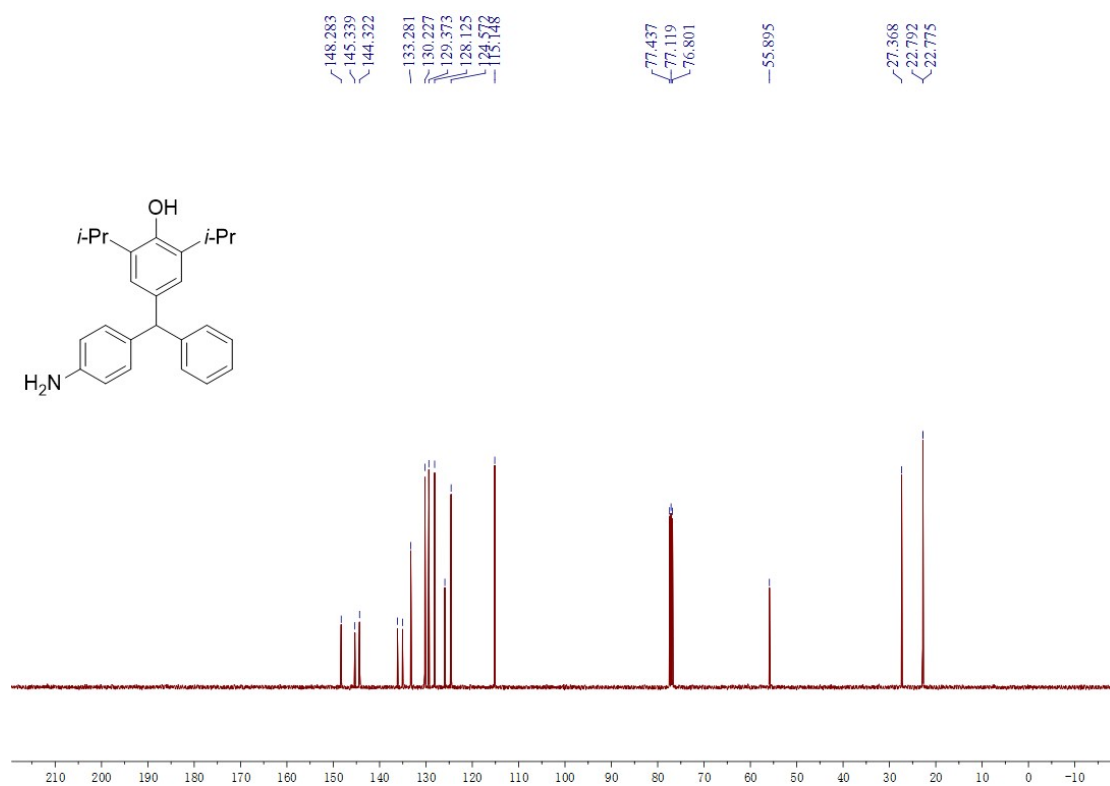
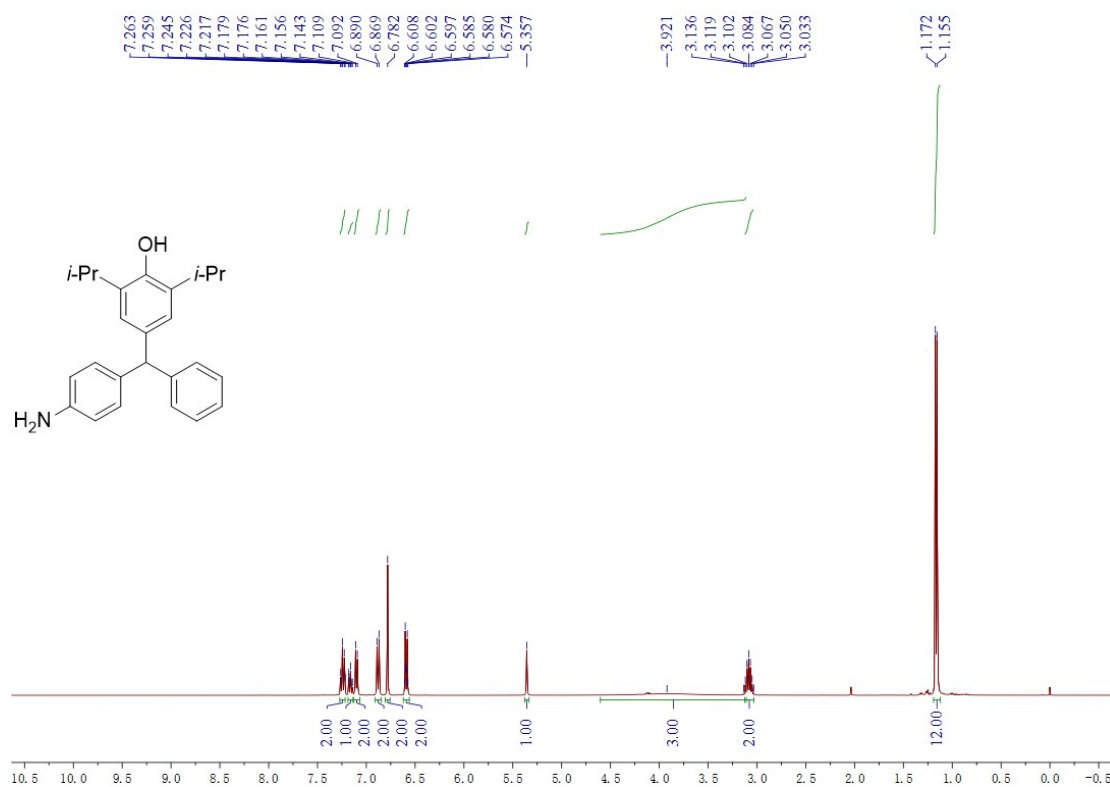
4-((4-Aminophenyl)(anthracen-9-yl)methyl)-2,6-di-tert-butylphenol (4aa)



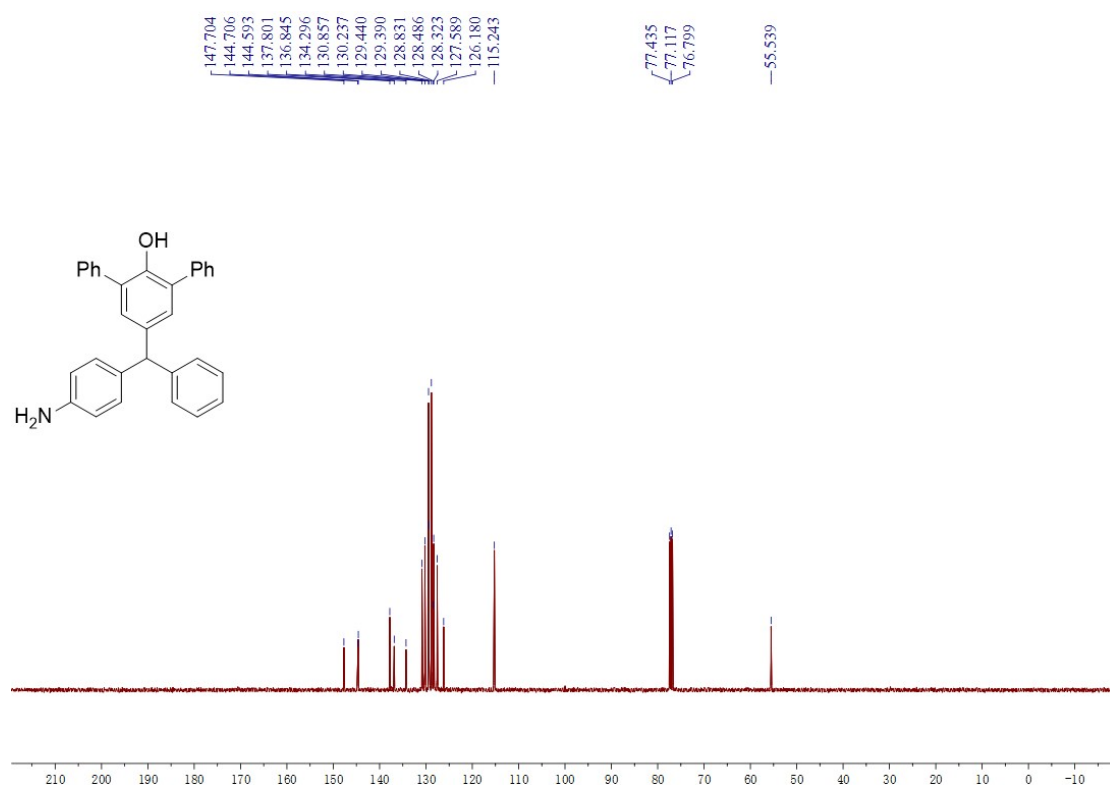
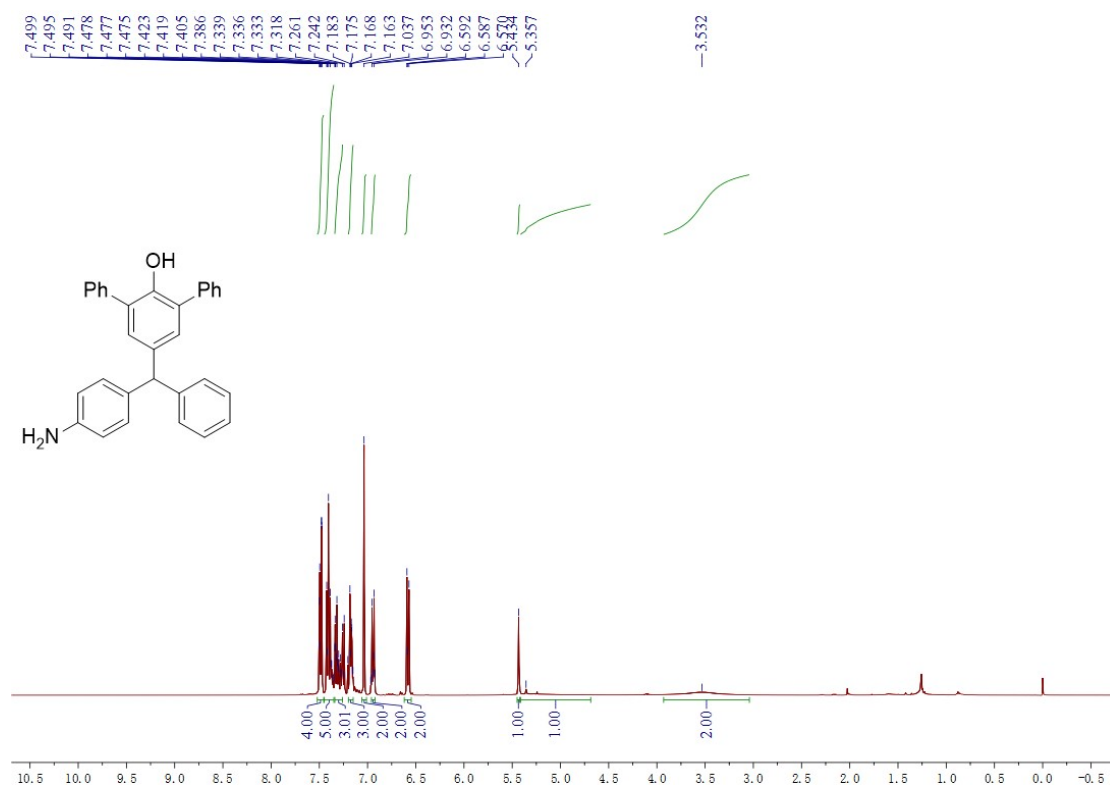
4-((4-Aminophenyl)(phenyl)methyl)-2,6-dimethylphenol (4ab)



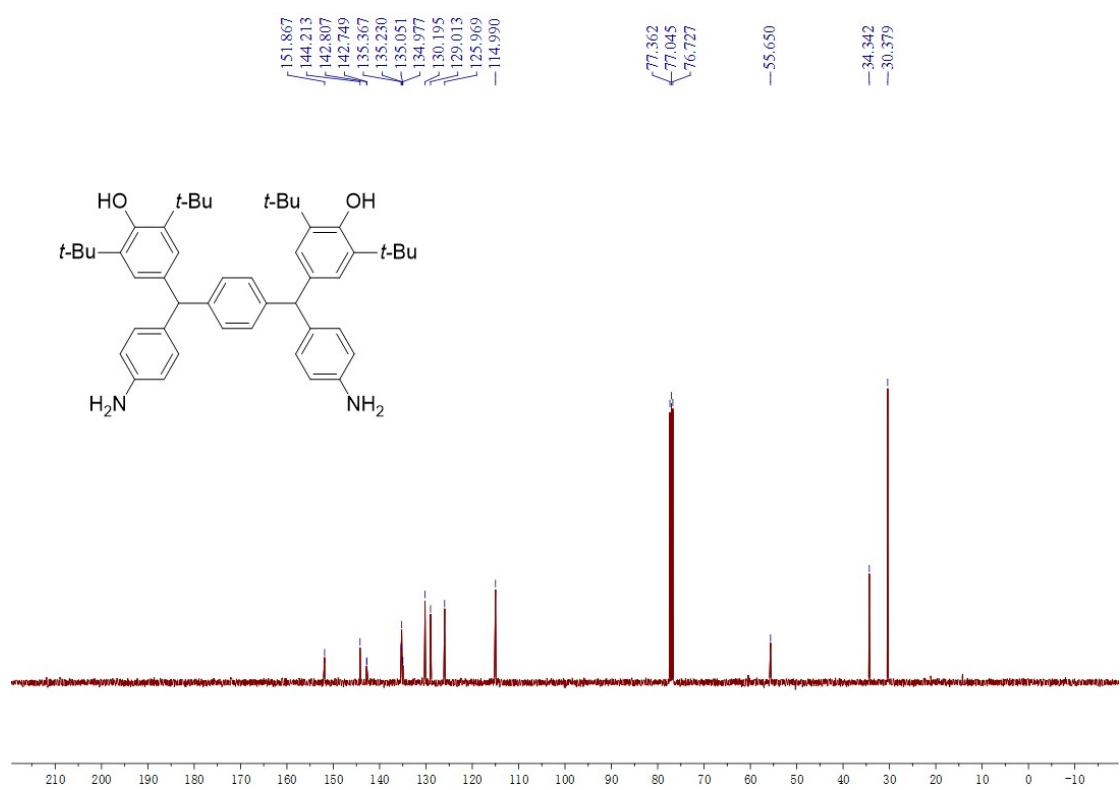
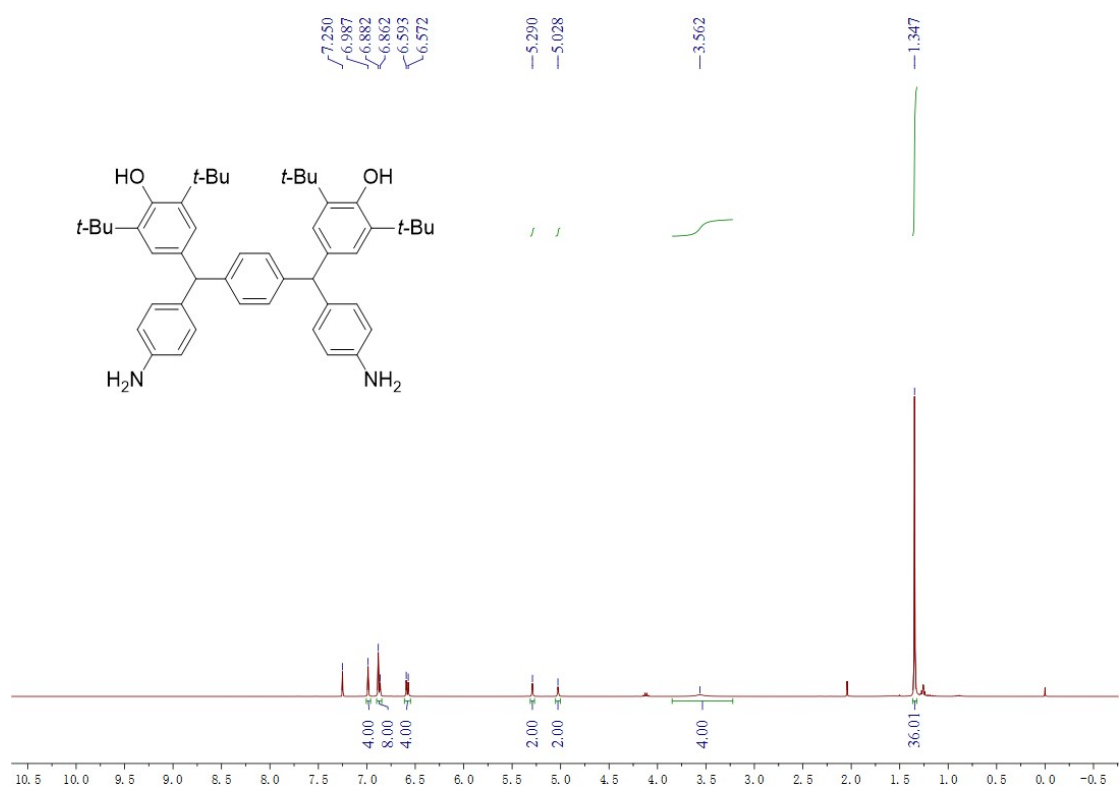
4-((4-Aminophenyl)(phenyl)methyl)-2,6-diisopropylphenol (4ac)



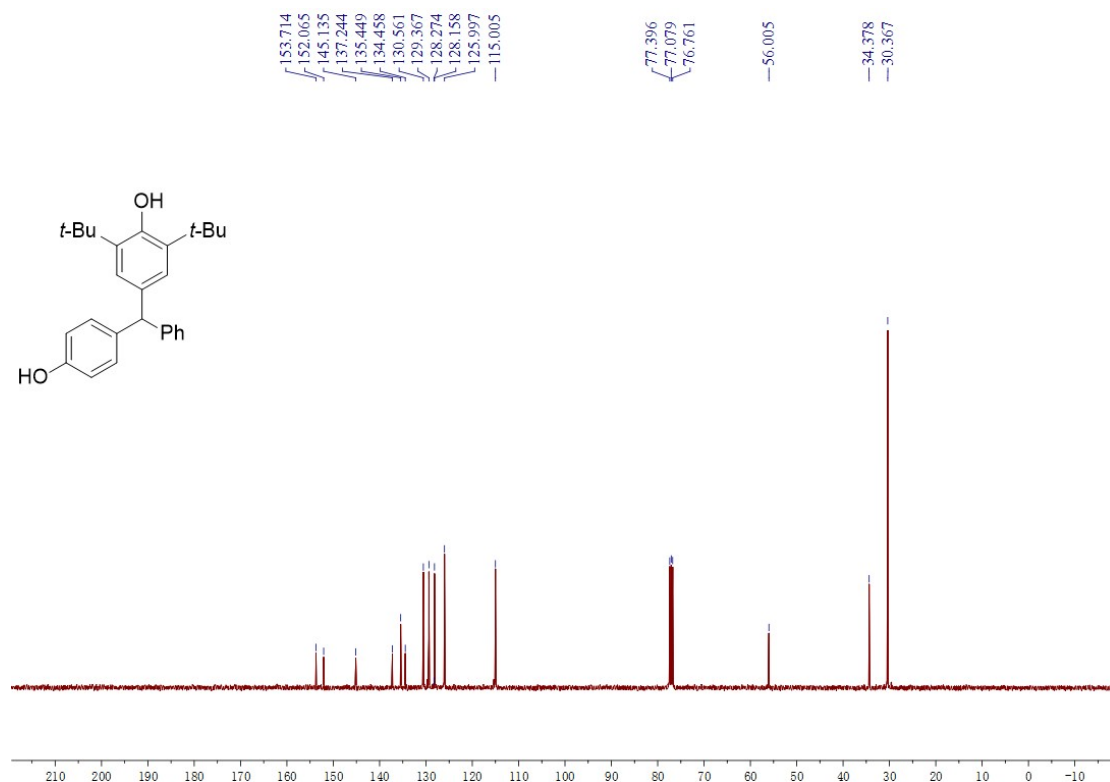
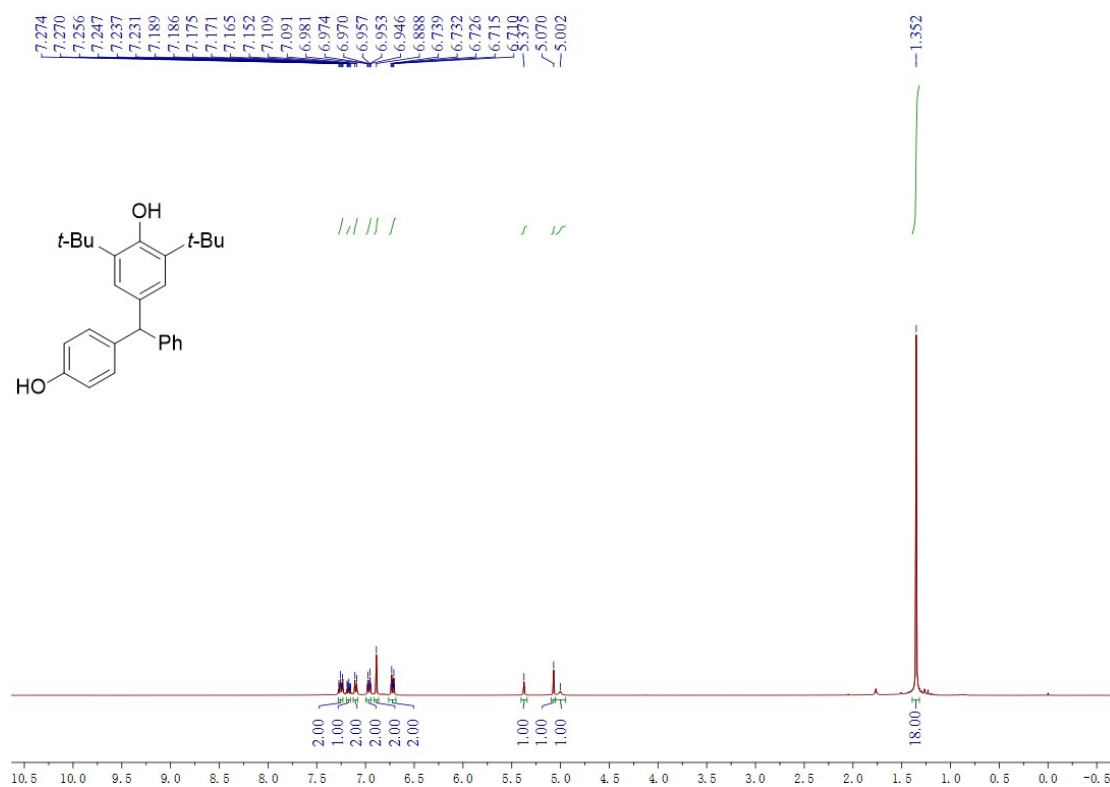
5'-((4-Aminophenyl)(phenyl)methyl)-[1,1':3',1''-terphenyl]-2'-ol (4ad)



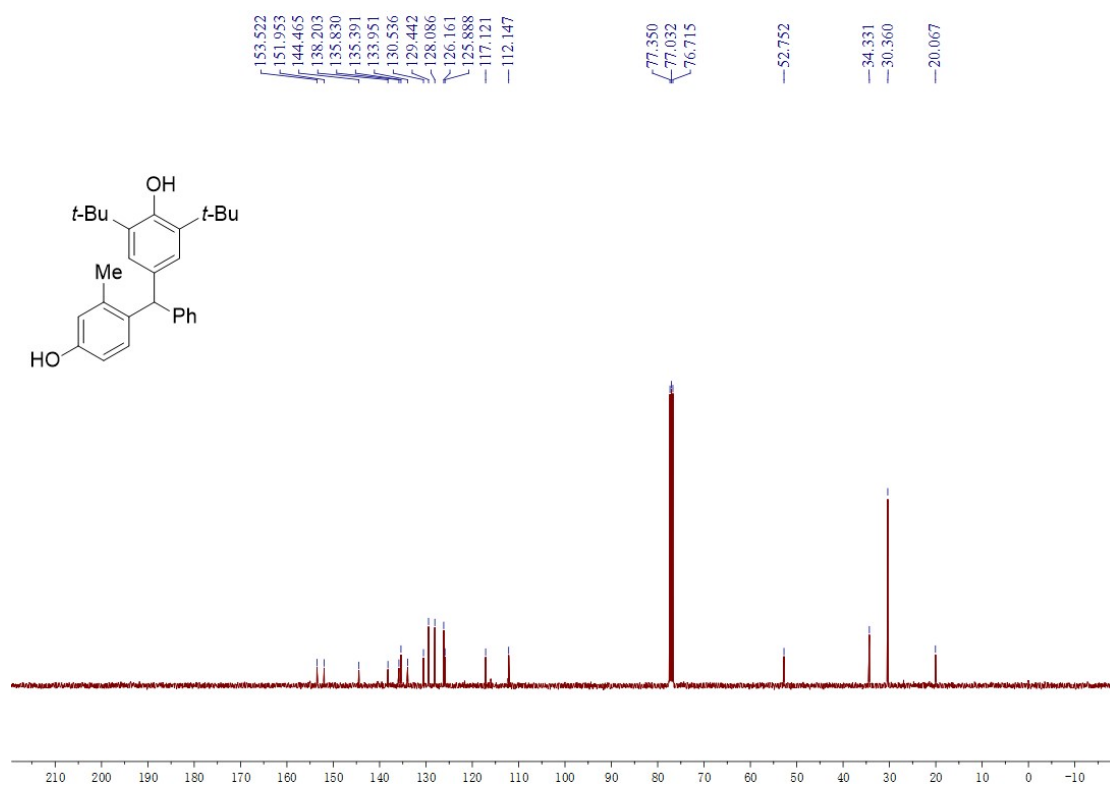
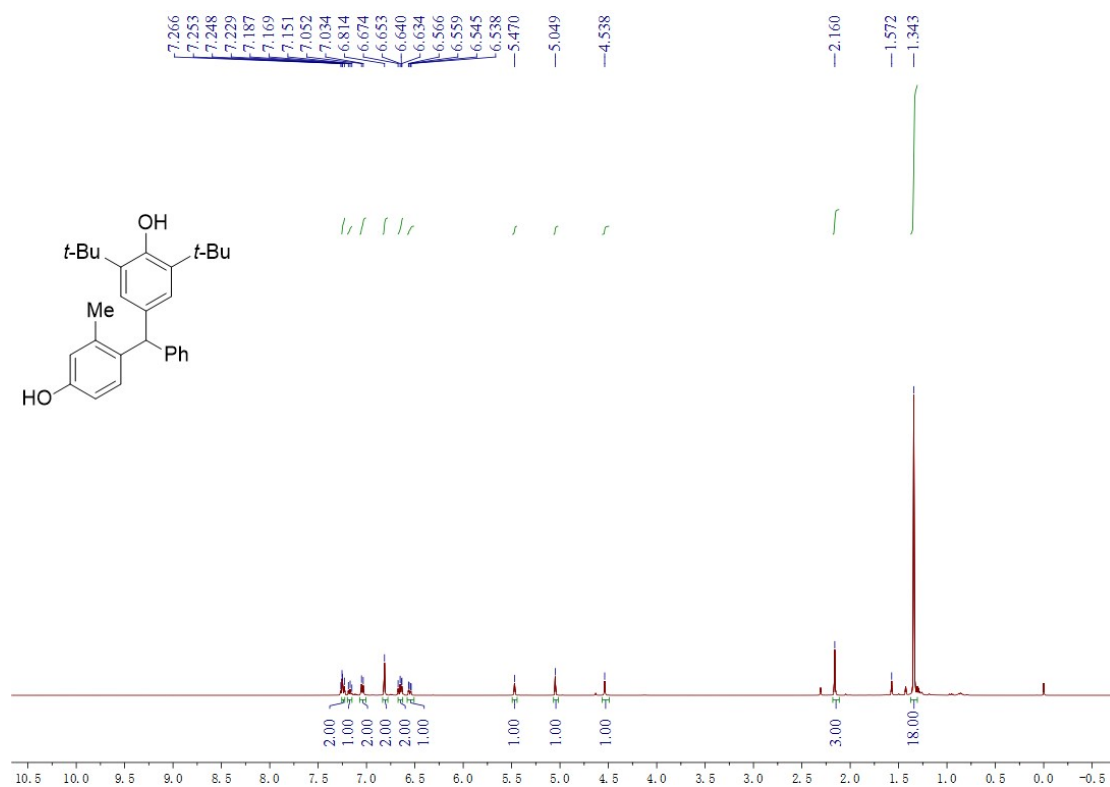
4,4'-(1,4-Phenylenebis((4-Aminophenyl)methylene))bis(2,6-di-tert-butylphenol) (4ae)



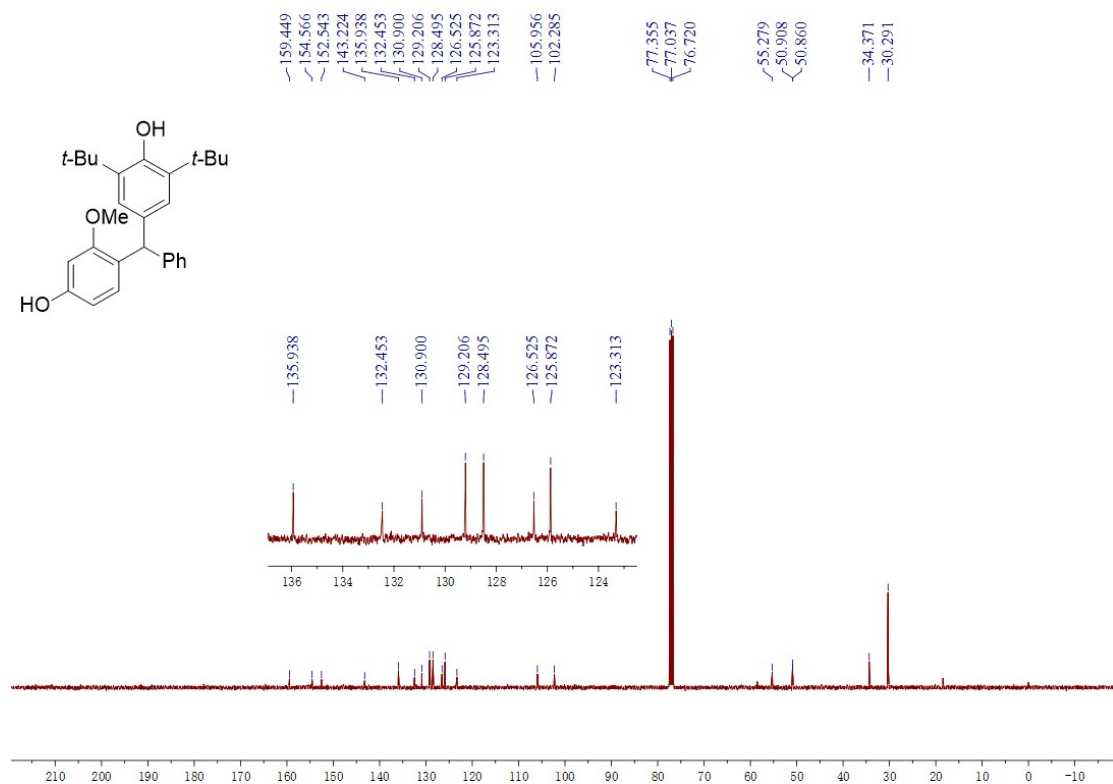
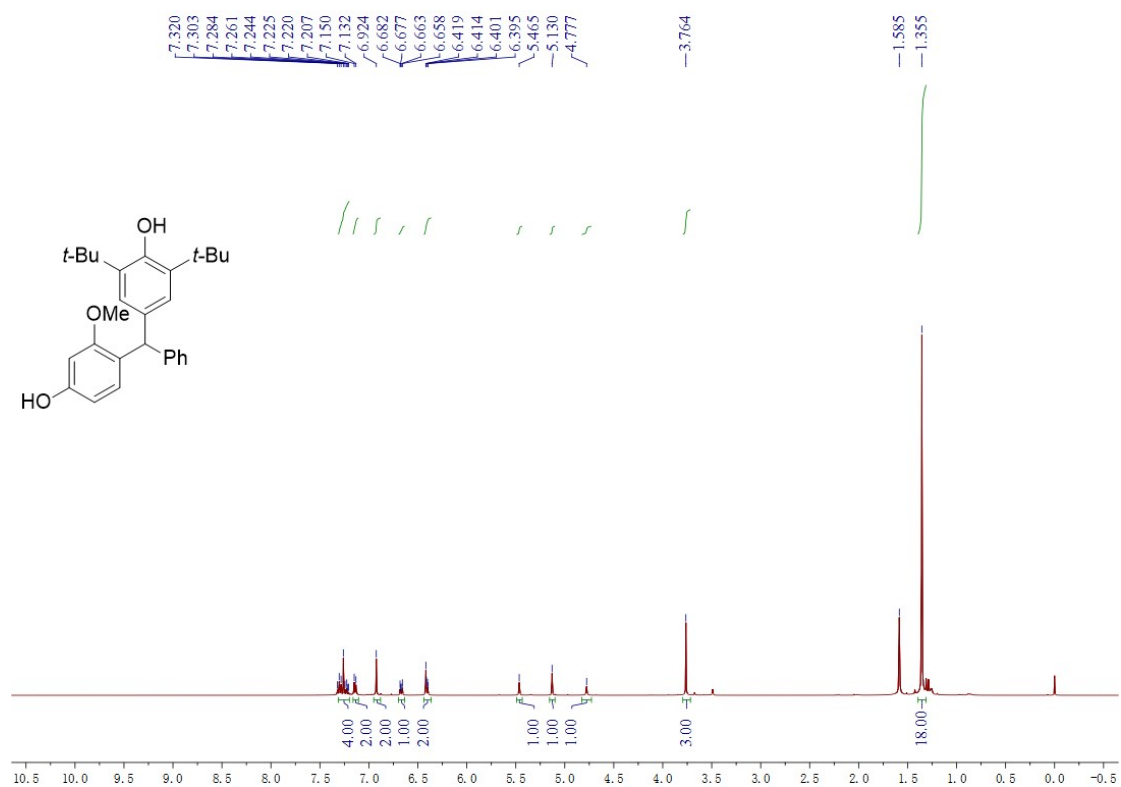
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(phenyl)methyl)phenol (6a)



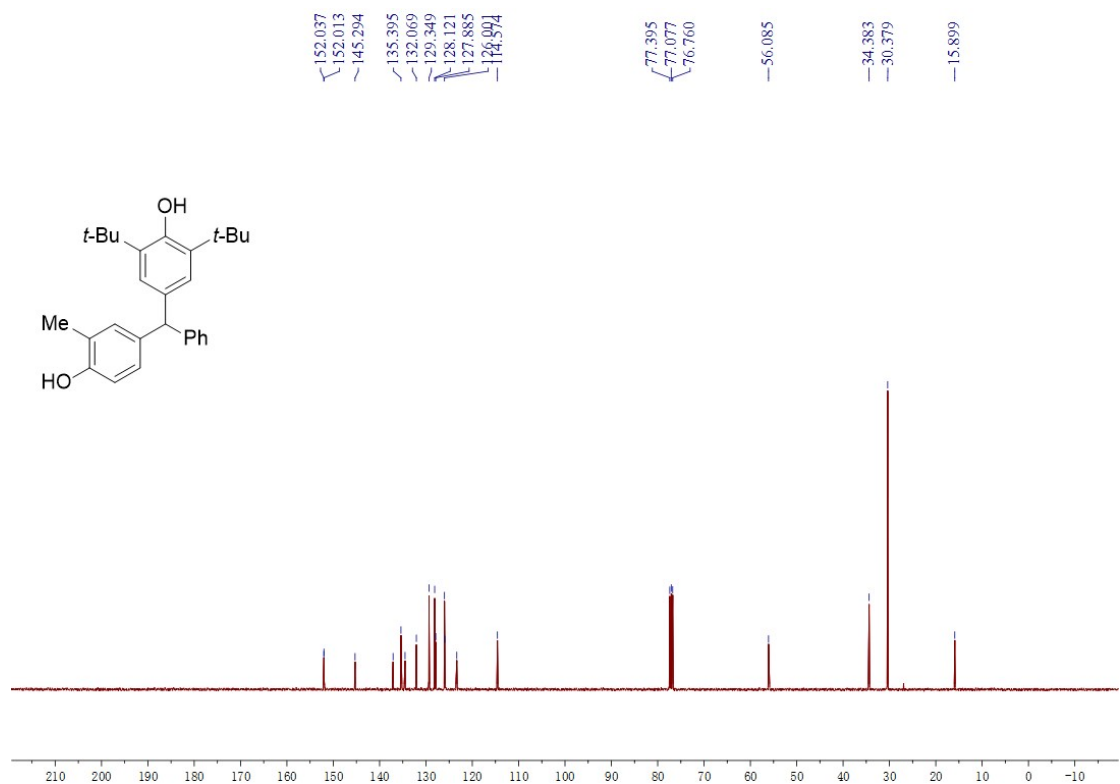
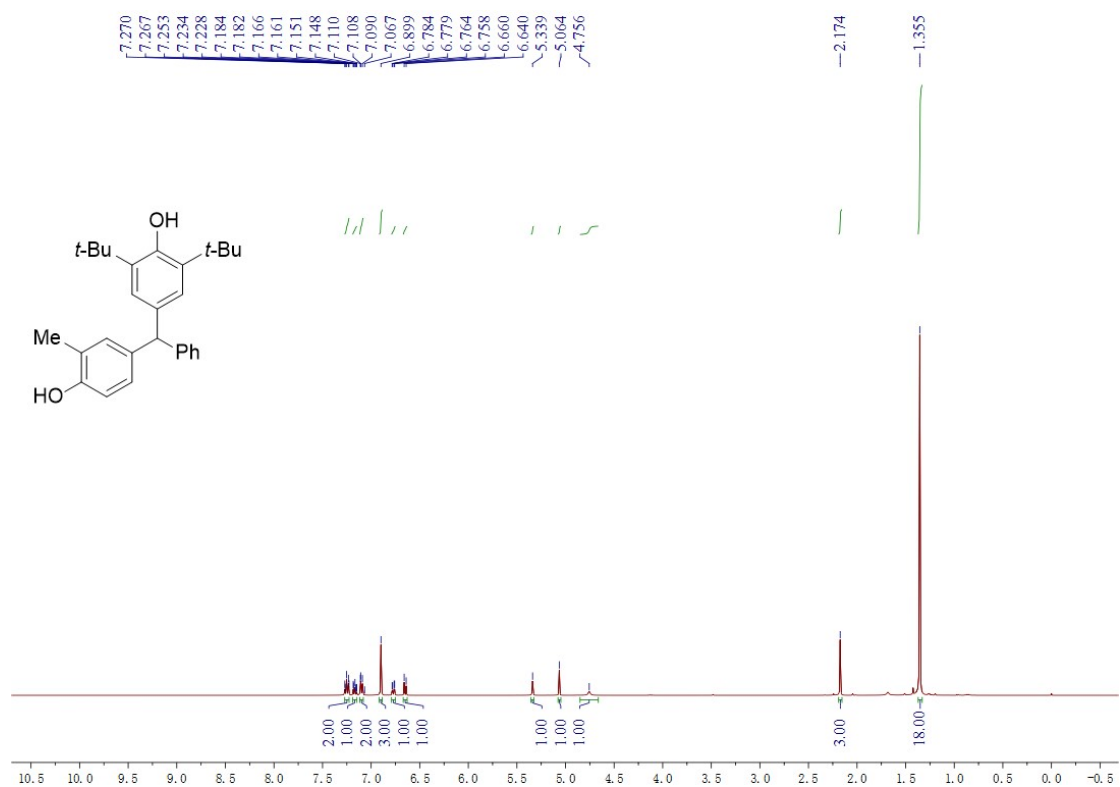
2,6-Di-*tert*-butyl-4-((4-hydroxy-2-methylphenyl)(phenyl)methyl)phenol (6b)



