

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071525\
 Data File : BM050460.D
 Acq On : 15 Jul 2025 18:11
 Operator : RC/JU
 Sample : Q2578-19
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-12-C

Quant Time: Jul 16 02:00:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 08 18:32:25 2025
 Response via : Initial Calibration

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

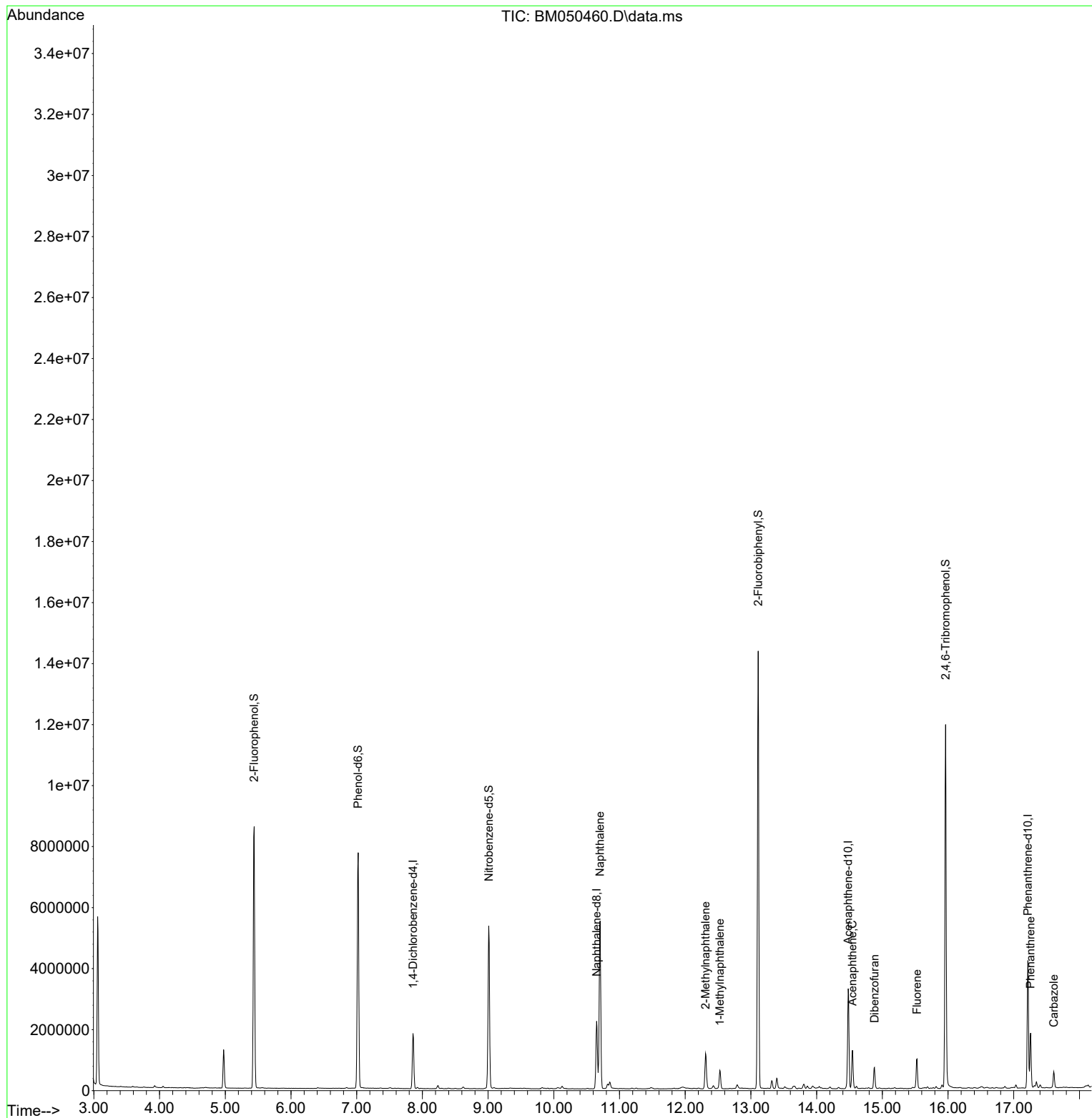
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	514304	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1853596	20.000	ng	-0.01
39) Acenaphthene-d10	14.480	164	1193542	20.000	ng	-0.01
64) Phenanthrene-d10	17.215	188	2415238	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	2529047	20.000	ng	-0.02
86) Perylene-d12	24.462	264	2680110	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3759347	125.821	ng	0.00
7) Phenol-d6	7.022	99	4413191	117.169	ng	0.00
23) Nitrobenzene-d5	9.010	82	2946564	81.101	ng	-0.01
42) 2,4,6-Tribromophenol	15.962	330	2287437	157.736	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	7492871	77.527	ng	-0.01
79) Terphenyl-d14	19.827	244	12275923	85.404	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	10.704	128	4503882	47.838	ng	100
37) 2-Methylnaphthalene	12.310	142	516159	8.683	ng	99
38) 1-Methylnaphthalene	12.527	142	289691	4.605	ng	99
52) Acenaphthene	14.545	154	461013	6.796	ng	99
55) Dibenzofuran	14.880	168	458114	4.383	ng	98
58) Fluorene	15.527	166	443103	5.147	ng	99
71) Phenanthrene	17.256	178	1121936	8.209	ng	99
73) Carbazole	17.609	167	349127	2.836	ng	99

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

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