

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050438.D
 Acq On : 14 Jul 2025 22:54
 Operator : RC/JU
 Sample : Q2578-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A5-01-C

Quant Time: Jul 15 02:05:29 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)

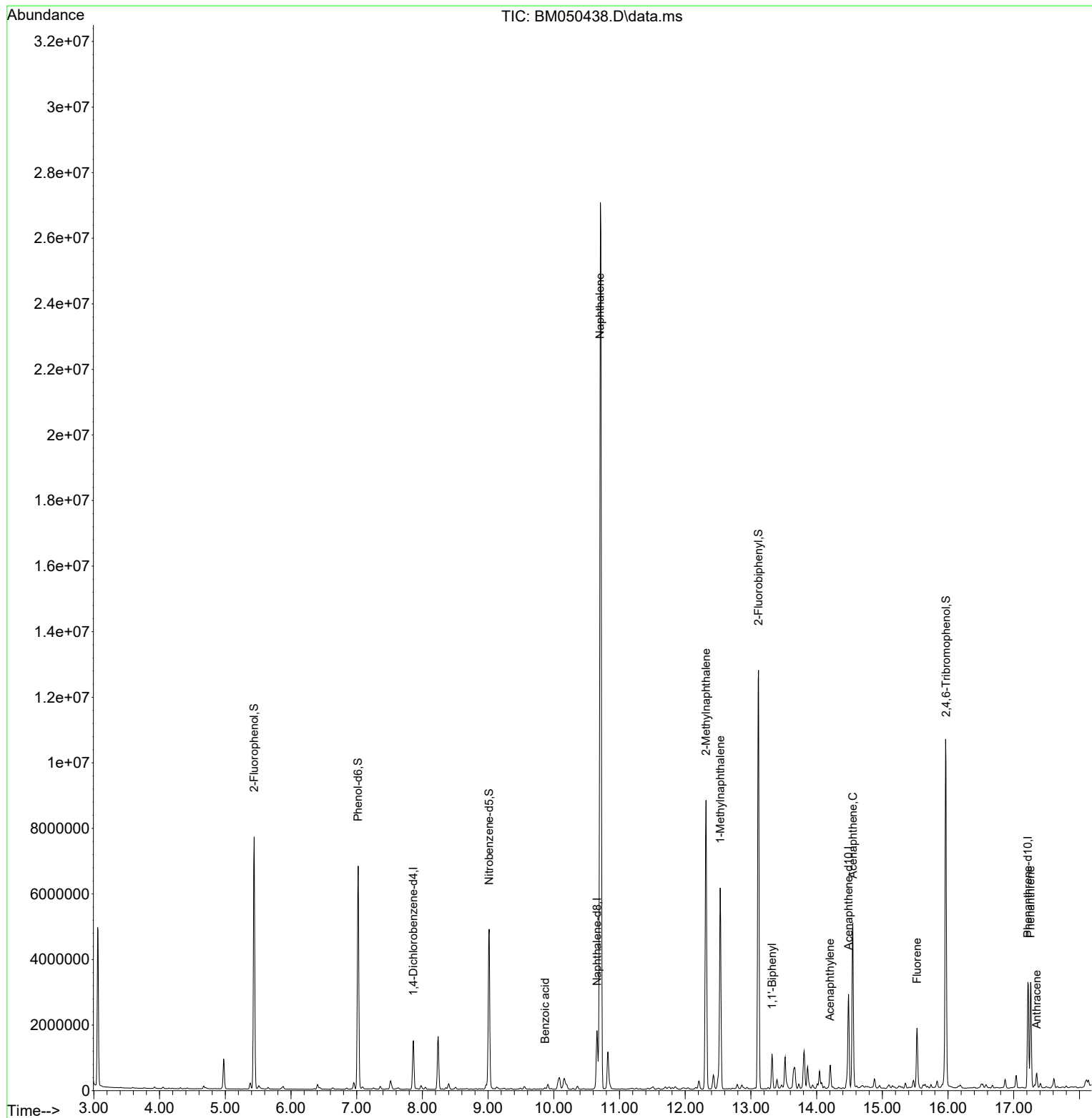
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	406303	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1510572	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	1001056	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	1993082	20.000	ng	-0.01
76) Chrysene-d12	21.439	240	1992538	20.000	ng	-0.02
86) Perylene-d12	24.468	264	2058330	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3156459	133.725	ng	0.00
7) Phenol-d6	7.022	99	3598339	120.930	ng	0.00
23) Nitrobenzene-d5	9.016	82	2652160	89.574	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1999830	164.420	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	6685623	82.476	ng	0.00
79) Terphenyl-d14	19.827	244	11365563	100.361	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	10.704	128	18532645	241.542	ng	# 91
32) Benzoic acid	9.869	122	15714	6.952	ng	# 81
37) 2-Methylnaphthalene	12.316	142	4048151	83.561	ng	99
38) 1-Methylnaphthalene	12.533	142	2728595	53.220	ng	99
46) 1,1'-Biphenyl	13.321	154	530987	7.149	ng	99
49) Acenaphthylene	14.204	152	301017	3.364	ng	# 95
52) Acenaphthene	14.551	154	1811059	31.830	ng	99
58) Fluorene	15.527	166	825826	11.437	ng	98
71) Phenanthrene	17.256	178	1968866	17.458	ng	100
72) Anthracene	17.351	178	277524	2.472	ng	99

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

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