

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050439.D
 Acq On : 14 Jul 2025 23:34
 Operator : RC/JU
 Sample : Q2578-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-09-C

Quant Time: Jul 15 02:05:48 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)

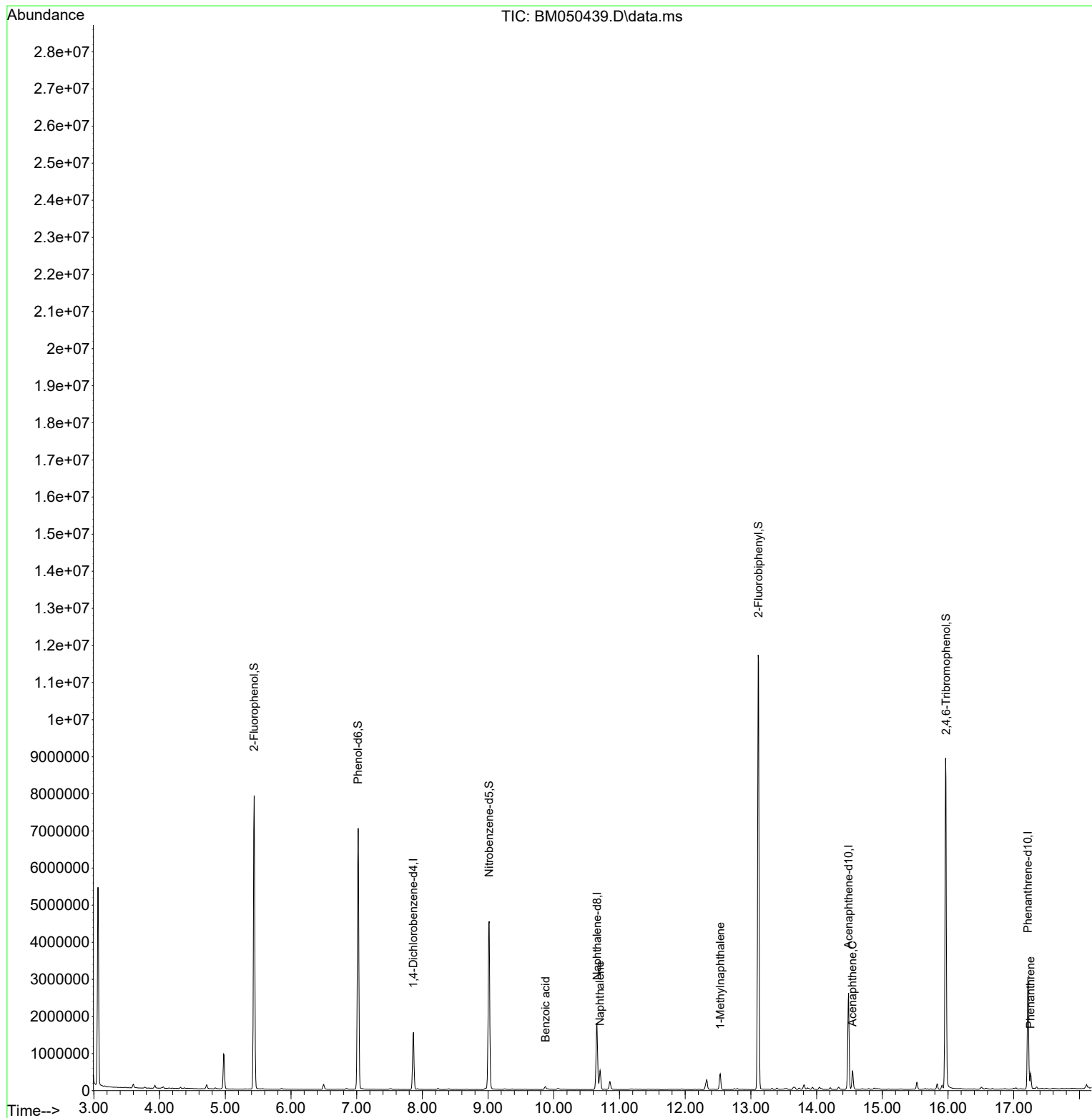
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	411418	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1469594	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	901307	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	1748222	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	1798338	20.000	ng	-0.02
86) Perylene-d12	24.468	264	1921480	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3287677	137.552	ng	0.00
7) Phenol-d6	7.022	99	3815235	126.625	ng	0.00
23) Nitrobenzene-d5	9.016	82	2553831	88.658	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1703848	155.589	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	6159994	84.402	ng	0.00
79) Terphenyl-d14	19.827	244	10338595	101.151	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	10.704	128	397005	5.319	ng	99
32) Benzoic acid	9.875	122	26605	7.677	ng	97
38) 1-Methylnaphthalene	12.533	142	193865	3.887	ng	97
52) Acenaphthene	14.545	154	171714	3.352	ng	98
71) Phenanthrene	17.257	178	243662	2.463	ng	98

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

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