

Stabilizing Pd on the surface of hollow magnetic mesoporous spheres: a highly active and recyclable catalyst for hydrogenation and Suzuki coupling reaction

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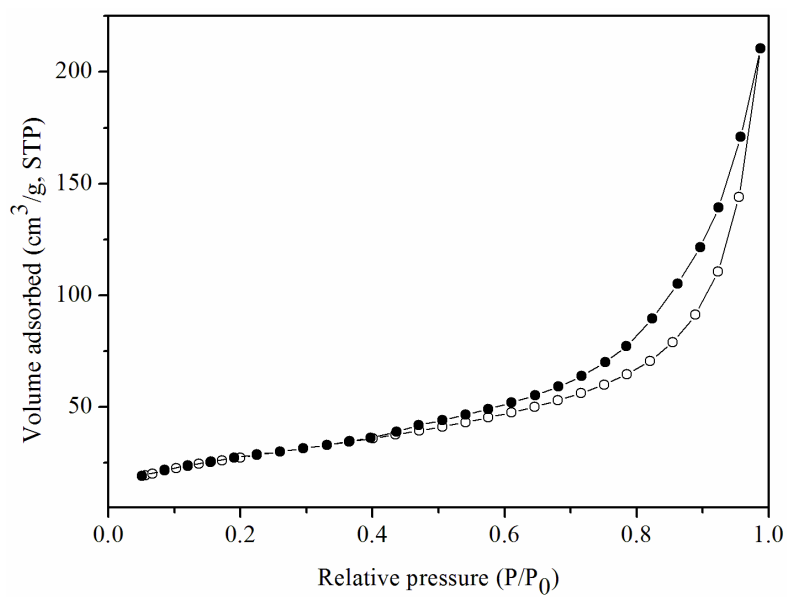


