

Supplementary Information

In situ MnO_x/N-doped carbon aerogels from cellulose as monolithic and highly efficient catalysts for upgrading of bioderived aldehydes

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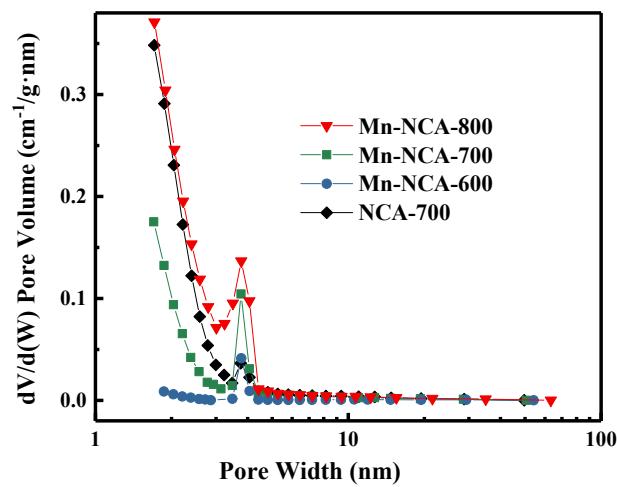


Fig. S1 Pore size distribution of Mn-NCA-T

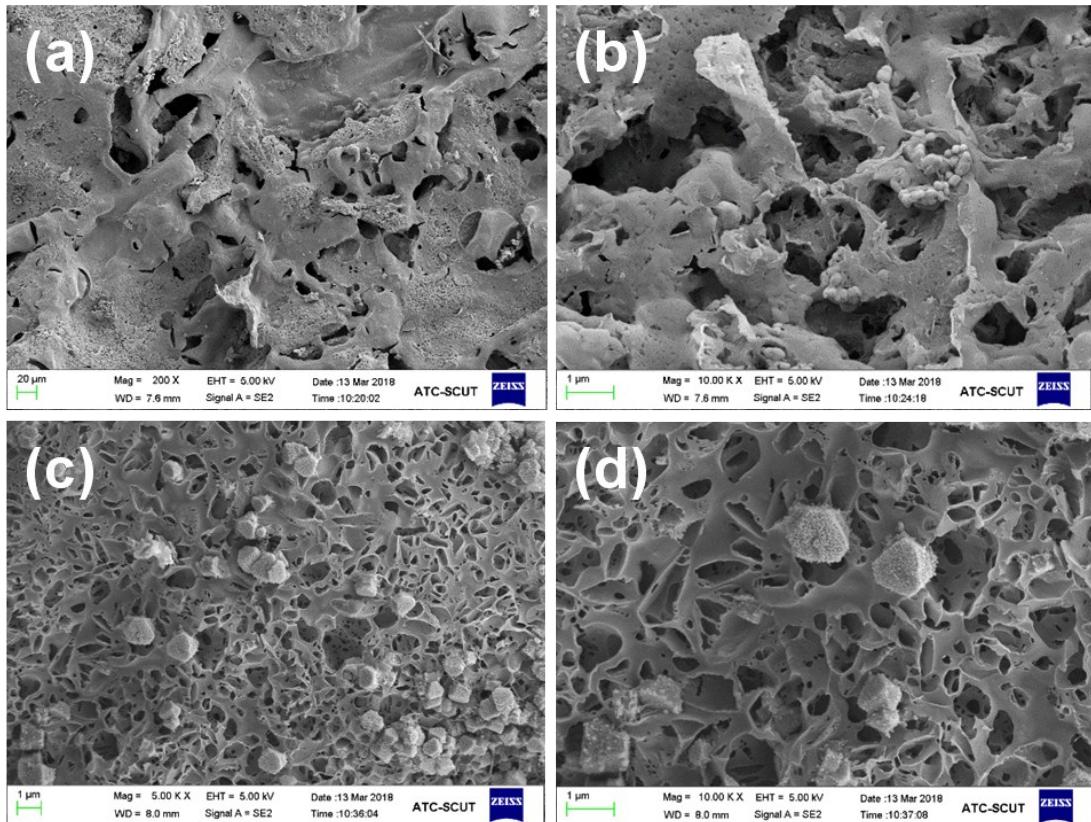


Fig. S2 SEM images of Mn-NCA-600 (a, b) and Mn-NCA-800 (c, d).

