

Table S1. Elemental compositions investigated by XPS.

	Graphene stacks	GO	PDA-GO_30	RGO_30
[N]:[C] _{spec}	–	0.009	0.096	0.035
[O]:[C] _{spec}	0.029	0.166	0.27	0.165
[F]:[C] _{spec}	–	–	–	0.005
[Na]:[C] _{spec}	0.001	–	0.0004	0.013
[Si]:[C] _{spec}	0.004	0.006	0.013	0.005
[S]:[C] _{spec}	–	0.012	0.002	0.002
[Cl]:[C] _{spec}	0.0003	0.0002	–	0.0002

	PDA-GO_90	PDA-GO_180
[N]:[C] _{spec}	0.0831	0.0843
[O]:[C] _{spec}	0.2376	0.261
[F]:[C] _{spec}	0.003	0.012
[Na]:[C] _{spec}	–	–
[Si]:[C] _{spec}	0.006	0.024

Table S2. Wide-scan spectrum of the sample recorded from the graphene stacks.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	RSF	Atomic Conc %	Mass Conc %
C 1s	284.0	0.58	64387	0.278	96.71	95.17
O 1s	532.7	2.65	5528	0.780	2.76	3.62
Na 1s	1070.6	2.50	9106	1.685	0.09	0.18
Si 2p	103.2	4.06	6433	0.328	0.41	0.94
Cl 2p	199.3	0.19	1377	0.891	0.03	0.09

Table S3. C 1s spectrum of the sample recorded from the graphene stacks.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
C 1s <i>Gr</i>	284.00	0.75	51530	71650	80.05	80.05
C 1s <i>C</i>	285.70	1.00	1840	1675	2.86	2.86
C 1s <i>Sh1</i>	286.60	2.34	5302	2067	8.23	8.23
C 1s <i>Sh2</i>	289.81	2.34	3389	1321	5.26	5.26
C 1s <i>Sh3</i>	291.77	2.34	1933	753	3.00	3.00
C 1s <i>Sh4</i>	294.36	2.34	396.0	154	0.61	0.61

C 1s *Gr*: AS(90,0.45).**Table S4.** Wide-scan spectrum of the GO sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	RSF	Atomic Conc %	Mass Conc %
C 1s	284.0	0.90	44545	0.278	83.79	78.23
N 1s	401.2	0.55	731	0.477	0.79	0.86
O 1s	531.6	2.80	22345	0.780	13.90	17.29
Si 2p	101.2	2.45	6503	0.328	0.51	1.12
S 2p	168.2	2.93	25088	0.668	0.98	2.44
Cl 2p	199.3	0.20	855	0.891	0.02	0.07

Table S5. C 1s spectrum of the GO sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
C 1s <i>Gr</i>	283.99	0.91	25608	27376	57.50	57.50
C 1s <i>Ph</i>	284.75	1.14	5083	4072	11.41	11.41
C 1s <i>B</i>	285.72	1.14	808	647	1.81	1.81
C 1s <i>C</i>	286.07	1.20	3498	2656	7.85	7.85
C 1s <i>D</i>	287.29	1.20	1144	869	2.57	2.57
C 1s <i>F</i>	288.36	1.20	1208	916	2.71	2.71
C 1s <i>Sh1</i>	286.36	2.36	3461	1335	7.77	7.77
C 1s <i>Sh2</i>	289.57	2.36	2212	853	4.96	4.96
C 1s <i>Sh3</i>	291.54	2.36	1262	487	2.83	2.83
C 1s <i>Sh4</i>	294.13	2.36	261	101	0.58	0.58

C 1s *Gr*: AS(60,1.0).**Table S6.** N 1s spectrum of the GO sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
N 1s <i>L</i>	399.46	1.83	191	95	26.11	26.11
N 1s <i>M</i>	401.15	1.83	540	268	73.89	73.89

Table S7. Wide-scan spectrum of the PDA-GO_30 sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	RSF	Atomic Conc %	Mass Conc %
C 1s	284.7	2.54	13568	0.278	72.31	66.18
N 1s	400.1	1.74	2299	0.477	6.95	7.42
O 1s	532.6	2.39	11051	0.780	19.57	23.86
Na 1s	1070.9	1.98	815	1.685	0.03	0.05
Si 2p	102.2	2.49	4280	0.328	0.96	2.06
S 2p	168.5	2.07	1584	0.668	0.18	0.43

Table S8. C 1s spectrum of the PDA-GO_30 sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
C 1s <i>Ph</i>	284.40	1.08	3729	3474	27.56	27.56
C 1s <i>A</i>	285.00	1.05	3022	2623	22.33	22.33
C 1s <i>B</i>	285.81	1.13	2600	2097	19.21	19.21
C 1s <i>C</i>	286.57	1.13	2600	2097	19.21	19.21
C 1s <i>C1</i>	286.57	1.13	0	0	0.00	0.00
C 1s <i>Sh1</i>	288.15	1.82	1327	666	9.80	9.80
C 1s <i>Sh2</i>	290.80	1.82	256	129	1.89	1.89