Fig.S1 N₂ sorption isotherms of HMMS-NH₂-Pd catalyst

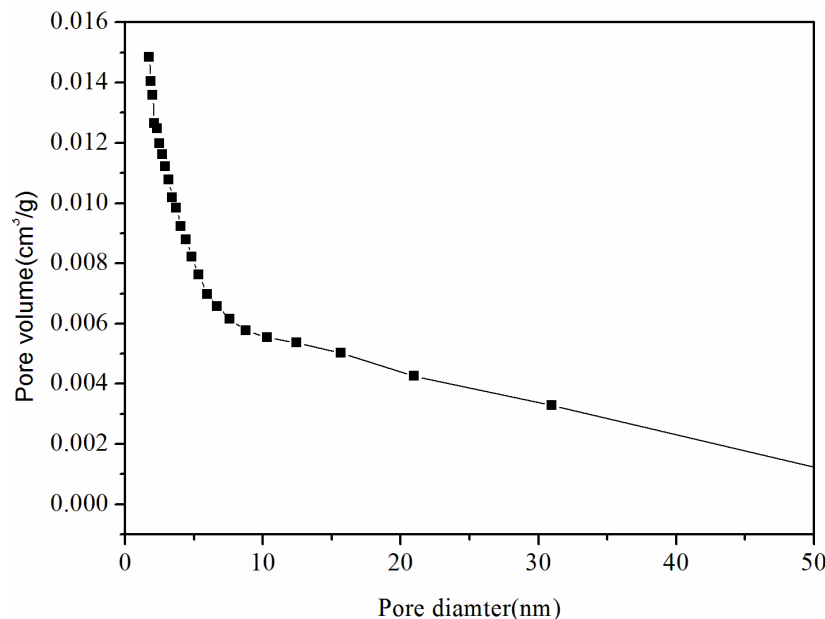


Fig. S2 The pore size distribution of HMMS-NH₂-Pd catalyst

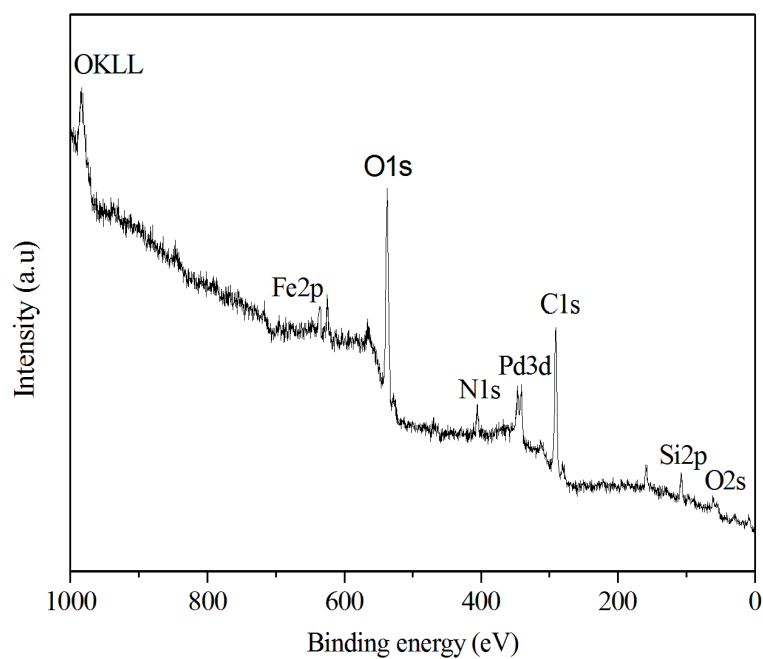


Fig. S3 XPS spectrum of the elemental survey scan of HMMS-NH₂-Pd

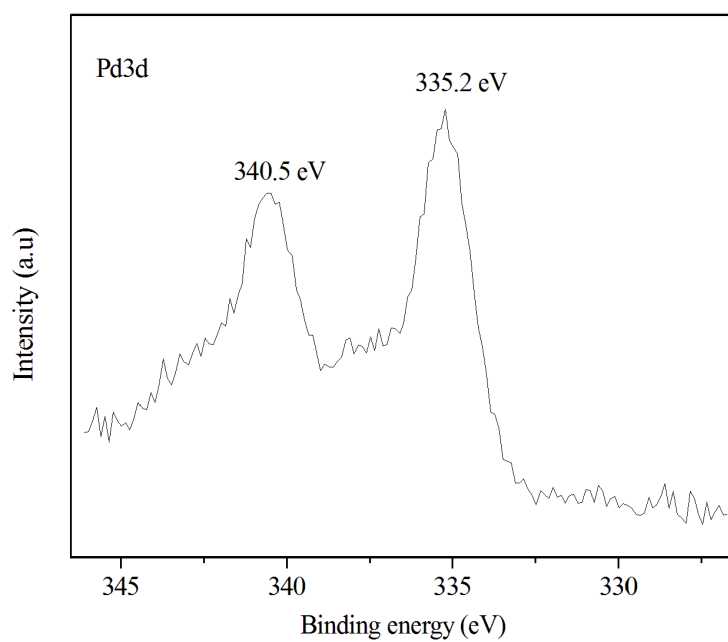


Fig. S4 XPS spectrum of the HMMS-NH₂-Pd showing Pd 3d_{5/2} and Pd 3d_{3/2} binding energies