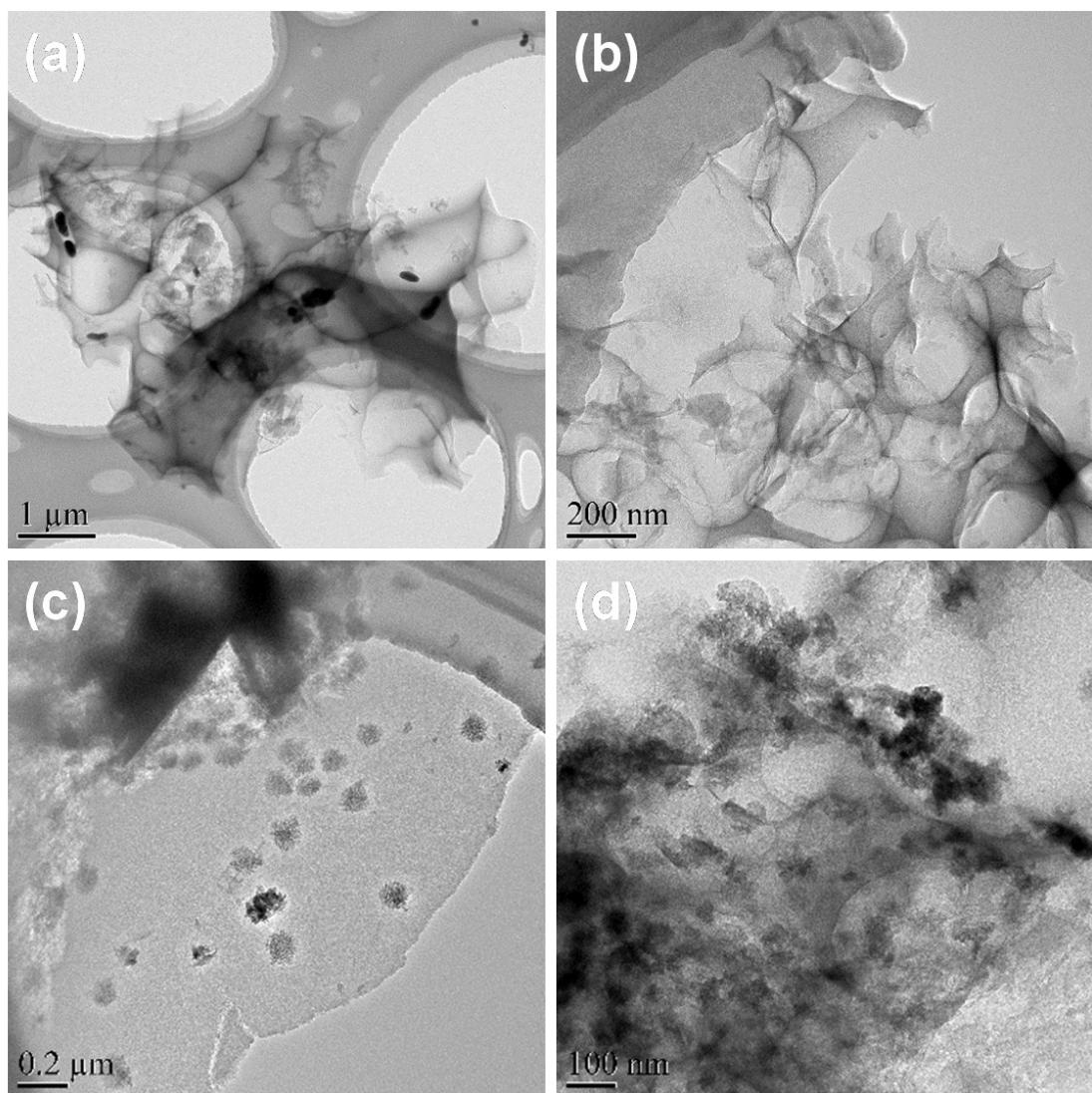


Fig. S3 TEM images of Mn-NCA-600 (a, b) and Mn-NCA-800 (c, d).

