

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071525\
 Data File : BM050461.D
 Acq On : 15 Jul 2025 18:51
 Operator : RC/JU
 Sample : Q2579-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-13-C

Quant Time: Jul 16 02:01:19 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)

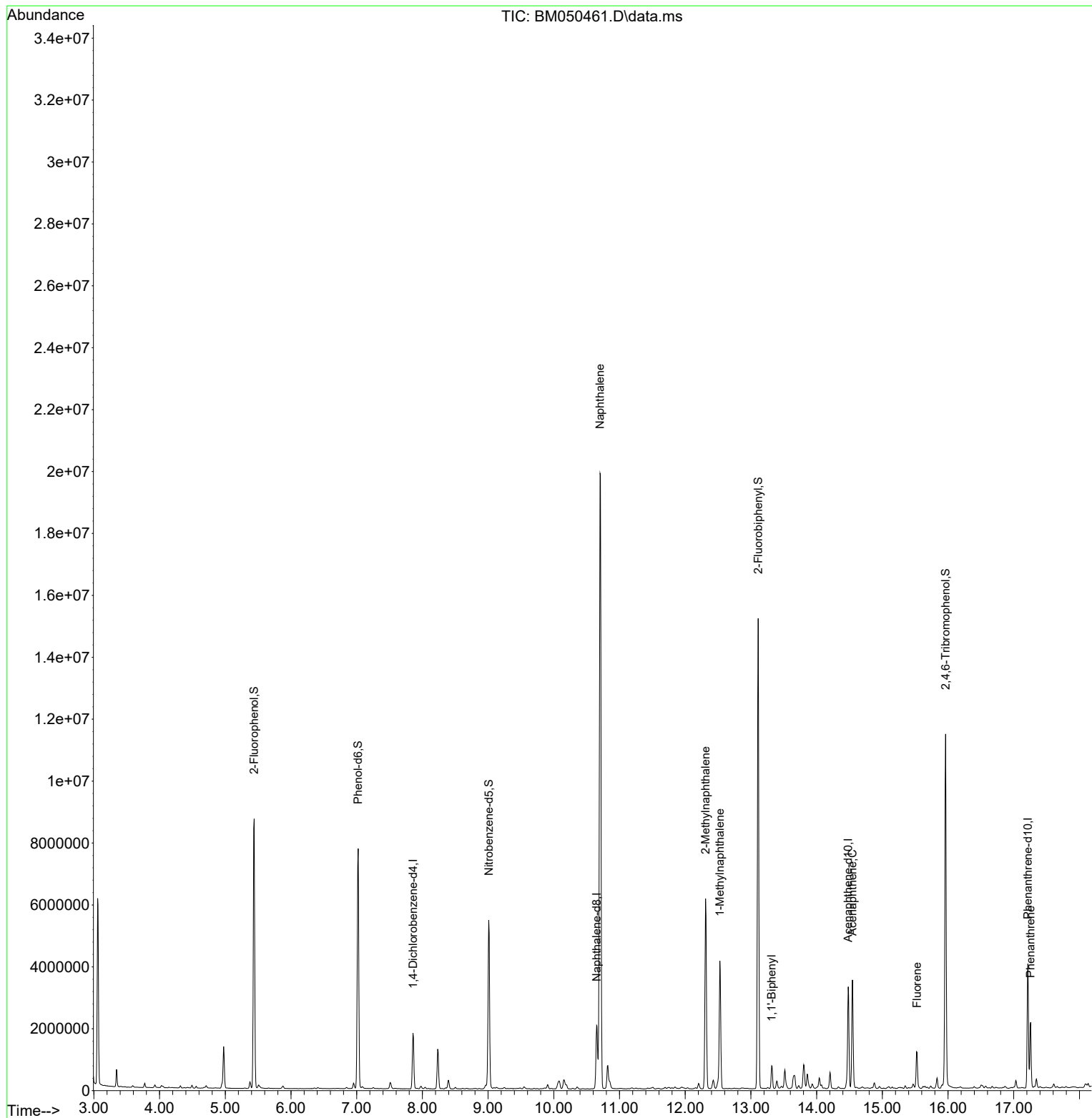
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	509697	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1822441	20.000	ng	-0.01
39) Acenaphthene-d10	14.480	164	1185090	20.000	ng	-0.01
64) Phenanthrene-d10	17.215	188	2353649	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	2415694	20.000	ng	-0.02
86) Perylene-d12	24.462	264	2549306	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3816425	128.886	ng	0.00
7) Phenol-d6	7.022	99	4413626	118.240	ng	0.00
23) Nitrobenzene-d5	9.010	82	3058960	85.634	ng	-0.01
42) 2,4,6-Tribromophenol	15.963	330	2256514	156.714	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	7894274	82.263	ng	-0.01
79) Terphenyl-d14	19.827	244	11949213	87.031	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	10.704	128	16057452	173.468	ng	# 94
37) 2-Methylnaphthalene	12.310	142	2855800	48.861	ng	98
38) 1-Methylnaphthalene	12.528	142	1925228	31.125	ng	99
46) 1,1'-Biphenyl	13.316	154	393195	4.472	ng	100
52) Acenaphthene	14.545	154	1273137	18.901	ng	99
58) Fluorene	15.527	166	566655	6.629	ng	99
71) Phenanthrene	17.257	178	1298144	9.747	ng	99

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

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