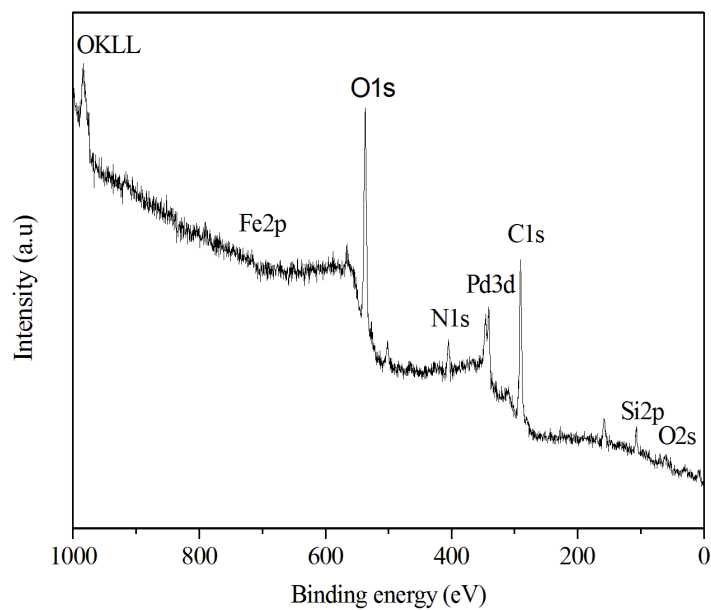


Fig. S5 XPS spectrum of the elemental survey scan of reused HMMS-NH₂-Pd

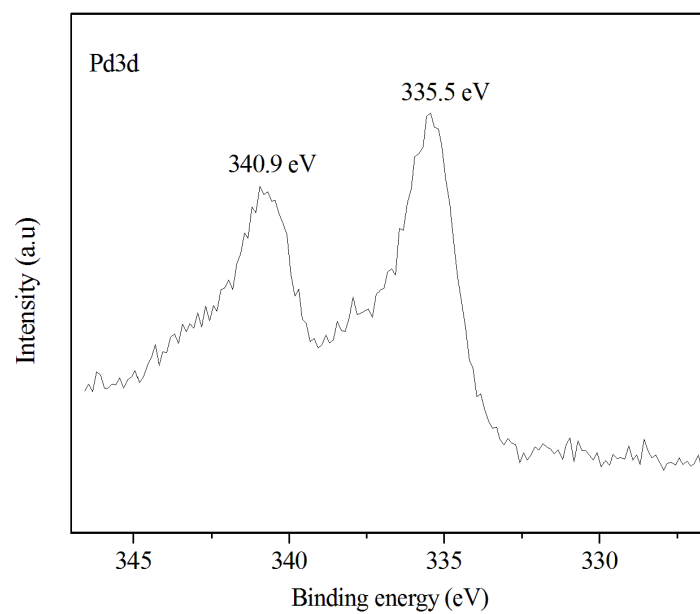
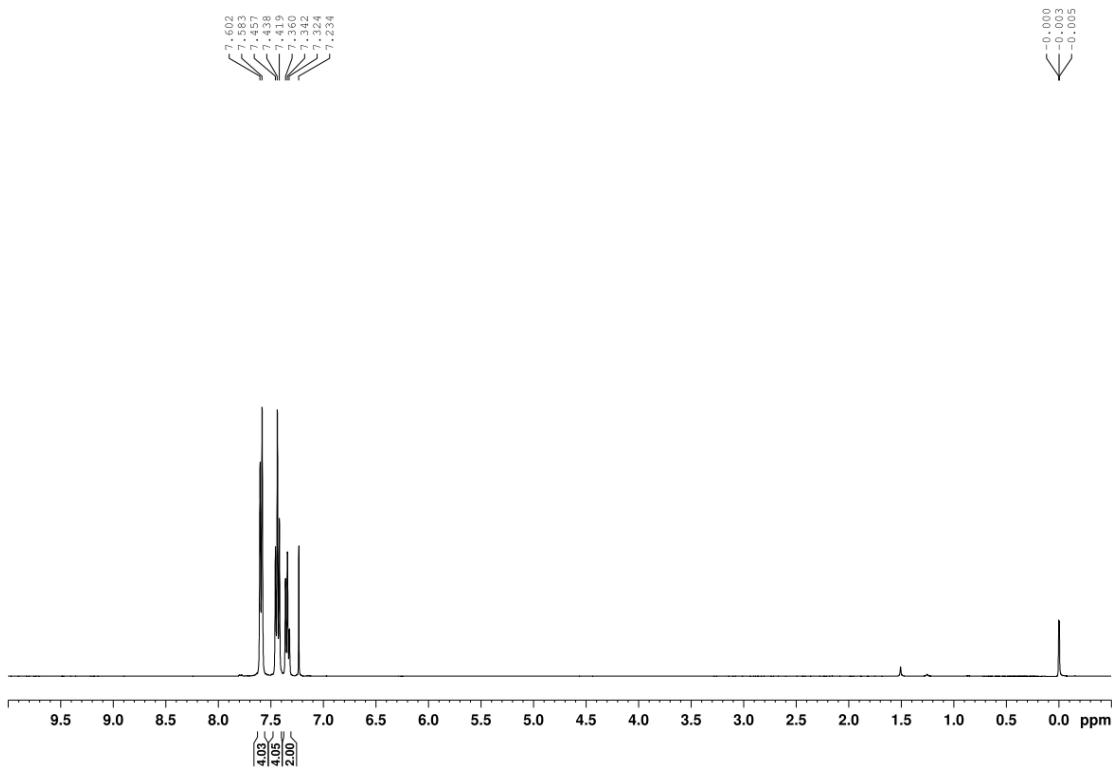
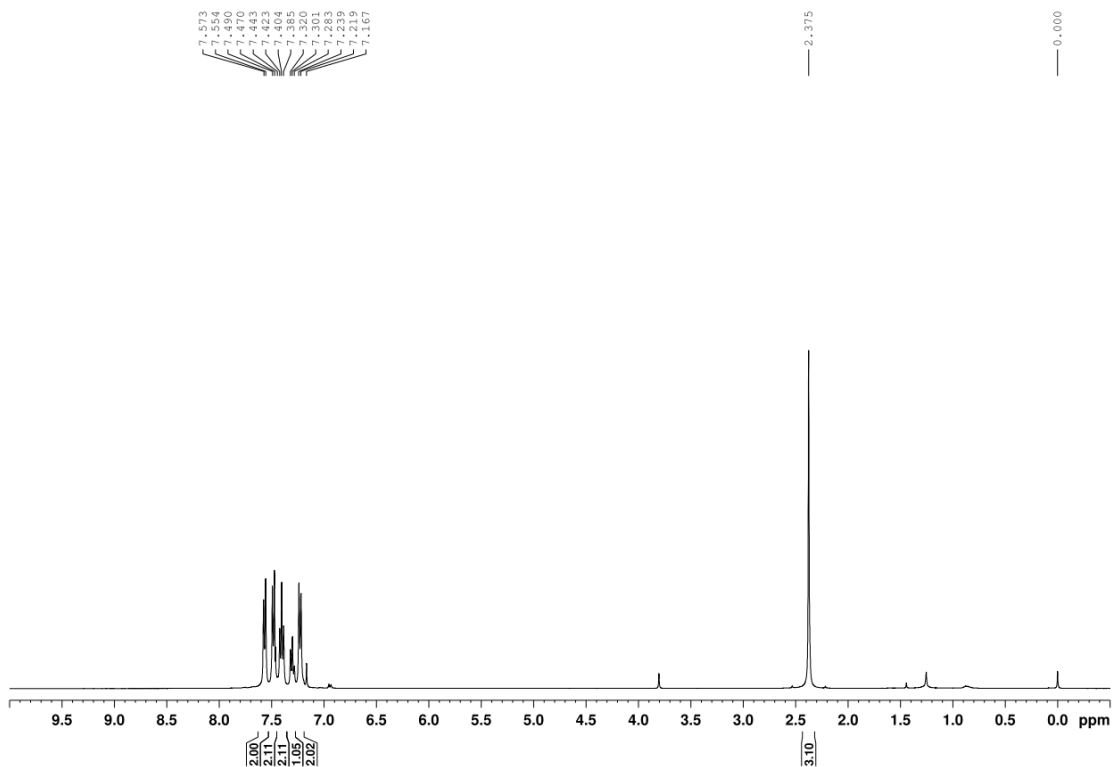