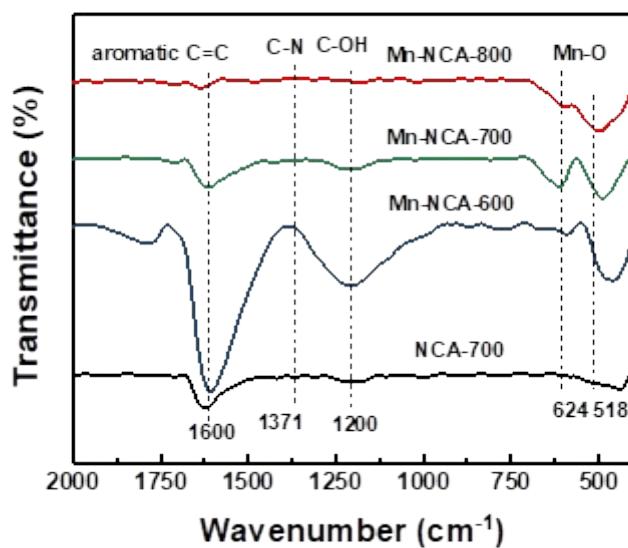


Fig. S4 FT-IR spectra of NCA-700 and Mn-NCA-T.

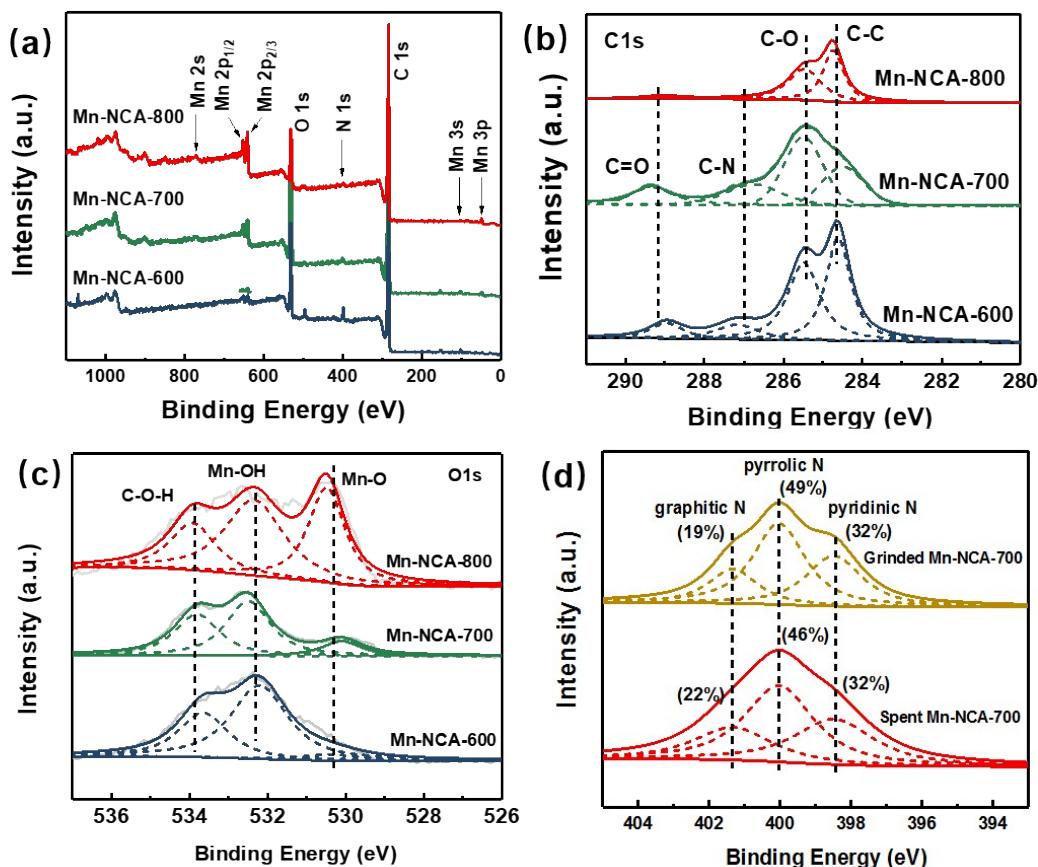


Fig. S5 XPS spectrum of survey scan (a), C 1s peaks (b), O 