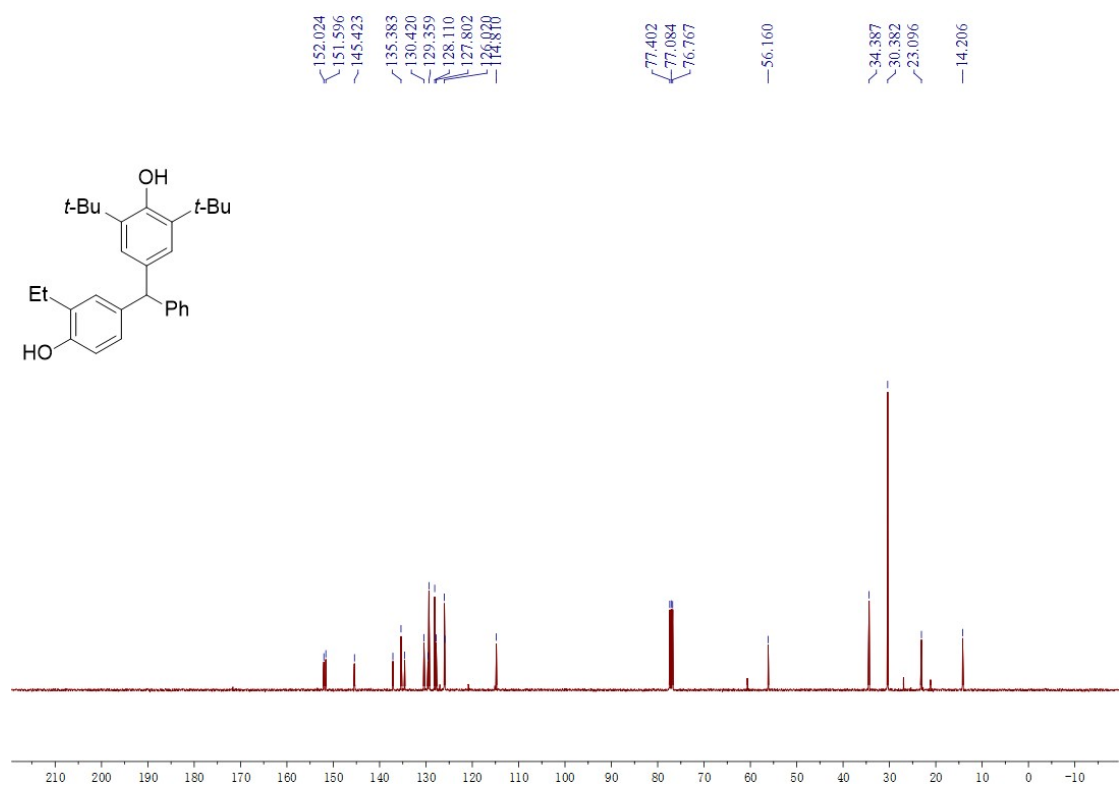
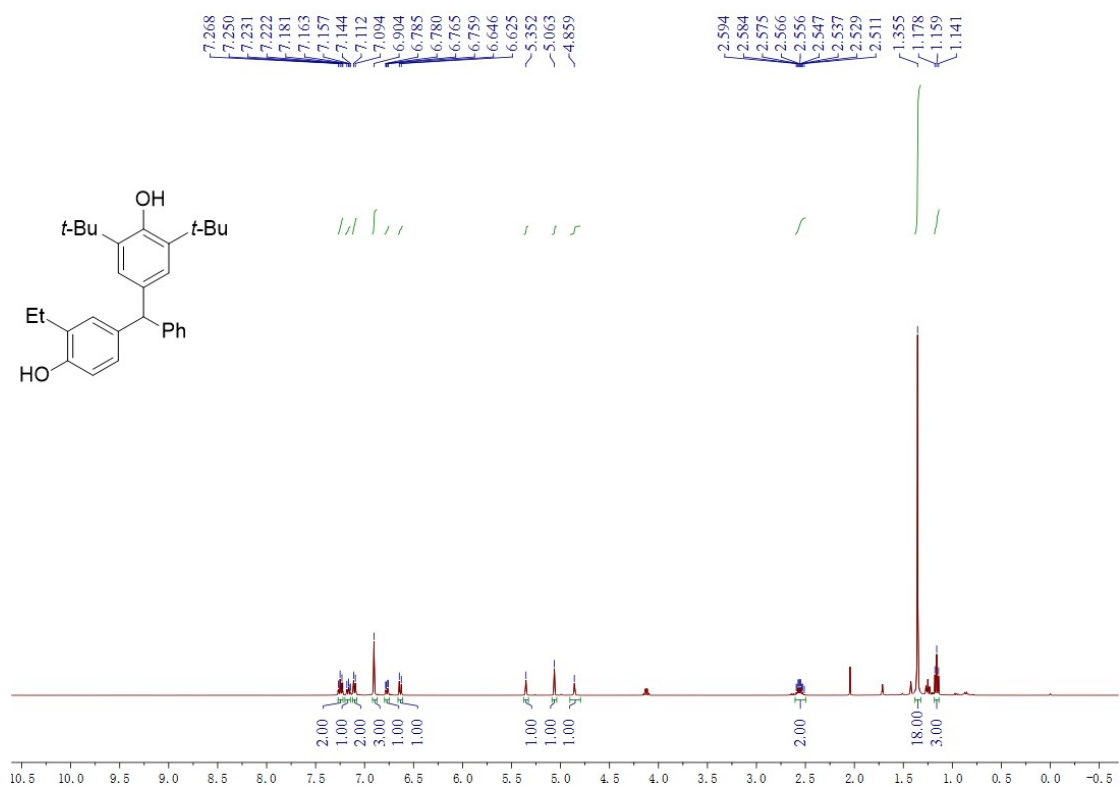
2,6-Di-tert-butyl-4-((4-hydroxy-2-methoxyphenyl)(phenyl)methyl)phenol (6c)



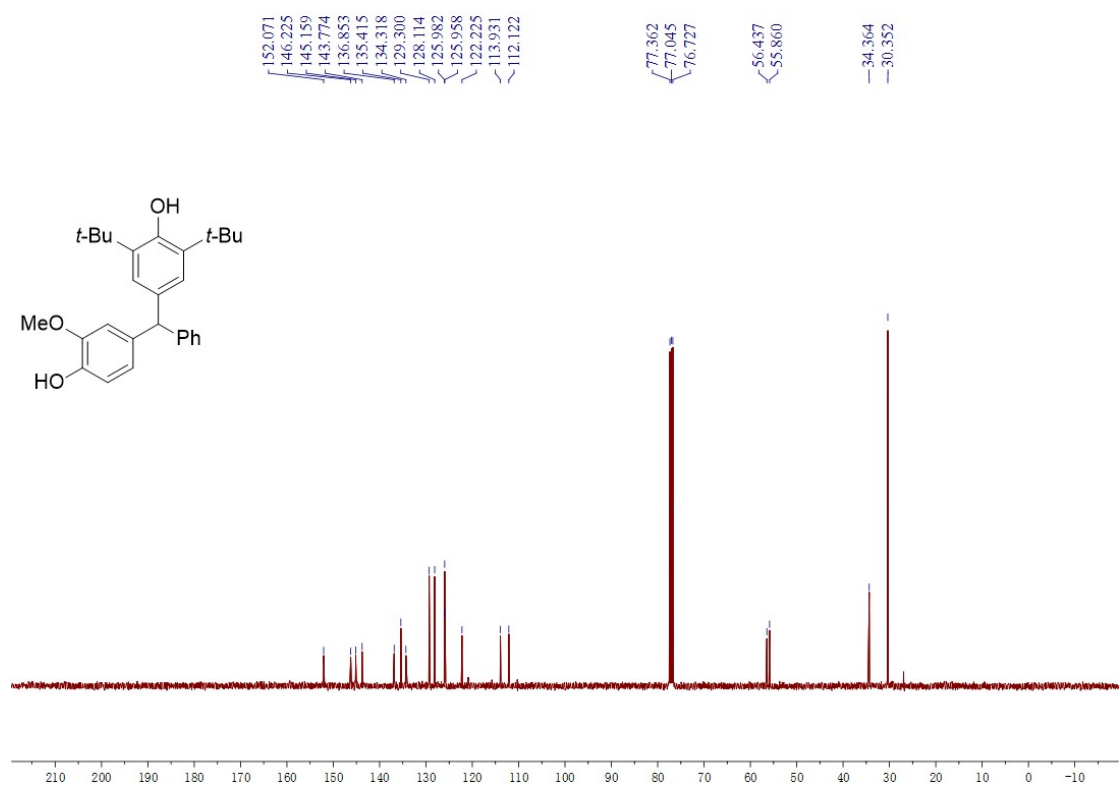
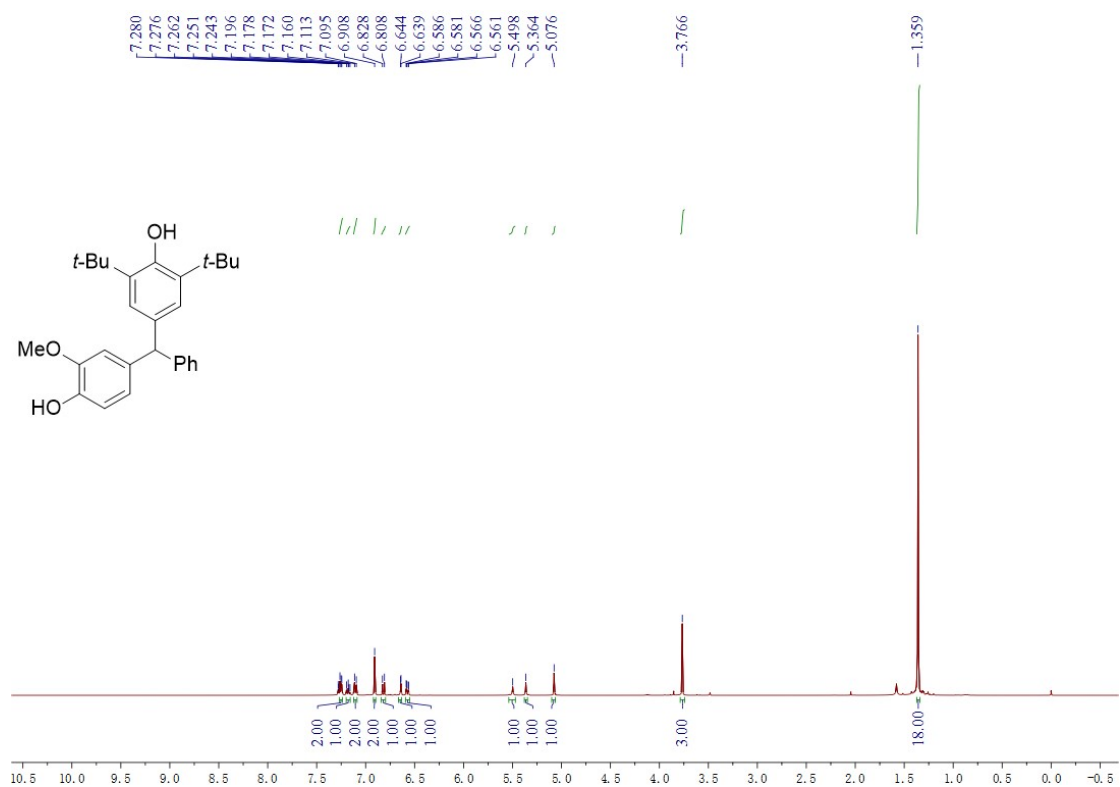
2,6-Di-tert-butyl-4-((4-hydroxy-3-methylphenyl)(phenyl)methyl)phenol (6d)



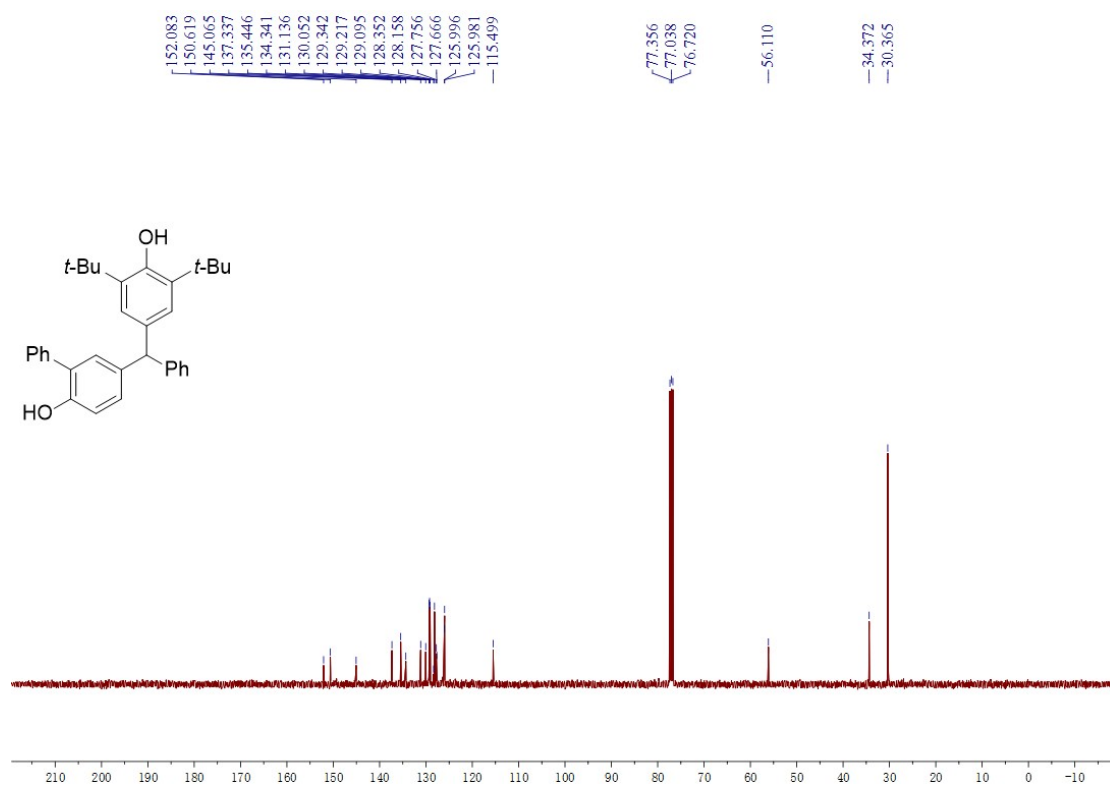
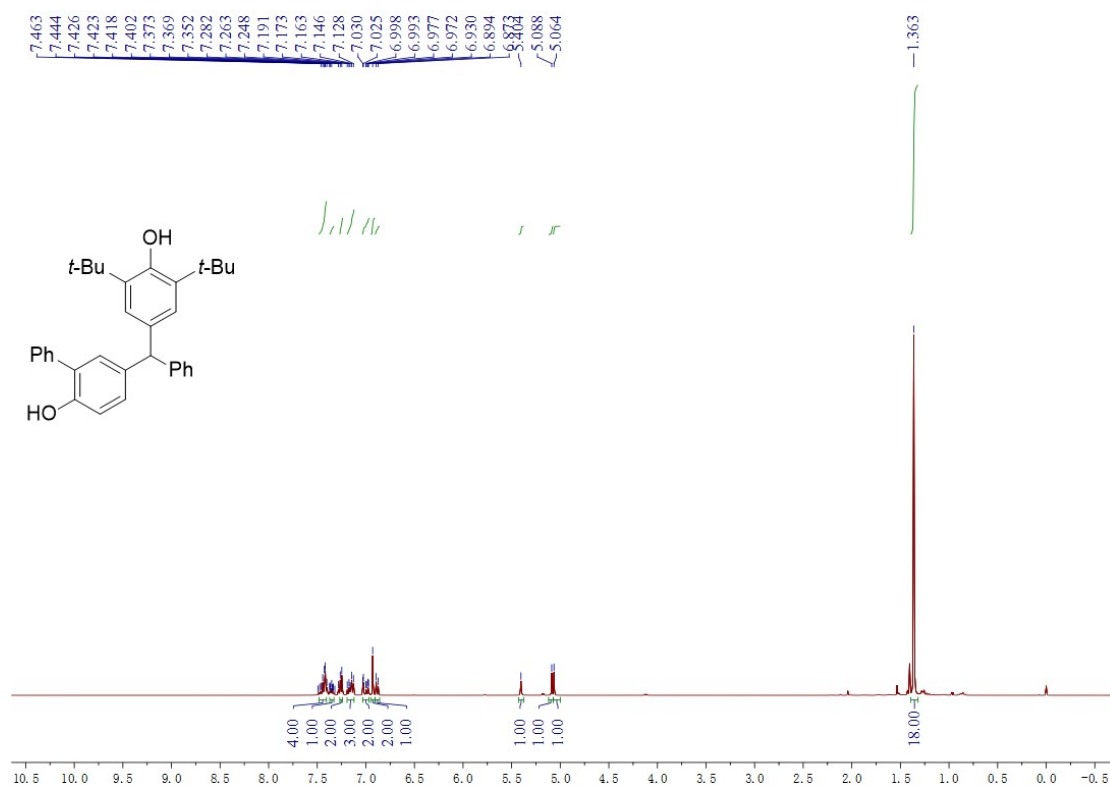
2,6-Di-tert-butyl-4-((3-ethyl-4-hydroxyphenyl)(phenyl)methyl)phenol (6e)



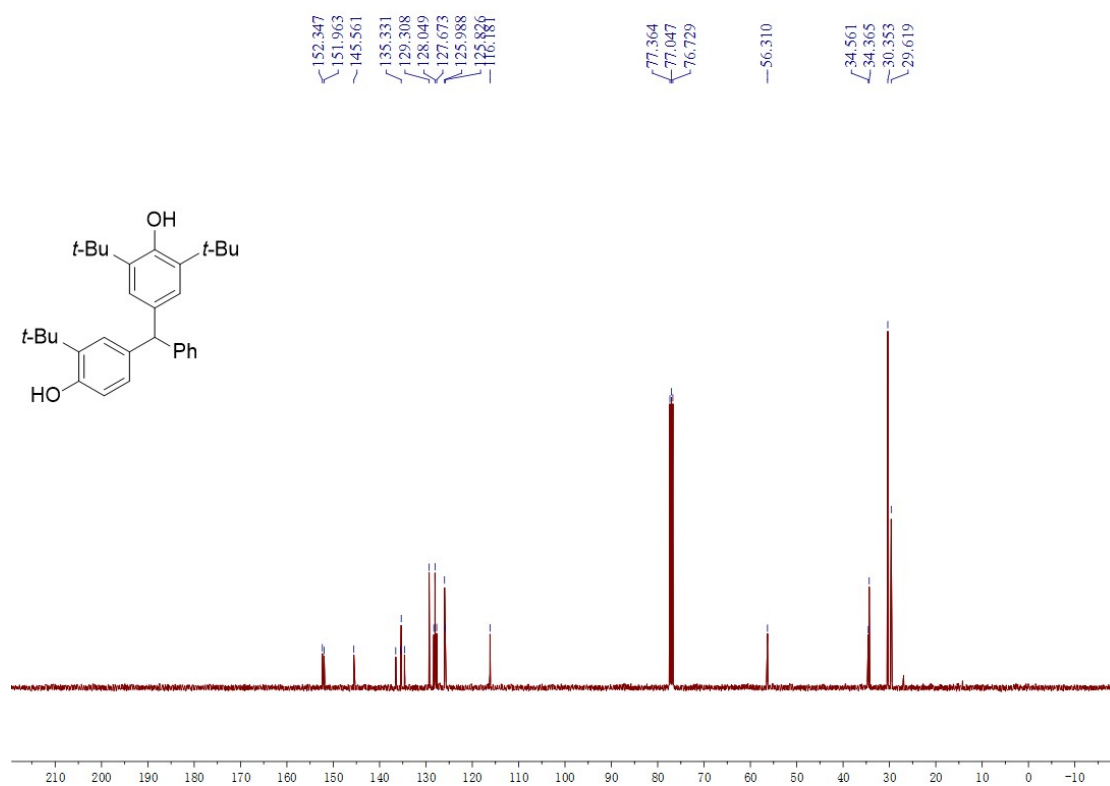
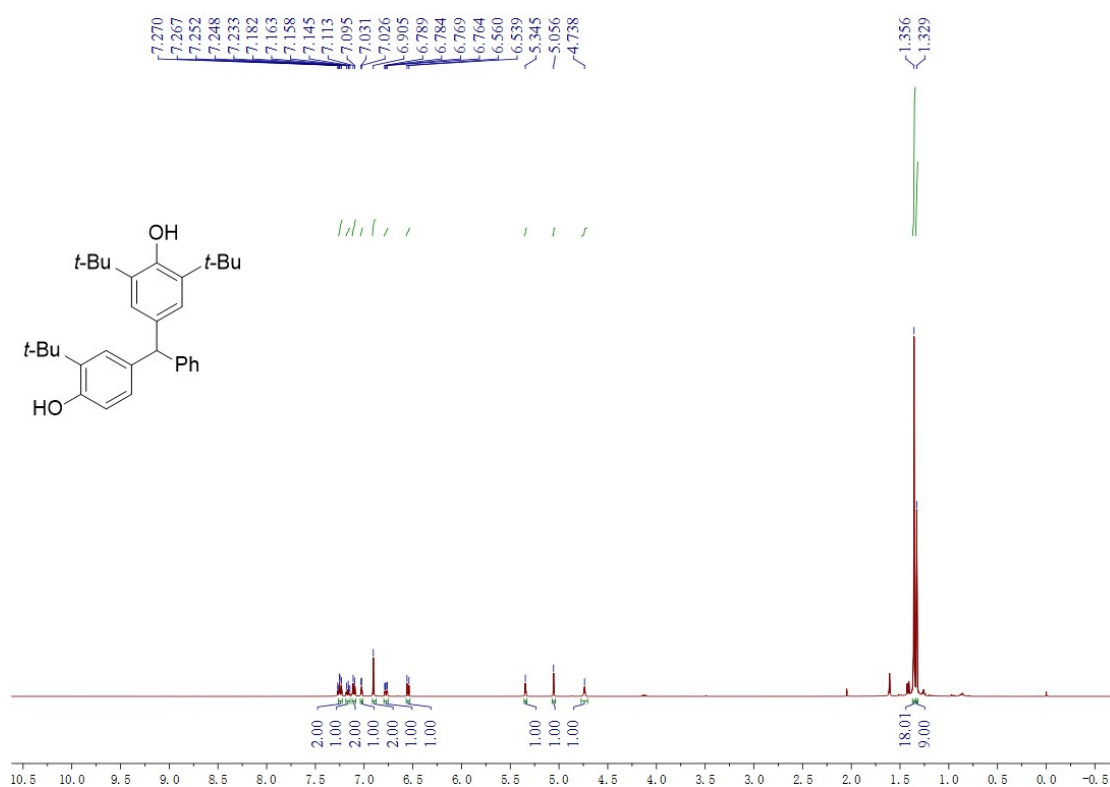
2,6-Di-tert-butyl-4-((4-hydroxy-3-methoxyphenyl)(phenyl)methyl)phenol (6f)



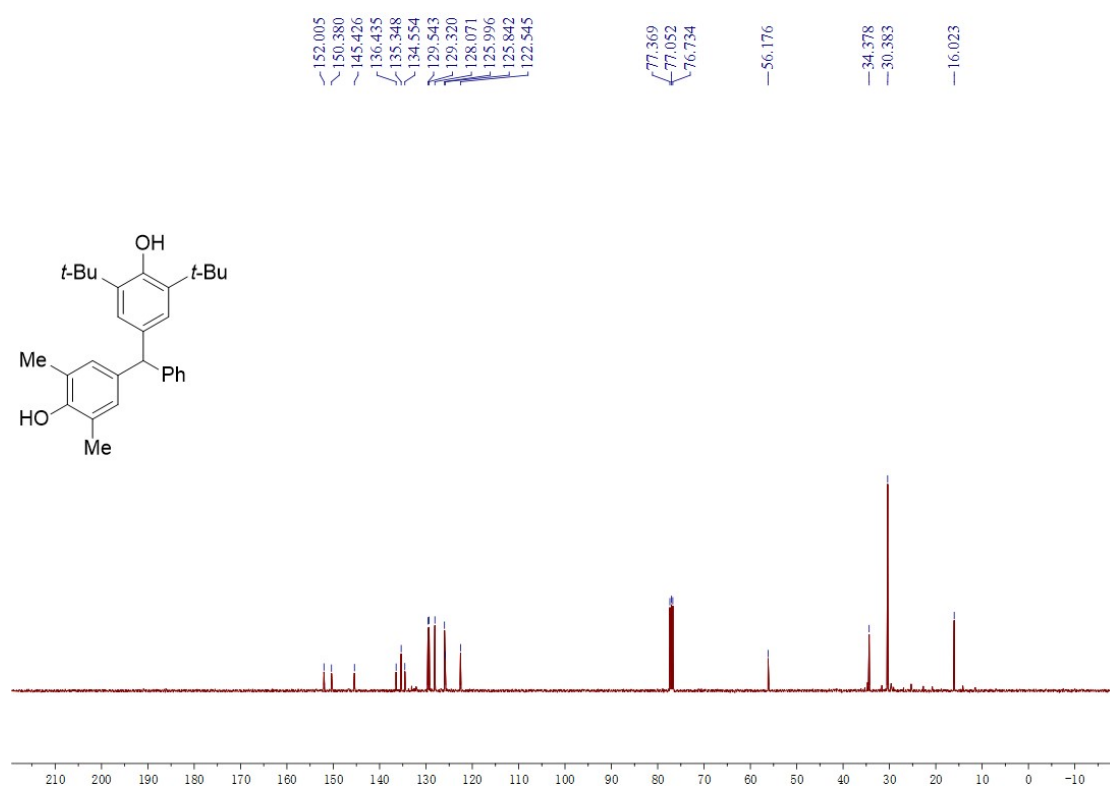
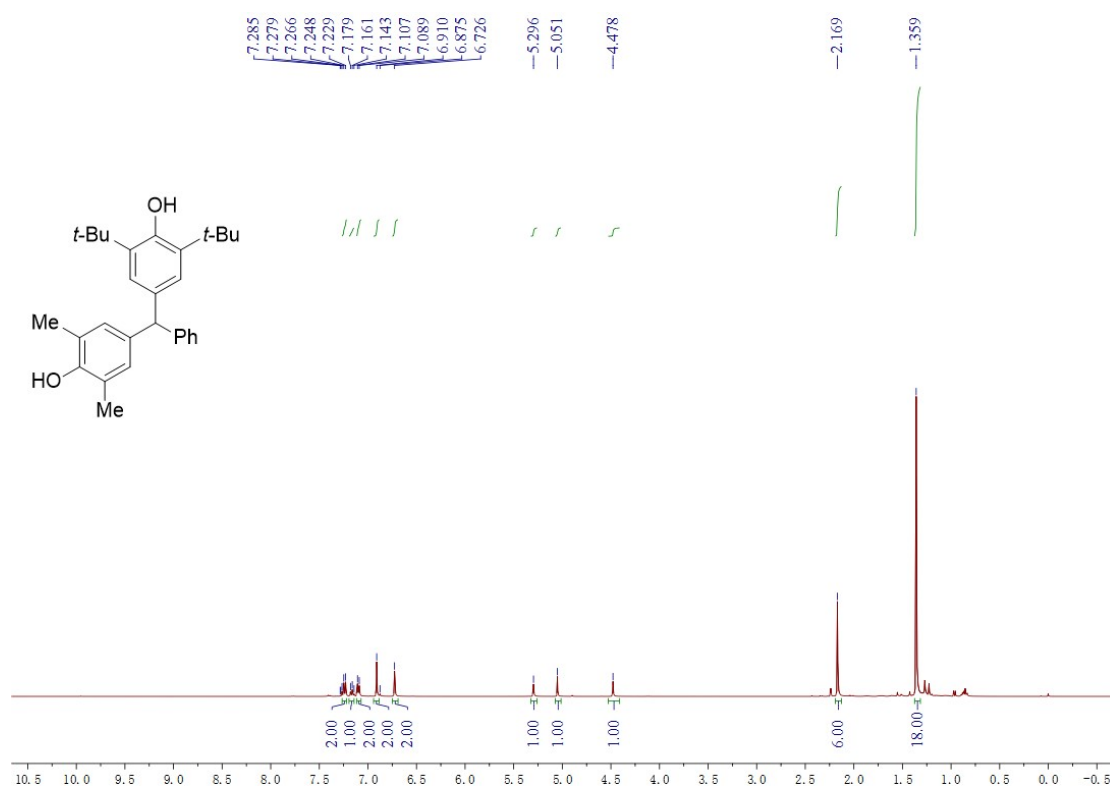
5-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-[1,1'-biphenyl]-2-ol (6g)



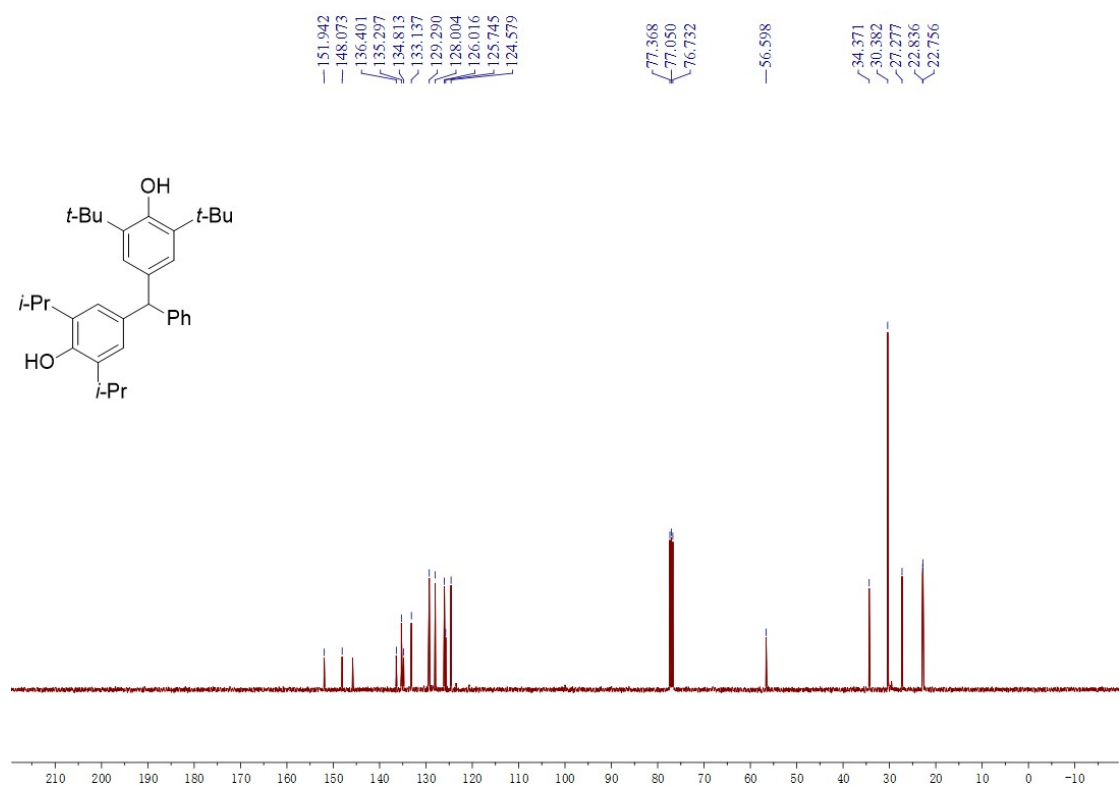
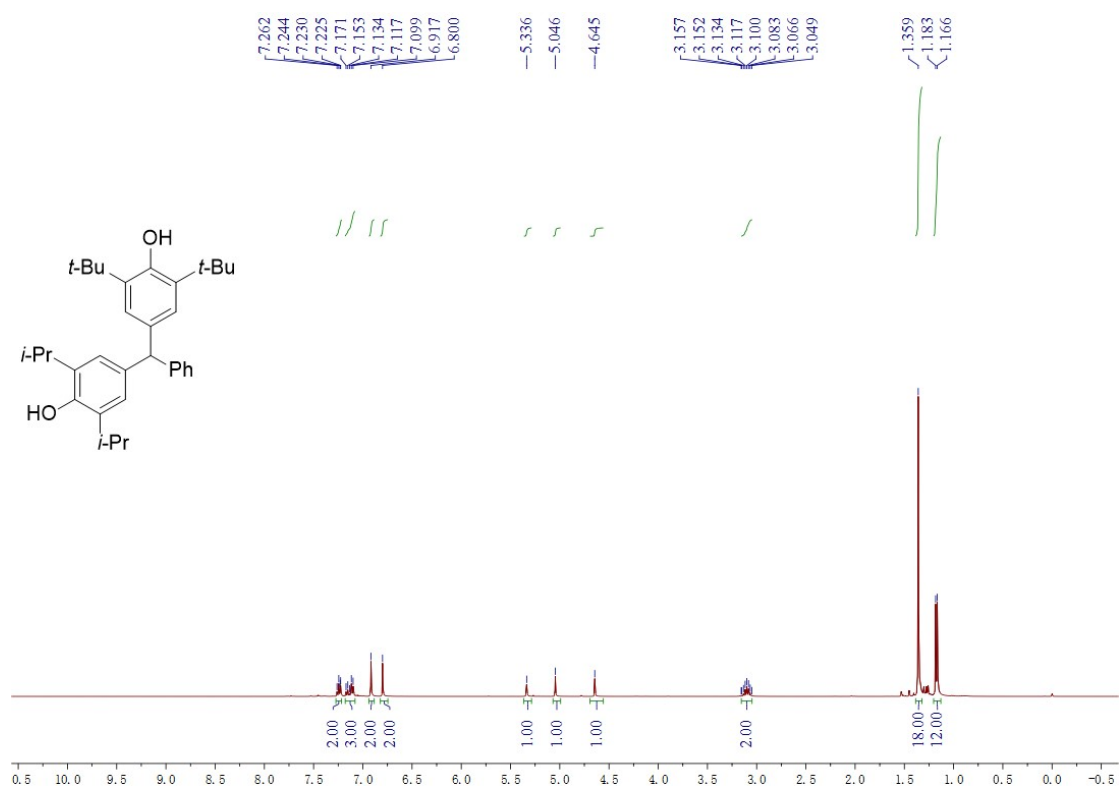
2,6-Di-tert-butyl-4-((4-hydroxy-3-isopropylphenyl)(phenyl)methyl)phenol (6h)



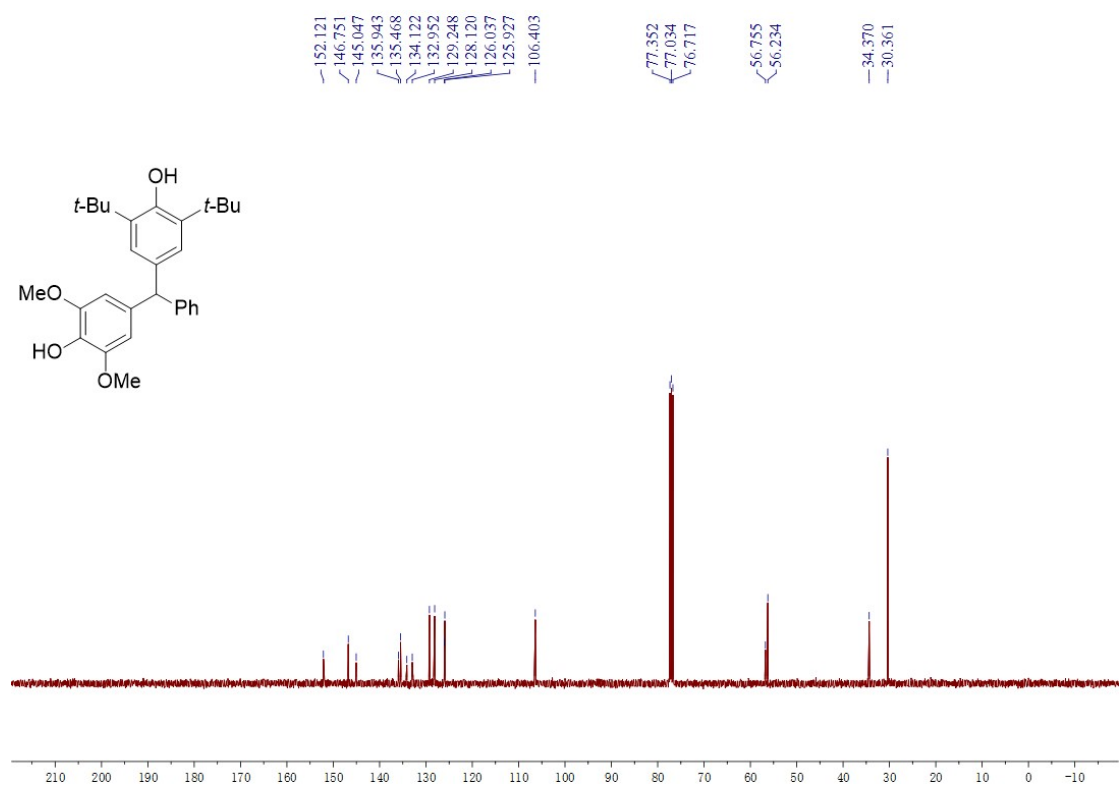
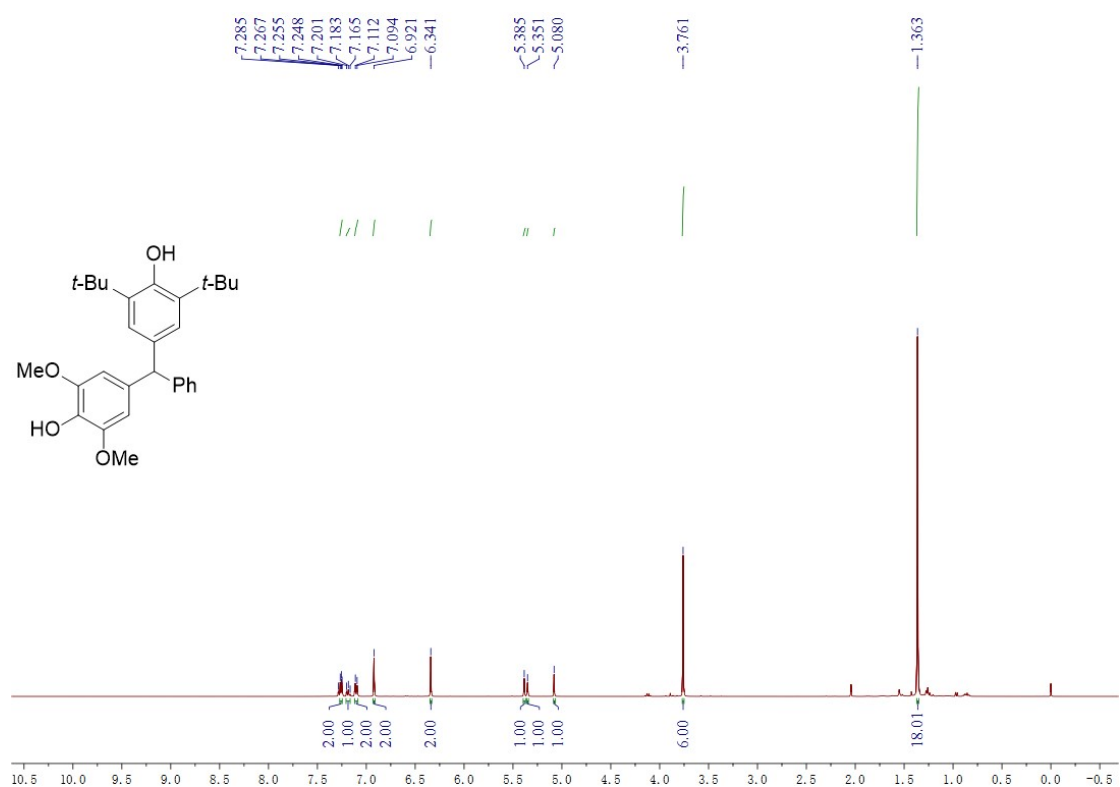
2,6-Di-tert-butyl-4-((4-hydroxy-3,5-dimethylphenyl)(phenyl)methyl)phenol (6j)