Fig. S6 XPS spectrum of the reused HMMS-NH₂-Pd showing Pd 3d_{5/2} and Pd 3d_{3/2} binding energies

1. Biphenyl. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 7.342 (t, 2H, $^3J_{\text{HH}}=7.2$ Hz), 7.438 (t, 4H, $^3J_{\text{HH}}=7.6$ Hz), 7.593 (d, 4H, $^2J_{\text{HH}}=4.0$ Hz)

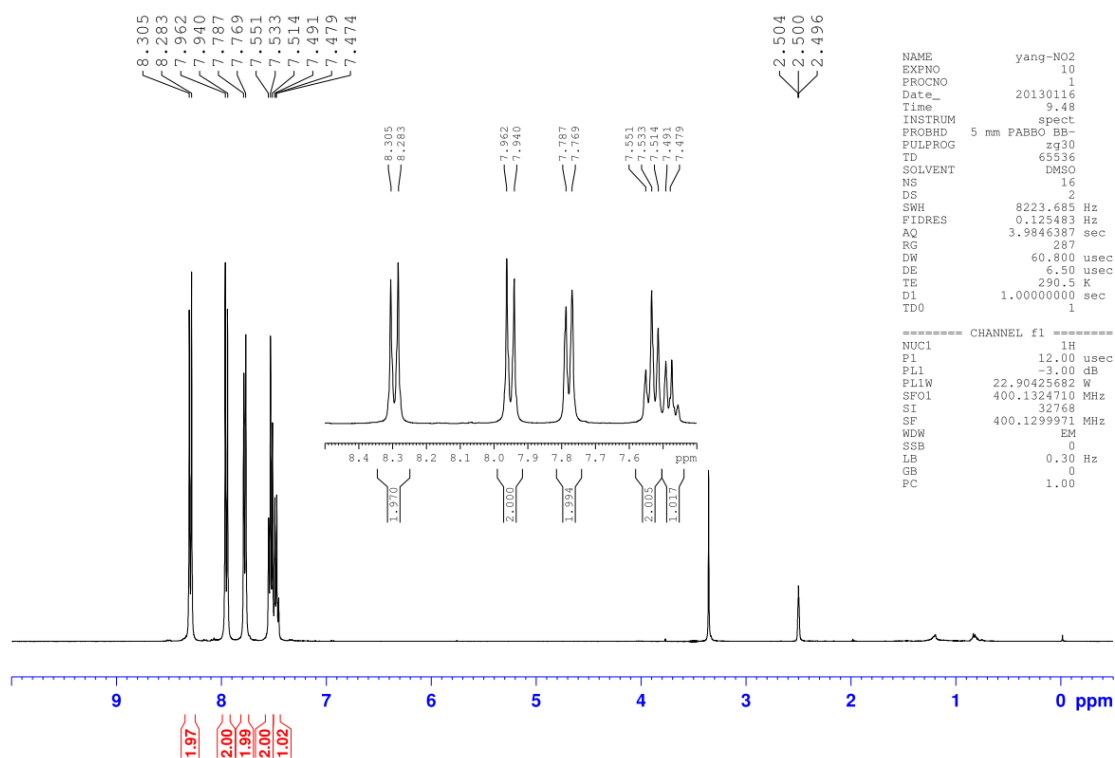


2. 4-methyl-biphenyl. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 2.375 (s, 3H, CH_3), 7.229 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz), 7.301 (t, 1H, $^3J_{\text{HH}}=7.2$ Hz), 7.404 (t, 2H, $^3J_{\text{HH}}=7.6$ Hz), 7.467 (t, 2H, $^3J_{\text{HH}}=9.6$ Hz), 7.563 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz)