1s peaks (c) of Mn-NCA-T and N 1s peaks (d) of grinded and spent Mn-NCA-700 catalysts.

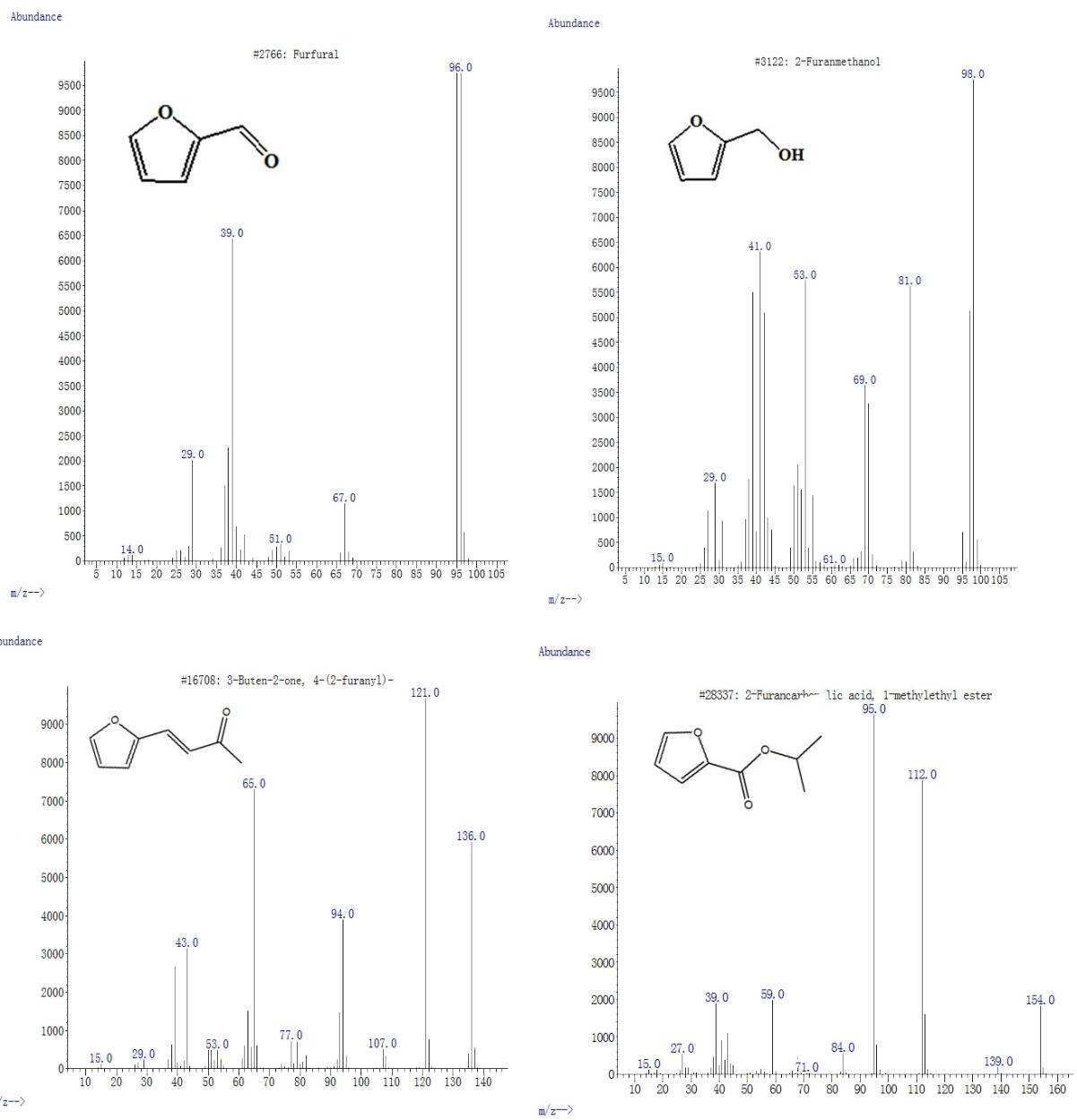


Fig. S6 Representative MS spectra of the reaction mixture.