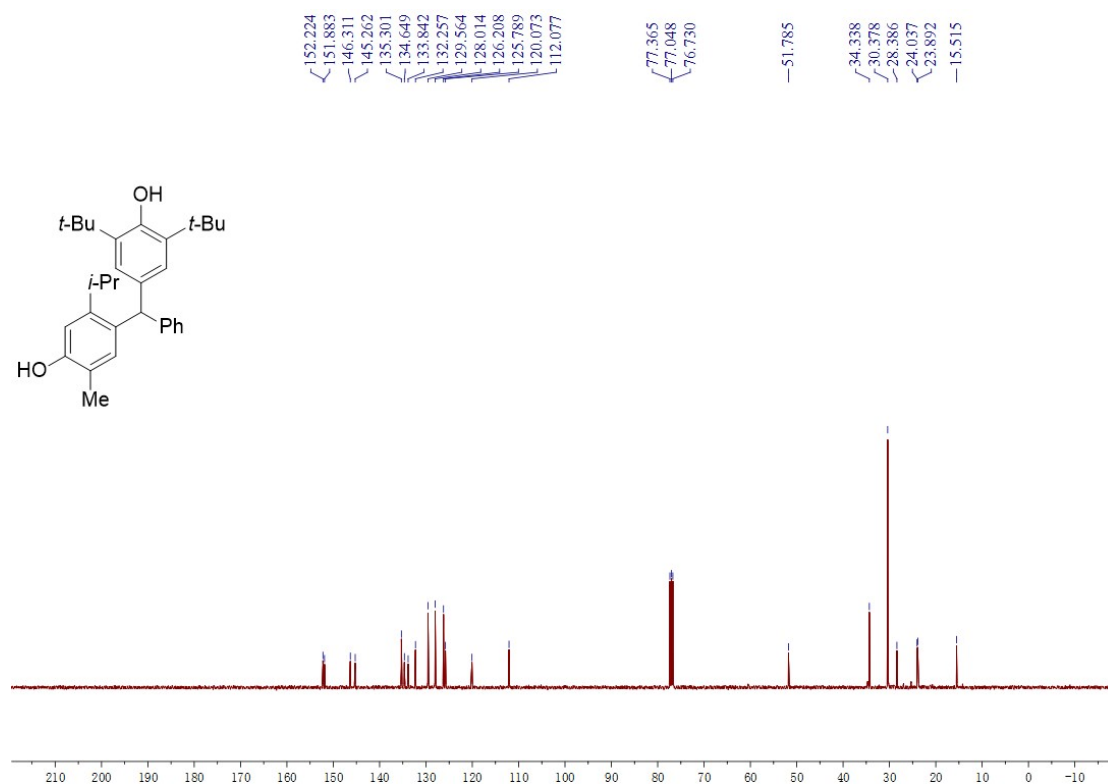
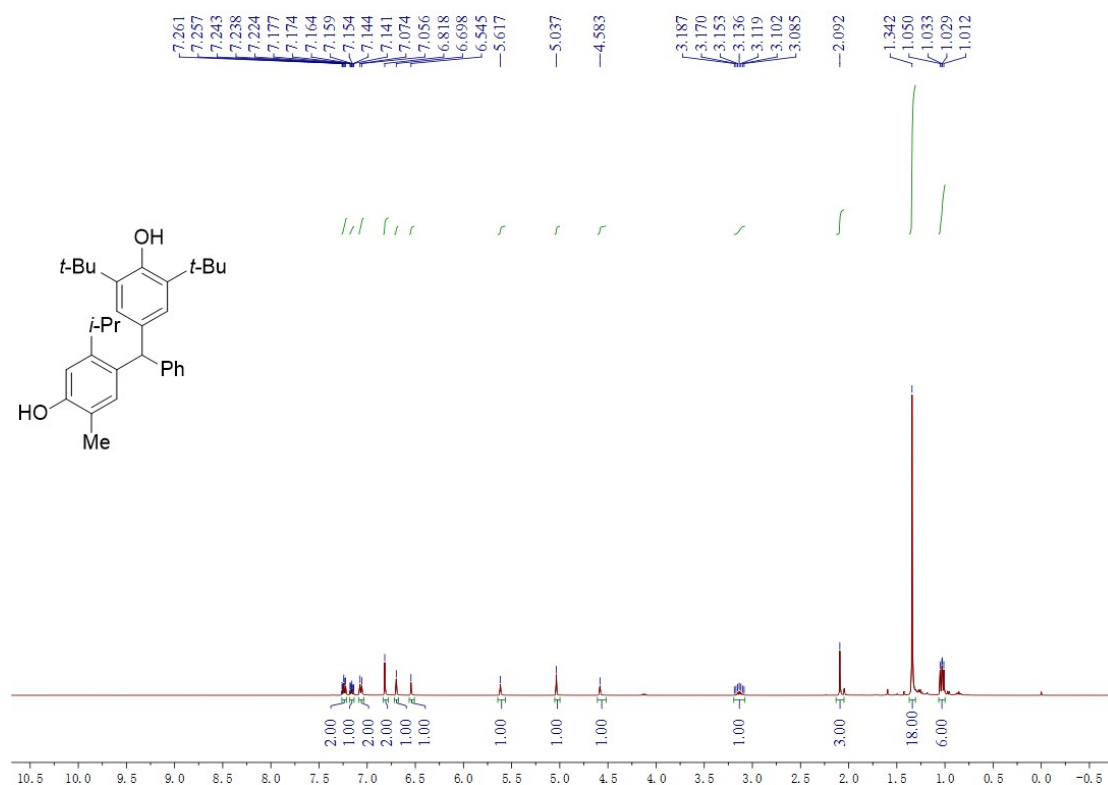
2,6-Di-tert-butyl-4-((4-hydroxy-3,5-diisopropylphenyl)(phenyl)methyl)phenol (6k)



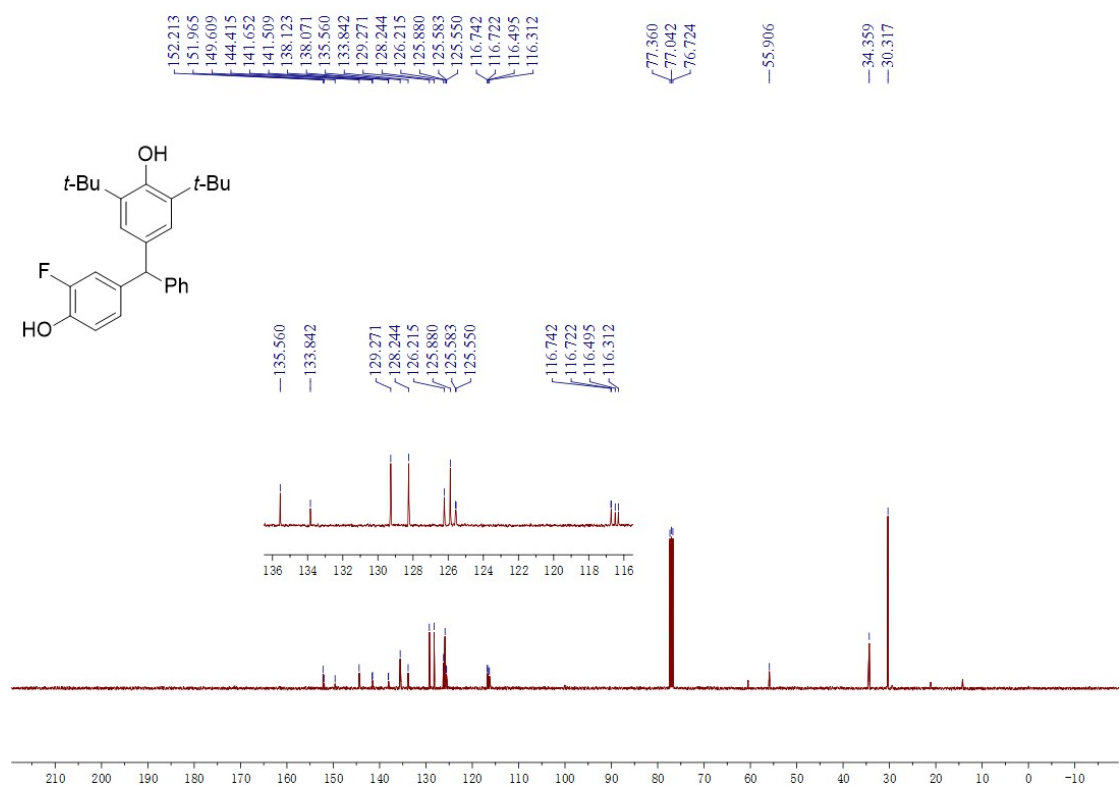
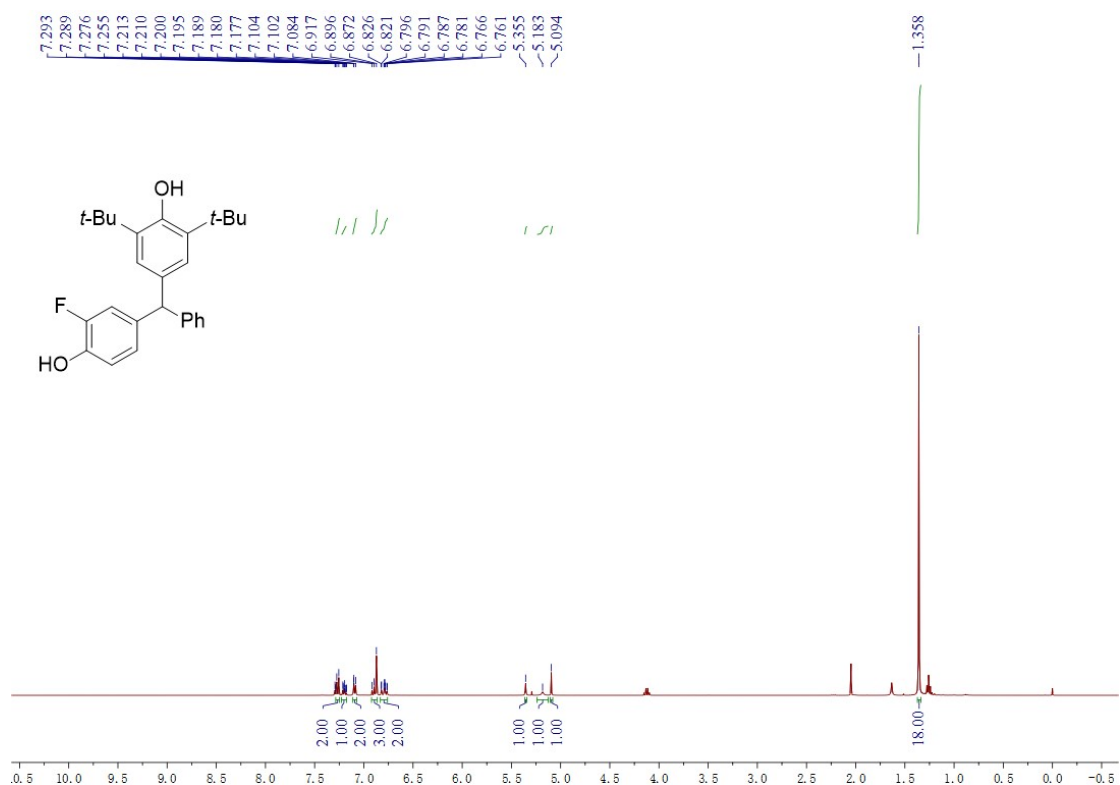
2,6-Di-tert-butyl-4-((4-hydroxy-3,5-dimethoxyphenyl)(phenyl)methyl)phenol (6l)

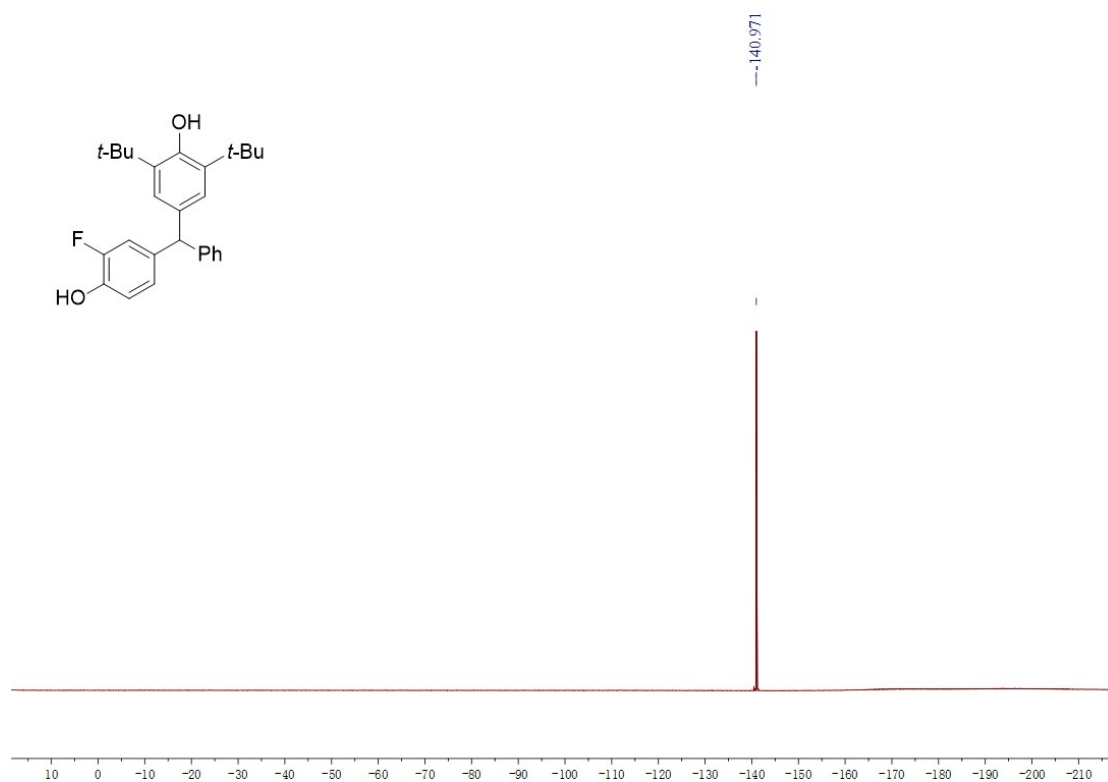


**4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-5-isopropyl-2-methylphenol
(6m)**

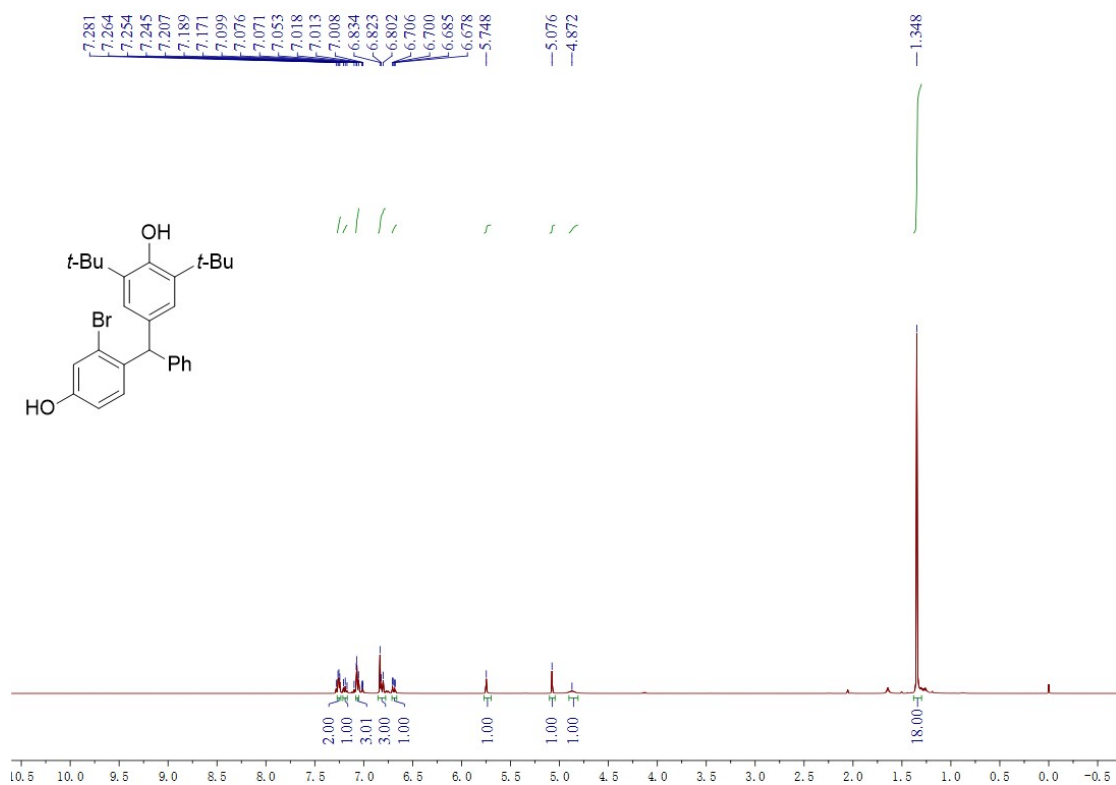


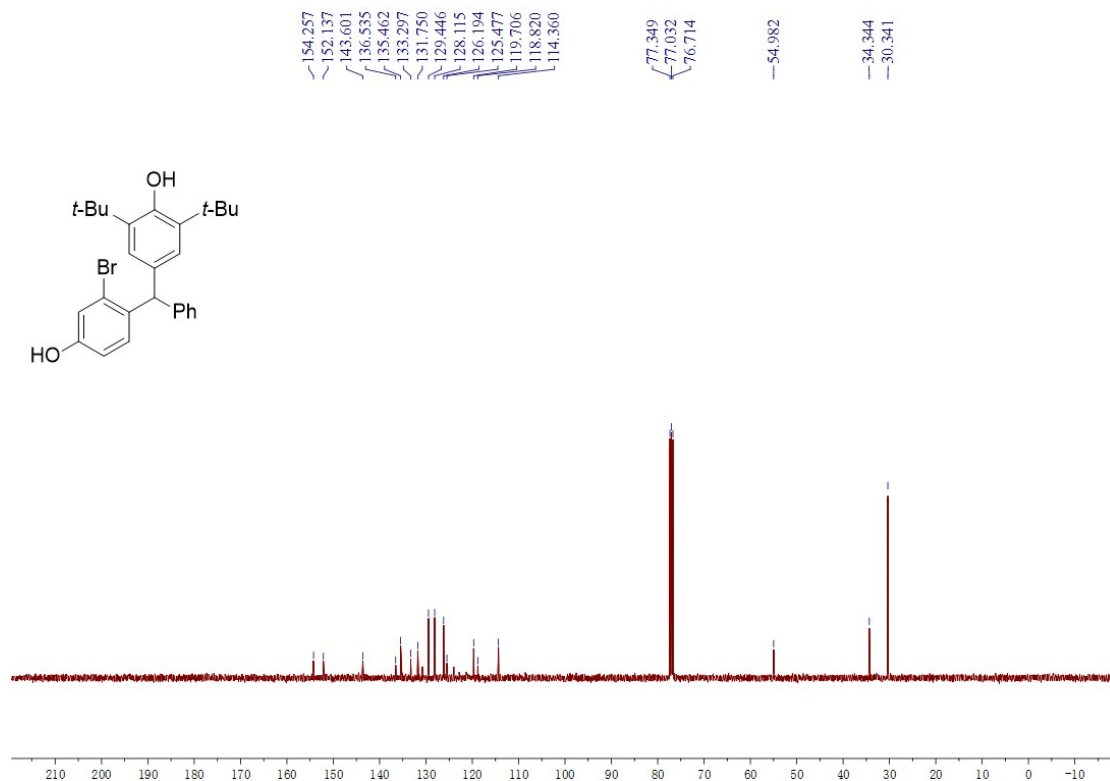
2,6-Di-*tert*-butyl-4-((3-fluoro-4-hydroxyphenyl)(phenyl)methyl)phenol (6n)



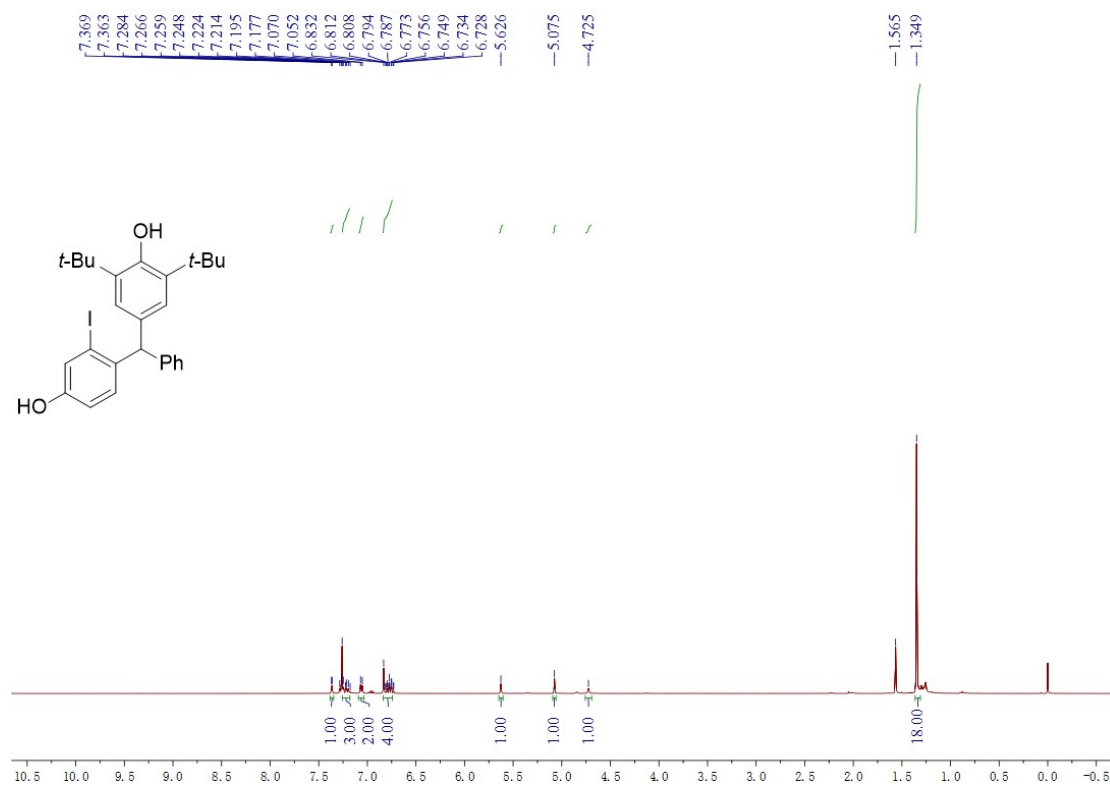


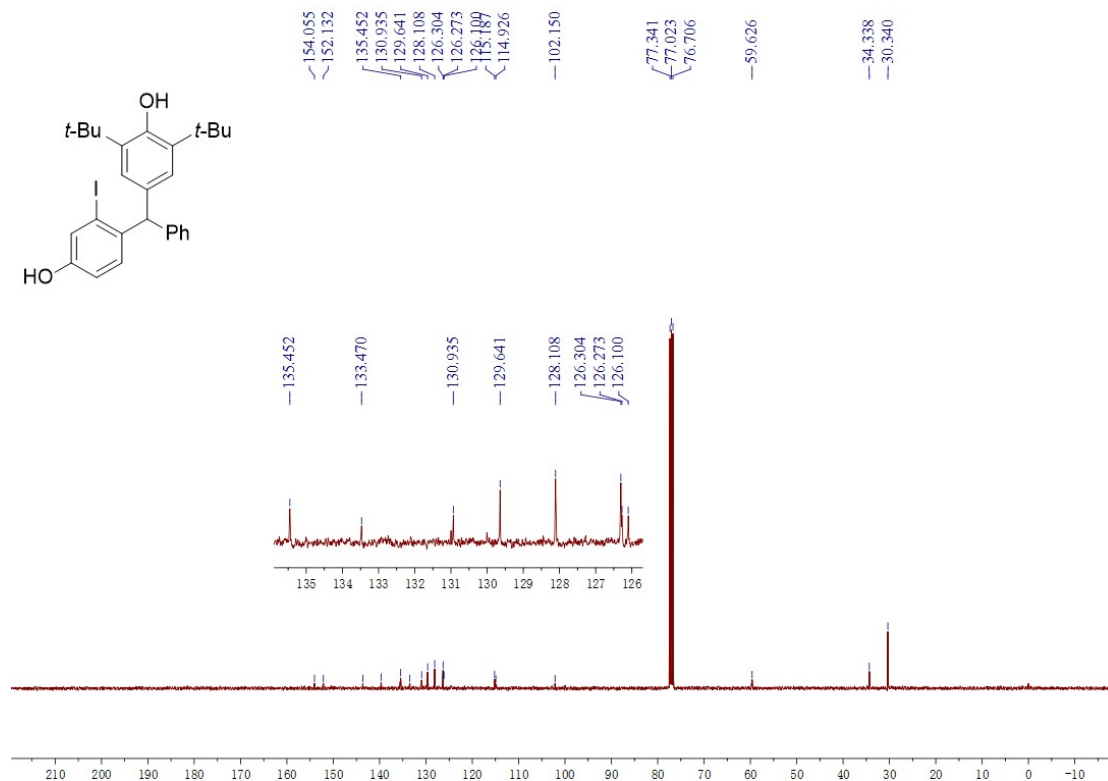
4-((2-Bromo-4-hydroxyphenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (60)



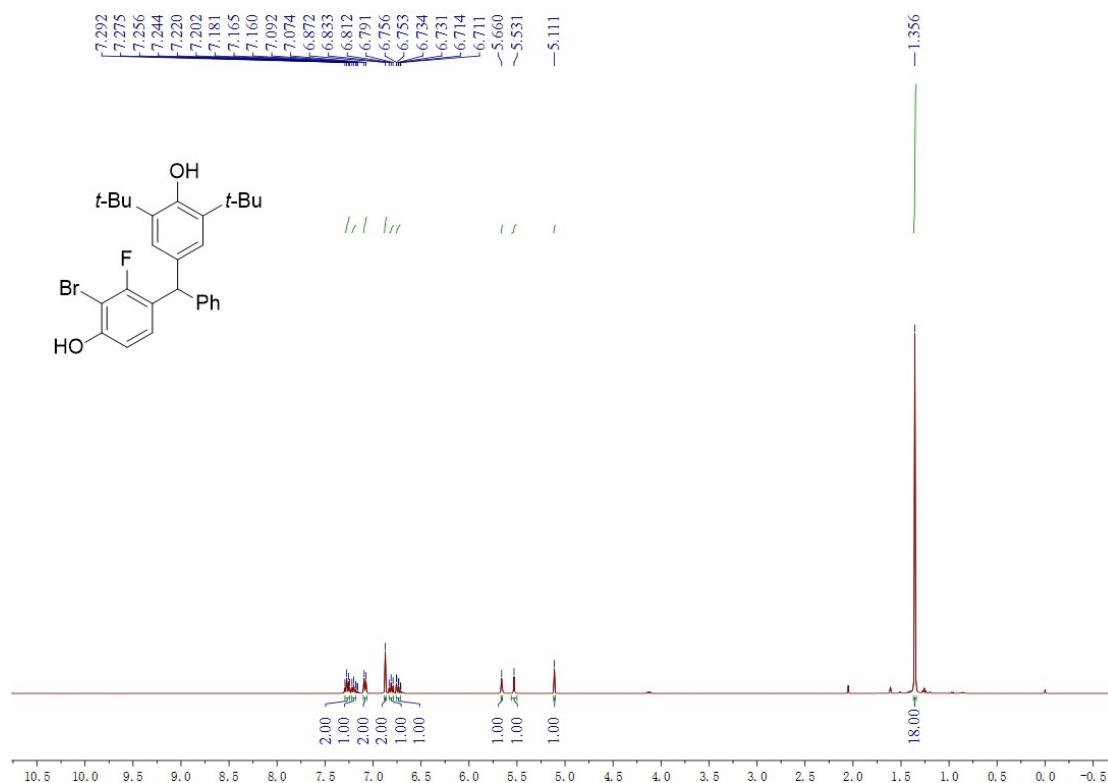


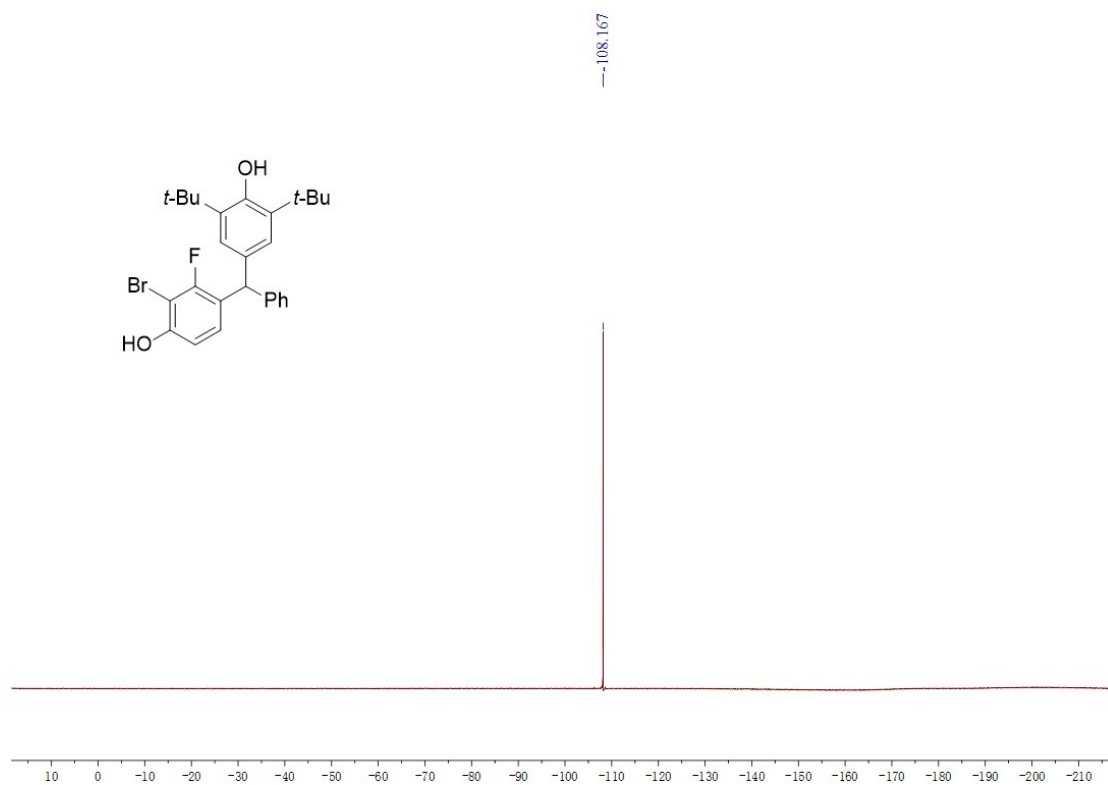
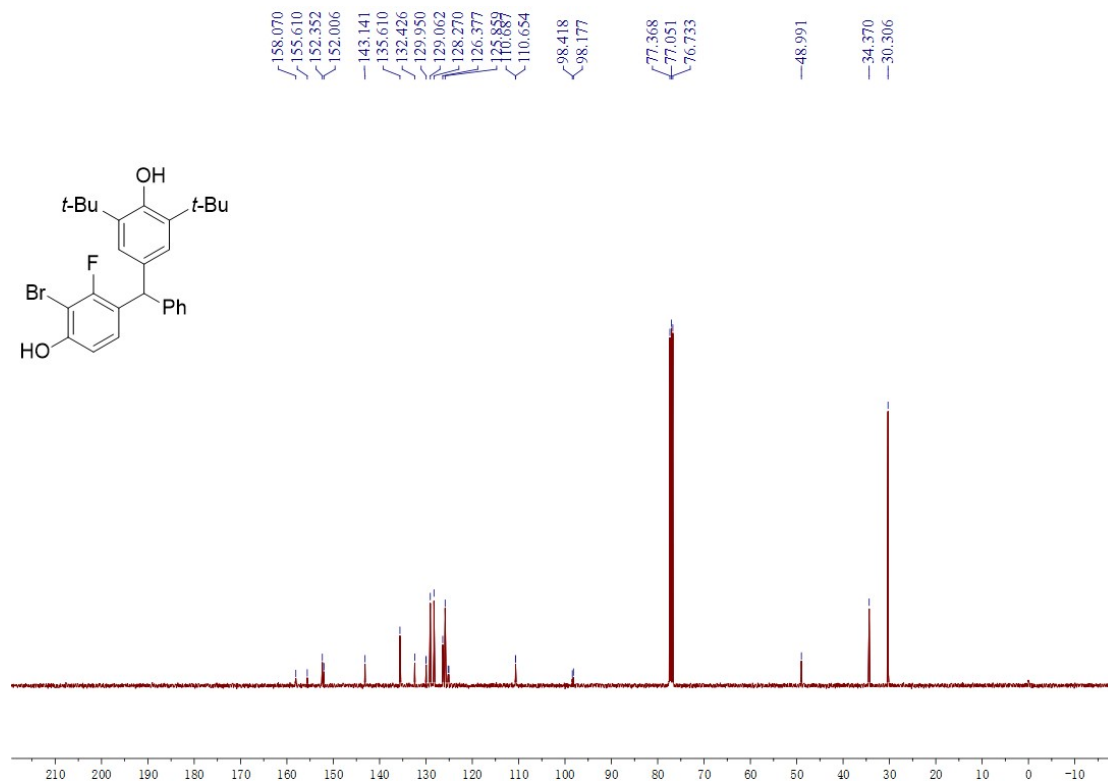
2,6-Di-tert-butyl-4-((4-hydroxy-2-iodophenyl)(phenyl)methyl)phenol (6p)



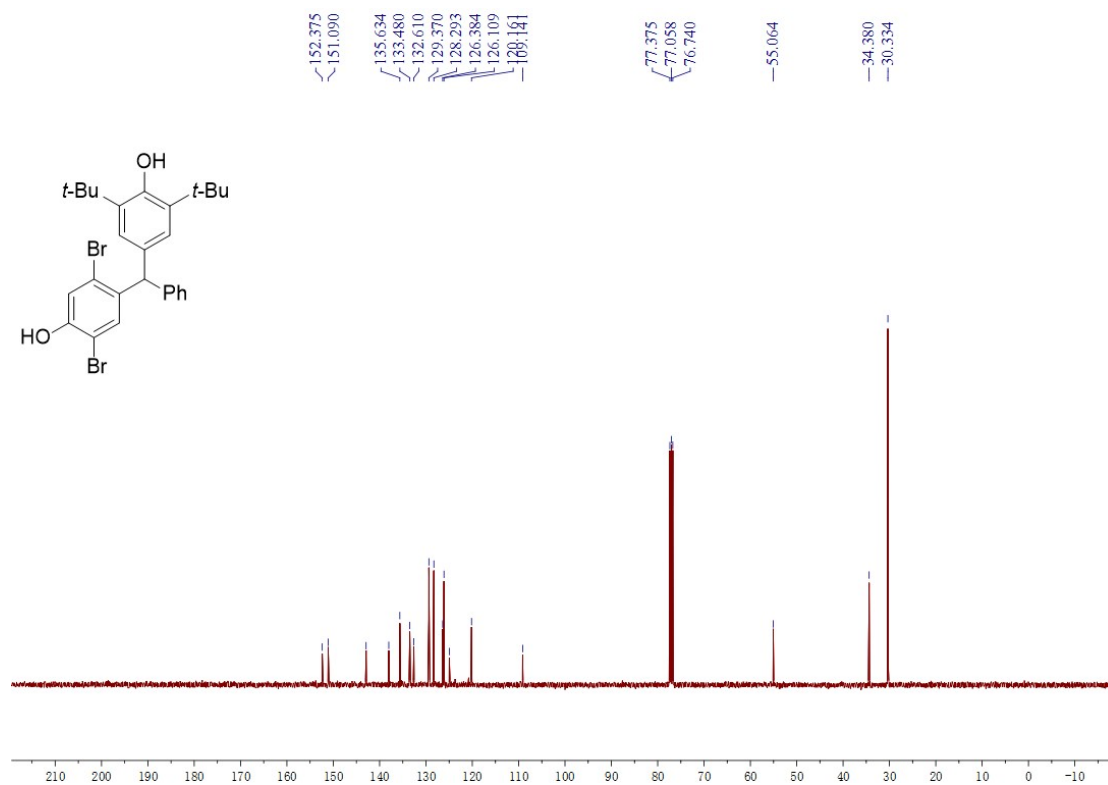
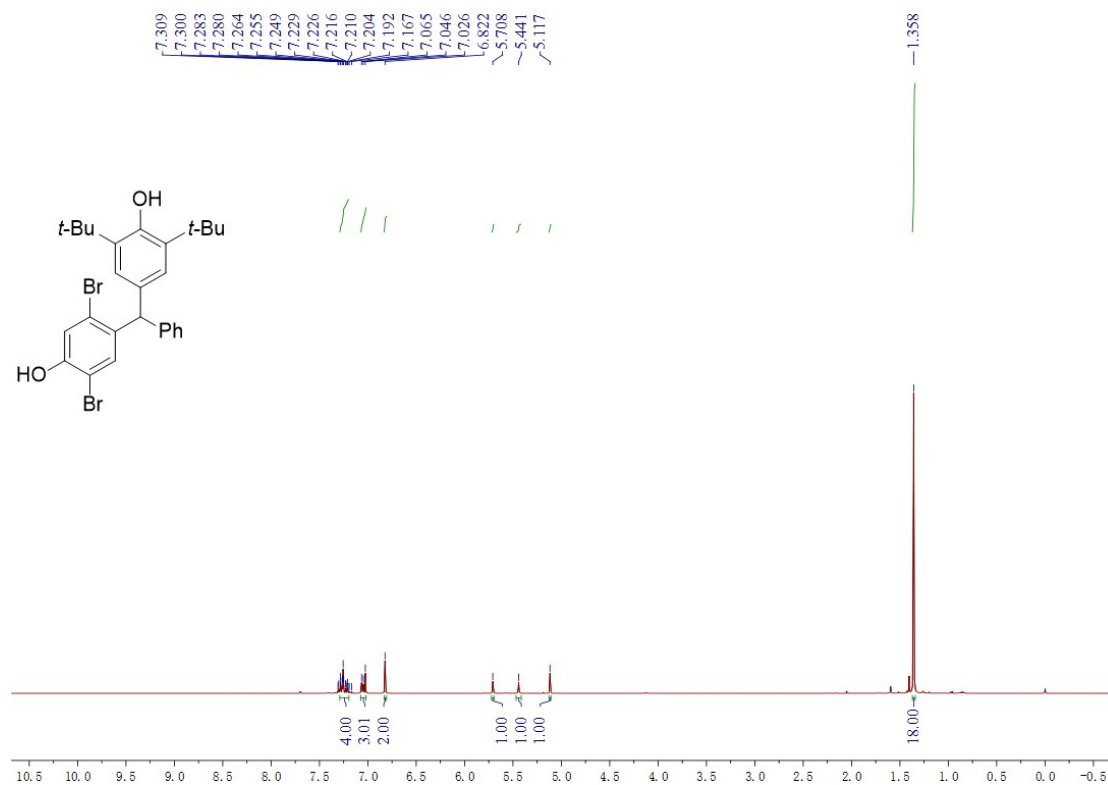