Table S9. N 1s spectrum of the PDA-GO_30 sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
N 1s <i>K</i>	398.61	1.60	247	141	10.75	10.75
N 1s <i>L</i>	400.06	1.60	1680	962	73.13	73.13
N 1s <i>M</i>	401.71	1.60	371	212	16.12	16.12

Table S10. Wide-scan spectrum of the RGO_30 sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	RSF	Atomic Conc %	Mass Conc %
C 1s	284.5	1.00	39897	0.278	81.66	76.34
N 1s	400.0	1.81	2430	0.477	2.82	3.08
O 1s	532.6	3.27	19847	0.780	13.48	16.79
F 1s	689.3	2.91	15673	1.000	0.39	0.58
Na 1s	1070.3	2.44	76371	1.685	1.08	1.94
Si 2p	102.1	2.34	4260	0.328	0.37	0.80
S 2p	168.7	3.29	3880	0.668	0.17	0.41
Cl 2p	199.3	0.59	622	0.891	0.02	0.05

Table S11. C 1s spectrum of the RGO_30 sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
C 1s <i>Ph</i>	284.46	1.13	24831	22184	62.03	62.03
C 1s <i>B</i>	285.34	1.46	2760	1720	6.89	6.89
C 1s <i>C</i>	286.29	1.46	2760	1720	6.89	6.89
C 1s <i>C'</i>	286.29	1.46	2441	1521	6.10	6.10
C 1s <i>F</i>	289.36	1.46	225	140	0.56	0.56
C 1s <i>Sh1</i>	287.92	2.44	4733	1769	11.81	11.81
C 1s <i>Sh2</i>	290.77	2.44	1812	677	4.52	4.52
C 1s <i>Sh3</i>	293.23	2.44	478	179	1.19	1.19

C 1s *Ph*: AS(80,0.55).**Table S12.** N 1s spectrum of the RGO_30 sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
N 1s <i>K</i>	398.85	2.09	373	162	15.73	15.73
N 1s <i>L</i>	400.07	2.09	1999	964	84.27	84.27

Table S13. Wide-scan spectrum of the PDA-GO_90 sample.

Peak.	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	RSF	Atomic Conc %	Mass Conc %
C 1s	284.8	2.38	12829	0.278	75.21	69.84
N 1s	400.1	1.67	1878	0.477	6.25	6.76
O 1s	533.0	2.58	9174	0.780	17.87	22.10
F 1s	689.9	2.90	3266	1.000	0.24	0.35
Si 2p	102.2	2.53	1771	0.328	0.44	0.95

Table S14. C 1s spectrum of the PDA-GO₉₀ sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
C 1s <i>Ph</i>	284.63	1.06	3203	2754	24.93	24.93
C 1s <i>A</i>	285.00	1.24	2909	2137	22.63	22.63
C 1s <i>B</i>	285.71	1.23	2135	1583	16.61	16.61
C 1s <i>C</i>	286.47	1.23	2135	1583	16.61	16.61
C 1s <i>C1</i>	286.47	1.23	434	323	3.38	3.38
C 1s <i>Sh1</i>	287.70	1.80	1269	644	9.86	9.86
C 1s <i>Sh2</i>	289.16	1.80	528	268	4.10	4.10
C 1s <i>Sh3</i>	291.31	1.80	241	123	1.87	1.87

Table S15. N 1s spectrum of the PDA-GO₉₀ sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
N 1s <i>K</i>	398.77	1.54	220	131	11.66	11.66
N 1s <i>L</i>	400.19	1.54	1453	862	76.87	76.87
N 1s <i>M</i>	401.72	1.54	217	129	11.48	11.48

Table S16. Wide-scan spectrum of the PDA-GO₁₈₀ sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	RSF	Atomic Conc %	Mass Conc %
C 1s	284.9	2.23	10051	0.278	72.40	65.76
N 1s	399.9	1.51	1492	0.477	6.10	6.46
O 1s	532.7	2.37	7891	0.780	18.89	22.85
F 1s	689.1	2.51	10076	1.000	0.89	1.28
Si 2p	102.5	2.63	5659	0.328	1.72	3.65

Table S17. C 1s spectrum of the PDA-GO₁₈₀ sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
C 1s <i>Ph</i>	284.35	1.23	2956	2422	29.49	29.49
C 1s <i>A</i>	285.00	1.08	2515	2125	25.08	25.08
C 1s <i>B</i>	285.63	1.22	1687	1259	16.82	16.82
C 1s <i>C</i>	286.37	1.22	1687	1259	16.82	16.82
C 1s <i>C1</i>	286.37	1.22	0	0	0.00	0.00
C 1s <i>Sh1</i>	287.99	2.31	1059	418	10.55	10.55
C 1s <i>Sh2</i>	290.64	2.31	125	49	1.24	1.24

Table S18. N 1s spectrum of the PDA-GO₁₈₀ sample.

Peak	Position BE (eV)	FWHM (eV)	Raw Area (CPS)	Height	Atomic Conc %	Mass Conc %
N 1s <i>K</i>	398.51	1.53	174	103	11.59	11.59
N 1s <i>L</i>	400.05	1.53	1188	707	79.22	79.22
N 1s <i>M</i>	401.50	1.53	138	82	9.19	9.19