

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050254.D
 Acq On : 09 Jun 2025 22:56
 Operator : RC/JU
 Sample : Q2236-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-05-C

Quant Time: Jun 10 01:28:25 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)

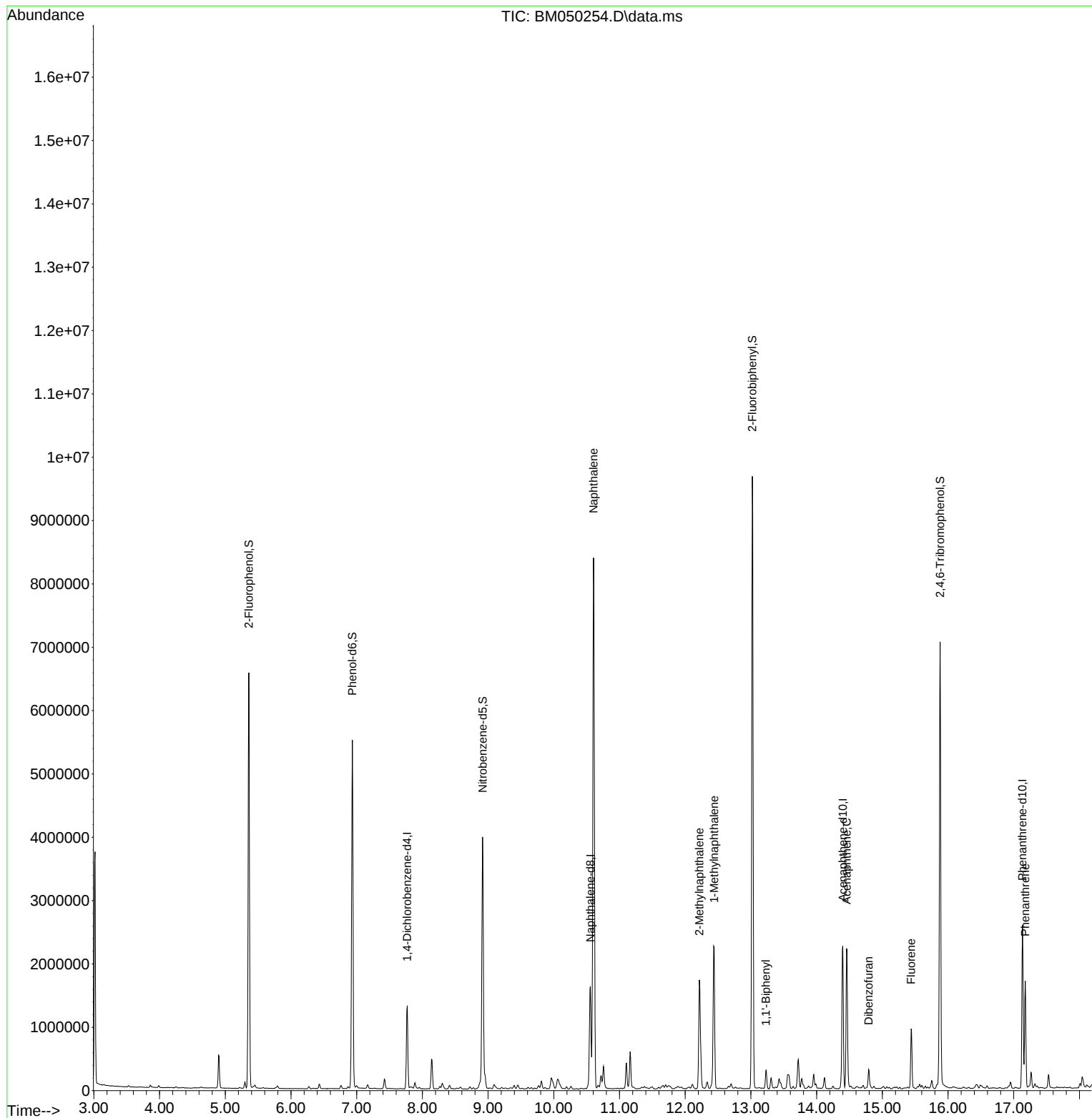
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	372657	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	1419602	20.000	ng	0.00
39) Acenaphthene-d10	14.398	164	799341	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1429372	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1316639	20.000	ng	-0.01
86) Perylene-d12	24.339	264	1352313	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	3073333	137.573	ng	0.00
7) Phenol-d6	6.934	99	3489531	118.659	ng	0.00
23) Nitrobenzene-d5	8.916	82	2529986	92.688	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1300437	143.032	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	5056880	85.649	ng	0.00
79) Terphenyl-d14	19.762	244	6346327	91.342	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.604	128	7308166	99.571	ng	100
37) 2-Methylnaphthalene	12.216	142	763922	17.253	ng	99
38) 1-Methylnaphthalene	12.439	142	1094227	23.284	ng	99
46) 1,1'-Biphenyl	13.228	154	154552	2.582	ng	98
52) Acenaphthene	14.457	154	801278	17.021	ng	98
55) Dibenzofuran	14.792	168	169828	2.466	ng	99
58) Fluorene	15.439	166	416718	7.900	ng	100
71) Phenanthrene	17.174	178	985702	12.072	ng	99

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

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