3-Bromo-4-((3,5-di-tert-butyl-4-hydroxyphenyl)(phenyl)methyl)-2-fluorophenol (6q)

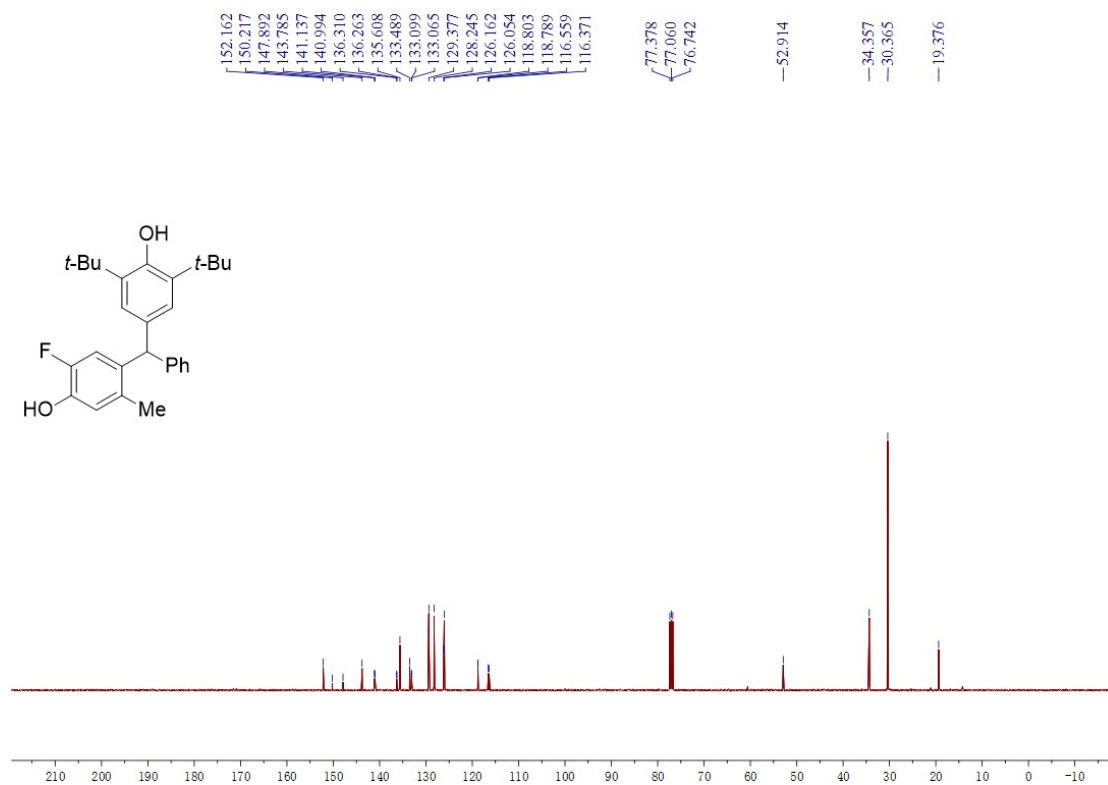
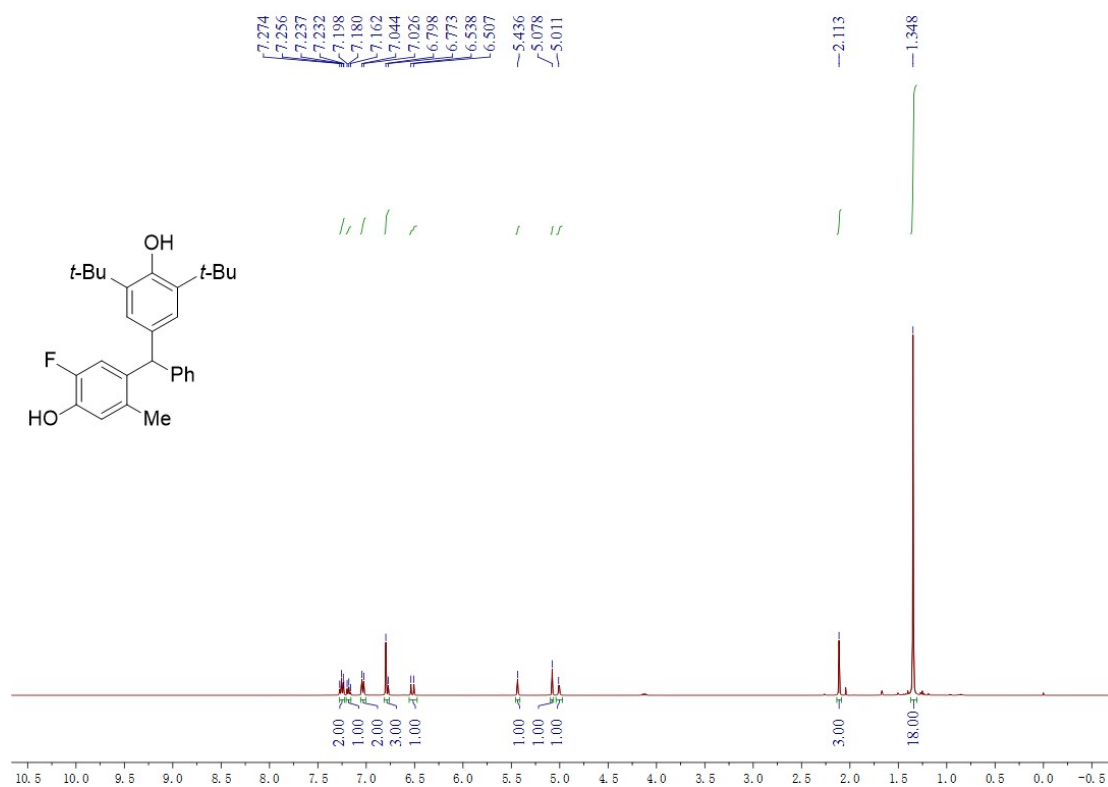


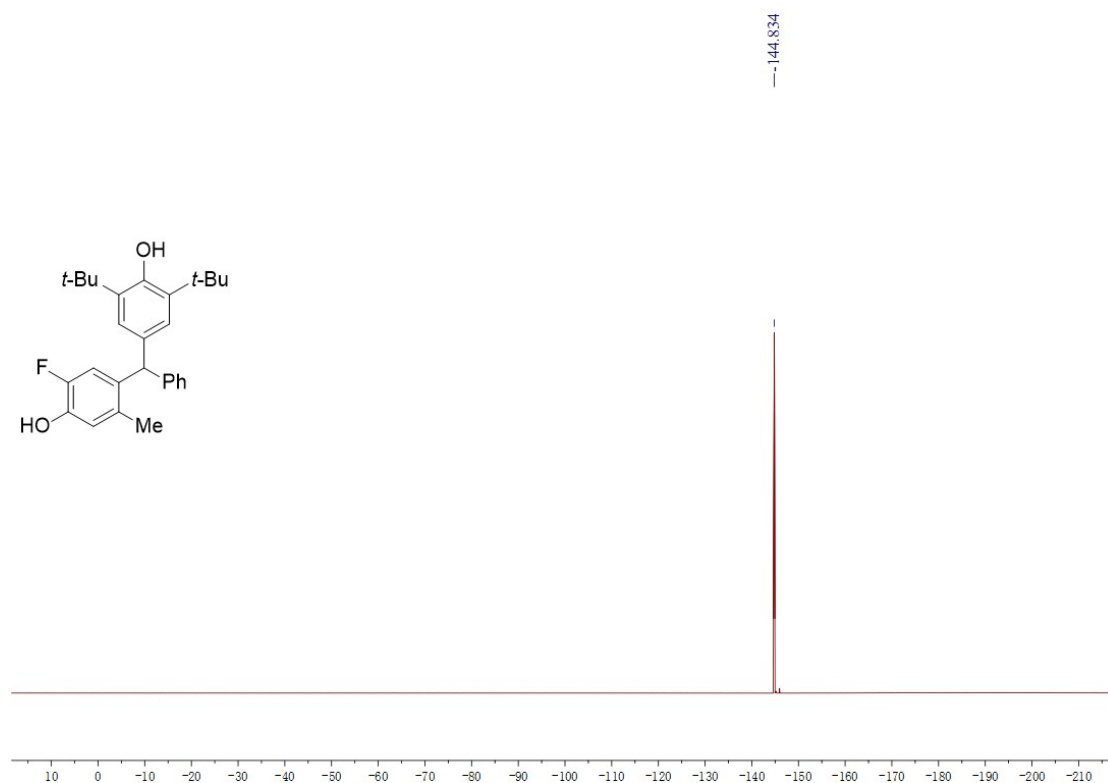


2,5-Dibromo-4-((3,5-di-tert-butyl-4-hydroxyphenyl)(phenyl)methyl)phenol (6r)

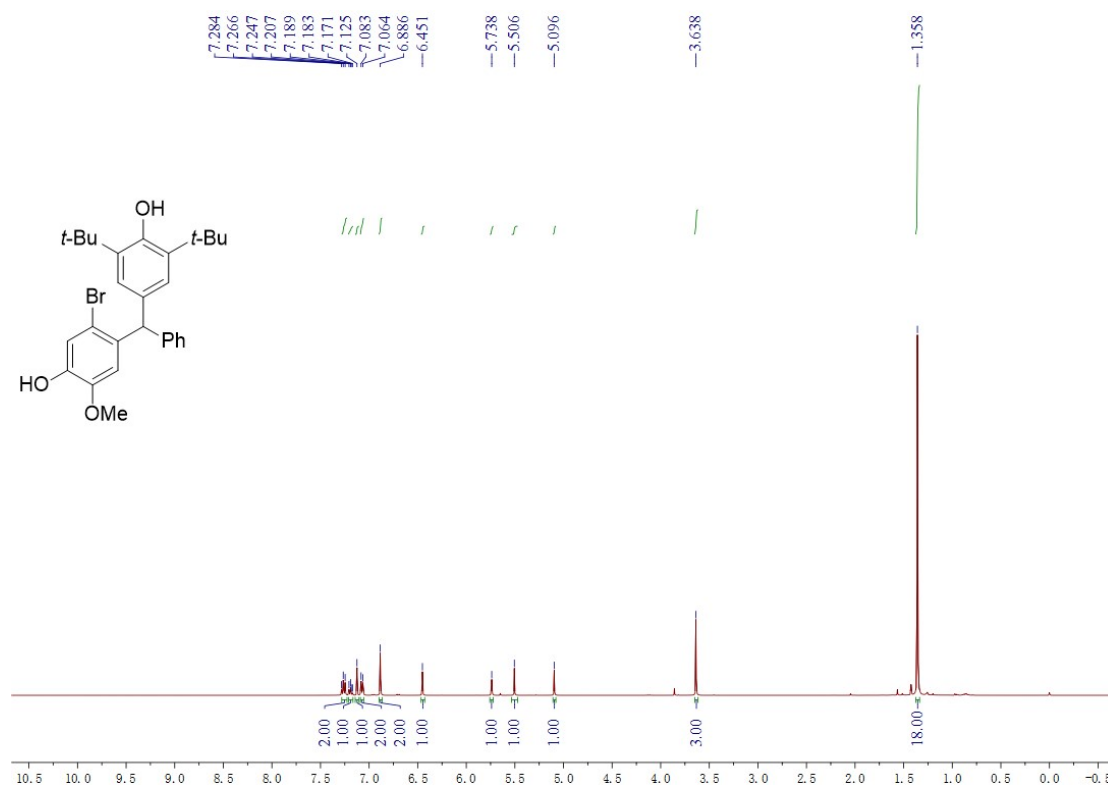


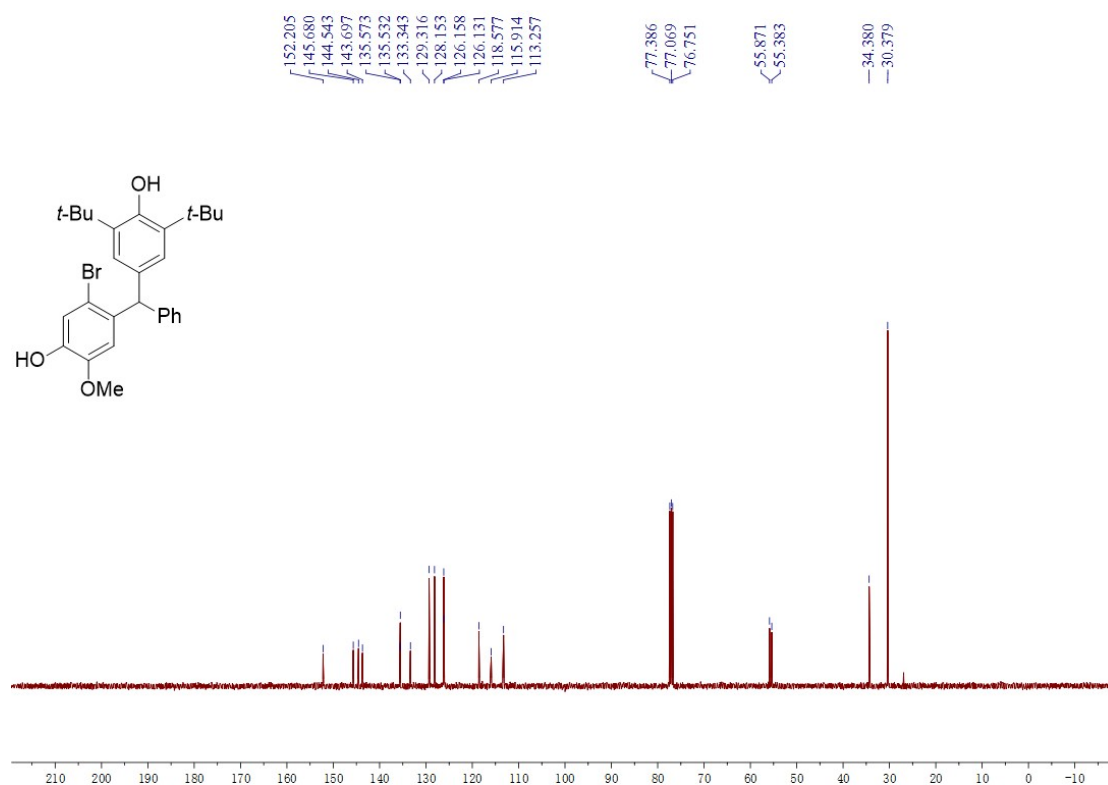
4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(phenyl)methyl)-2-fluoro-5-methylphenol (6s)



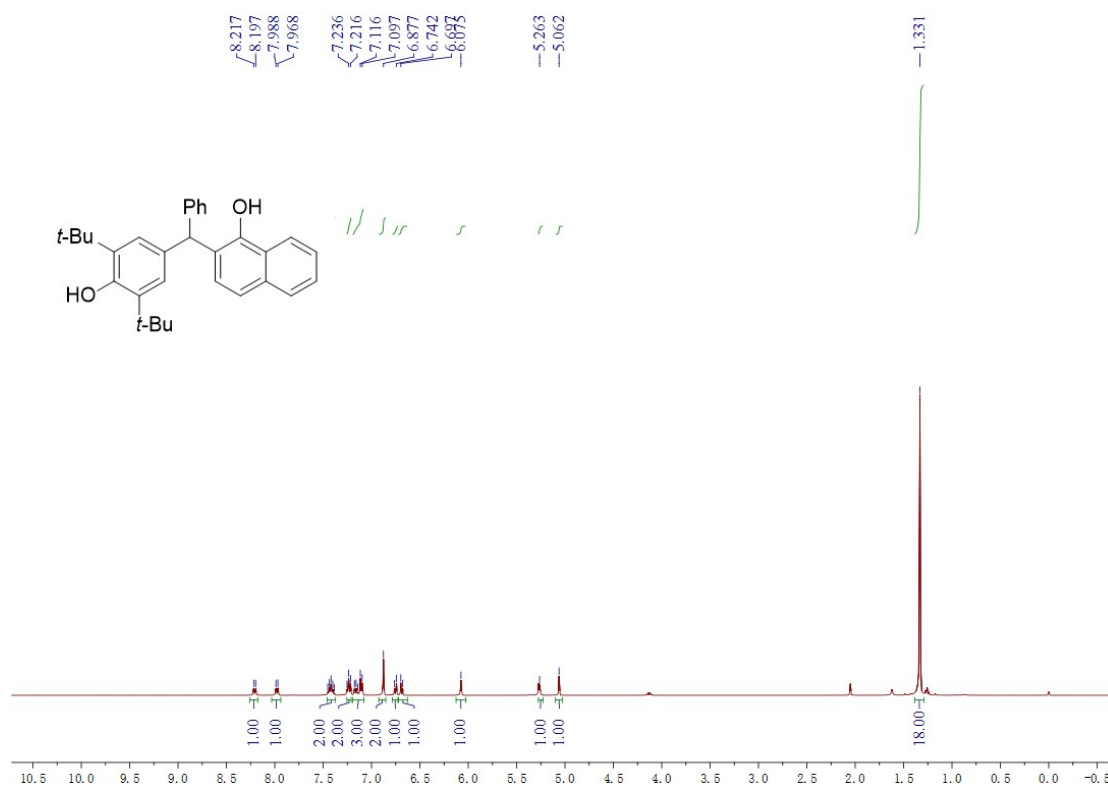


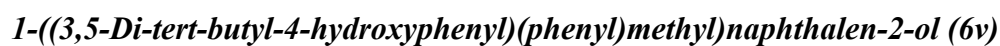
5-Bromo-4-((3,5-di-tert-butyl-4-hydroxyphenyl)(phenyl)methyl)-2-methoxyphenol (6t)

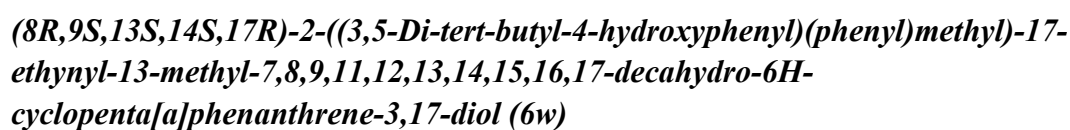


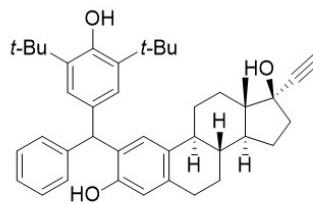


2-((3,5-Di-tert-butyl-4-hydroxyphenyl)(phenyl)methyl)naphthalen-1-ol (6u)

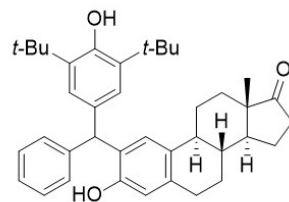








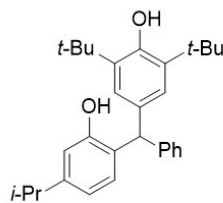
Chemical structure of compound 10 is shown. The ¹H NMR spectrum (CDCl₃) displays peaks from 0 to 10.5 ppm. Key features include a broad singlet at ~7.1 ppm (OH), aromatic signals between 6.5-7.5 ppm, aliphatic signals between 1.0-2.5 ppm, and a large solvent peak at 7.26 ppm. Integration values are provided below the baseline.

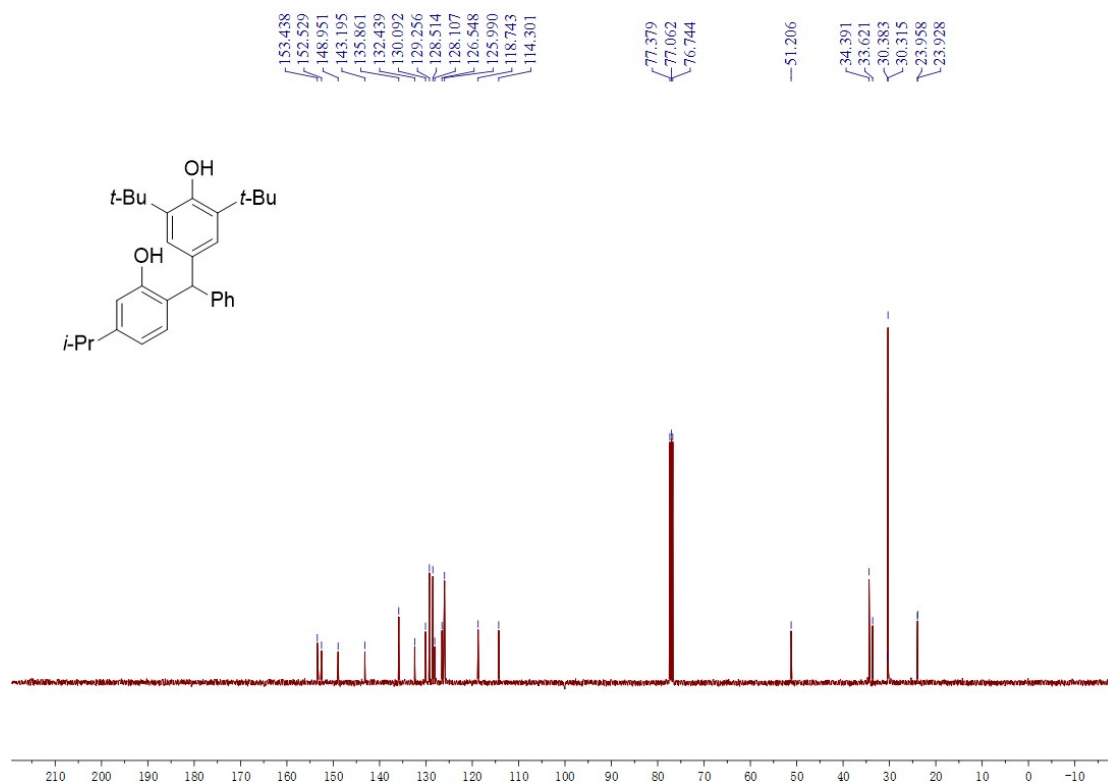


Chemical structure of 2,6-di-*t*-butyl-4-(1-(4-*i*-propylphenyl)-2-hydroxyethyl)phenol (1) is shown. The structure features a central biphenyl core with two hydroxyl groups at the 2 and 6 positions of the central ring, and two *t*-butyl groups at the 4 and 4' positions. The 1-position of the central ring is substituted with a 1-(4-*i*-propylphenyl)-2-hydroxyethyl group. The chemical shift values (ppm) for the protons are listed below the structure, grouped by color-coded brackets: blue for aromatic protons, green for aliphatic protons, and red for the *i*-propyl group protons.

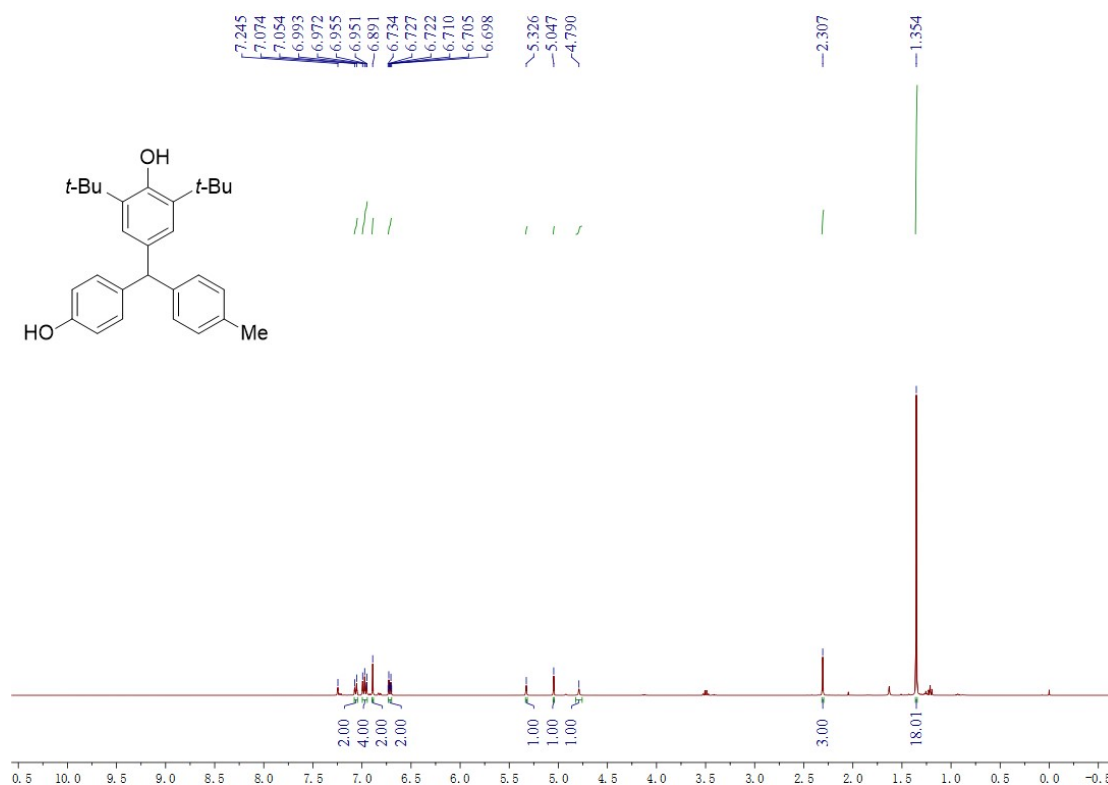
Chemical shift values (ppm):

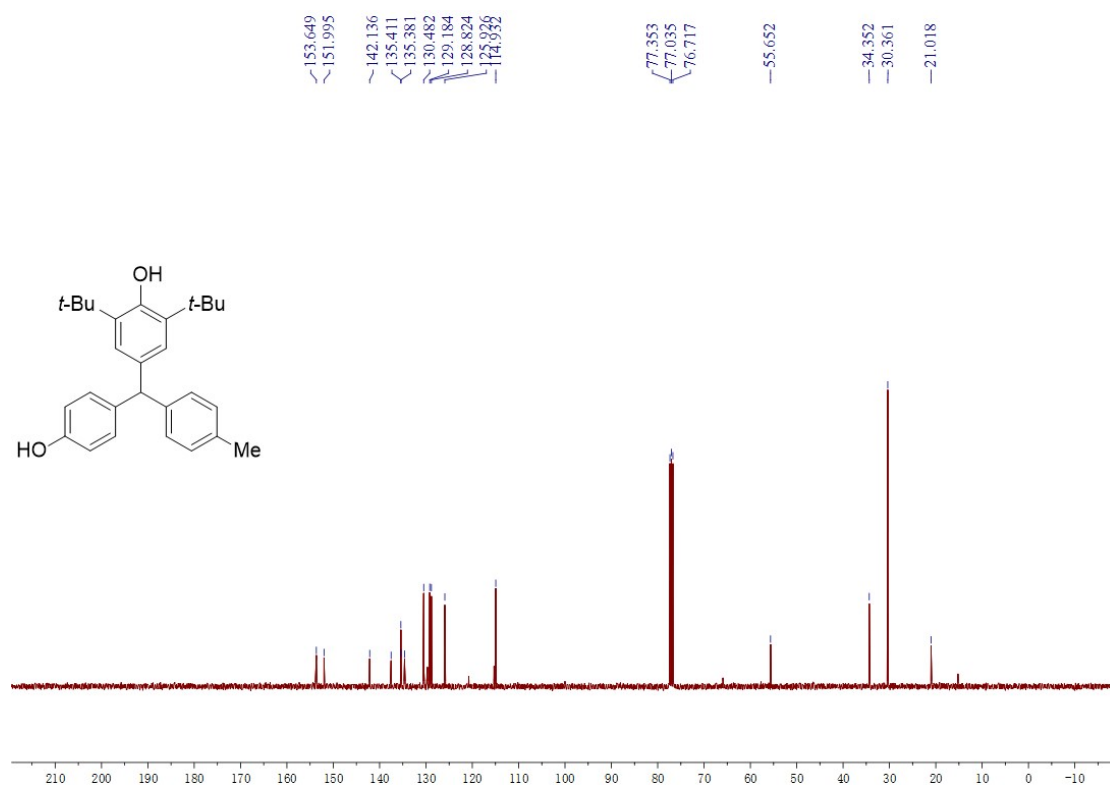
- Aromatic protons (blue): 7.314, 7.296, 7.277, 7.245, 7.236, 7.222, 7.217, 7.212, 7.199, 7.186, 7.166, 7.163, 7.145, 7.130, 6.930, 6.732, 6.723, 6.720, 6.710, 6.700, 6.693, 5.564, 5.121, 4.632.
- Aliphatic protons (green): 2.883, 2.866, 2.849, 2.832, 2.814, 2.797, 2.780.
- i*-Propyl group protons (red): 1.354, 1.225, 1.208.



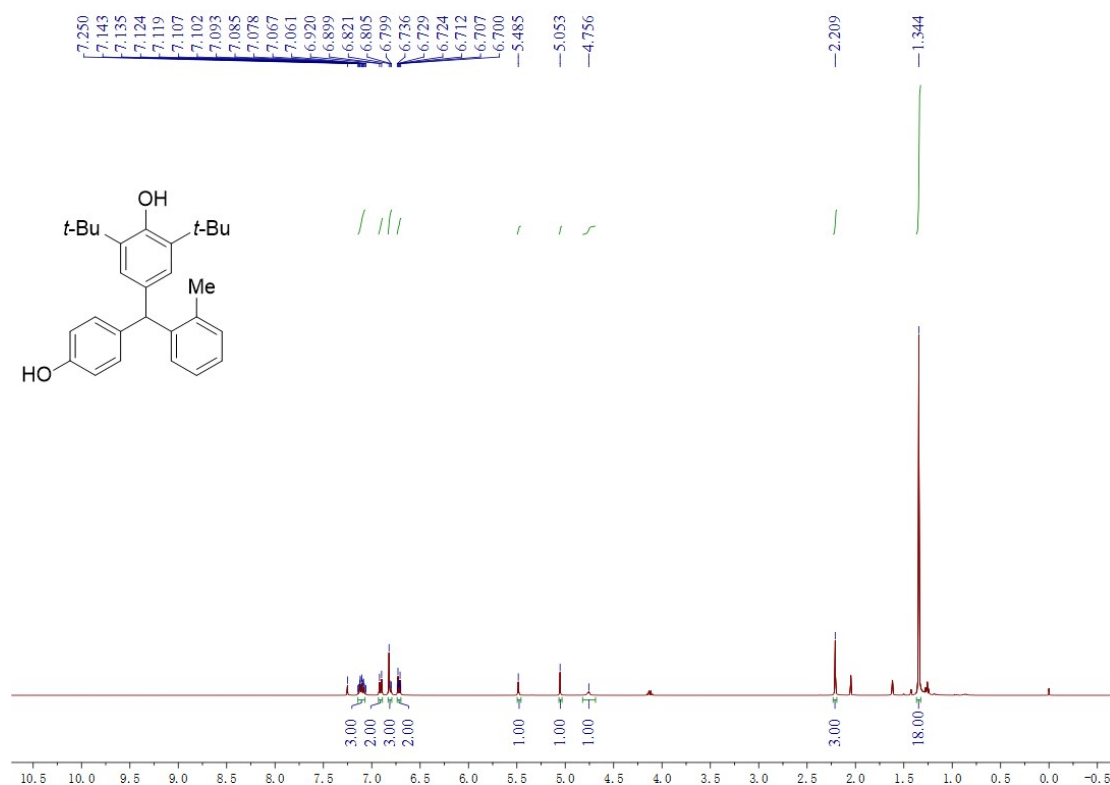


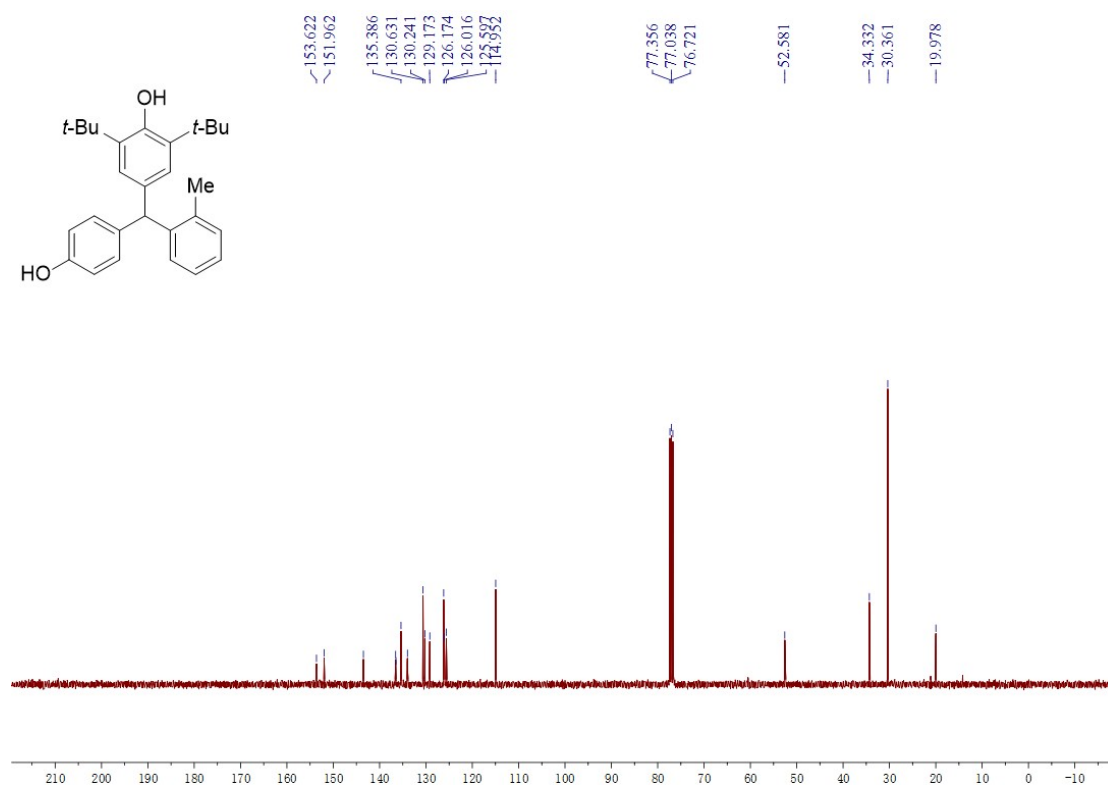
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(p-tolyl)methyl)phenol (7a)



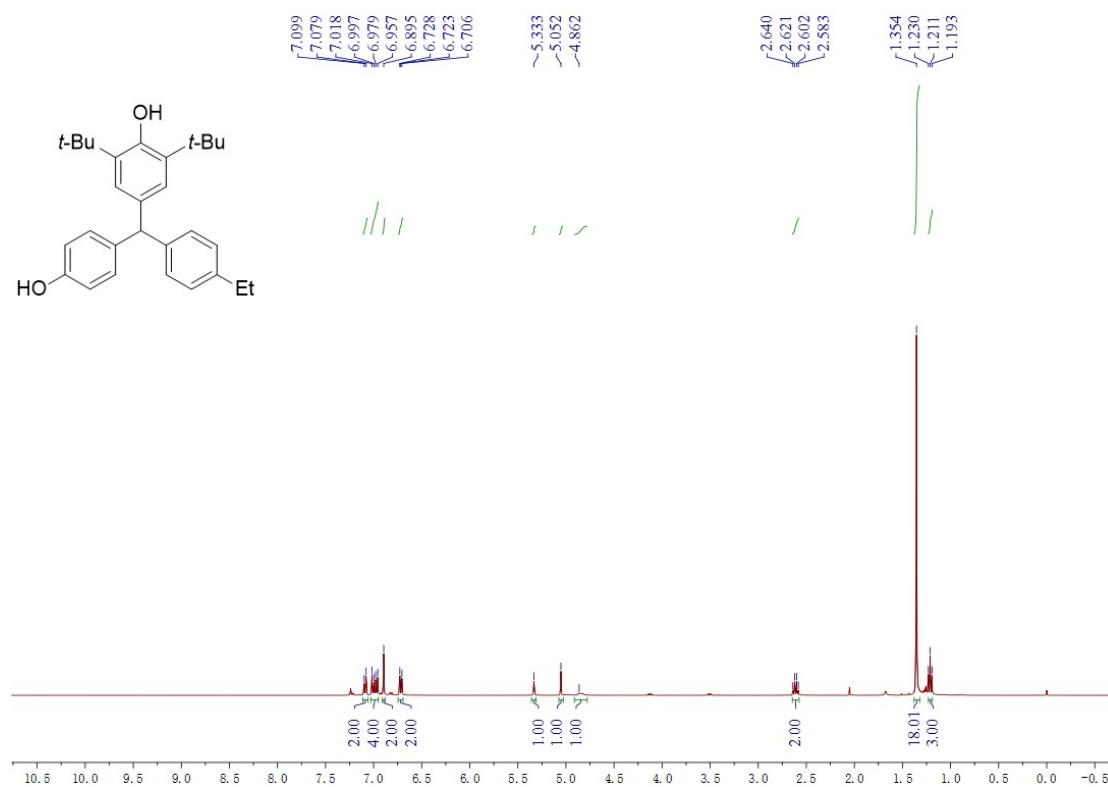


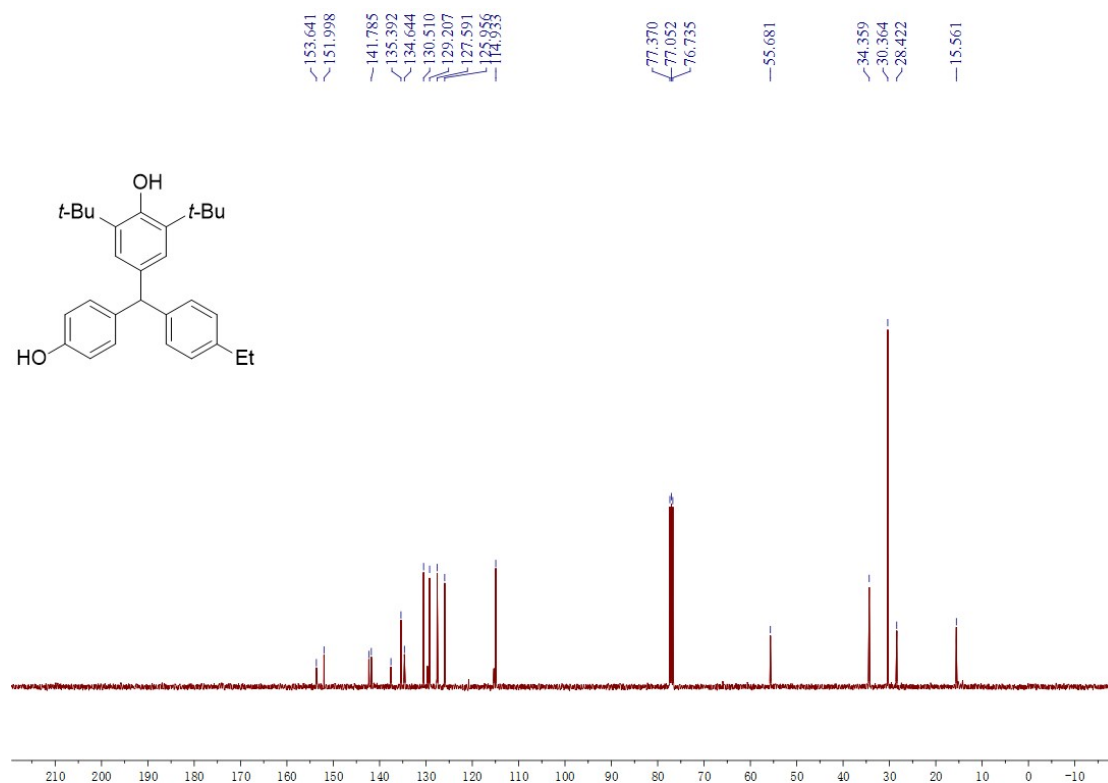
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(o-tolyl)methyl)phenol (7b)



