

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071525\
 Data File : BM050462.D
 Acq On : 15 Jul 2025 19:31
 Operator : RC/JU
 Sample : Q2579-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-14-C

Quant Time: Jul 16 02:01:39 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)

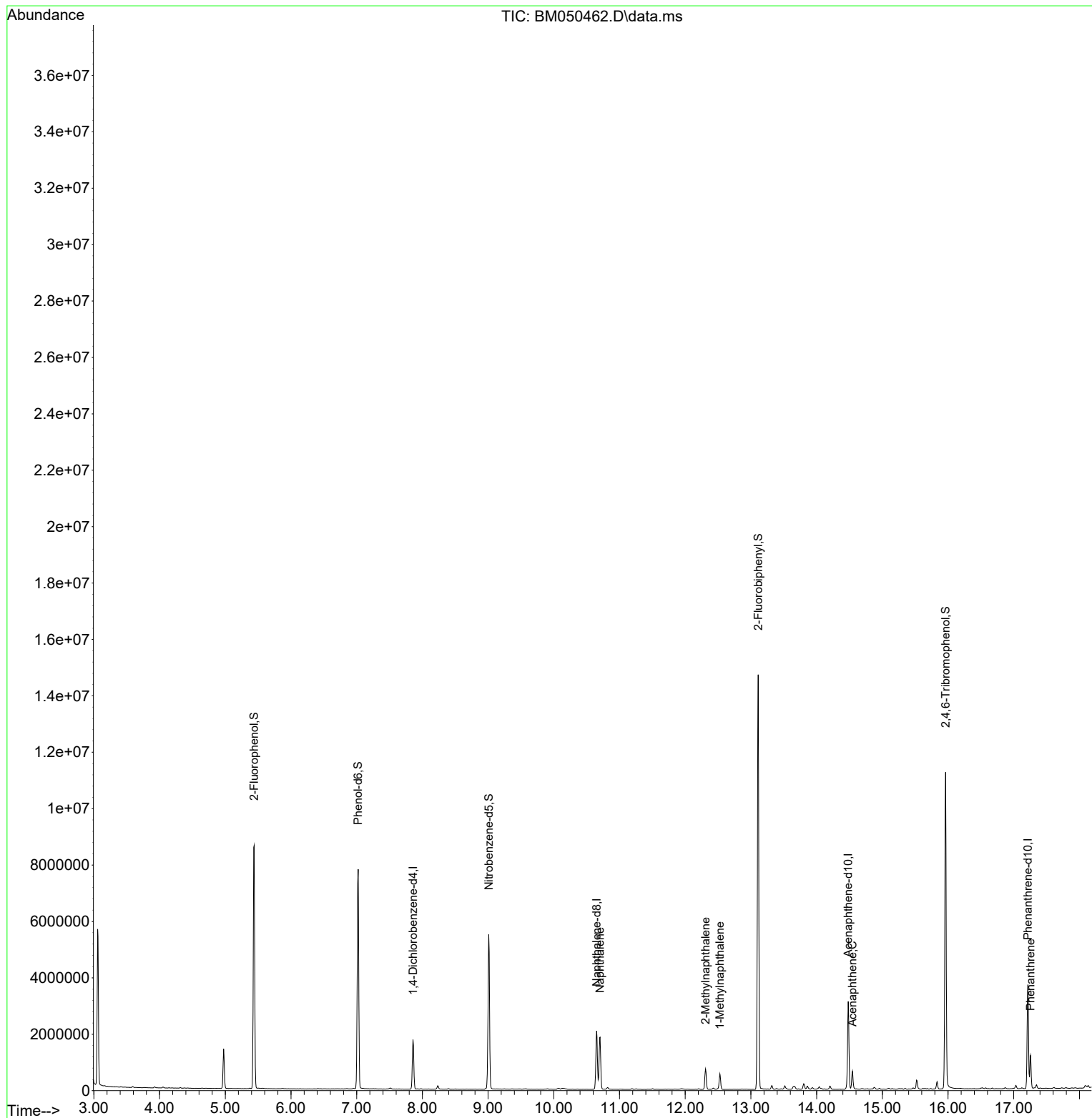
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	490534	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1731974	20.000	ng	-0.01
39) Acenaphthene-d10	14.480	164	1104765	20.000	ng	-0.01
64) Phenanthrene-d10	17.215	188	2272075	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	2341414	20.000	ng	-0.02
86) Perylene-d12	24.462	264	2438835	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3818830	134.006	ng	0.00
7) Phenol-d6	7.022	99	4431503	123.357	ng	0.00
23) Nitrobenzene-d5	9.010	82	2993647	88.183	ng	-0.01
42) 2,4,6-Tribromophenol	15.963	330	2248808	167.534	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	7678842	85.836	ng	-0.01
79) Terphenyl-d14	19.821	244	12141866	91.240	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	10.704	128	1566713	17.809	ng	100
37) 2-Methylnaphthalene	12.310	142	345027	6.212	ng	97
38) 1-Methylnaphthalene	12.528	142	259006	4.406	ng	98
52) Acenaphthene	14.545	154	225160	3.586	ng	99
71) Phenanthrene	17.257	178	730132	5.679	ng	99

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

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