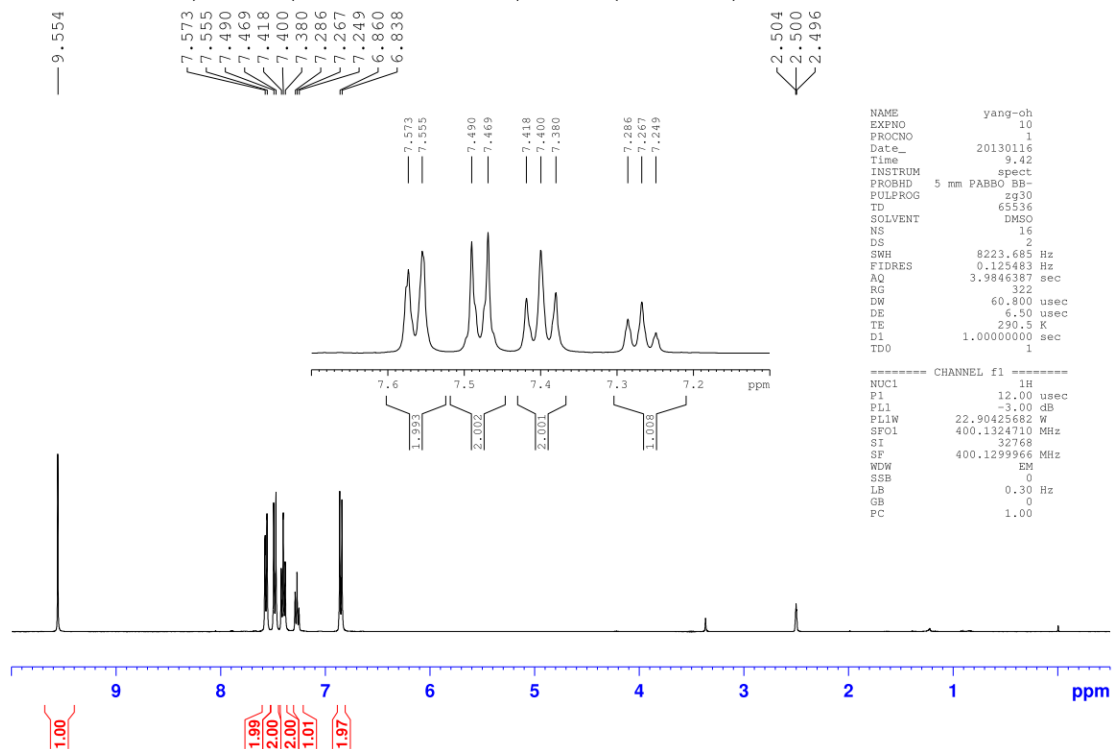


3. 4-Nitro-biphenyl. ^1H NMR ($(\text{CD}_3)_2\text{SO}$, 400 MHz, 25°C)

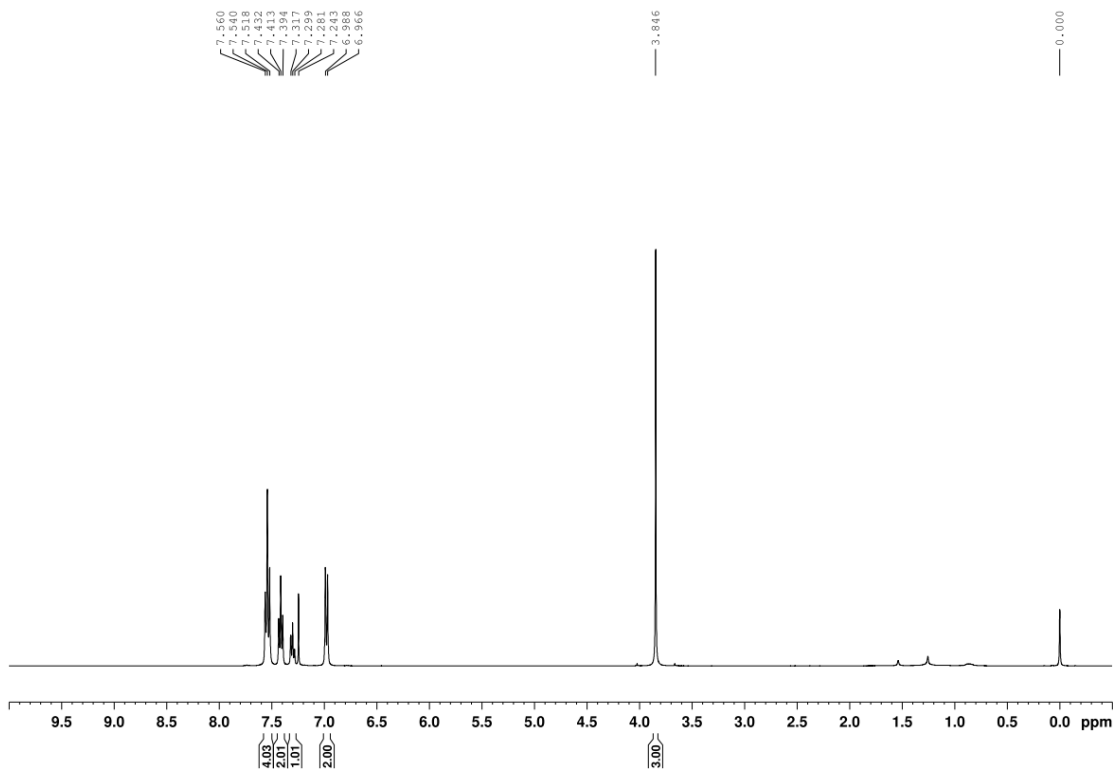
δ 6.849 (d, 1H, $^2J_{\text{HH}}=8.8$ Hz), 7.267 (t, 2H, $^3J_{\text{HH}}=7.6$ Hz), 7.400 (t, 2H, $^3J_{\text{HH}}=7.8$ Hz), 7.479 (d, 2H, $^2J_{\text{HH}}=8.4$ Hz), 7.564 (d, 2H, $^2J_{\text{HH}}=7.6$ Hz)



