

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071525\
 Data File : BM050451.D
 Acq On : 15 Jul 2025 12:10
 Operator : RC/JU
 Sample : Q2489-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G2(6-12)

Quant Time: Jul 15 12:55:46 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)

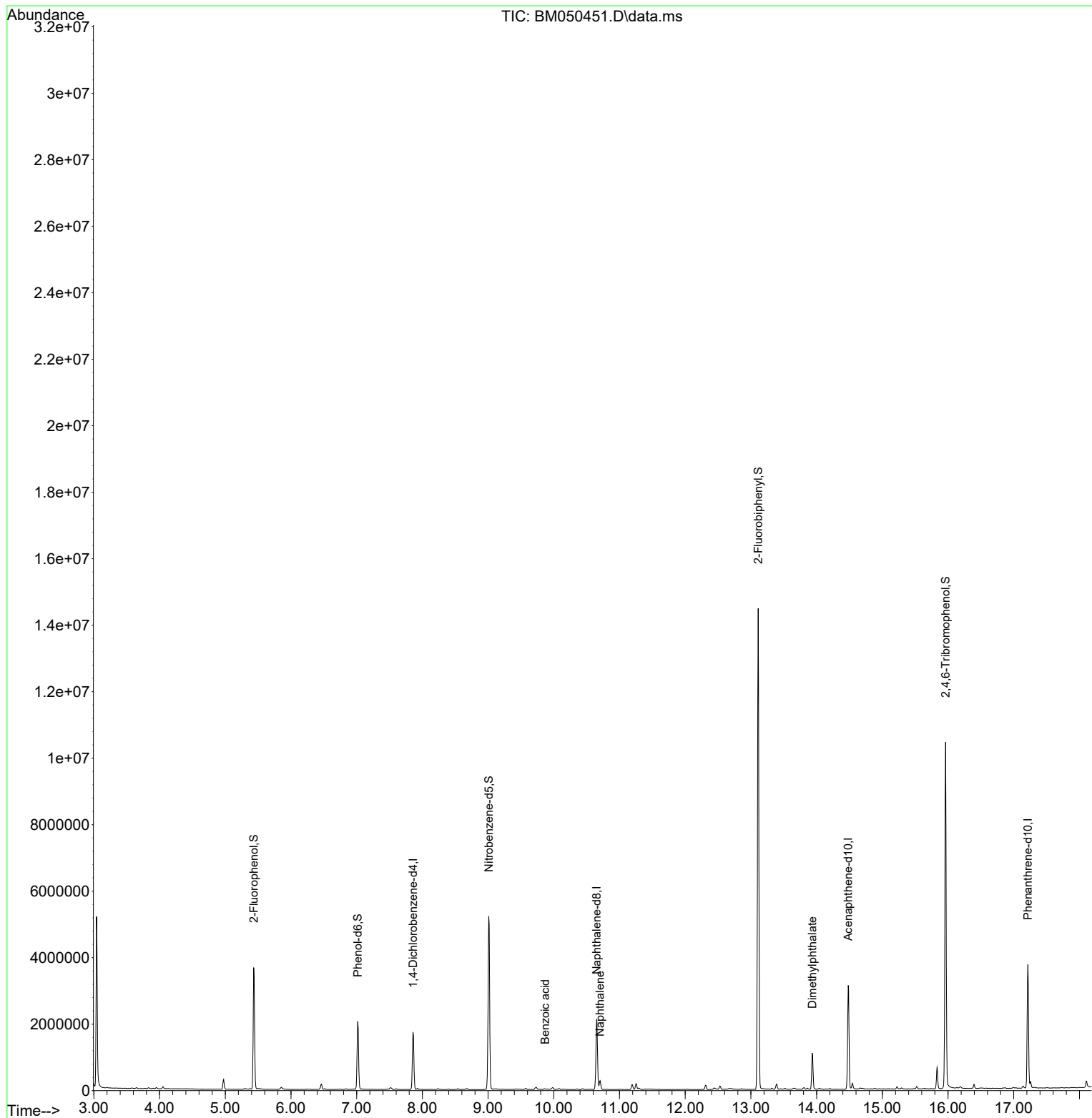
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	491189	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1785204	20.000	ng	-0.01
39) Acenaphthene-d10	14.480	164	1129360	20.000	ng	-0.01
64) Phenanthrene-d10	17.215	188	2268318	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	2383036	20.000	ng	-0.02
86) Perylene-d12	24.462	264	2558932	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1560429	54.684	ng	0.00
7) Phenol-d6	7.016	99	1125435	31.286	ng	0.00
23) Nitrobenzene-d5	9.010	82	2957069	84.508	ng	-0.01
42) 2,4,6-Tribromophenol	15.962	330	1987550	144.846	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	7510632	82.127	ng	-0.01
79) Terphenyl-d14	19.827	244	11933627	88.109	ng	-0.02
Target Compounds						
31) Naphthalene	10.704	128	205983	2.272	ng	99
32) Benzoic acid	9.869	122	8654m	6.430	ng	
50) Dimethylphthalate	13.933	163	726536	9.344	ng	99

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

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