

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071025\
 Data File : BM050397.D
 Acq On : 10 Jul 2025 19:59
 Operator : RC/JU
 Sample : Q2489-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G4(0-6)

Quant Time: Jul 11 01:26:54 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)

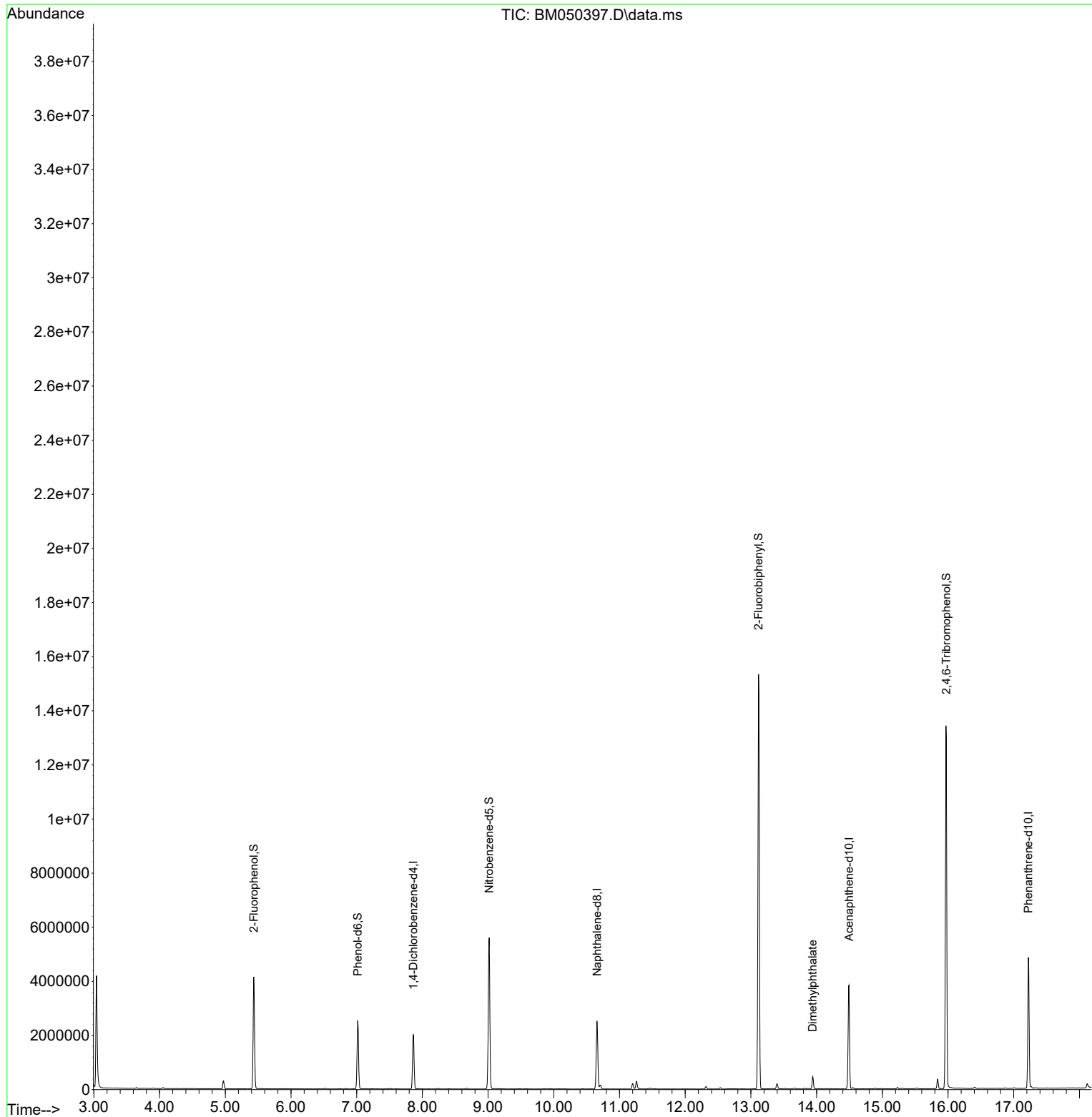
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	561976	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	2091010	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	1369169	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	2769354	20.000	ng	0.00
76) Chrysene-d12	21.451	240	2918200	20.000	ng	0.00
86) Perylene-d12	24.486	264	2873438	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1770139	54.219	ng	0.00
7) Phenol-d6	7.016	99	1368150	33.243	ng	0.00
23) Nitrobenzene-d5	9.016	82	3124285	76.229	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	2667817	160.369	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	8273364	74.622	ng	0.00
79) Terphenyl-d14	19.839	244	13018069	78.489	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.939	163	289254	3.069	ng	Qvalue 99

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

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