

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050255.D
 Acq On : 09 Jun 2025 23:35
 Operator : RC/JU
 Sample : Q2236-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-06-C

Quant Time: Jun 10 01:28:43 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 09 11:54:38 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

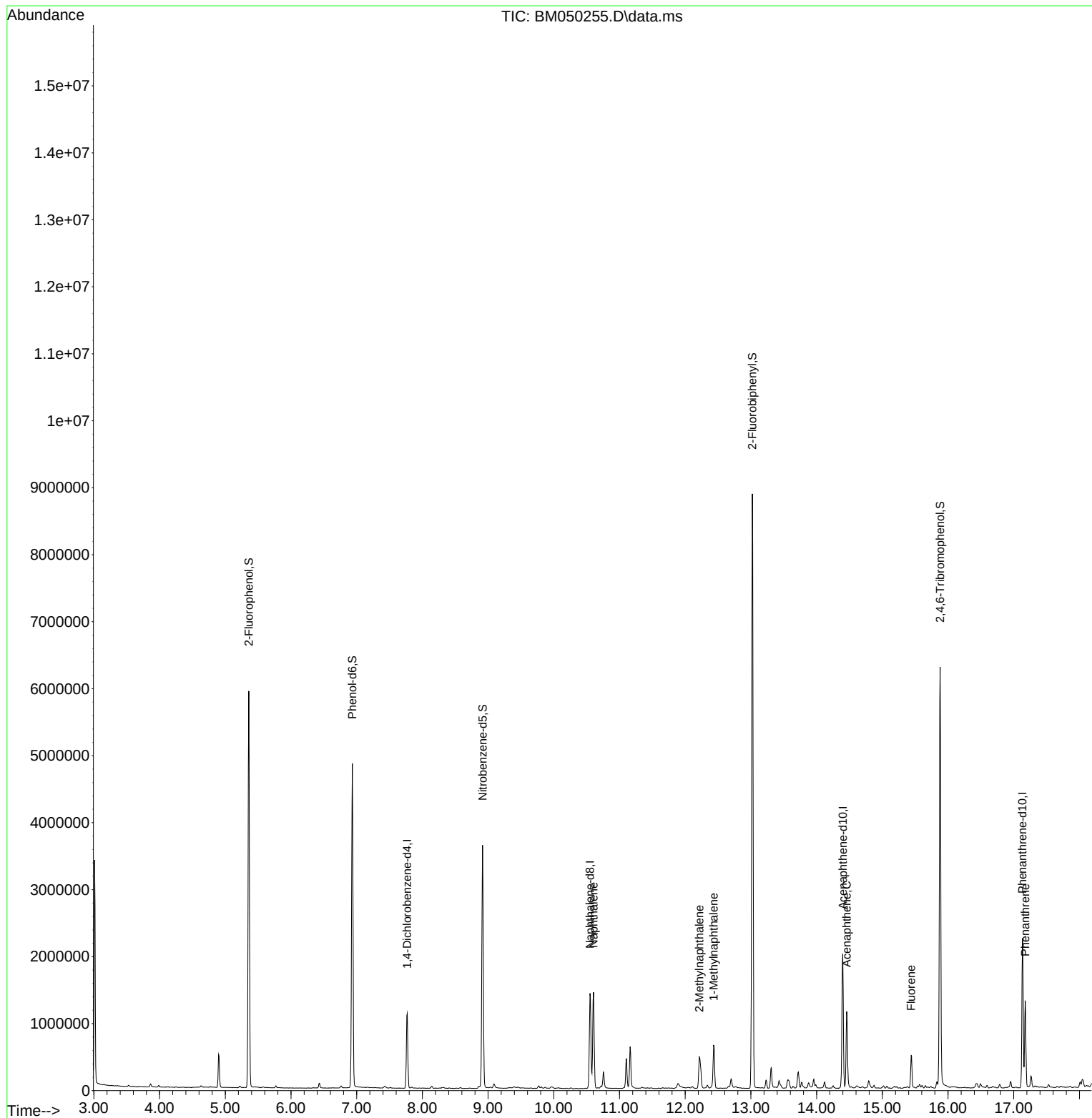
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	313628	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1214917	20.000	ng	0.00
39) Acenaphthene-d10	14.398	164	691337	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1261240	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1230719	20.000	ng	-0.01
86) Perylene-d12	24.338	264	1312536	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2745871	146.049	ng	0.00
7) Phenol-d6	6.934	99	3109740	125.647	ng	0.00
23) Nitrobenzene-d5	8.916	82	2265879	96.998	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1196426	152.150	ng	0.00
45) 2-Fluorobiphenyl	13.021	172	4612245	90.322	ng	0.00
79) Terphenyl-d14	19.762	244	6079199	93.606	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.604	128	1202299	19.141	ng	100
37) 2-Methylnaphthalene	12.216	142	204859	5.406	ng	99
38) 1-Methylnaphthalene	12.433	142	309802	7.703	ng	99
52) Acenaphthene	14.457	154	408261	10.027	ng	99
58) Fluorene	15.439	166	201762	4.422	ng	97
71) Phenanthrene	17.174	178	760701	10.558	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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