

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050440.D
 Acq On : 15 Jul 2025 00:13
 Operator : RC/JU
 Sample : Q2578-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-10-C

Quant Time: Jul 15 02:06:04 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)

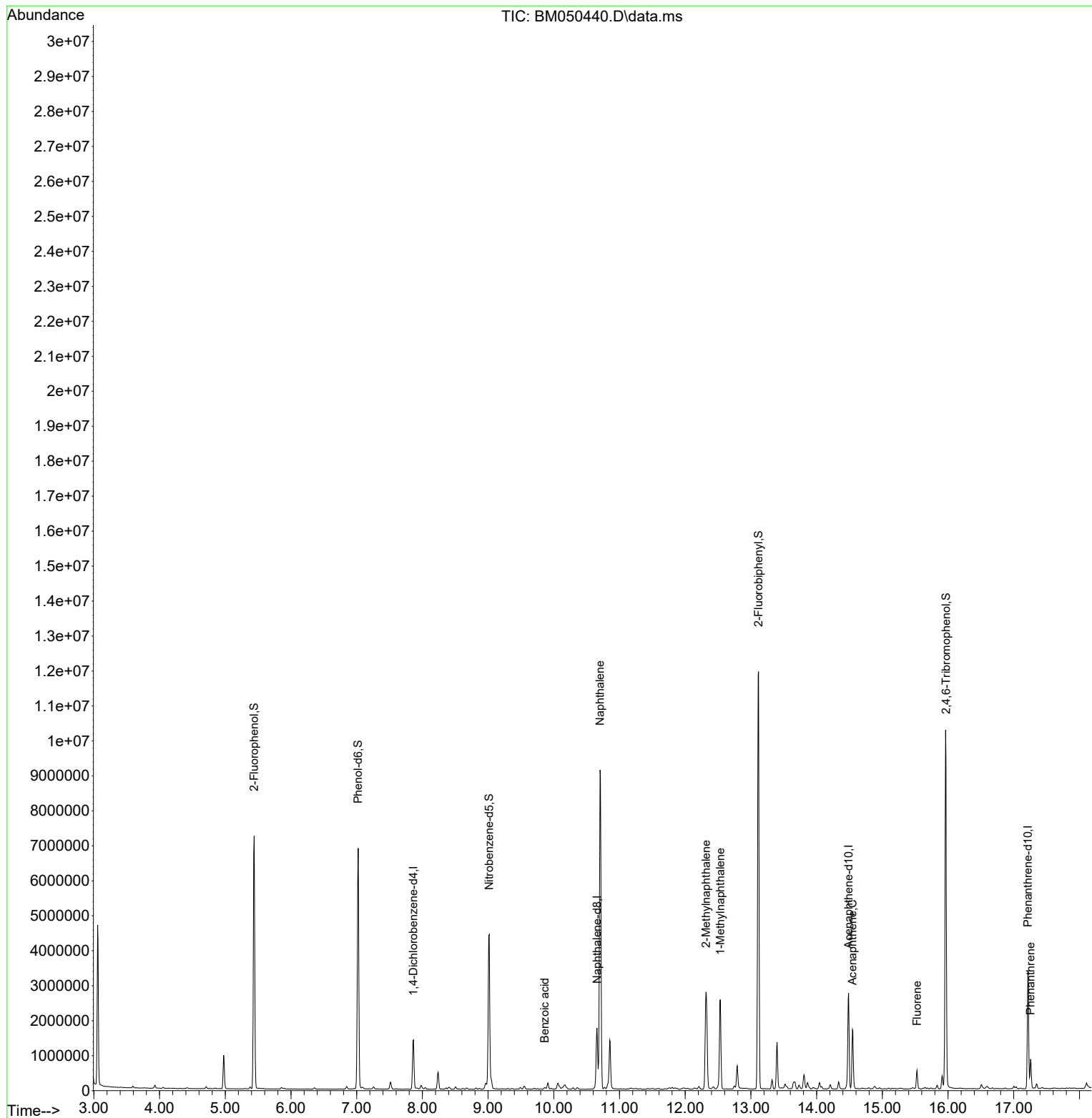
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	396203	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1451303	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	964408	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	1929252	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	1878020	20.000	ng	-0.02
86) Perylene-d12	24.462	264	1921664	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3062881	133.068	ng	0.00
7) Phenol-d6	7.022	99	3653974	125.930	ng	0.00
23) Nitrobenzene-d5	9.016	82	2454199	86.273	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1875599	160.066	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	6349448	81.305	ng	0.00
79) Terphenyl-d14	19.827	244	10799854	101.180	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	10.704	128	7482278	101.502	ng	100
32) Benzoic acid	9.863	122	13757	6.866	ng	83
37) 2-Methylnaphthalene	12.316	142	1075246	23.101	ng	99
38) 1-Methylnaphthalene	12.533	142	1189713	24.153	ng	98
52) Acenaphthene	14.545	154	600751	10.960	ng	98
58) Fluorene	15.527	166	221384	3.182	ng	98
71) Phenanthrene	17.257	178	496422	4.547	ng	100

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

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