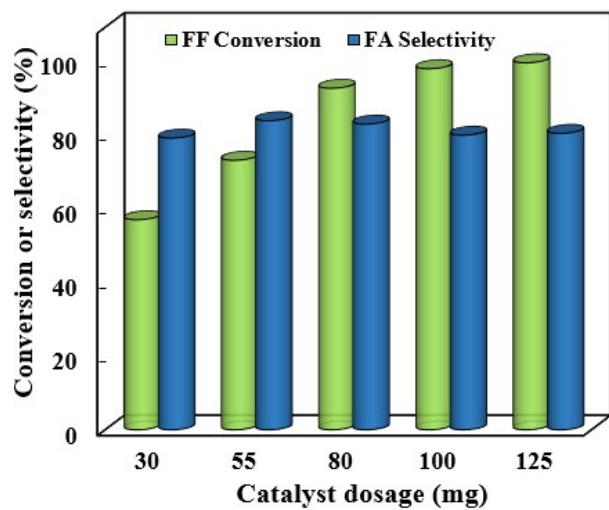


Fig. S7 Effect of the Mn-NCA-700 catalyst dosage on the conversion of FF to FA;

Reaction conditions: 0.05mmol furfural in 5ml isopropanol, 160 °C, 30 min.

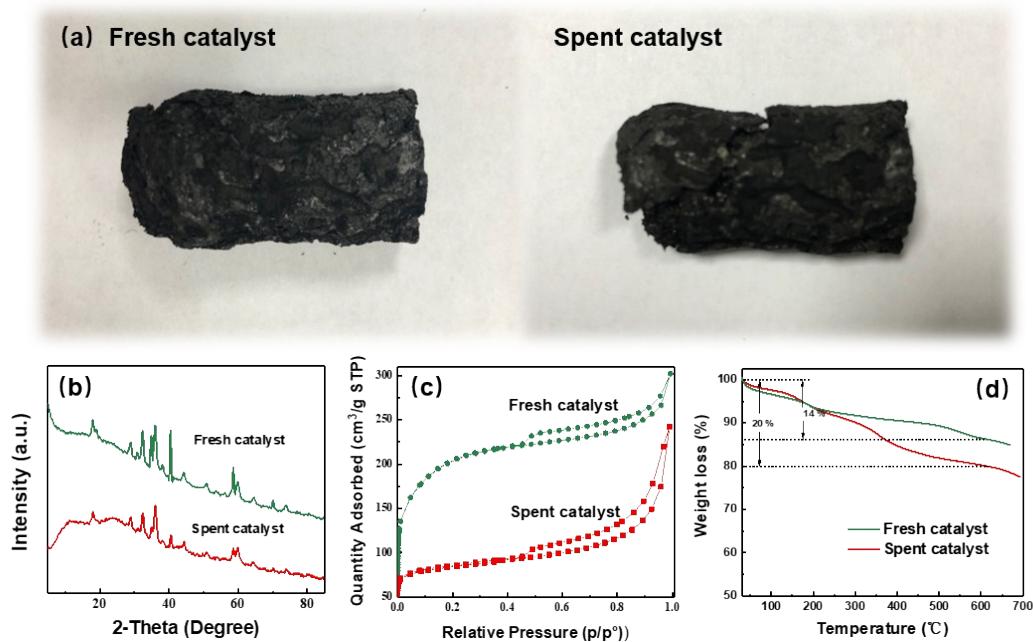


Fig. S8 Analyses of fresh and recovered Mn-NCA-700 catalyst after five cycles:

photograph (a), XRD patterns (b), N₂ adsorption-desorption isotherms (c), TG analyses (d).