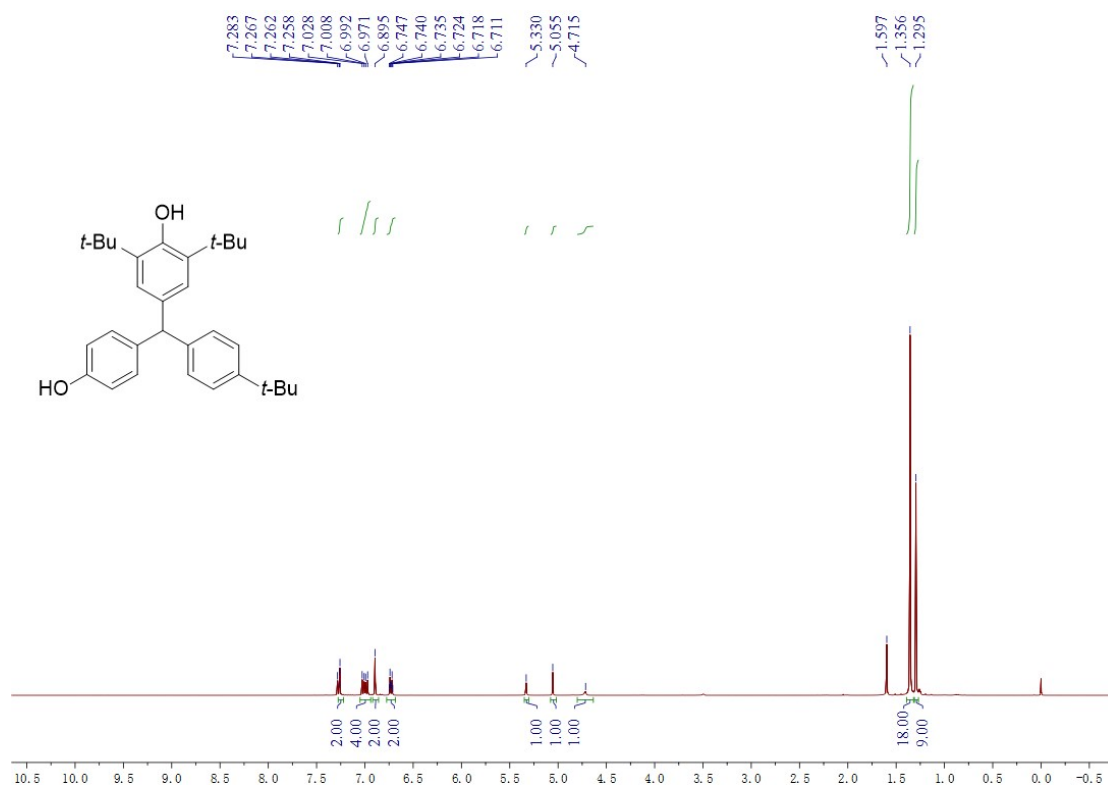


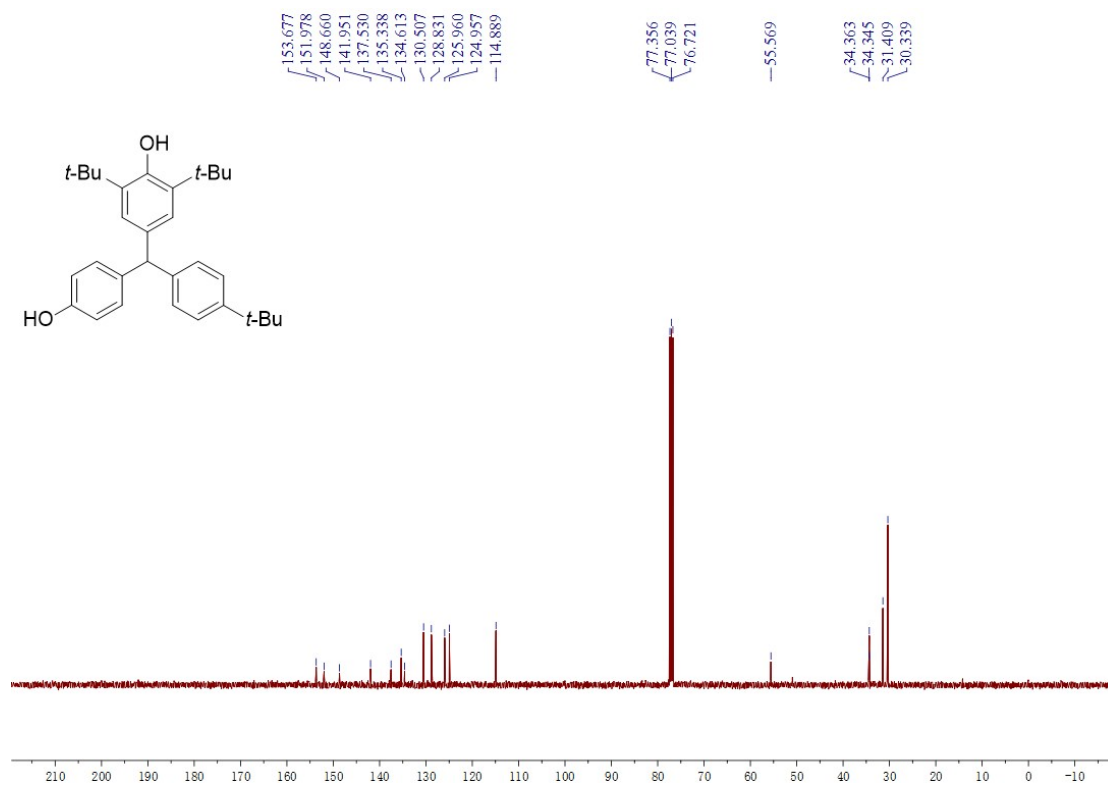
2,6-Di-tert-butyl-4-((4-ethylphenyl)(4-hydroxyphenyl)methyl)phenol (7c)



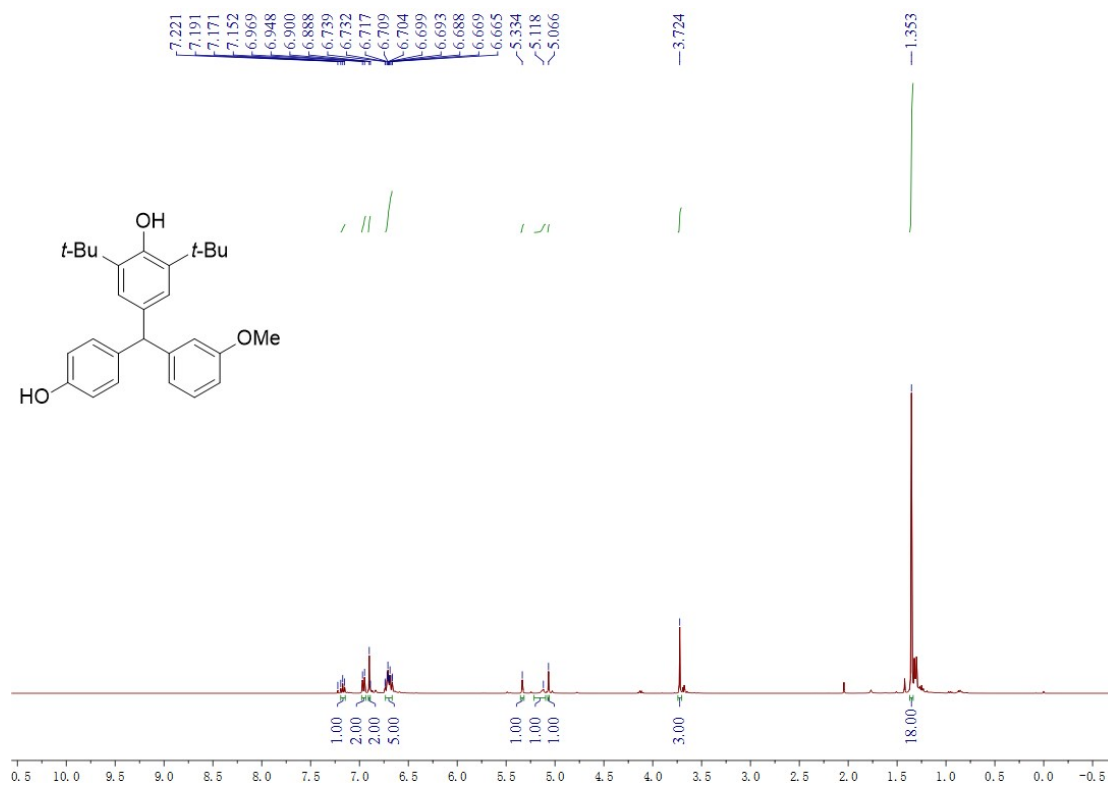


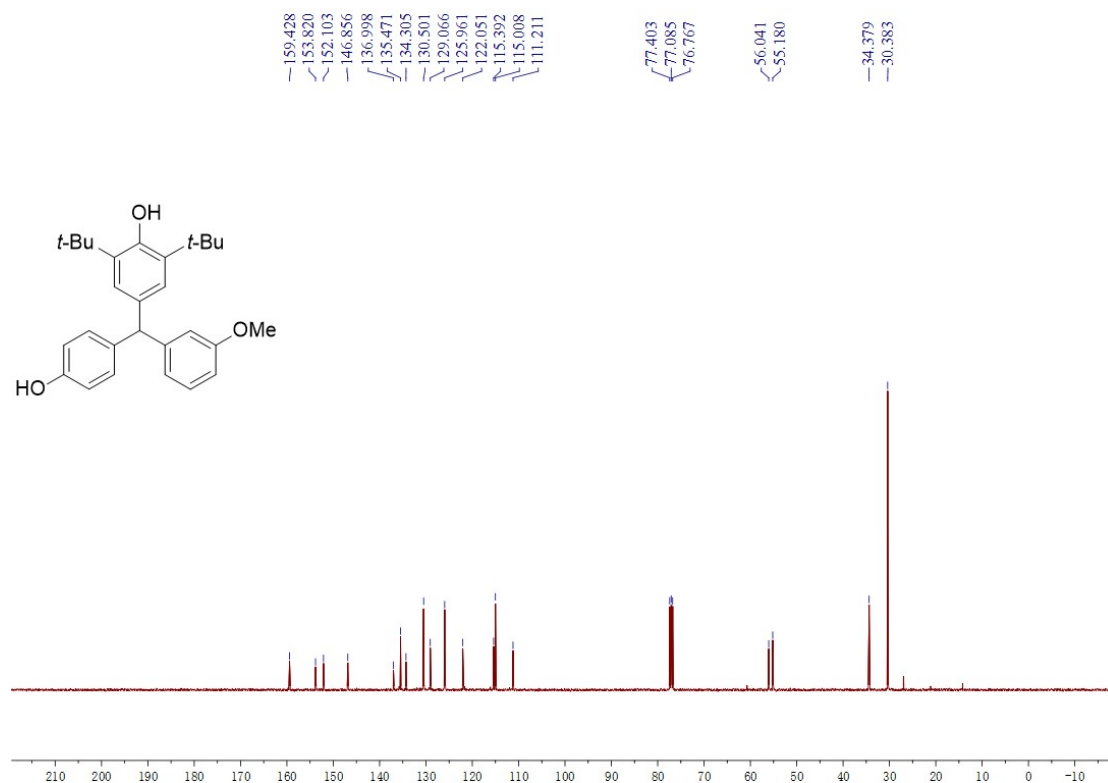
2,6-Di-tert-butyl-4-((4-(tert-butyl)phenyl)(4-hydroxyphenyl)methyl)phenol (7d)



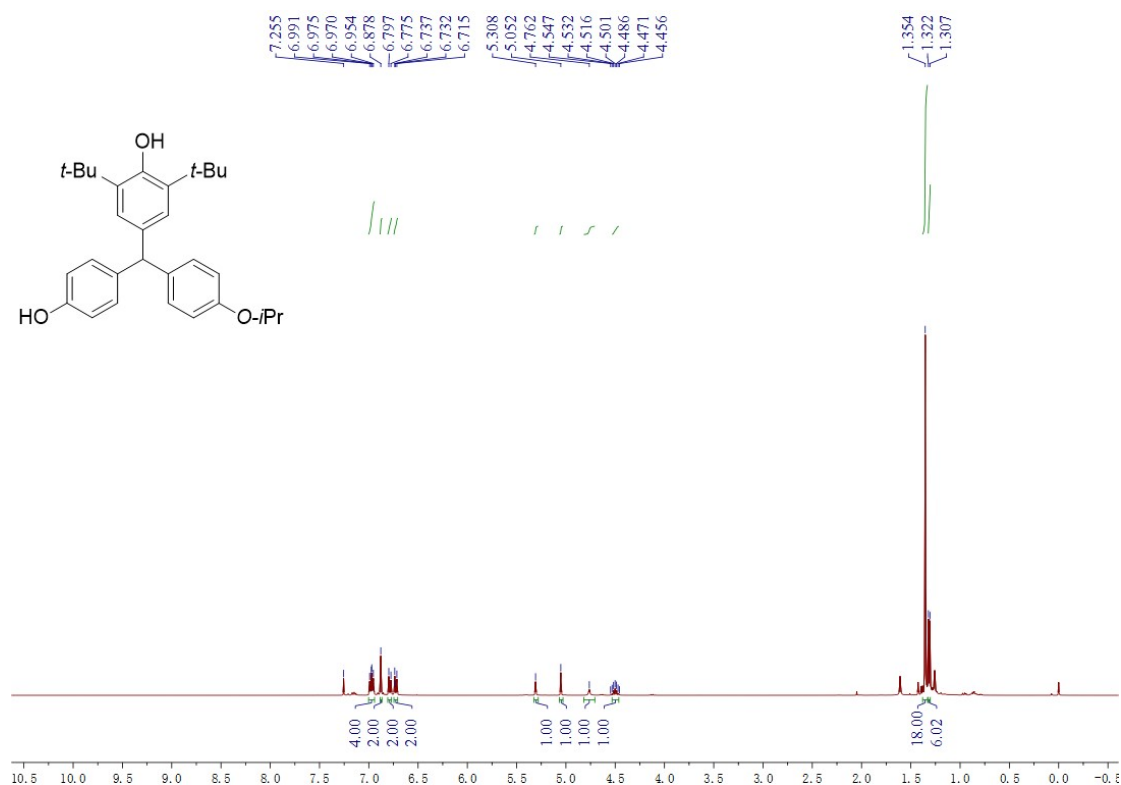


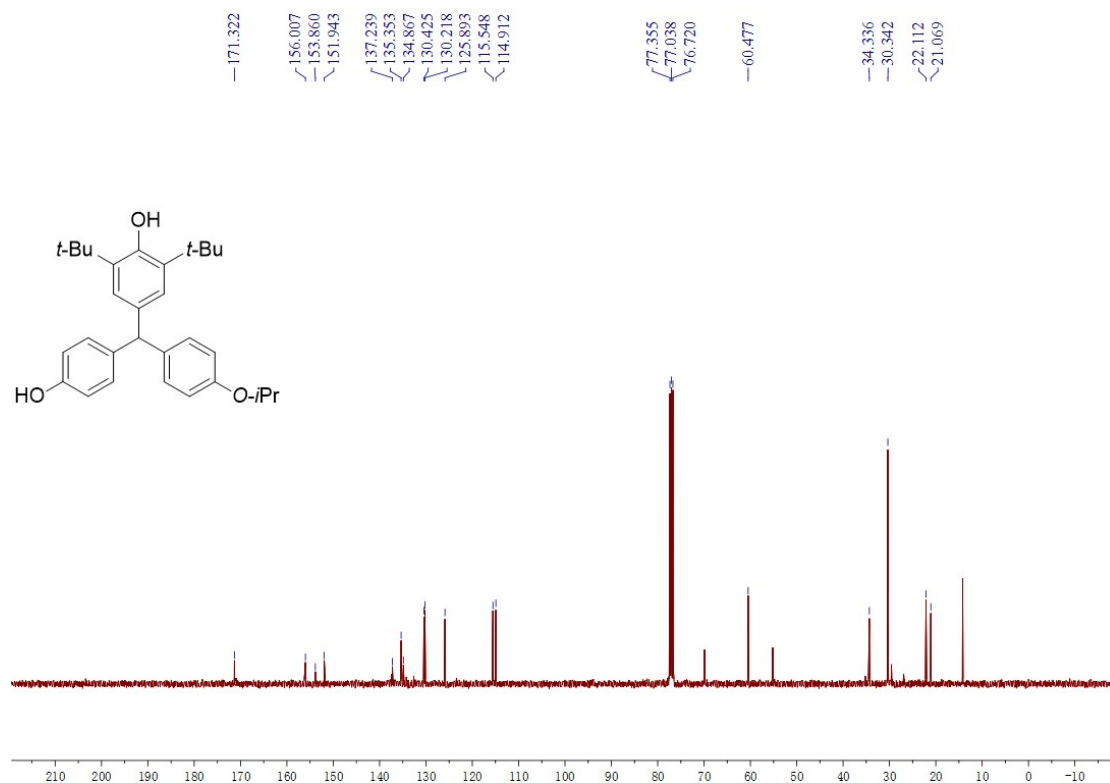
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(3-methoxyphenyl)methyl)phenol (7e)



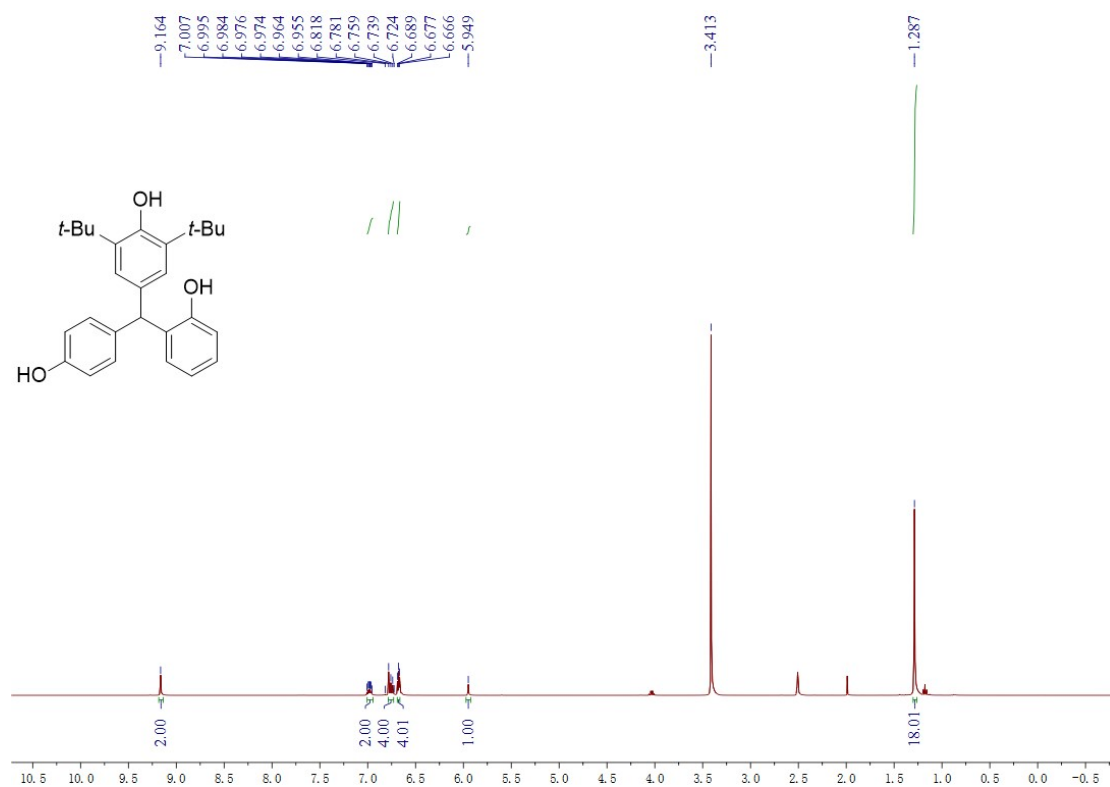


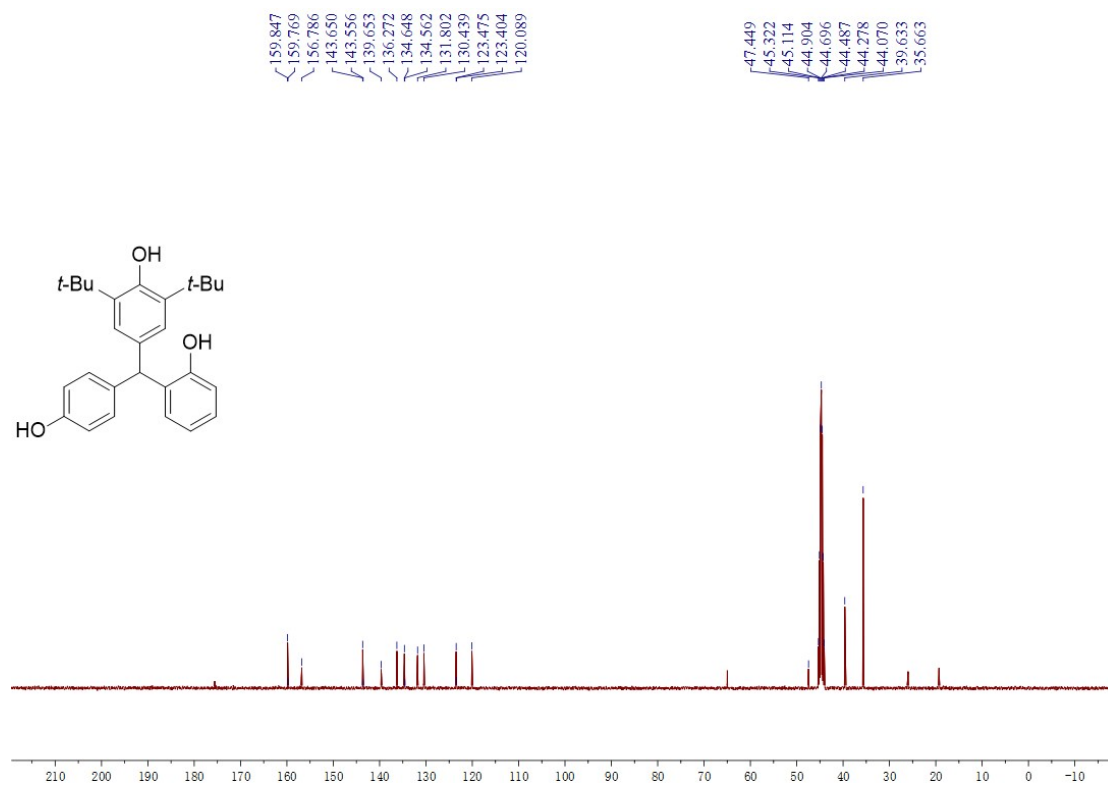
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(4-isopropoxyphenyl)methyl)phenol (7f)



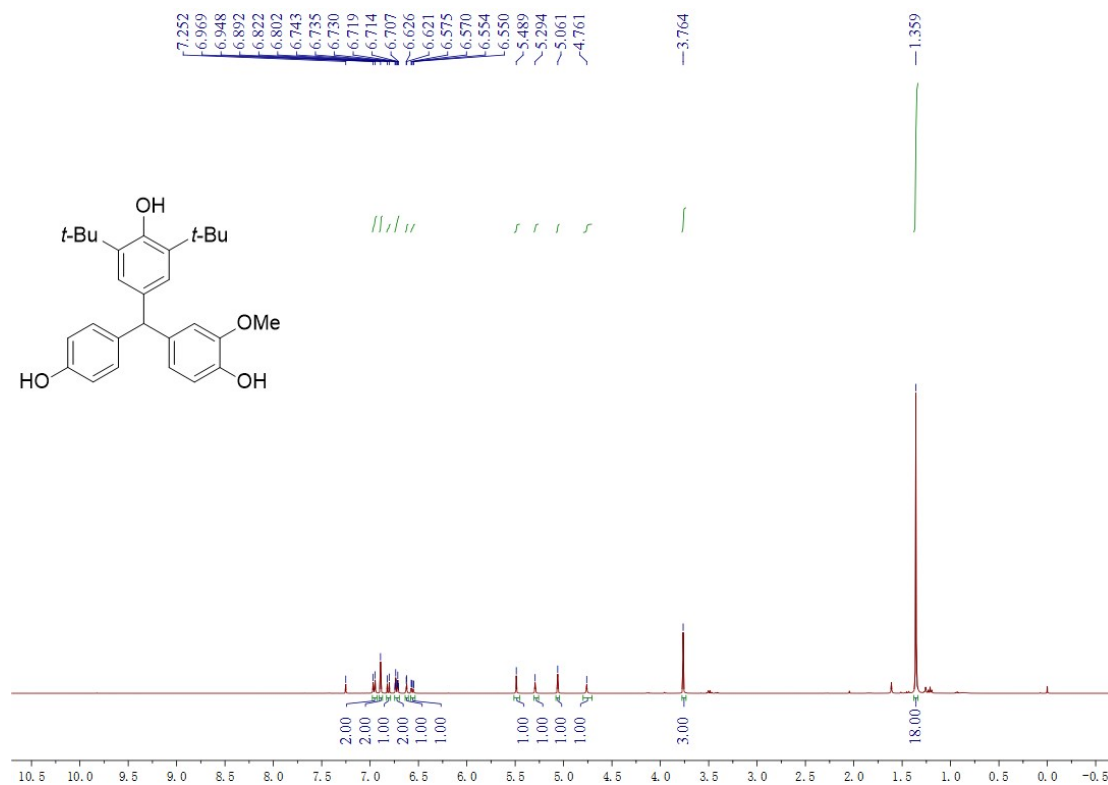


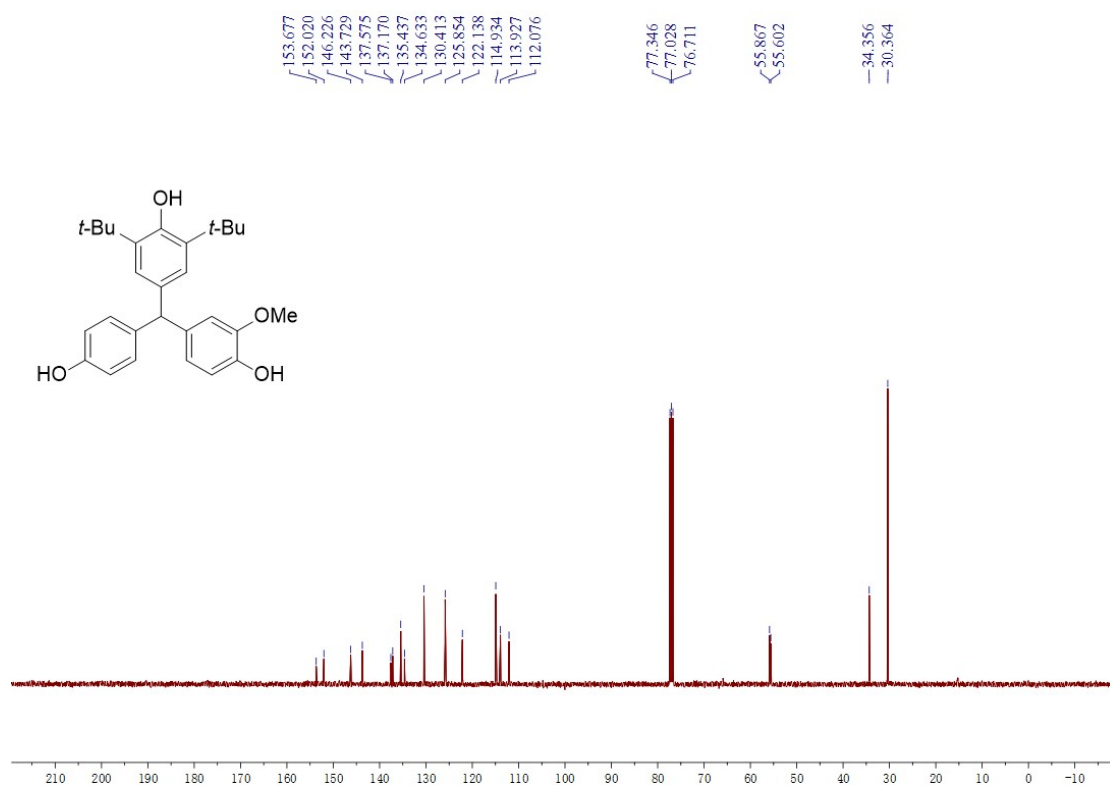
2,6-Di-tert-butyl-4-((2-hydroxyphenyl)(4-hydroxyphenyl)methyl)phenol (7h)



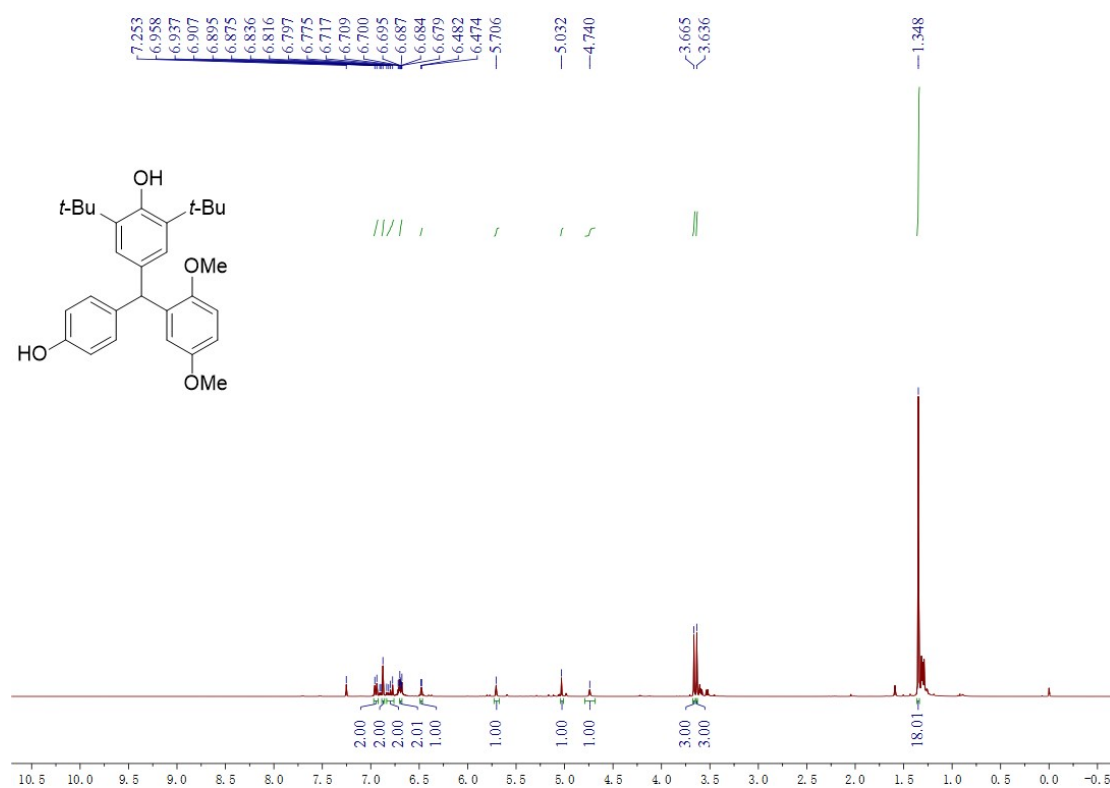


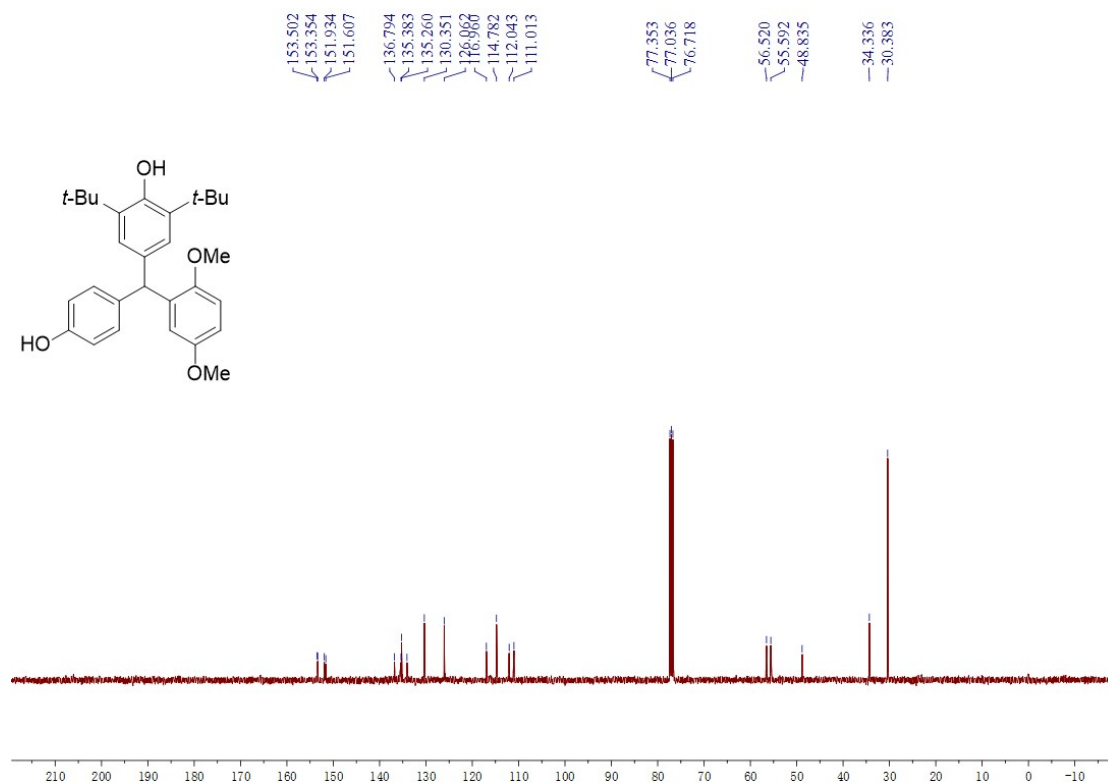
2,6-Di-tert-butyl-4-((4-hydroxy-3-methoxyphenyl)(4-hydroxyphenyl)methyl)phenol (7i)



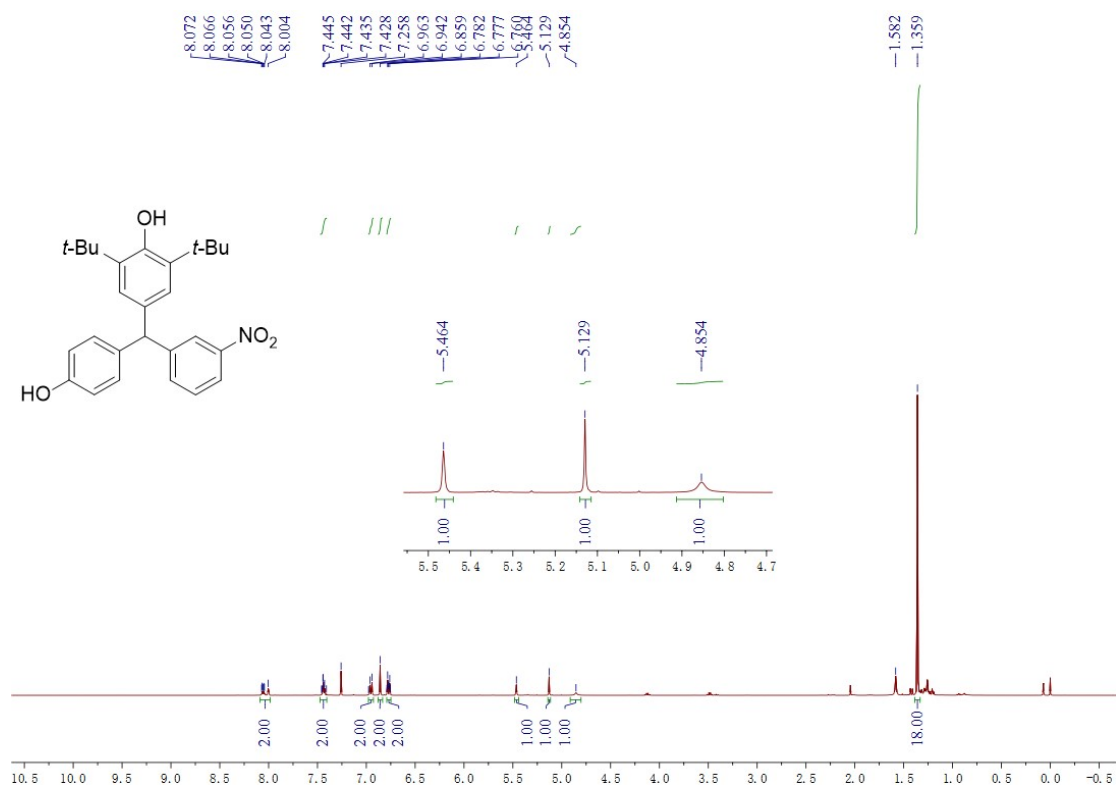


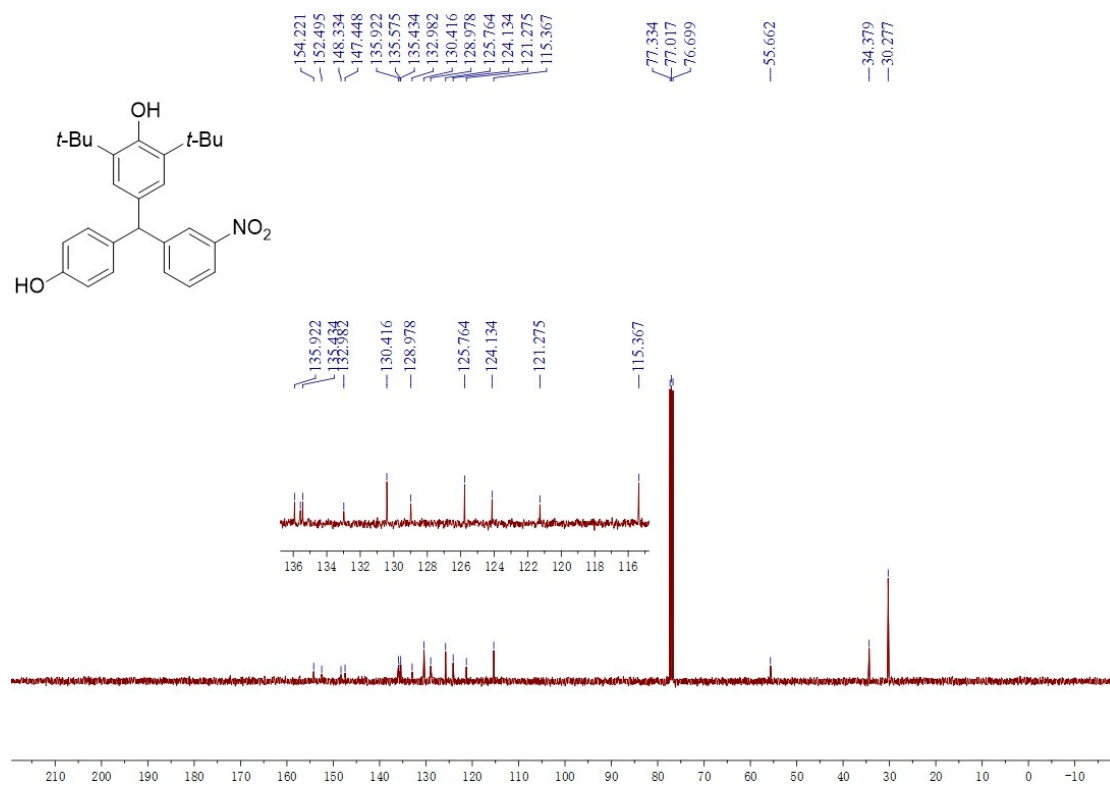
2,6-Di-tert-butyl-4-((2,5-dimethoxyphenyl)(4-hydroxyphenyl)methyl)phenol (7j)