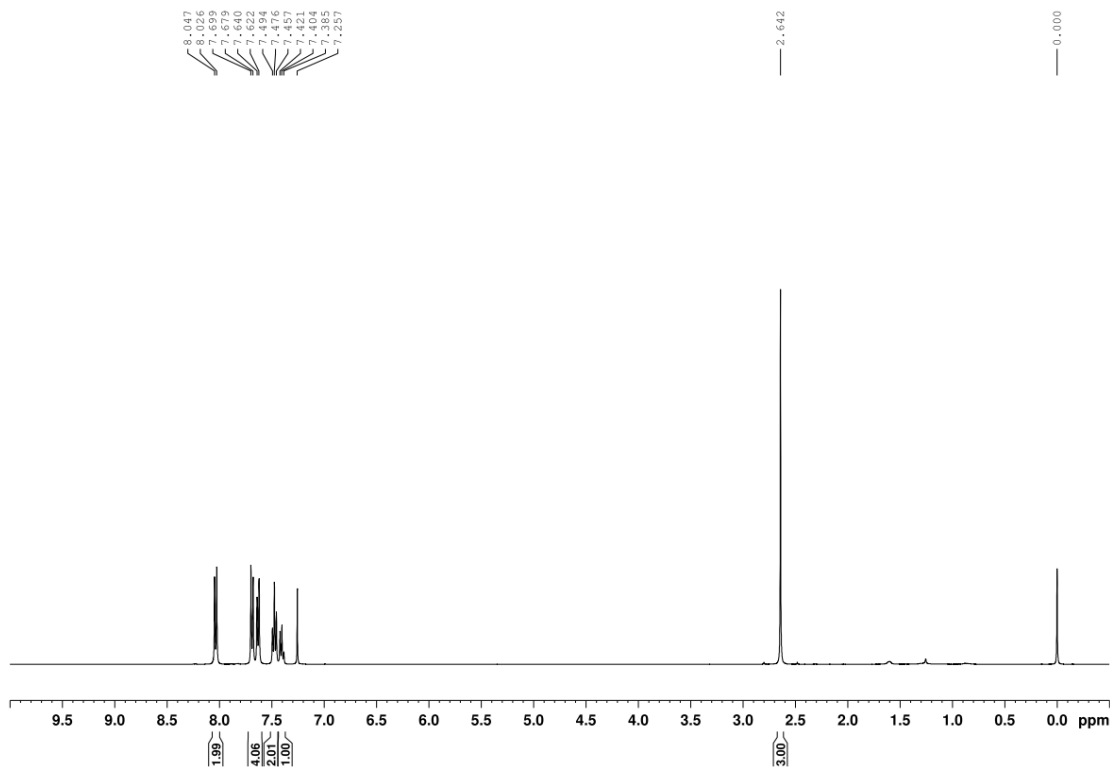
4. Biphenyl-4-ol. ^1H NMR ($(\text{CD}_3)_2\text{SO}$, 400 MHz, 25°C)
 δ 6.845 (d, 2H, $^2J_{\text{HH}}=4.8$ Hz), 7.267 (t, 1H, $^3J_{\text{HH}}=7.6$ Hz), 7.400 (d, 2H, $^2J_{\text{HH}}=7.2$ Hz), 7.469 (d, 2H, $^2J_{\text{HH}}=8.8$ Hz), 7.555 (d, 2H, $^2J_{\text{HH}}=8.8$ Hz), 9.554 (s, 1H, OH)



