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Supporting information

Effect of solvent in the hydrogenation of acetophenone catalyzed by Pd/S-DVB.

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Data Type: Counts Mag: 210 Acc. Voltage: 30.0 kV

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Project: POL 2

P. [bar]

11.16

11.05

10.95

10.85



Fig. S1 Elemental mapping (SEM) for Pd/S-DVB catalyst.

Reaction time	0	15	30	45	60	75	90	105	120	135	150	165	180	195	210	225	240
Pressure [bar]	11.30	10.90	10.40	9.90	9.80	9.70	9.60	9.60	9.60	9.50	9.50	9.50	9.50	9.50	9.50	9.50	9.50

Tab. S1.	Pressure	changes in	time fo	or hydrogei	nation o	f APh	in MeOH
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Reaction time	0	15	30	45	60	75	90	105	120	135	150	165	180	195	210	225	240
Pressure [bar]	11.10	10.70	10.20	9.90	9.20	9.00	8.90	8.80	8.75	8.70	8.70	8.70	8.60	8.60	8.60	8.50	8.50

R. time	0	15	30	45	60	75	90	105	120	x _i H ₂ O
P. [bar]	11.30	10.60	9.90	9.50	9.20	9.00	8.90	8.70	8.70	0.00
P. [bar]	11.40	11.30	11.00	10.70	10.60	10.40	10.30	10.20	10.10	0.28
P. [bar]	11.90	11.60	11.60	11.50	11.40	11.20	11.10	11.00	10.90	0.46
P. [bar]	12.80	12.70	12.60	12.60	12.50	12.50	12.40	12.40	12.30	0.81

10.80

10.68

10.60

10.30

10.27

1.00

Tab. S2. Pressure changes in time for hydrogenation of APh in BuOH

R. time	135	150	165	180	195	210	225	240	x _i H ₂ O
P. [bar]	8.60	8.60	8.50	8.50	8.50	8.50	8.50	8.40	0.00
P. [bar]	10.00	9.90	9.80	9.80	9.70	9.70	9.70	9.70	0.28
P. [bar]	10.80	10.70	10.60	10.50	10.40	10.30	10.20	10.20	0.46
P. [bar]	12.30	12.20	12.20	12.10	12.10	12.10	12.10	12.10	0.81
P. [bar]	10.17	10.05	9.95	9.85	9.76	9.71	9.67	9.54	1.00

Tab. S3. Pressure changes in time for hydrogenation of APh in i-PrOH







Fig. S2 TEM micrographs of Pd/S-DVB after catalytic reaction



Fig. S3 XPS P 2p core level spectra for sample before reaction in 'as received' form a), for sample before reaction after soft cleaning 'in situ' b) for sample after reaction c)

GC-MS data





Fig. S4 GC-MS after APh hydrogenation (10 bar, 80°C, 4h)











Fig. S7 MS of 1-phenylethanol







Fig. S9. MS of 1-(2,5-dimethylphenyl)ethan-1-ol



Fig. S10. MS of 1-ethyl-2,5-dimethylbenzene



Fig. S11 GC-MS after 3-methylacetophenone hydrogenation (10 bar, 80°C, 4h, 3ml MeOH)



80

<<Target >> Line#:4 R.Time:3.090(Scan#:219) MassPeaks:313 RawMode:Averaged 3.085-3.095(218-220) BasePeak:105.10(6865394) BG Mode:Cale. from Peak Group 1 - Event 1 Scan 100 1 185



Fig. S14 GC-MS after 2-methylacetophenone hydrogenation (10 bar, 80°C, 4h)

Fig. S12 MS of 1-(3-tolyl)ethan-1-ol

40

60

20



100 120 140 160 180 200 220 240 260 280 300 320 340 360 380 400 420 440 460 480



Fig. S15. MS of 1-(2-tolyl)ethan-1-ol



Fig. S16.GC-MS 3-aminoacetophenone hydrogenation (10 bar, 80°C, 4h)



Fig. S17. MS of 3-ethylaniline



Fig. S18. MS of 1-(3-aminophenyl)ethan-1-ol