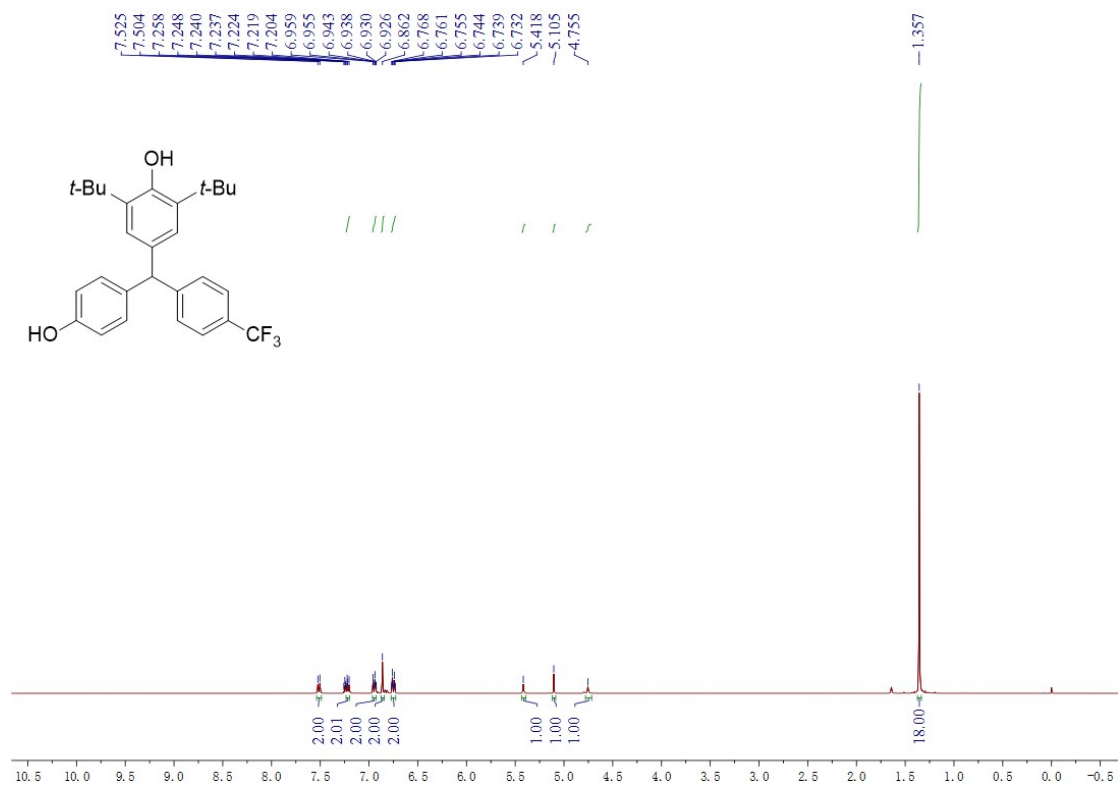


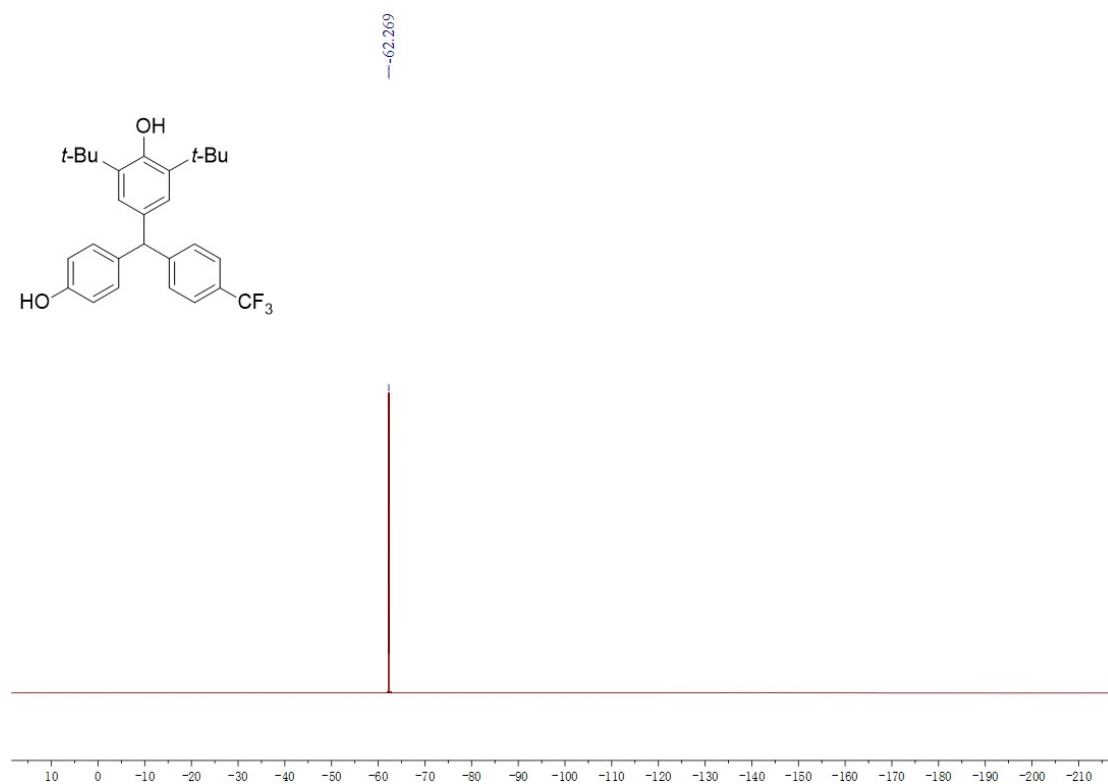
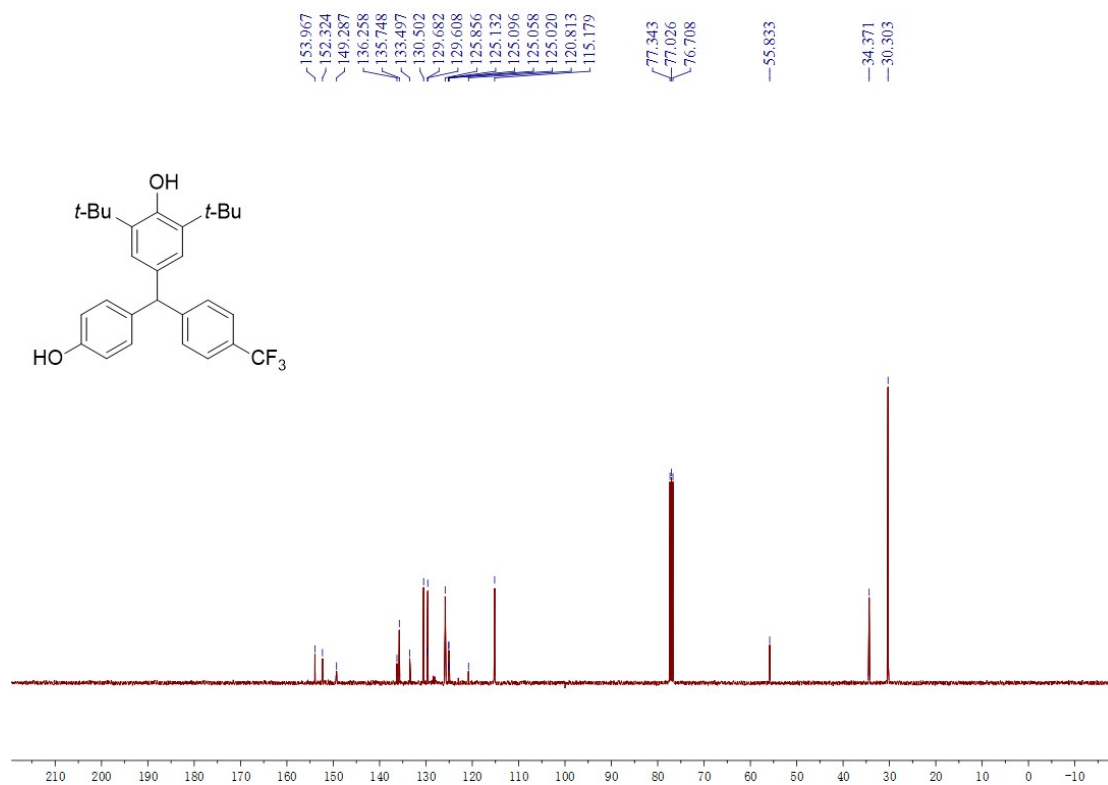
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(3-nitrophenyl)methyl)phenol (7k)



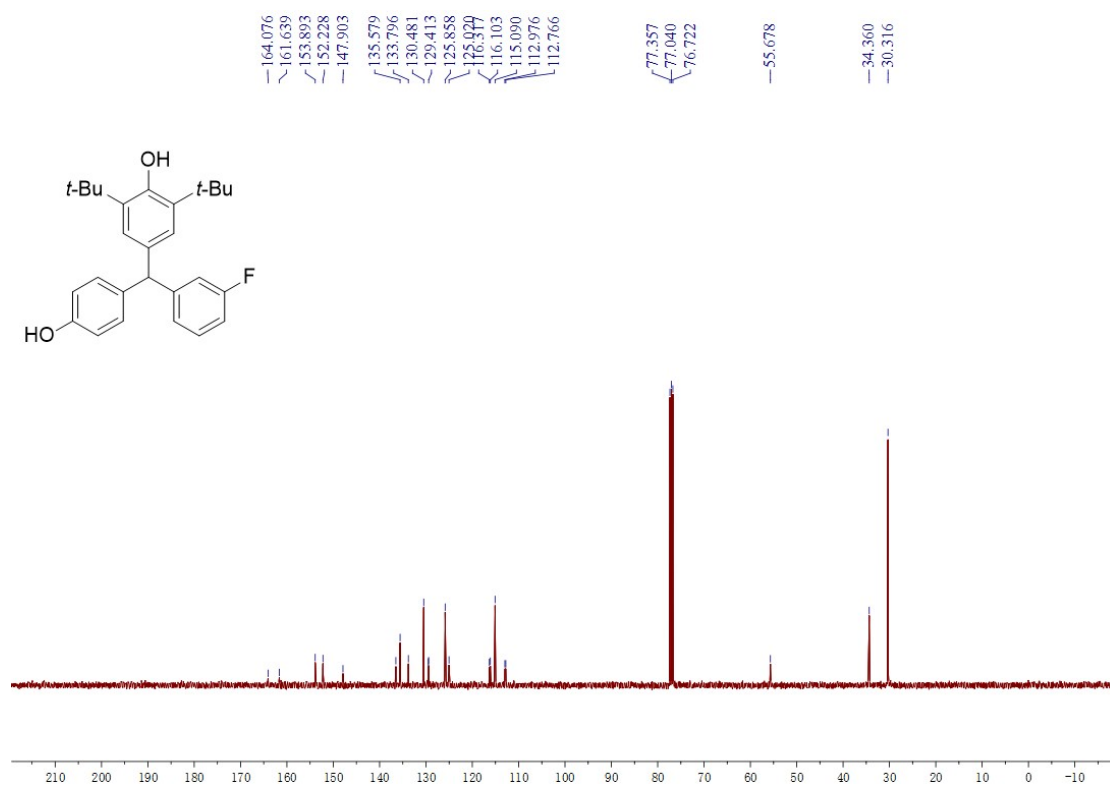
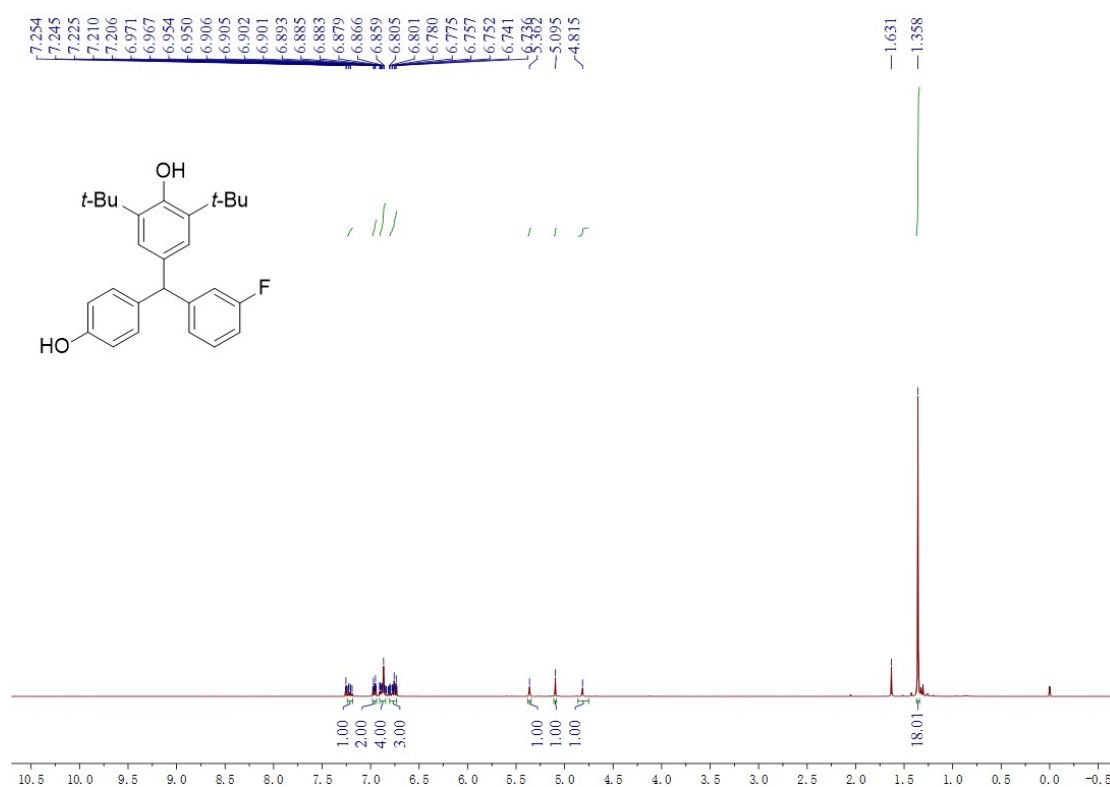


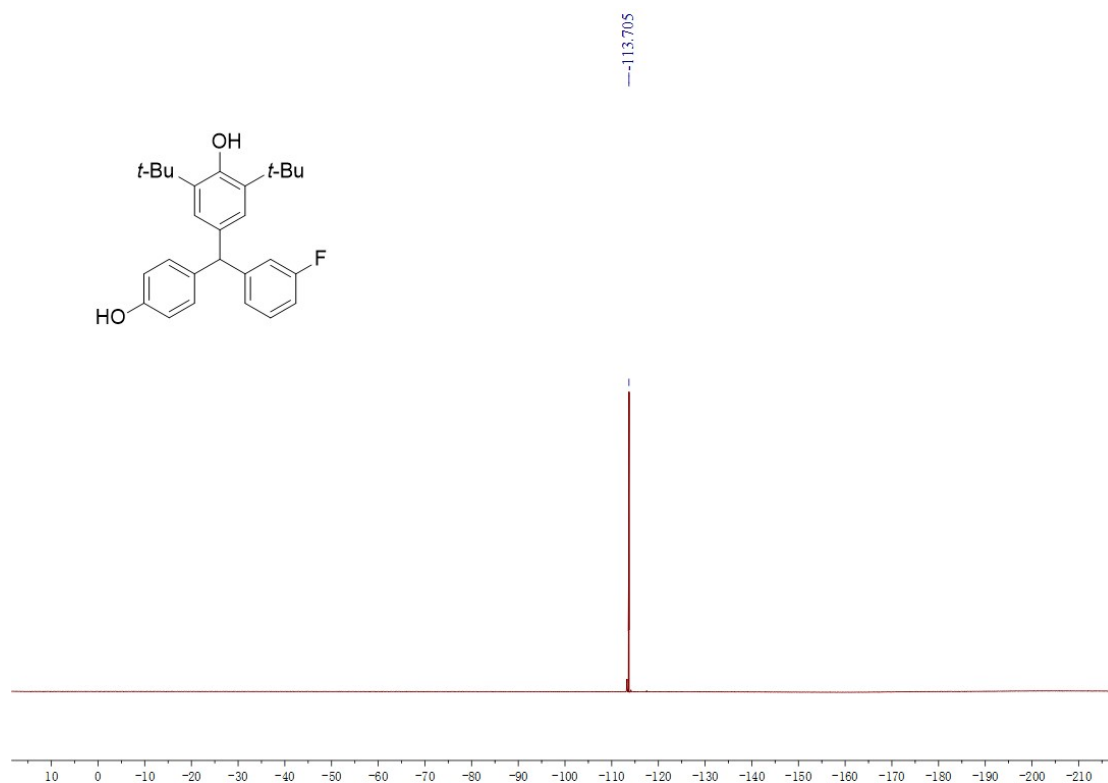
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(4-(trifluoromethyl)phenyl)methyl)phenol (7l)



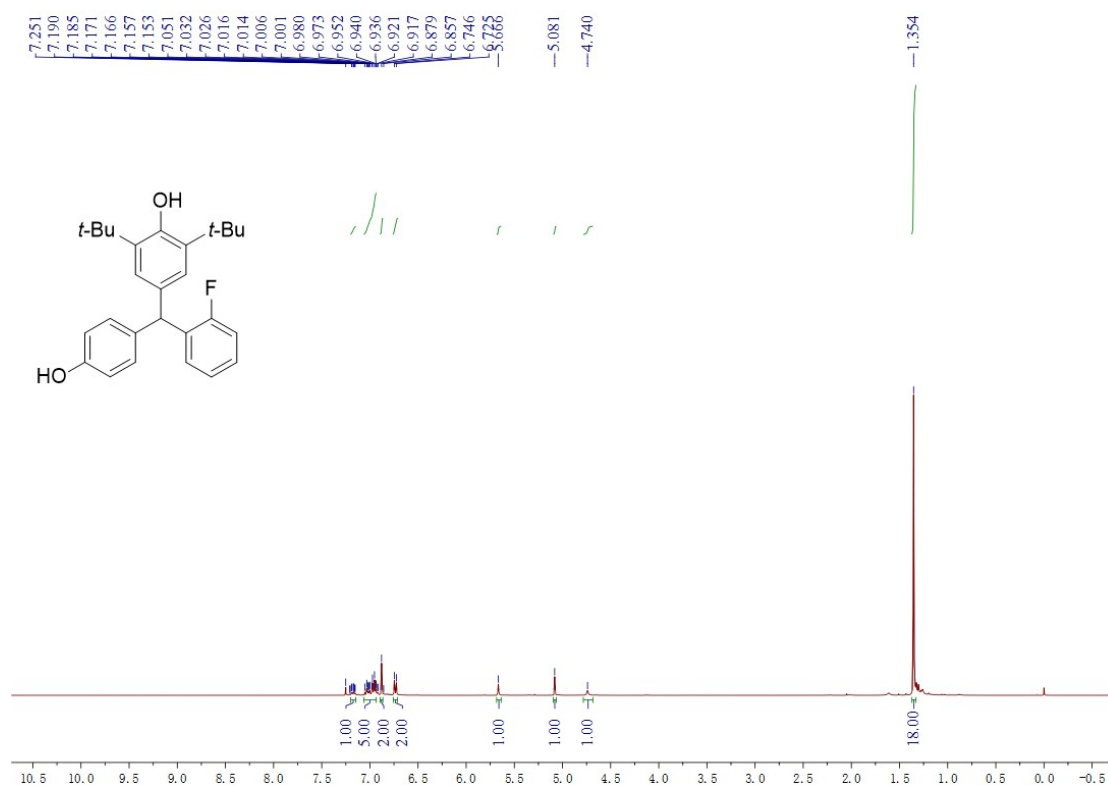


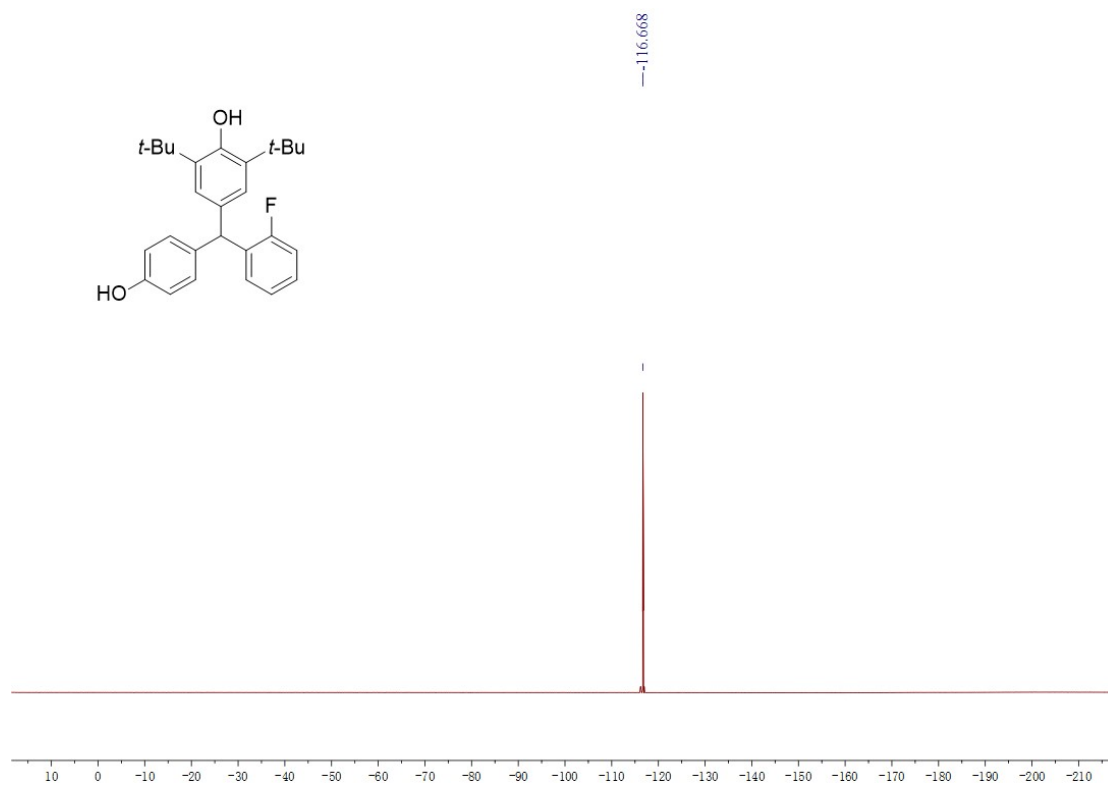
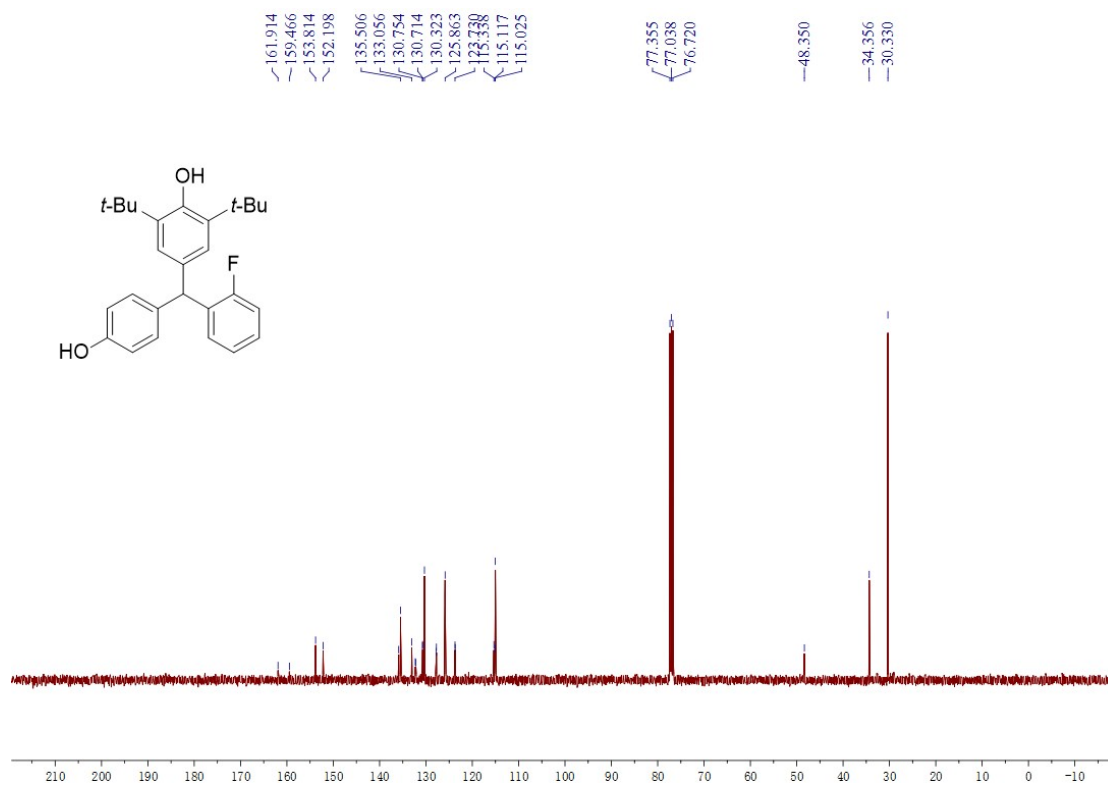
2,6-Di-tert-butyl-4-((3-fluorophenyl)(4-hydroxyphenyl)methyl)phenol (7m)



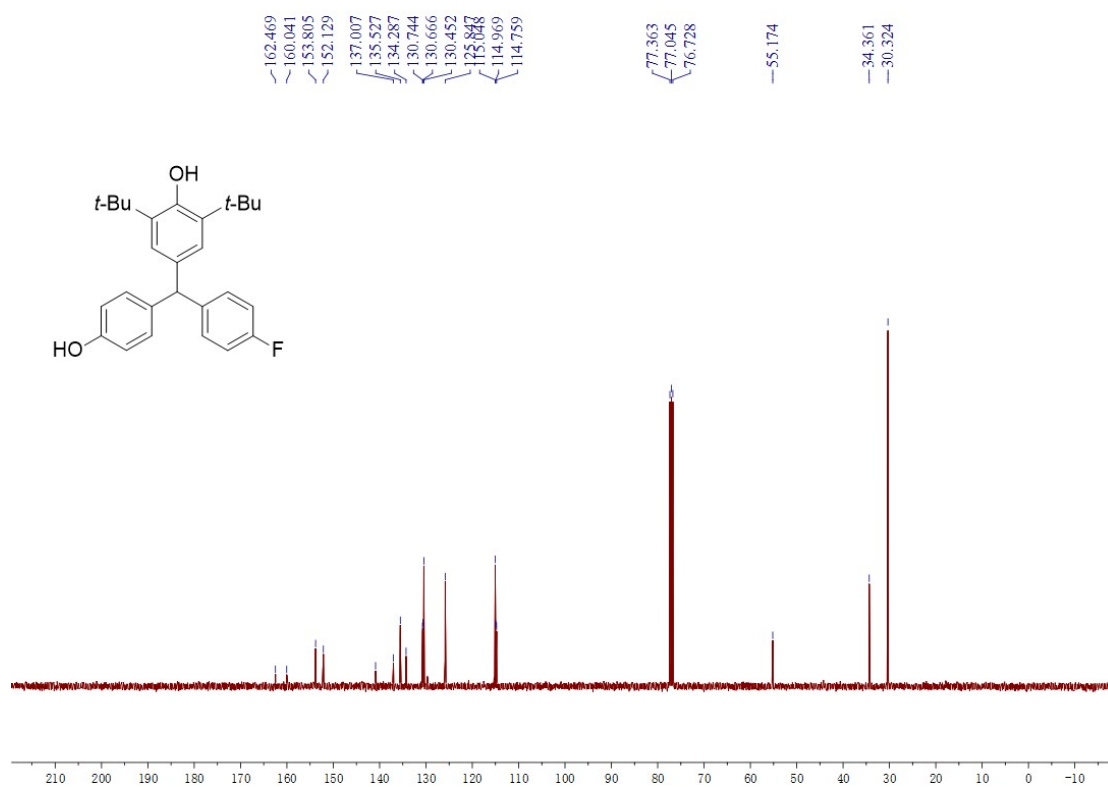
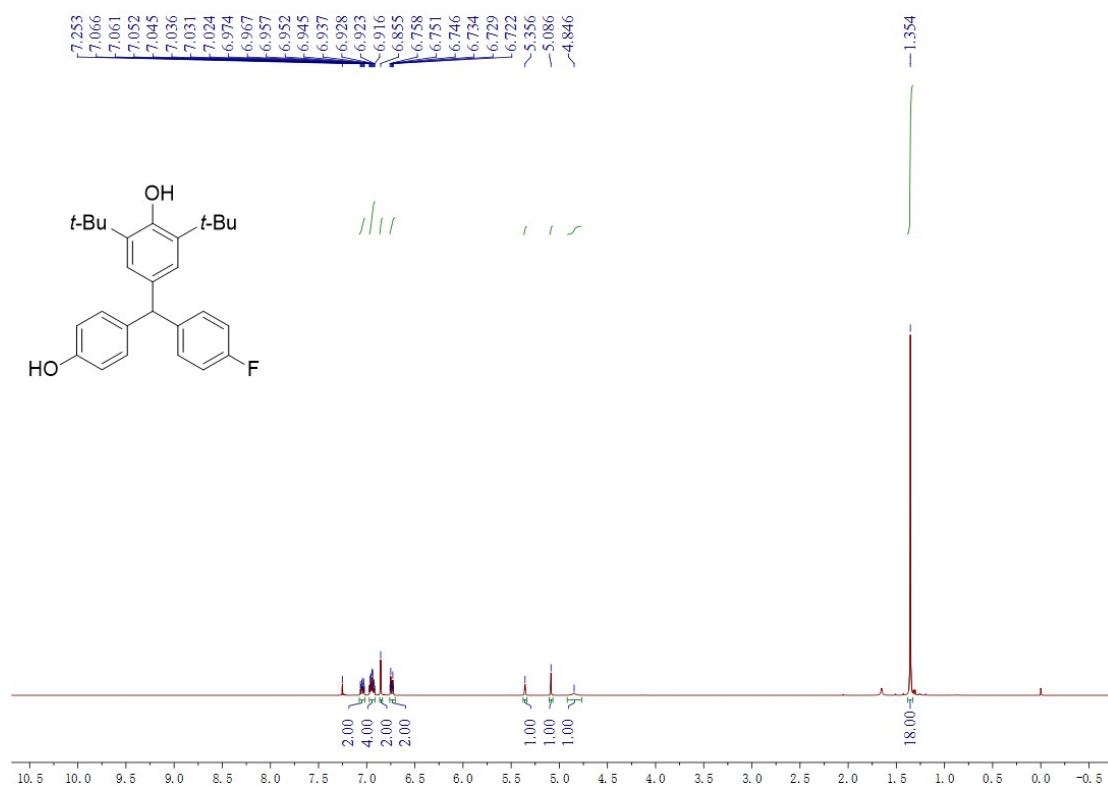


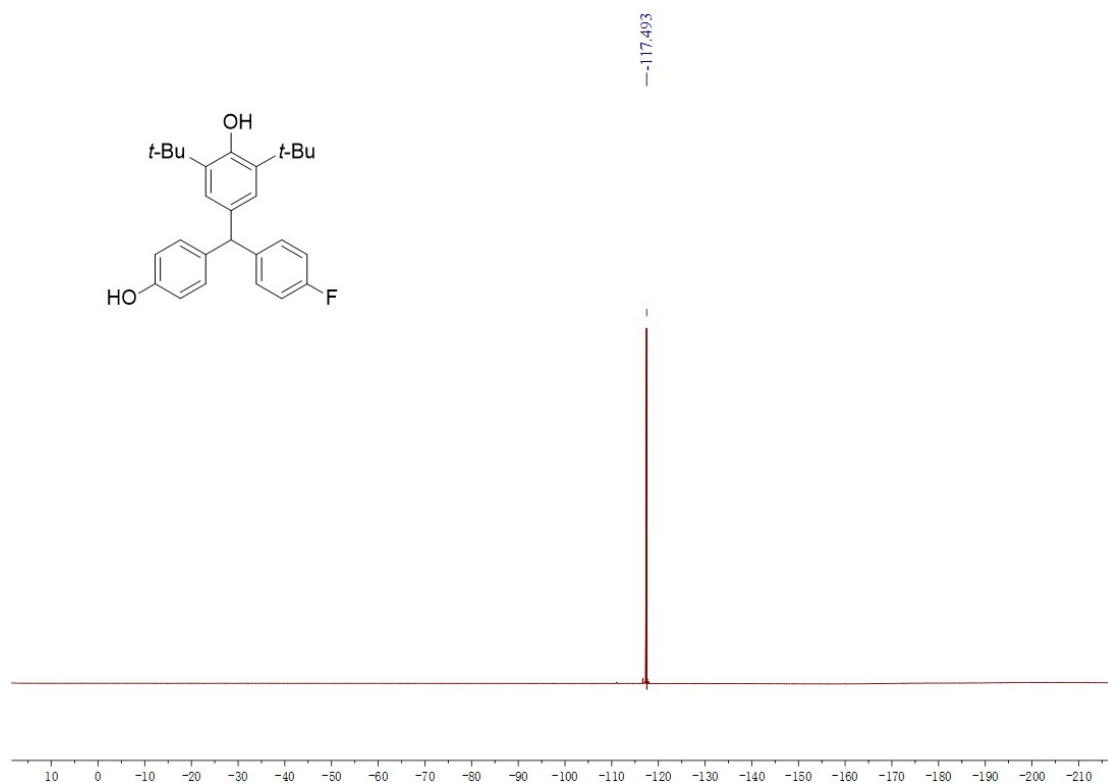
2,6-Di-tert-butyl-4-((2-fluorophenyl)(4-hydroxyphenyl)methyl)phenol (7n)



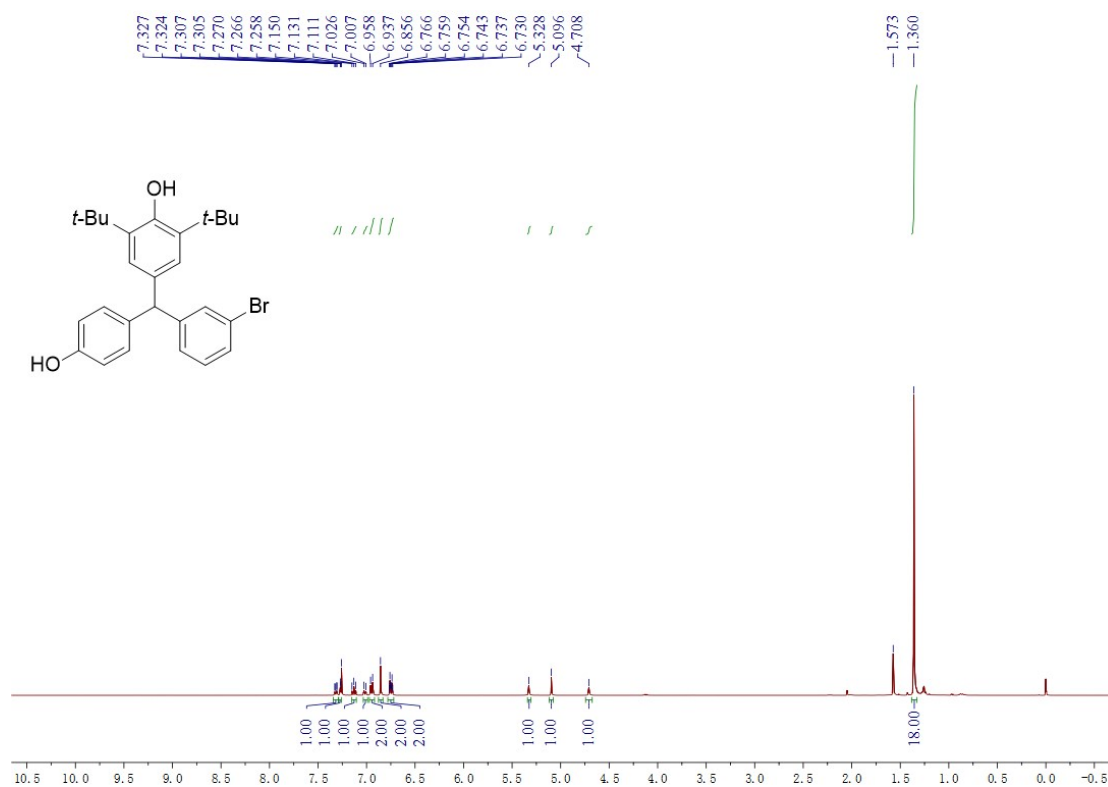


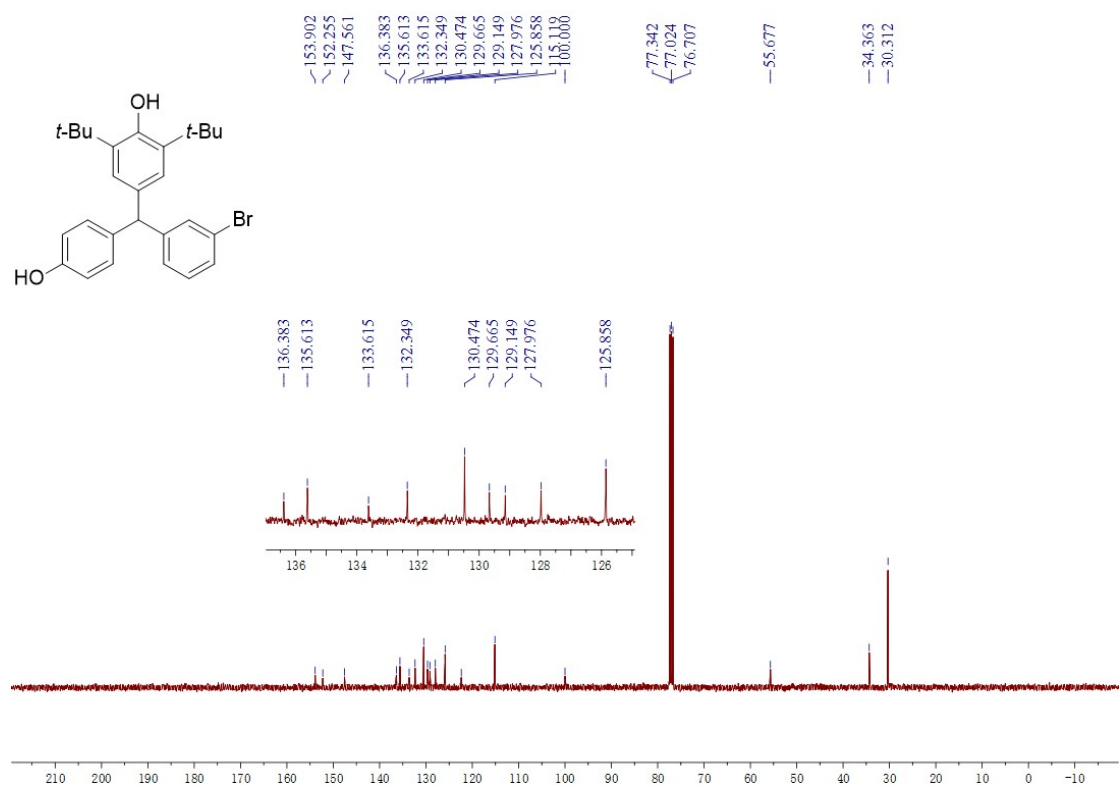
2,6-Di-*tert*-butyl-4-((4-fluorophenyl)(4-hydroxyphenyl)methyl)phenol (7o)