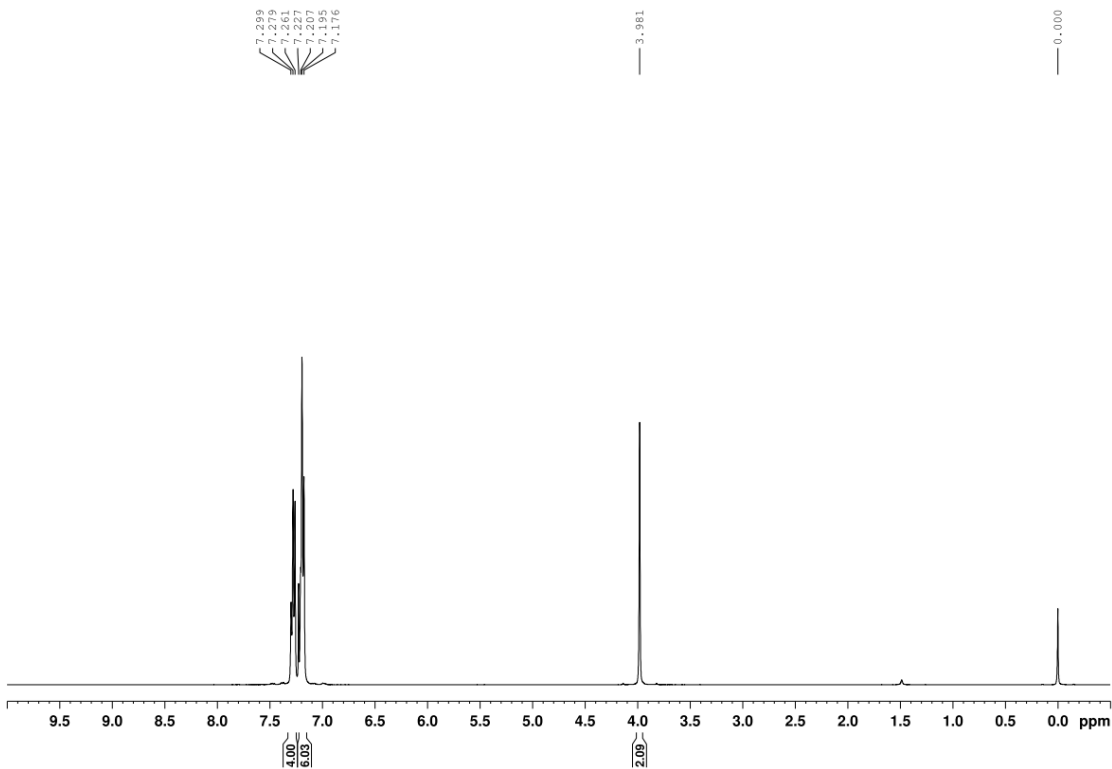
5. 4-Methoxy-biphenyl. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 3.846 (s, 3H, OCH_3), 6.977 (d, 2H, $^2J_{\text{HH}}=4.4$ Hz), 7.299 (t, 1H, $^3J_{\text{HH}}=7.2$ Hz), 7.413 (t, 2H, $^3J_{\text{HH}}=7.6$ Hz), 7.541 (t, 4H, $^3J_{\text{HH}}=7.6$ Hz)



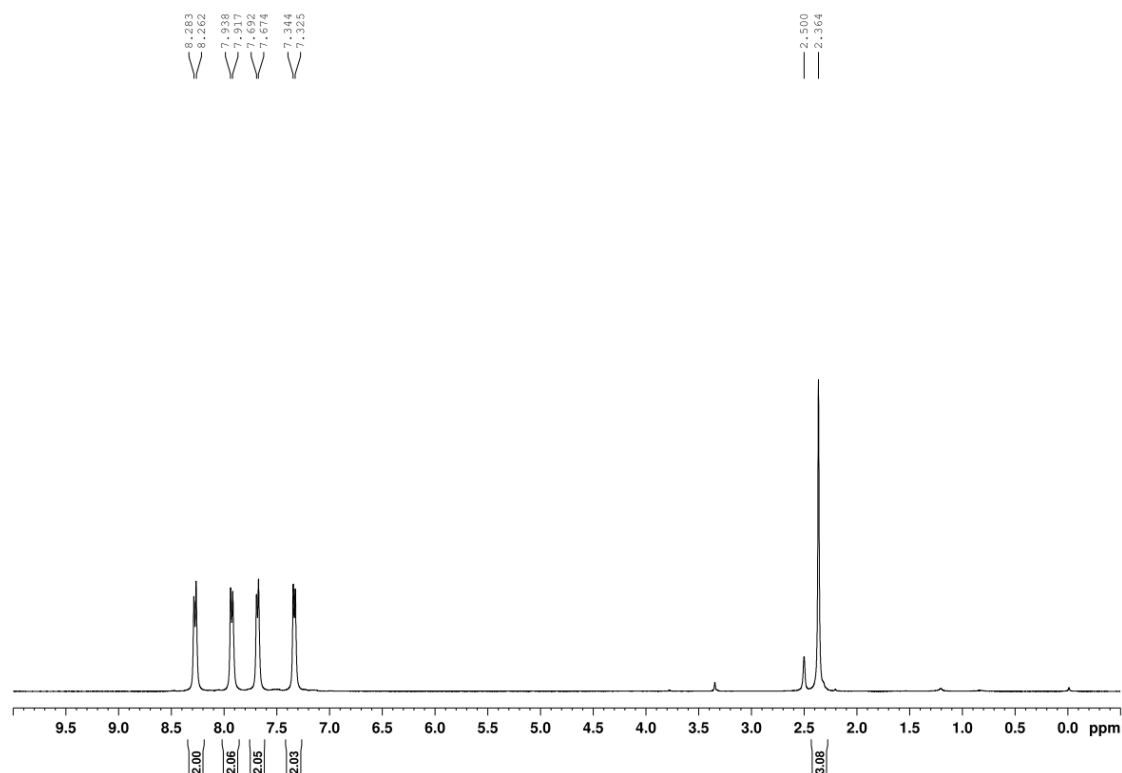
6. 1-biphenyl-4-yl-ethanone. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 2.642 (s, 3H, COCH_3), 7.403 (t, 1H, $^3J_{\text{HH}}=7.2$ Hz), 7.476 (t, 2H, $^3J_{\text{HH}}=7.6$ Hz), 7.631 (d, 2H, $^2J_{\text{HH}}=3.6$ Hz), 7.689 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz), 8.037 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz)



