Enhancement of peroxidase-like activity of N-doped graphene assembled with iron-tetrapyridylporphyrin

L Magerusan¹, C Socaci^{1*}, F Pogacean¹, M-C Rosu, A R Biris¹, M Coros¹, A. Turza¹, V Floare-Avram¹, G Katona², S Pruneanu^{1*}



Figure S1 HR-TEM images of porphyrin functionalized N-doped graphenes

Sample	Peak position	Interplanar	Crystallite size	Number of
	2 0 °	spacing (Å)	(Å)	layers
NGr	24.62	3.610	111	30.7
	26.11	3.408	51	14.96
NGr-TPyP	24.62	3.610	113	31.30
	26.16	3.401	58	17.05
TPyP-NGr	23.65	3.757	53	14.10
	26.05	3.432	32	9.32
NGr-FeTPyP	24.49	3.631	99	27.26

Table S1. XRD data for the porphyrin-functionalized N-doped graphenes

	25.93	3.430	59	17.20
FeTPyP-NGr	23.77	3.736	35	9.36
	26.11	3.414	35	10.25

Table S2. Assignments of each deconvoluted peak based on their binding energies (BE) andatomic concentrations [AC, %] for NGr-FeTPyP and FeTPyP-NGr

Flomont	Binding energy	Assignments		
Liement	NGr-(Fe)TPyP	(Fe)TPyP-NGr	Assignments	
C 1s	284.27 eV [39.73 %]	284.31 eV [42.33 %]	<i>sp</i> ² C=C / CH _n	
	284.88 eV [28.05 %]	284.86 eV [23.71 %]	sp^3 C-C / CH _n	
	286.13 eV [13.01 %]	285.73 eV [11.64 %]	C-N / C-O	
	287.92 eV [10.68 %]	287.47 eV [13.38 %]	C=O	
	289.83 eV [6.11 %]	289.11 eV [5.86 %]	OH-C=O / COOH	
	292.23 eV [2.42 %]	291.83 eV [3.08 %]	$\pi \rightarrow \pi^*$ shake up satellite	
N 1s	398.34 eV [53.66 %]	398.21 eV [31.35 %]	pyridinic N	
	400.17 eV [36.22 %]	400.10 eV [59.03 %]	pyrrolic N	
	402.52 eV [10.12 %]	402.90 eV [9.62 %]	quaternary N	
0 1s	529.64 [8.57 %]	-	Fe-O	
	531.30 [38.75 %]	530.84 [37.20 %]	C-O	
	533.14 [38.82%]	532.83 [52.97 %]	OH-C=O / COOH	
	535.37 [13.86 %]	535.75 [9.83 %]	adsobed H ₂ O	



Figure S2: Element mapping images of FeTPyP-NGr (a), revealing the distribution of C (Red), O (green), Fe(purple) and N(cyan) elements and the corresponding EDX spectrum (b)