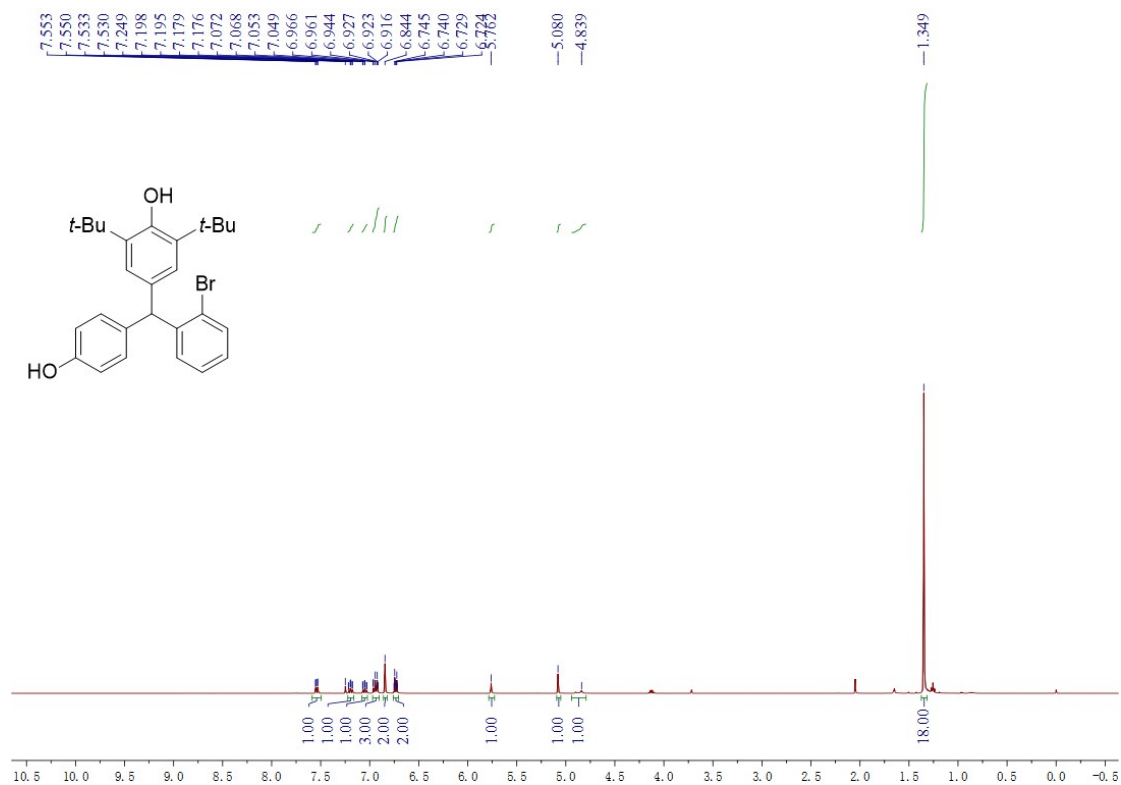


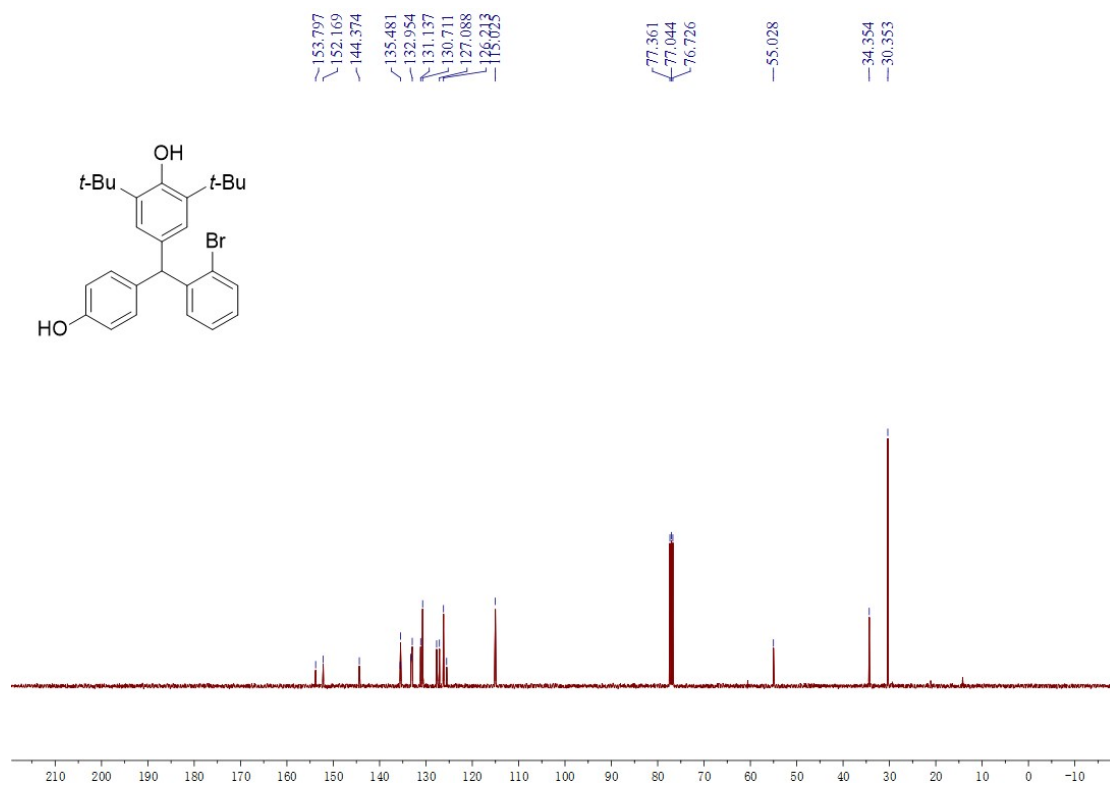
4-((3-Bromophenyl)(4-hydroxyphenyl)methyl)-2,6-di-tert-butylphenol (7p)



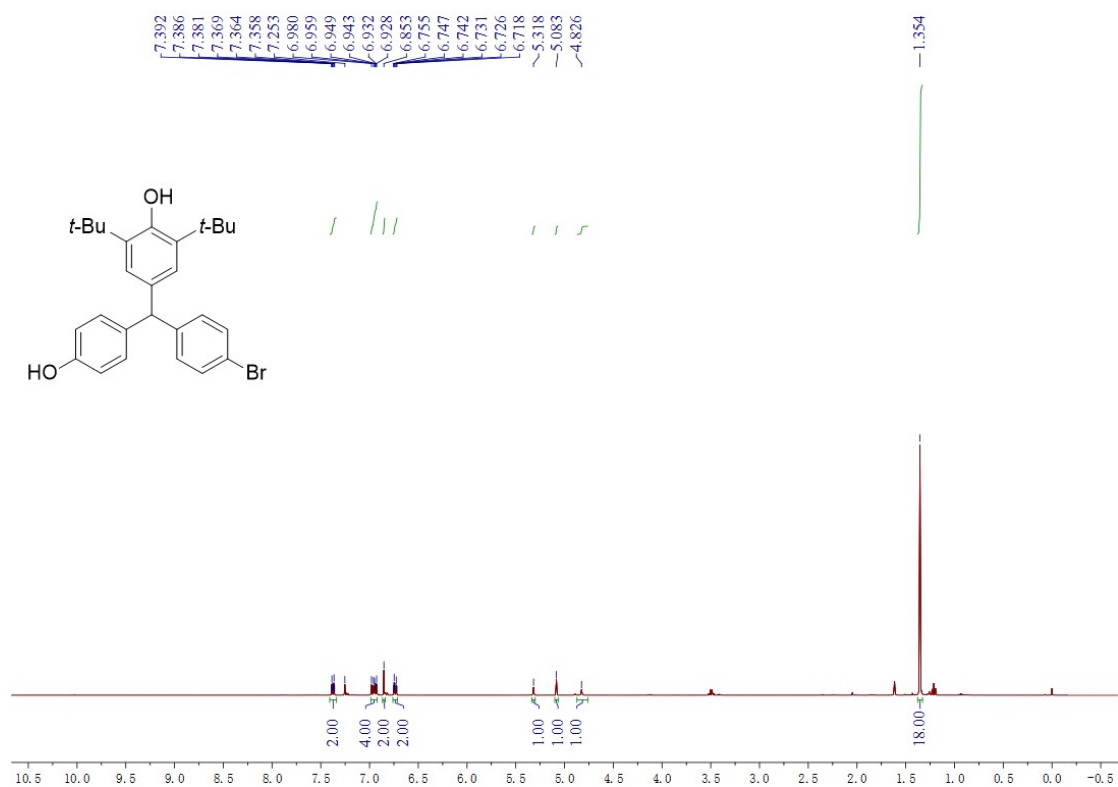


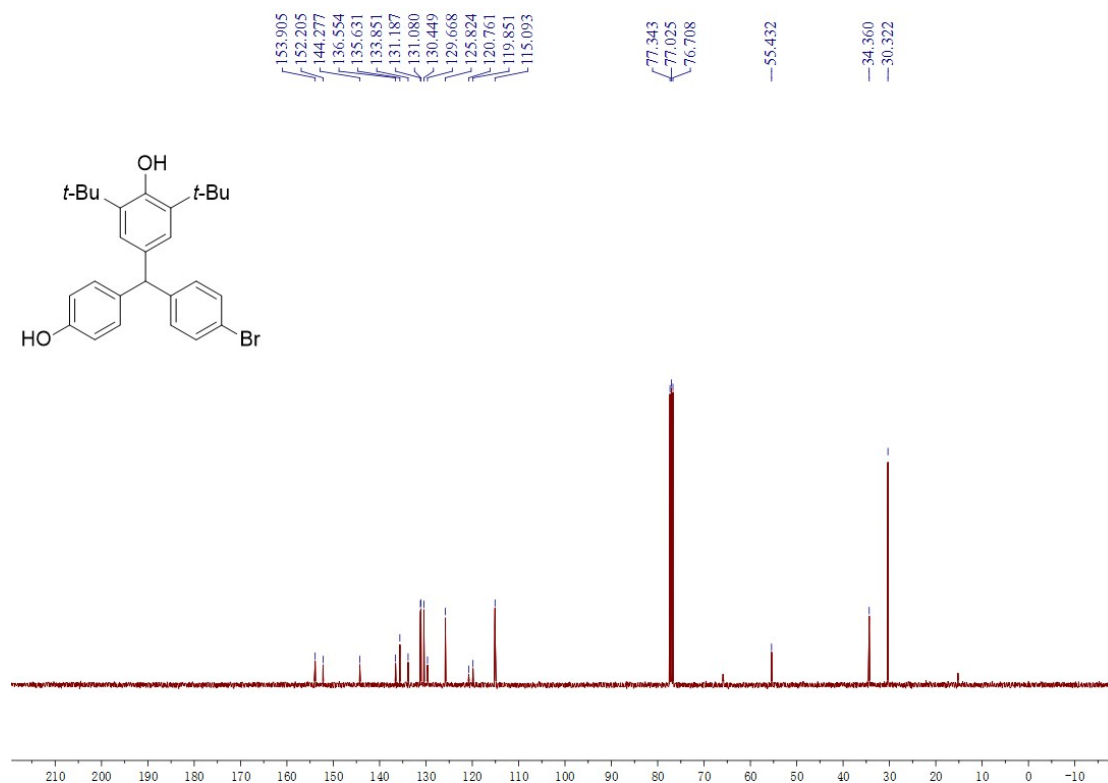
4-((2-Bromophenyl)(4-hydroxyphenyl)methyl)-2,6-di-tert-butylphenol (7q)



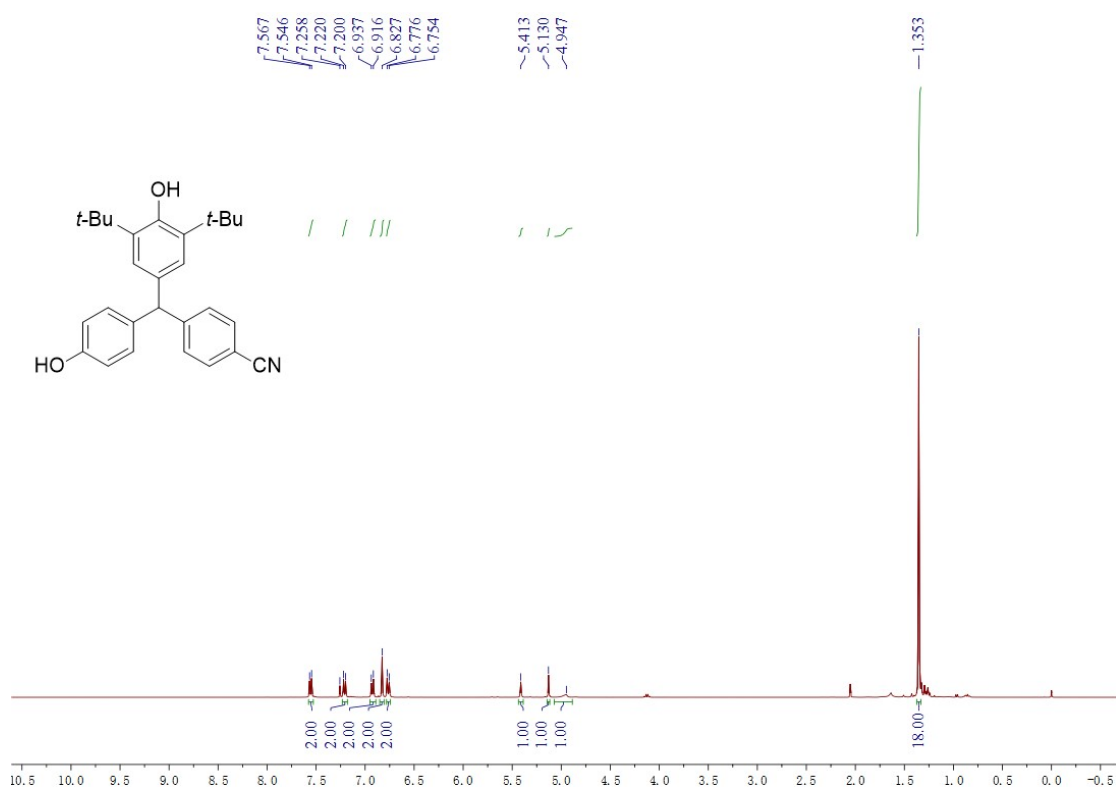


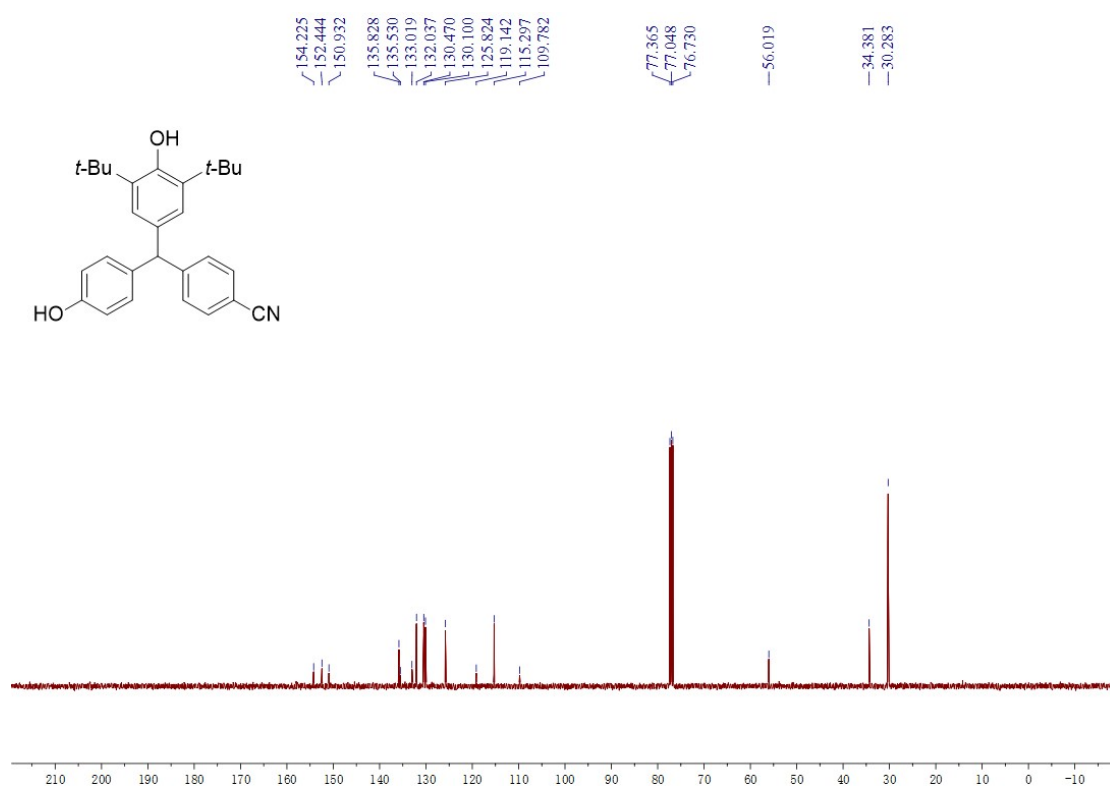
4-((4-Bromophenyl)(4-hydroxyphenyl)methyl)-2,6-di-tert-butylphenol (7r)



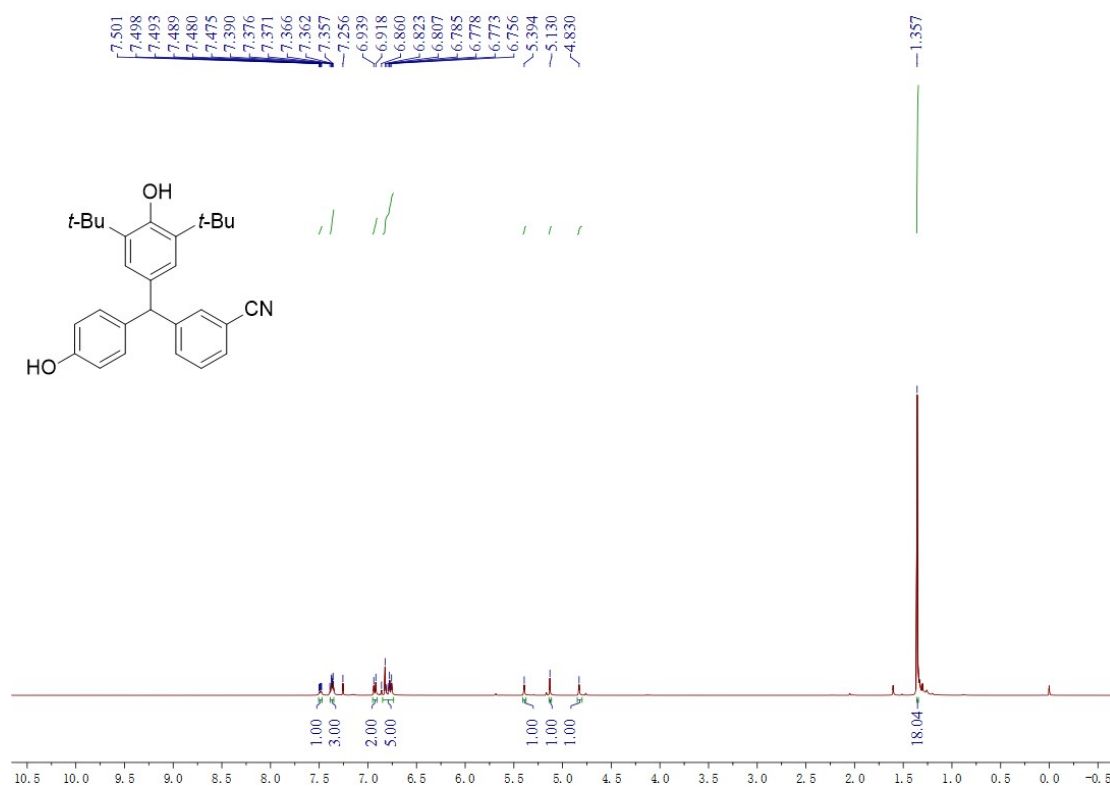


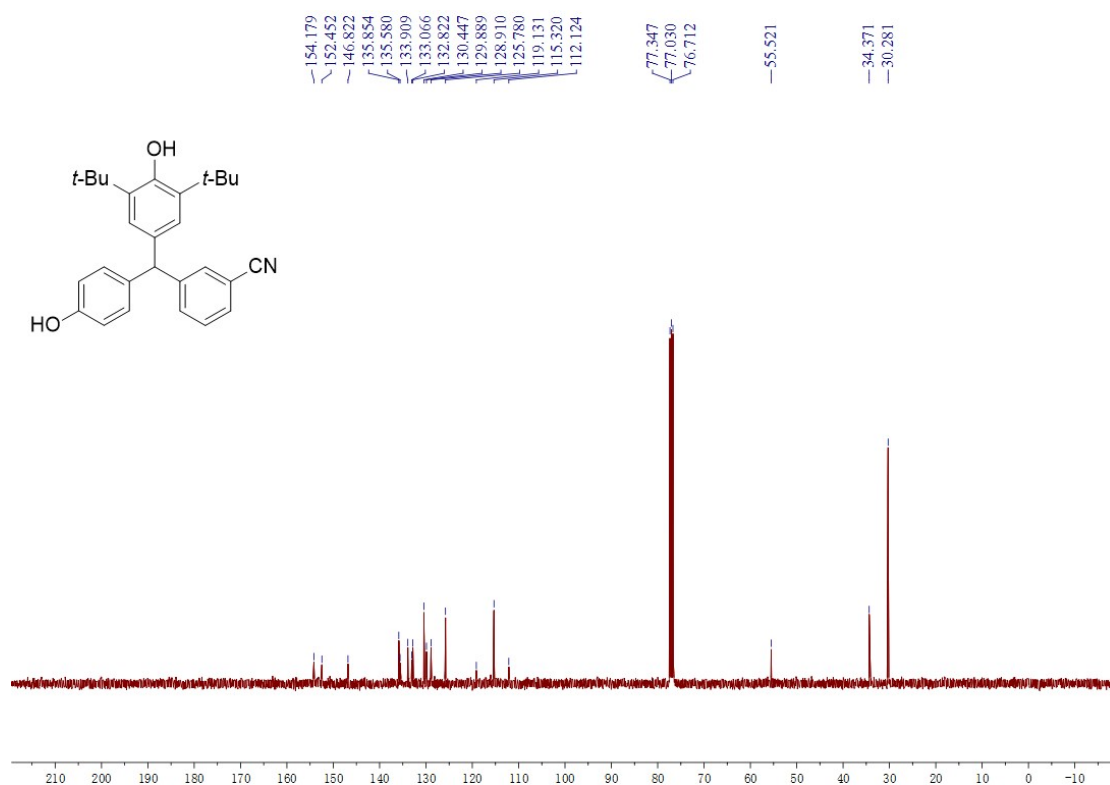
4-((3,5-Di-tert-butyl-4-hydroxyphenyl)(4-hydroxyphenyl)methyl)benzonitrile (7s)



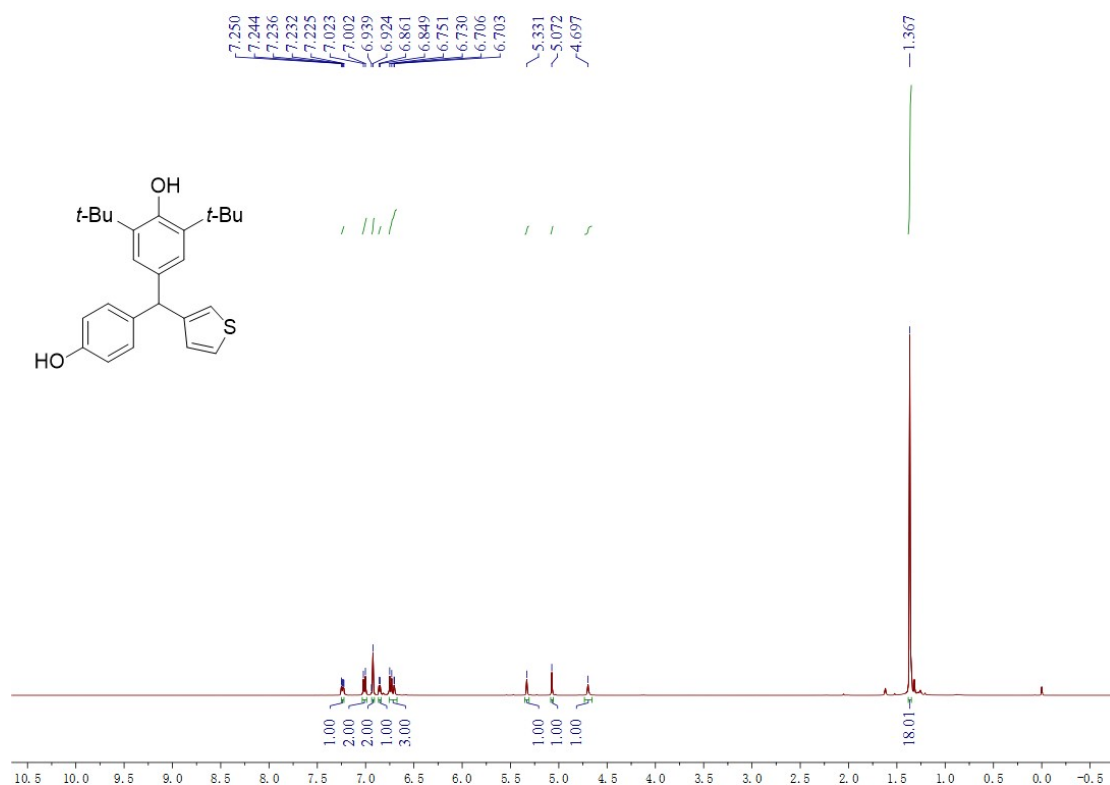


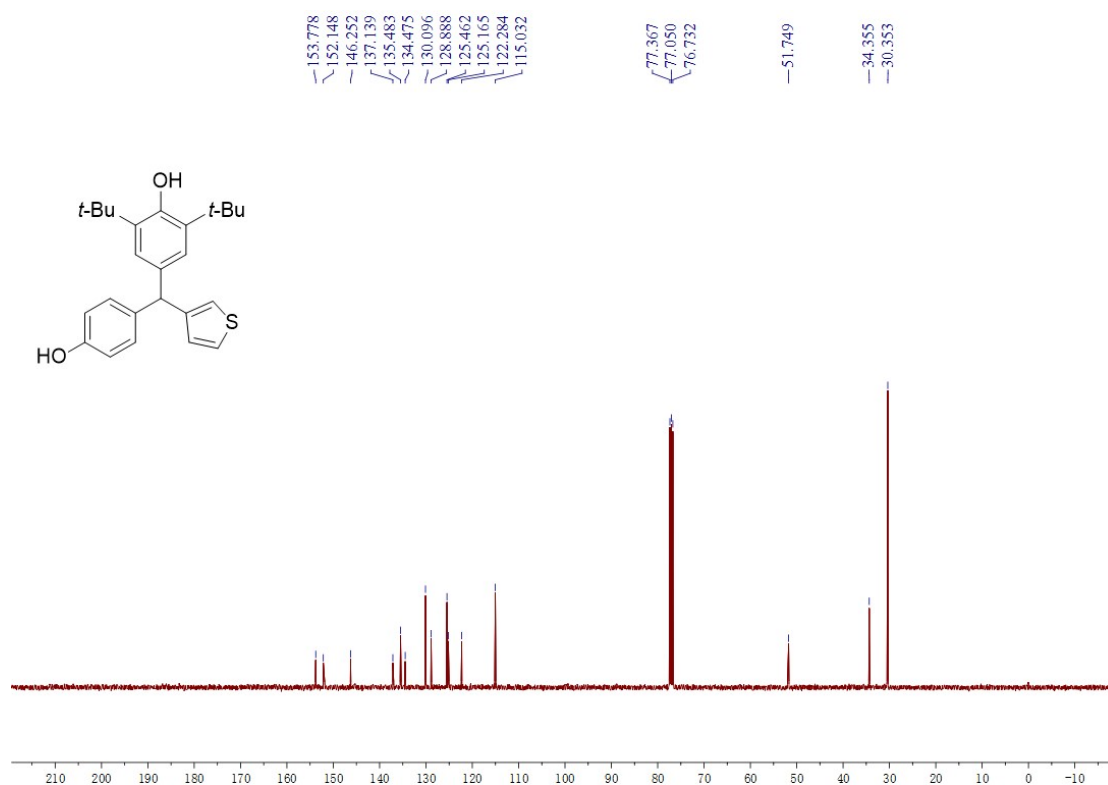
3-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(4-hydroxyphenyl)methyl)benzonitrile (7t)



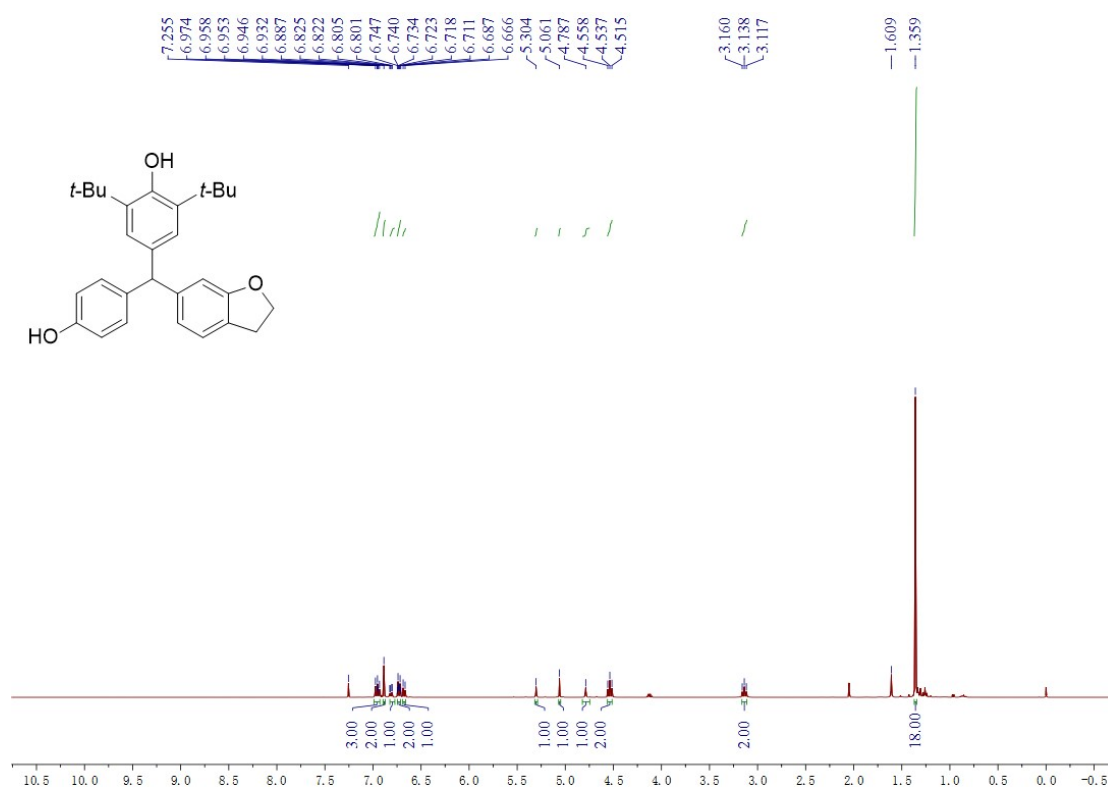


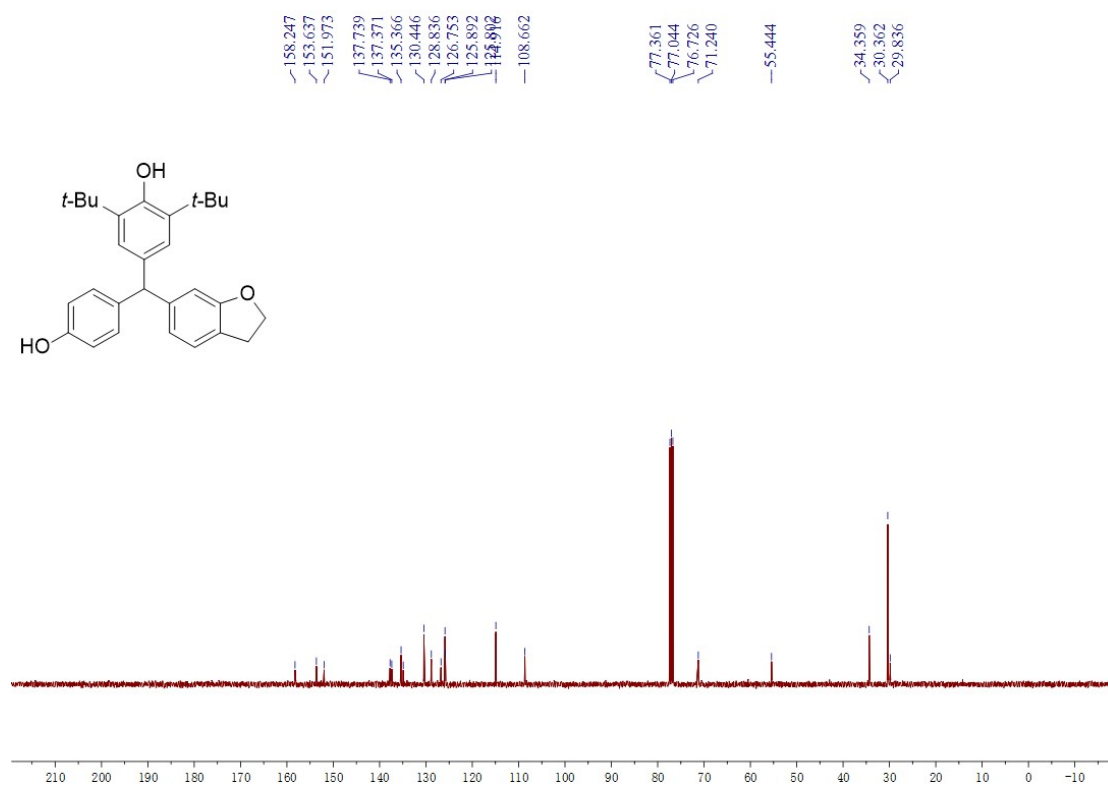
2,6-Di-tert-butyl-4-((4-hydroxyphenyl)(thiophen-3-yl)methyl)phenol (7u)



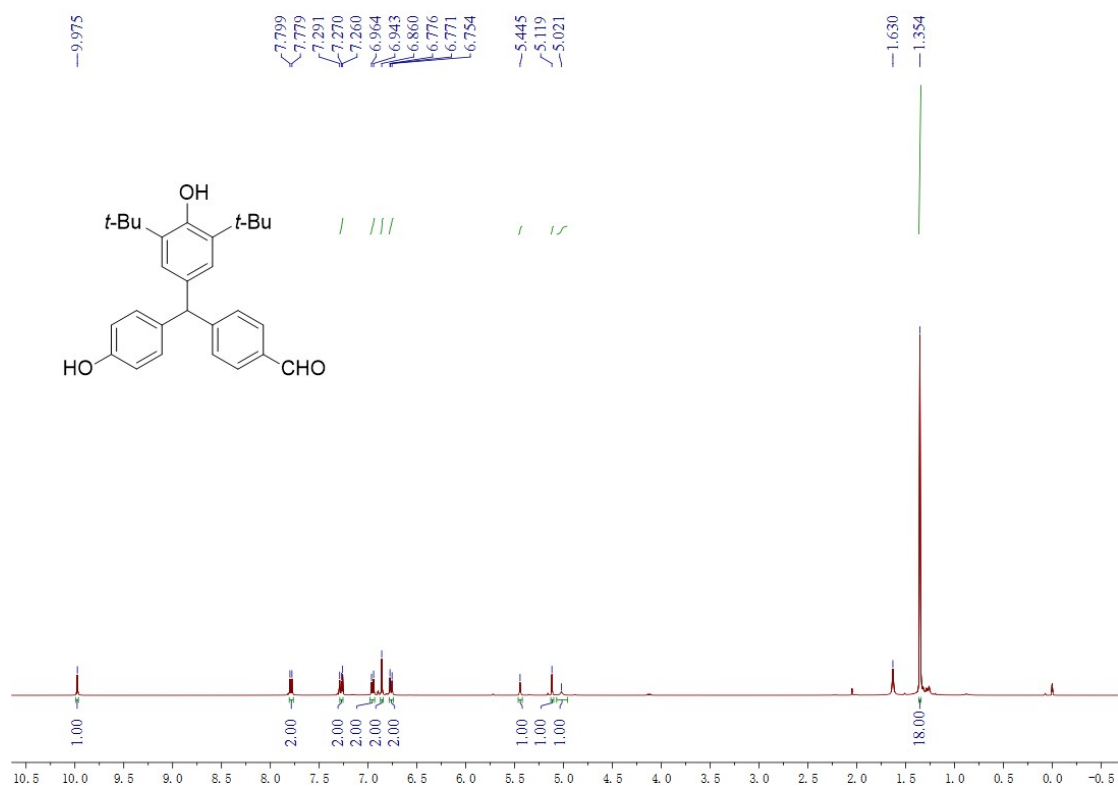


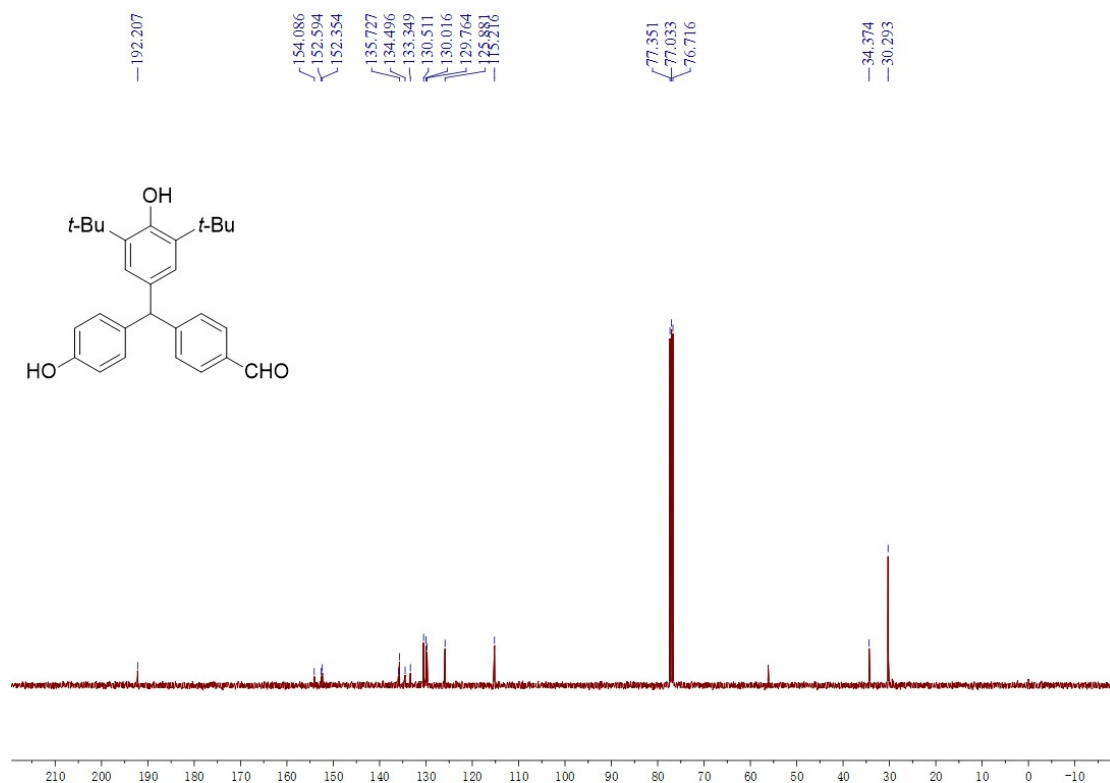
2,6-Di-tert-butyl-4-((2,3-dihydrobenzofuran-6-yl)(4-hydroxyphenyl)methyl)phenol (7x)