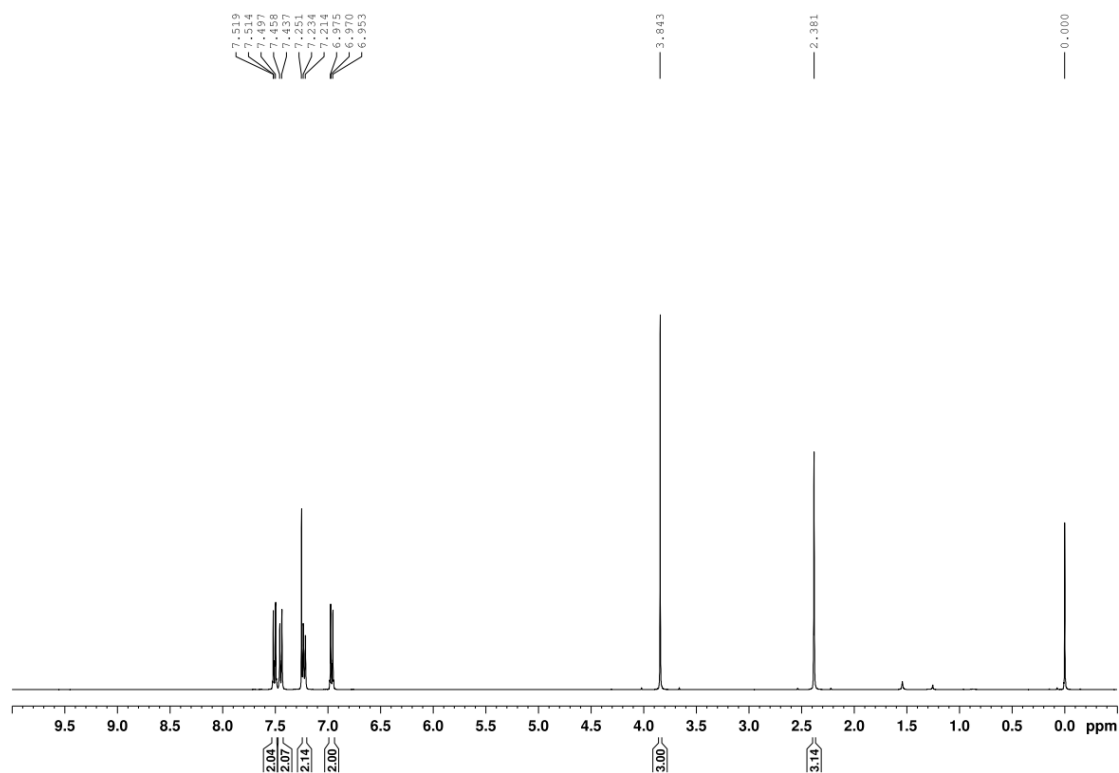
7. Diphenylmethane. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 3.981 (s, 2H, CH_2), 7.201 (m, 6H), 7.279 (t, 4H, $^3J_{\text{HH}}=7.6$ Hz)



8. 4-methyl-4'-nitro-biphenyl. ^1H NMR ($(\text{CD}_3)_2\text{SO}$, 400 MHz, 25°C)
 δ 2.364 (s, 3H, CH_3), 7.334 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz), 7.683 (d, 2H, $^2J_{\text{HH}}=3.6$ Hz), 7.927 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz), 8.273 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz)



9. 4'-methoxy-4-methyl-biphenyl. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 2.381 (s, 3H, CH_3), 3.843 (s, 3H, OCH_3), 6.966 (t, 2H, $^3J_{\text{HH}}=5.2$ Hz), 7.232 (t, 2H, $^3J_{\text{HH}}=7.6$ Hz), 7.448 (d, 2H, $^2J_{\text{HH}}=4.0$ Hz), 7.510 (t, 2H, $^3J_{\text{HH}}=5.2$ Hz)



10. 4,4'-dimethyl-biphenyl. ^1H NMR (CDCl_3 , 400 MHz, 25°C)
 δ 2.380 (s, 6H, CH_3), 7.229 (d, 4H, $^2J_{\text{HH}}=7.6$ Hz), 7.472 (d, 4H, $^2J_{\text{HH}}=8.0$ Hz)