Table S1 The physical properties and Raman results of various catalysts

Catalysts	Surface area (m ² ·g ⁻¹)	Pore volume (cm ³ /g)	Average pore diameter (nm)	I_D/I_G	I_{2D}/I_G
NCA-700	1083.14	0.61	3.46	0.9852	0.3014
Mn-NCA-600	308.49	0.18	8.27	1.053	0.4942
Mn-NCA-700	676.95	0.41	4.89	1.006	0.6690
Mn-NCA-800	909.83	0.52	2.42	1.010	-

Table S2 The compositions of various catalysts

Catalysts	Content (wt.%)					
	Metal ^a	Na ^a	N ^b	C ^b	H ^b	O (Calculated)
MnO _x	63.87	-	-	-	-	36.13
CA-700	-	0.11	0.52	77.35	2.3	19.68
NCA-700	-	0.293	2.66	65.66	5.7	25.65
Mn-CA-700	13.14	0.286	1.4	62.38	2.3	20.44
Mn-NCA-600	15.07	0.088	6.98	47.98	2.5	27.28
Mn-NCA-700	19.97	0.290	2.27	46.72	3.8	26.86
Mn-NCA-800	30.82	0.103	0.45	46.88	4.0	17.74
Grinded Mn-NCA-700	19.97	0.290	2.27	46.72	3.8	26.86
Co-NCA-700	17.11	0.197	0.6	65.26	2.7	14.12
Fe-NCA-700	20.23	0.345	0.76	67.31	1.9	9.38
Cu-NCA-700	31.12	0.538	3.47	35.64	4.0	25.21
Ni-NCA-700	15.50	0.477	0.88	72.42	1.5	9.20
Recycled Mn-NCA-700	18.46	0.174	2.09	49.95	2.8	26.51

^a Determined by ICP-OES.^b Determined by elemental analysis.

Table S3 The comparison of CTH of FF with H₂ or H-donor reaction systems with previously reported catalysts.

Entry	Catalyst	H-donor	Reaction conditions	FF Conv. (%)	FA Sel. (%)	Ref.
1	2%Pt-1%Re/TiO ₂ -ZrO ₂	5 MPa H ₂	130 °C, 8 h	100	95.7	[1]
2	Co/SBA-15	2 MPa H ₂	150 °C, 1 h	80	96	[2]
3	Cu-Fe(7:1)/SiO ₂	H ₂	-	-	-	[3]
4	Cu:Zn:Cr:Zr(3:2:1:3)	2 MPa H ₂	170 °C, 3.5 h	100	96	[4]
5	Fe-L4(L5)/C-800	2-propanol	160 °C, 15h	91.6	83	[5]
6	Pd/Fe ₂ O ₃	2-propanol	180 °C, 7.5h	100	34	[6]
7	Cu/MgO-Al ₂ O ₃	2-propanol	210 °C, 1h,	100	89	[7]
8	DyCl ₃	2-propanol	180 °C, 3h, (2Mpa N ₂)	98	97	[8]
9	Al ₇ Zr ₃ @Fe ₃ O ₄	2-propanol	180 °C, 4h	99.1	90.5	[9]
10	Fe ₂ O ₃ @HAP	2-propanol	180 °C, 10h	96.2	95.3	[10]
11	NiFe ₂ O ₄	2-propanol	180 °C, 6h	99	95	[11]
12	Cu ₂ Al	methanol	200 °C, 2.5h, (1Mpa N ₂)	100	94	[12]
13	Al ₂ O ₃ -S	2-propanol	130 °C, 6h	98	97	[13]
14	γ-Al ₂ O ₃	2-propanol	150 °C, 6h	100	90	[14]
15	Mn-NCA-700	2-propanol	200 °C, 0.5h	99	89	This work
16	Mn-NCA-700	2-propanol	160 °C, 1h	99	85	This work
17	Mn-NCA-700	2-propanol	140 °C, 1.2h	95	80	This work

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