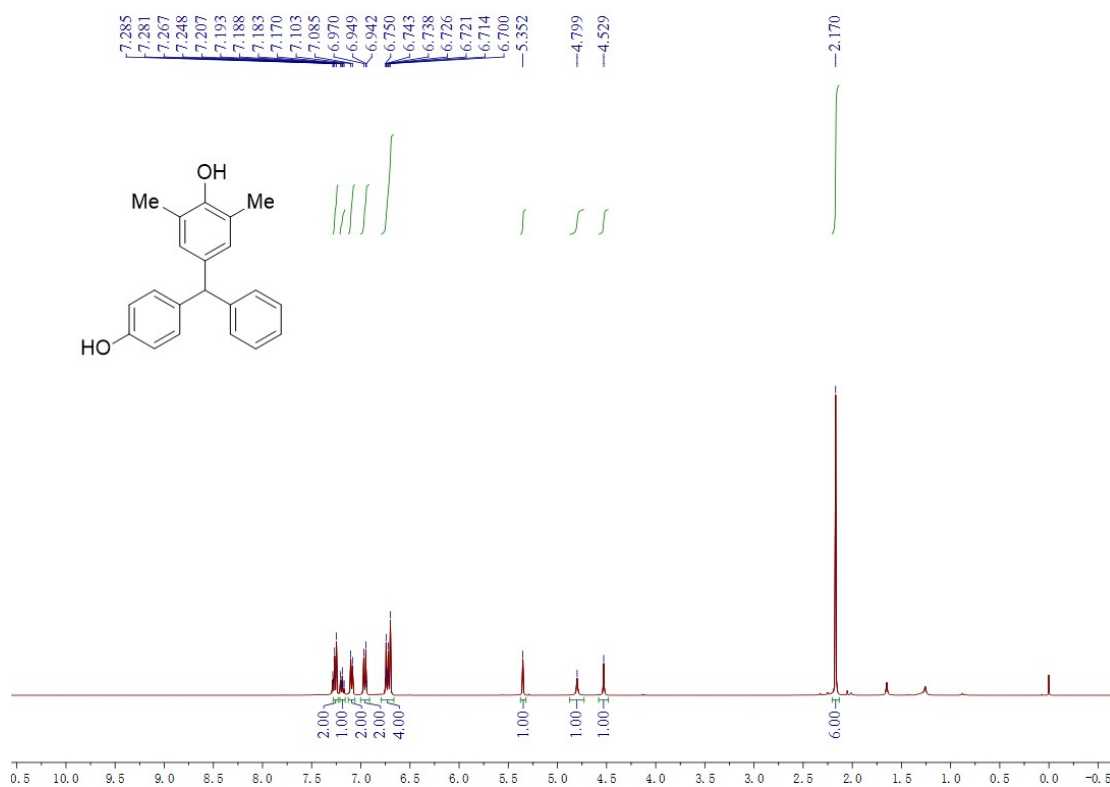


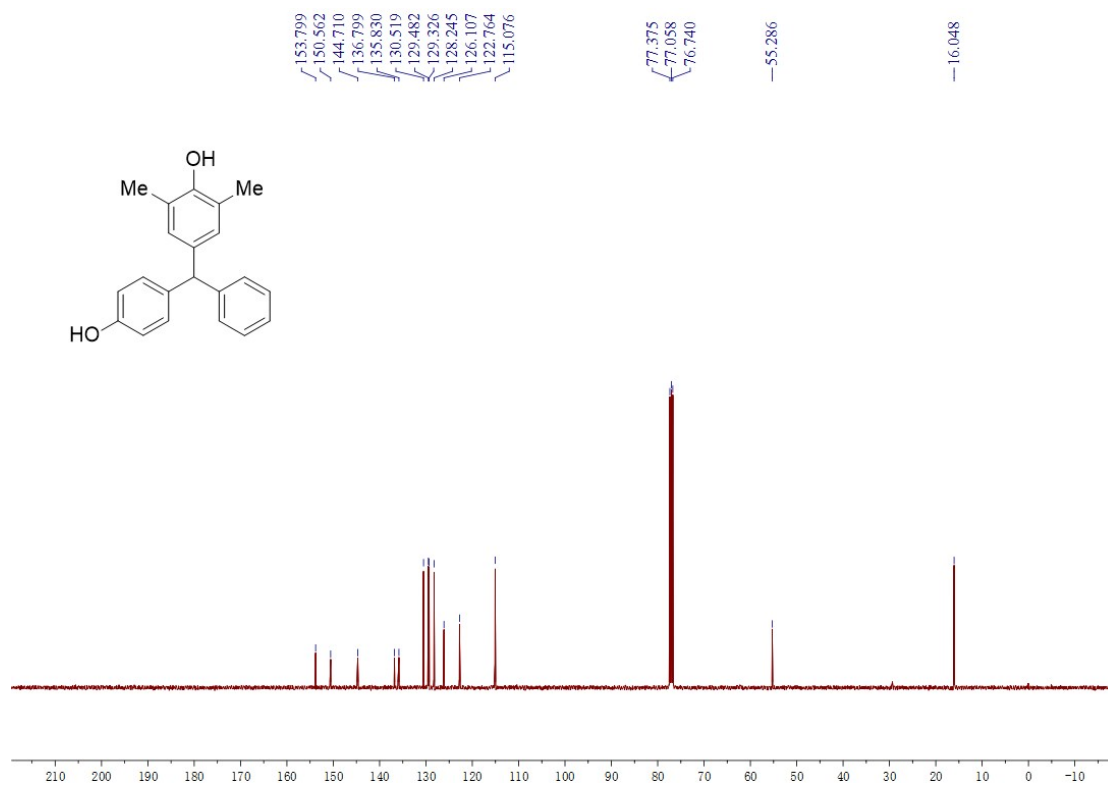
4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(4-hydroxyphenyl)methyl)benzaldehyde (7z)



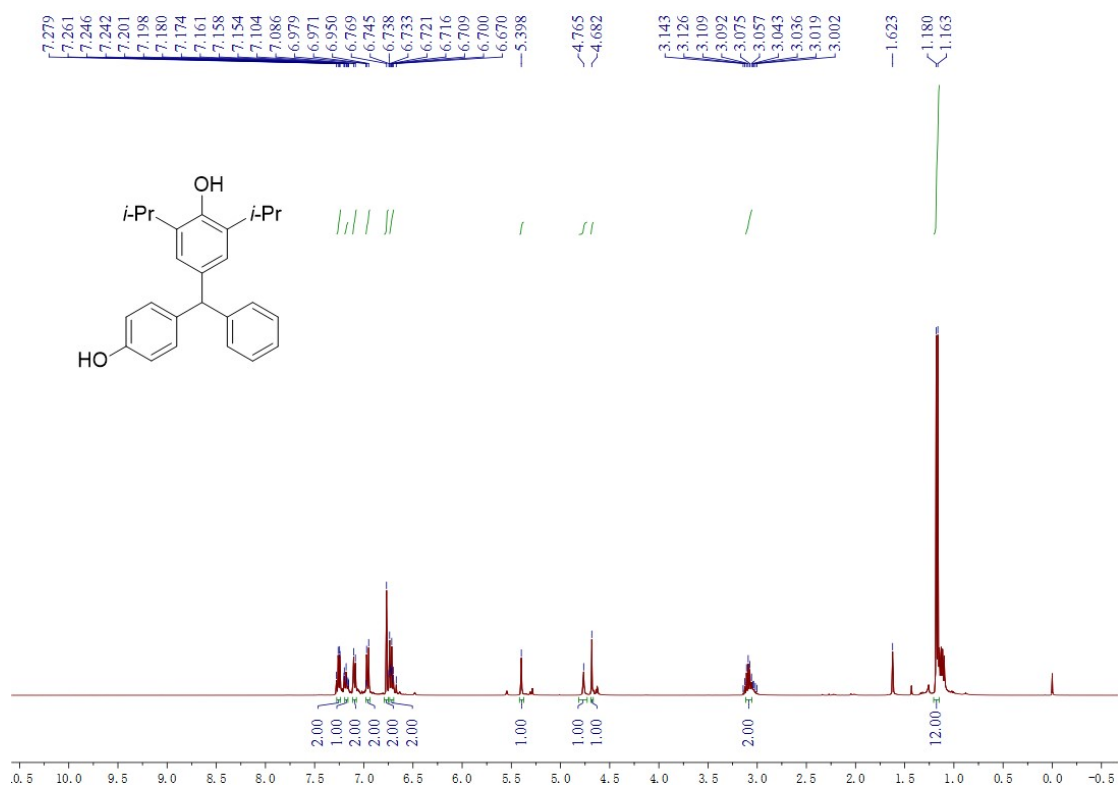


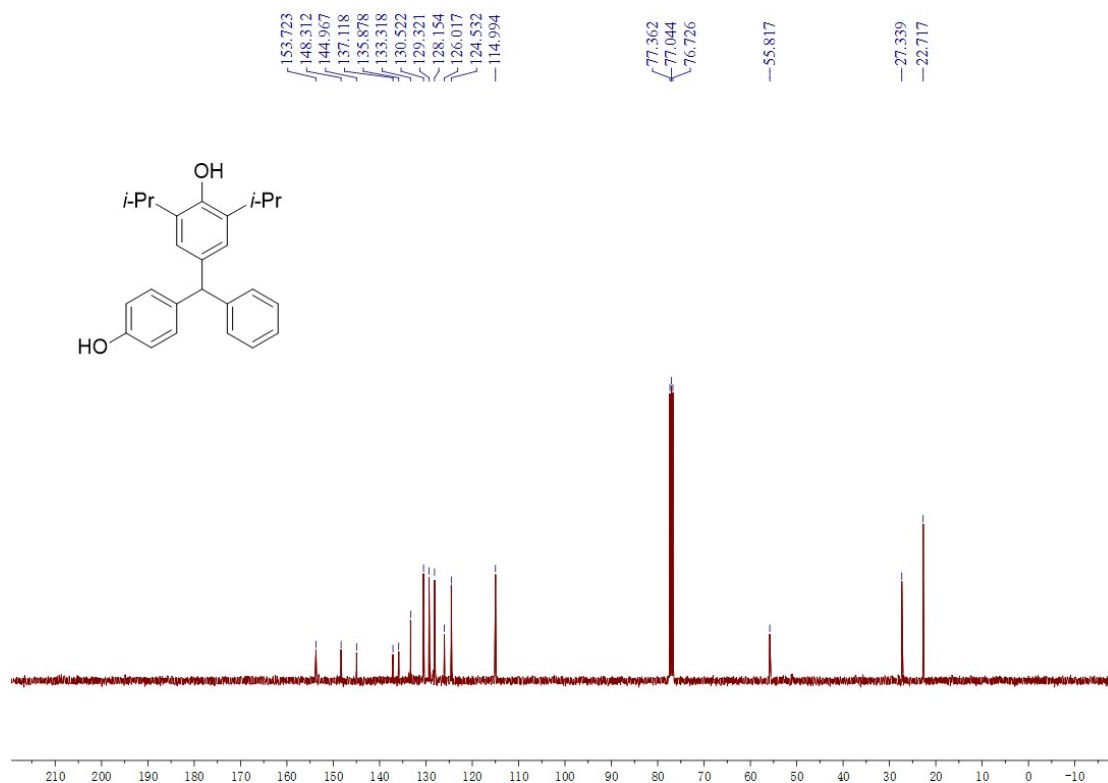
4-((4-Hydroxyphenyl)(phenyl)methyl)-2,6-dimethylphenol (7ab)



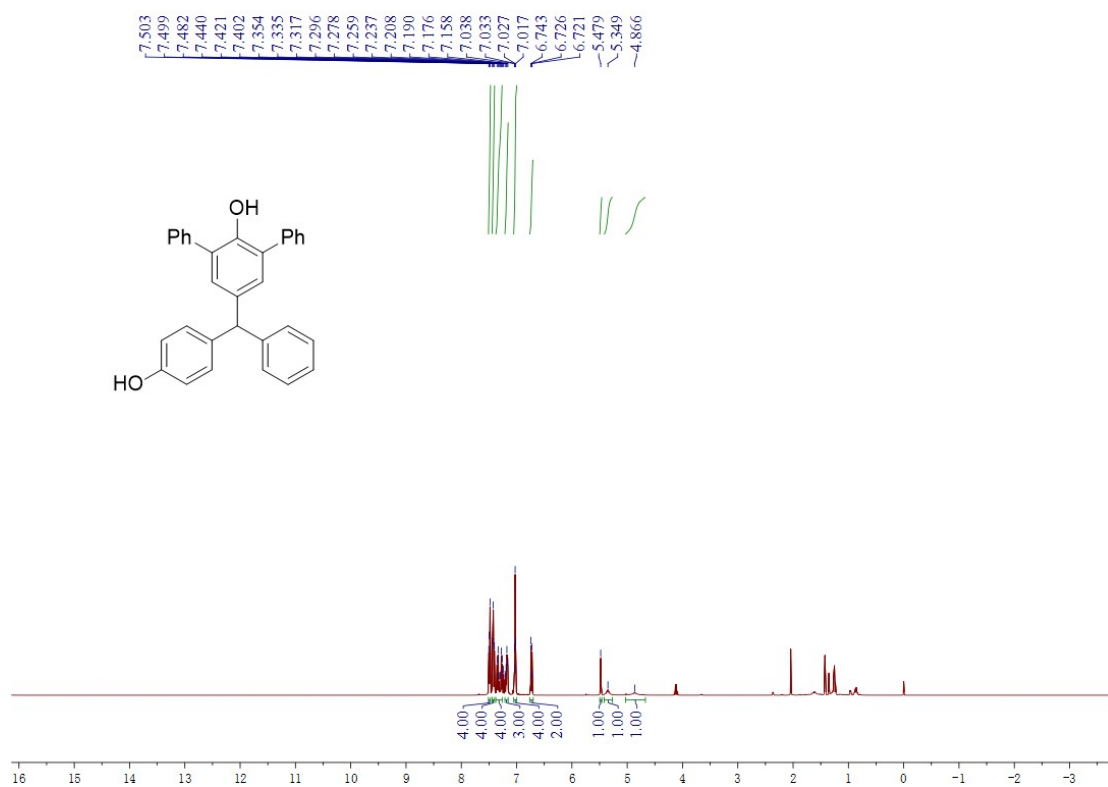


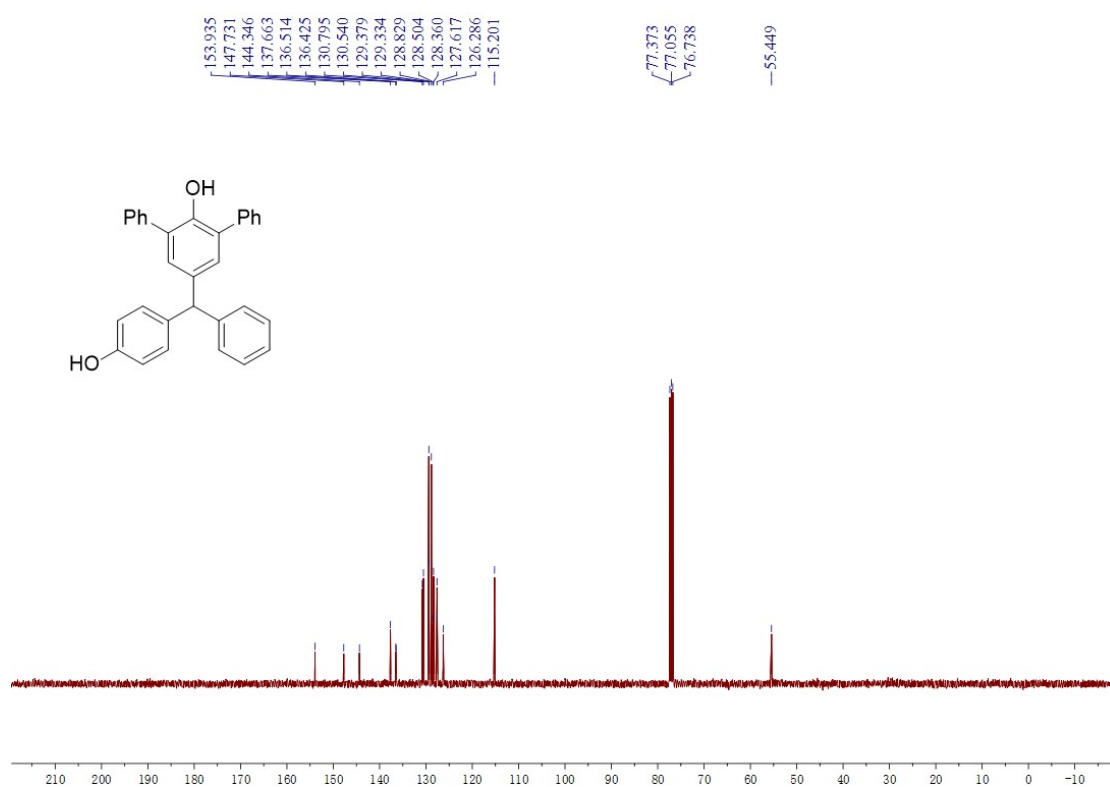
4-((4-Hydroxyphenyl)(phenyl)methyl)-2,6-diisopropylphenol (7ac)



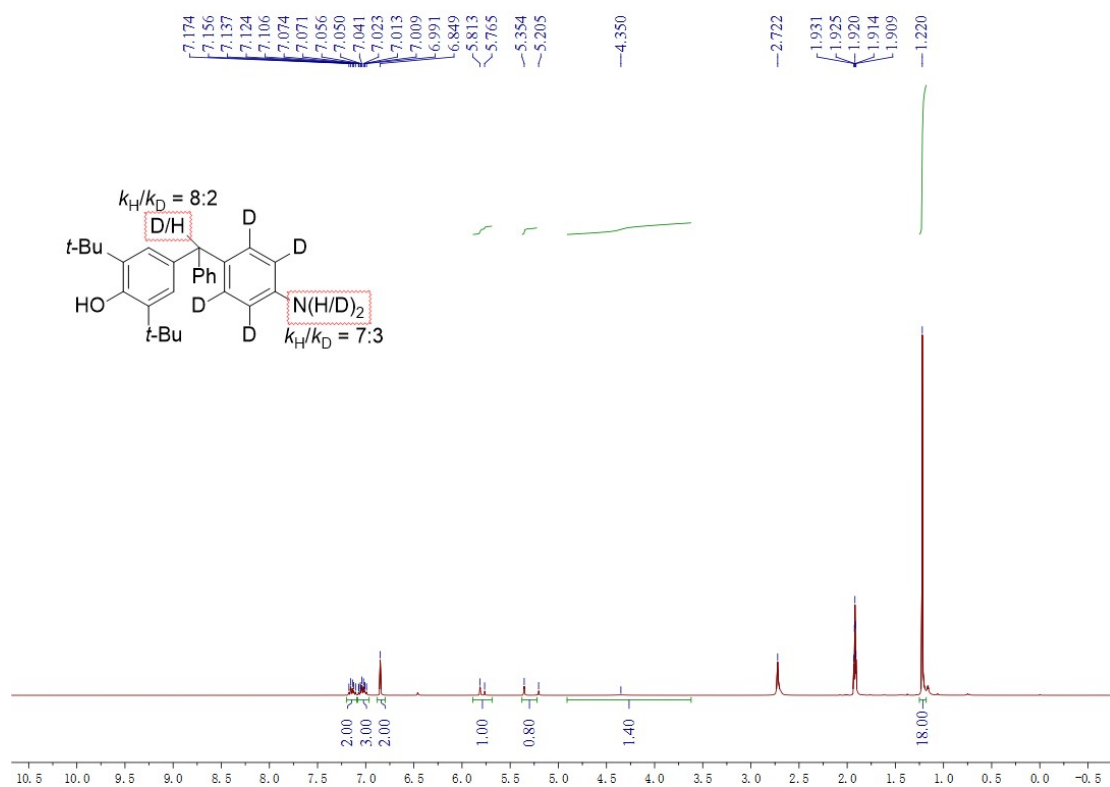


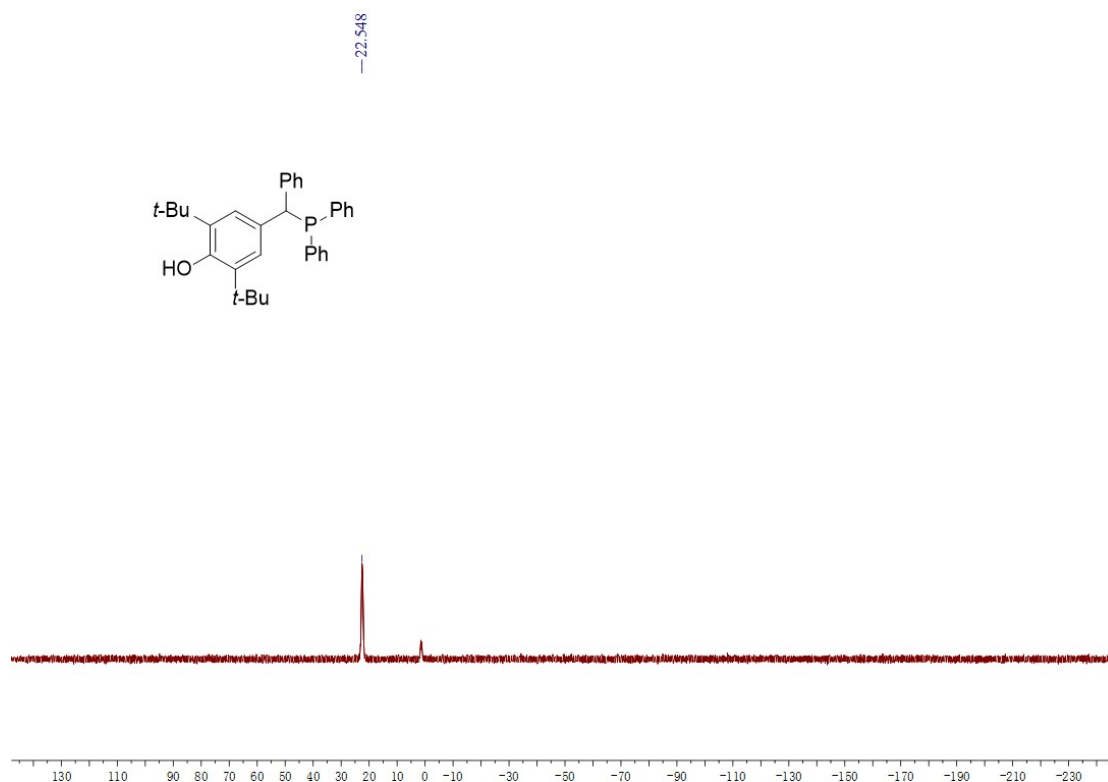
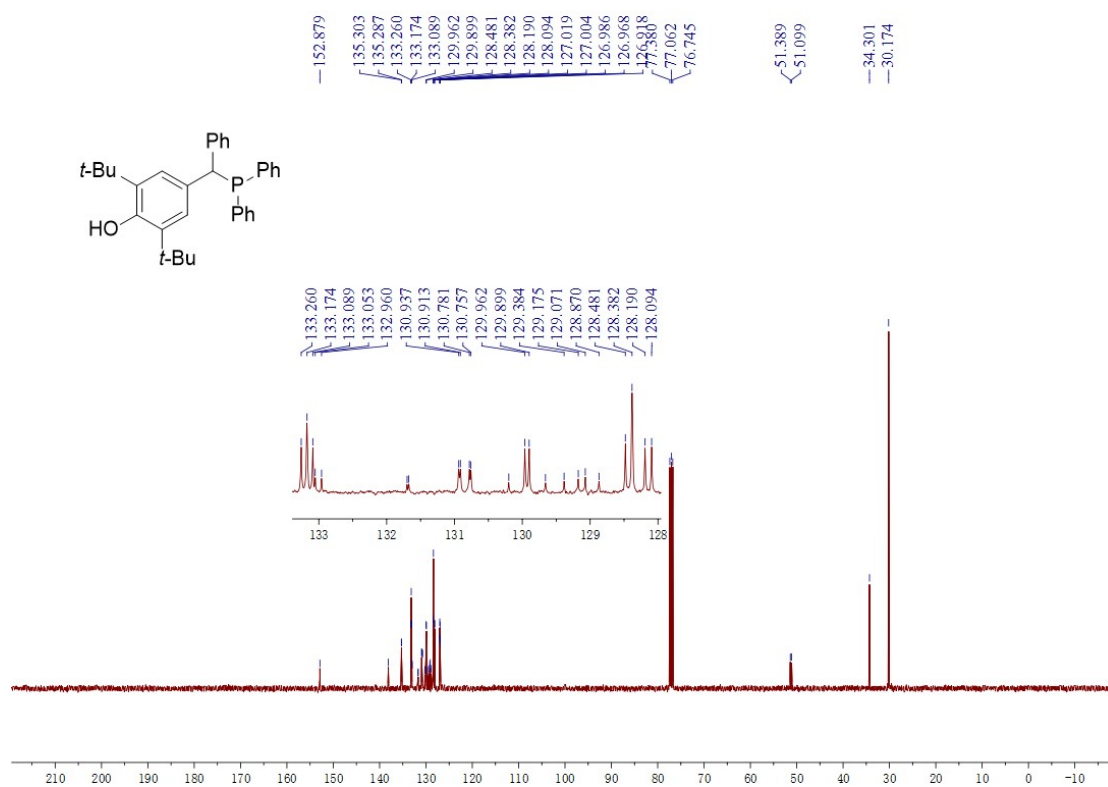
5'-((4-Hydroxyphenyl)(phenyl)methyl)-[1,1':3',1''-terphenyl]-2'-ol (7ad)



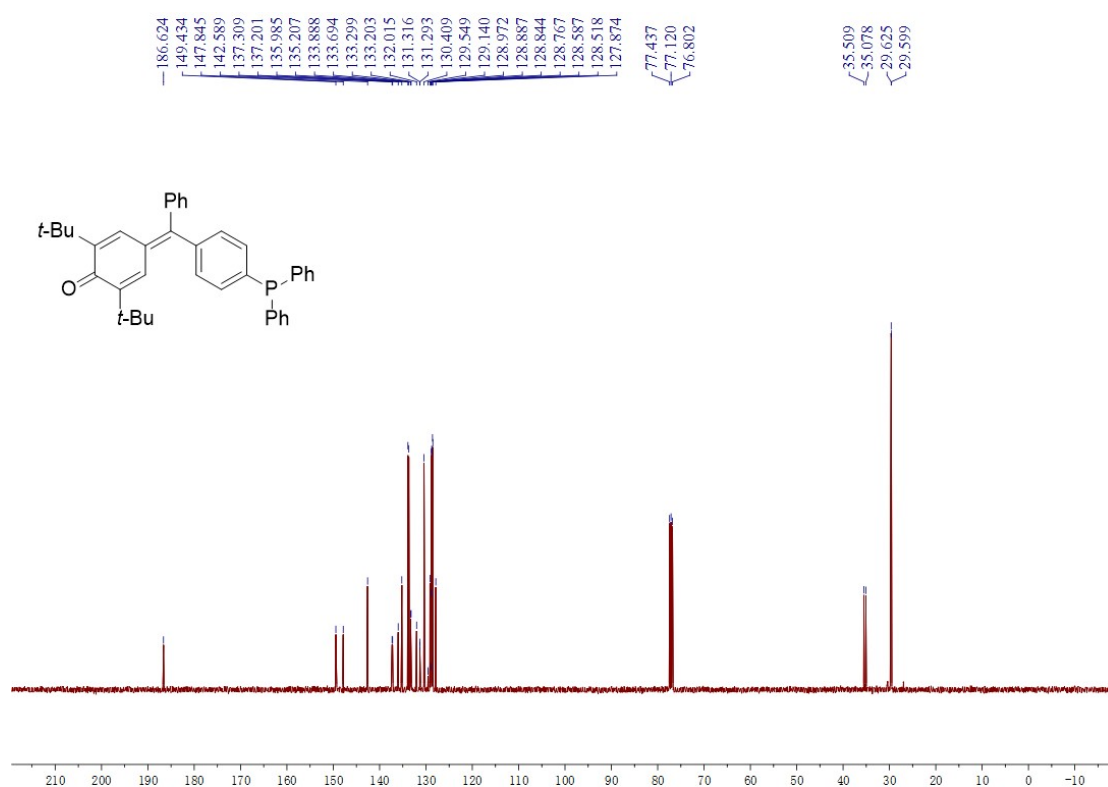
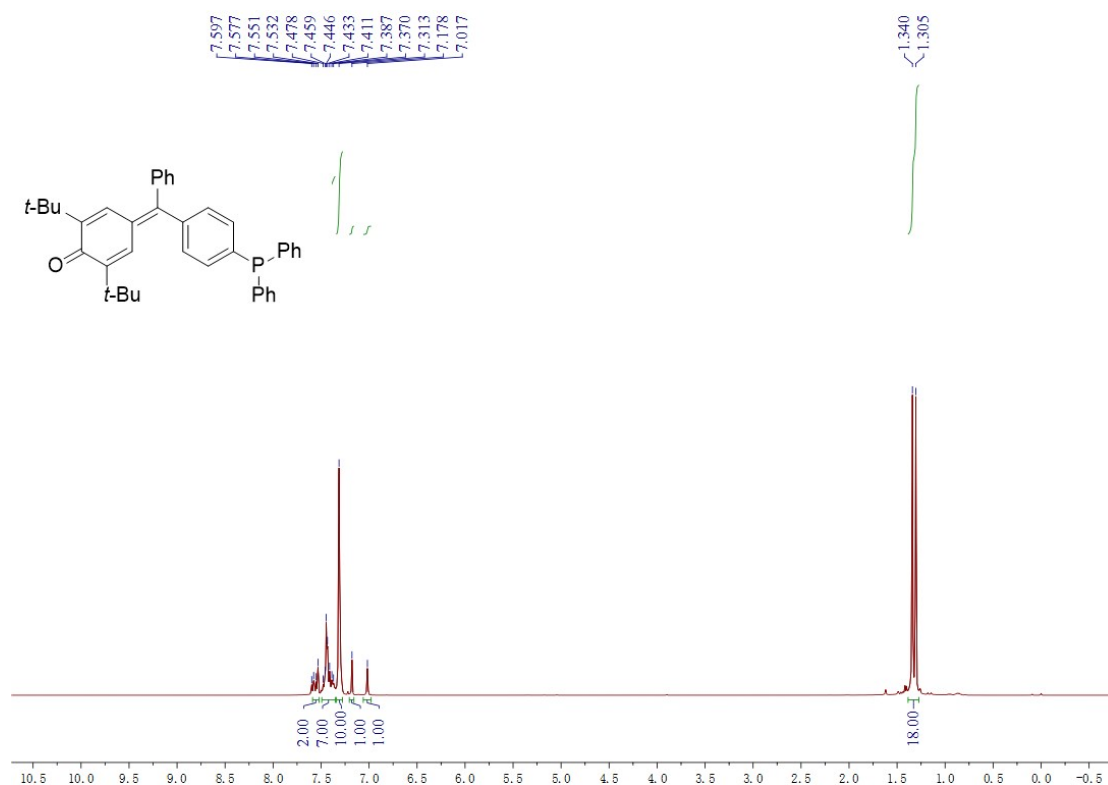


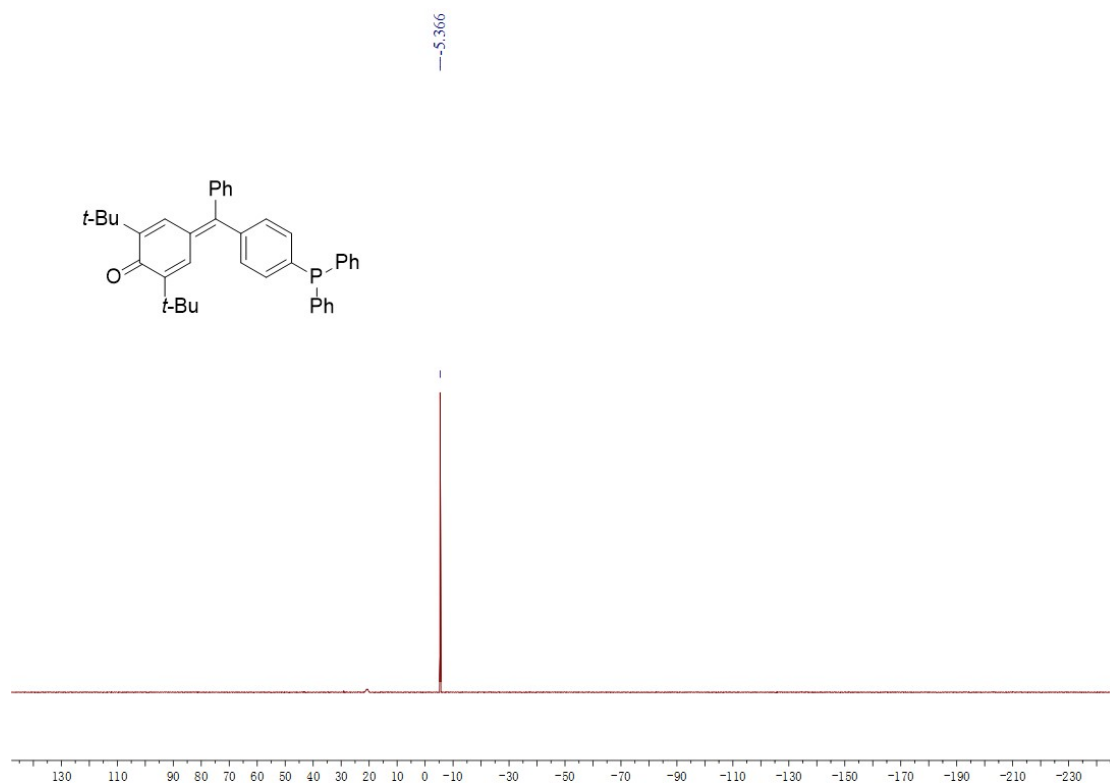
4-((4-Aminophenyl)(phenyl)methyl)-2,6-di-tert-butylphenol (3a')



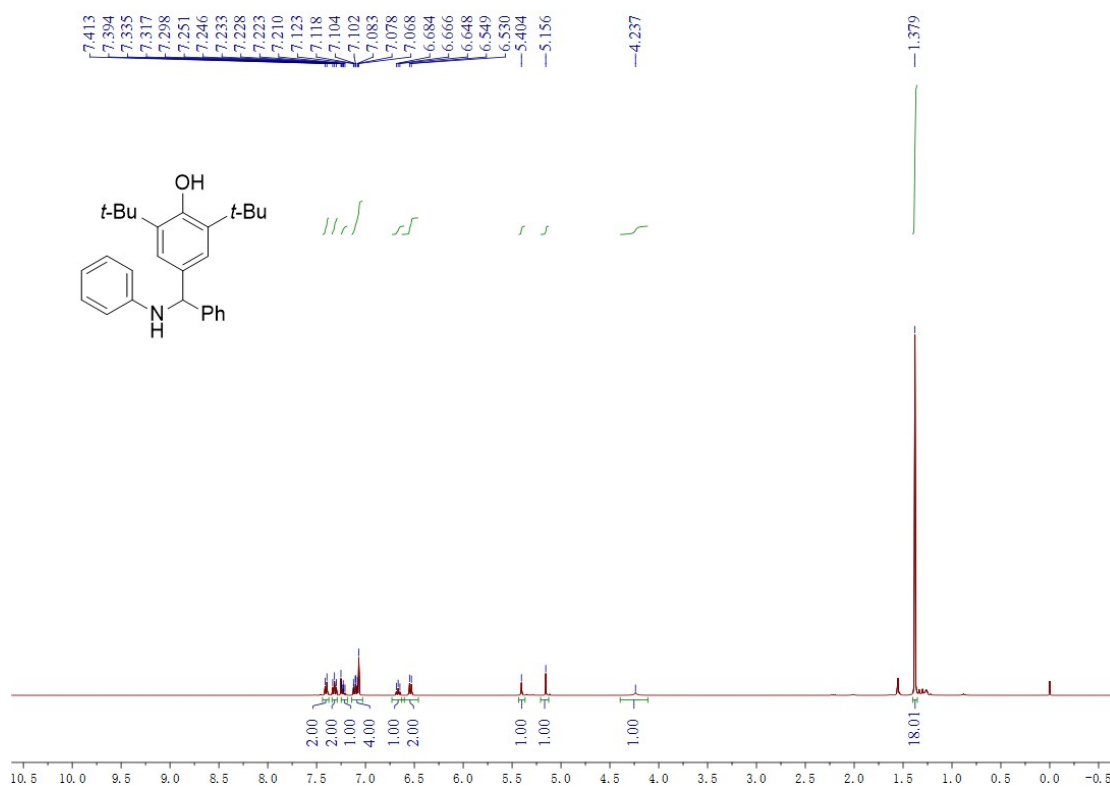


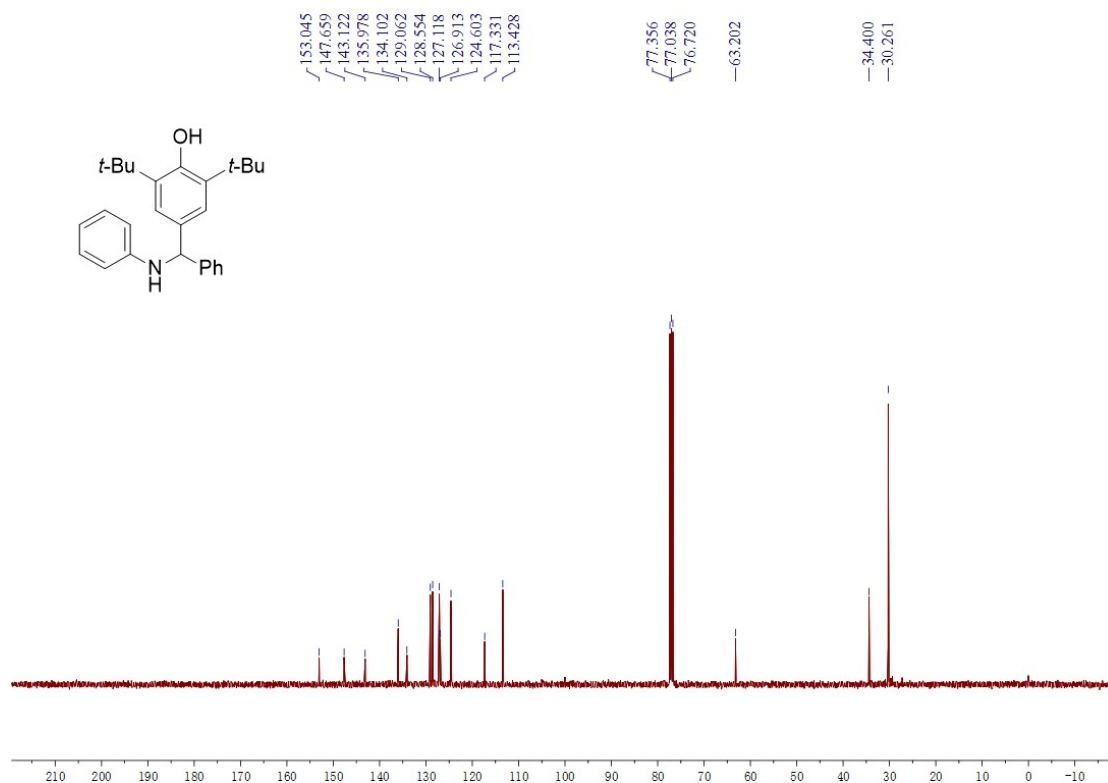
2,6-Di-*tert*-butyl-4-((4-(diphenylphosphanyl)phenyl)(phenyl)methylene)cyclohexa-2,5-dien-1-one (9b)





2,6-Di-tert-butyl-4-(phenyl(phenylamino)methyl)phenol (3a'')





Copies of HPLC spectra

