

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071725\
 Data File : BM050477.D
 Acq On : 17 Jul 2025 16:37
 Operator : RC/JU
 Sample : Q2605-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 V908

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 07/18/2025
 Supervised By :Jagrut Upadhyay 07/18/2025

Quant Time: Jul 17 17:07:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 17 16:20:15 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	491498	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	1719925	20.000	ng	0.00
39) Acenaphthene-d10	14.474	164	1091025	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	2296820	20.000	ng	0.00
76) Chrysene-d12	21.433	240	2469819	20.000	ng	0.00
86) Perylene-d12	24.456	264	2717634	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	3316543	116.152	ng	0.00
7) Phenol-d6	7.010	99	3782401	105.082	ng	0.00
23) Nitrobenzene-d5	8.998	82	2694093	79.915	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	2120838	159.990	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	6790256	76.859	ng	0.00
79) Terphenyl-d14	19.815	244	11967090	85.252	ng	0.00
Target Compounds						
32) Benzoic acid	9.851	122	16044m	6.851	ng	Qvalue

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

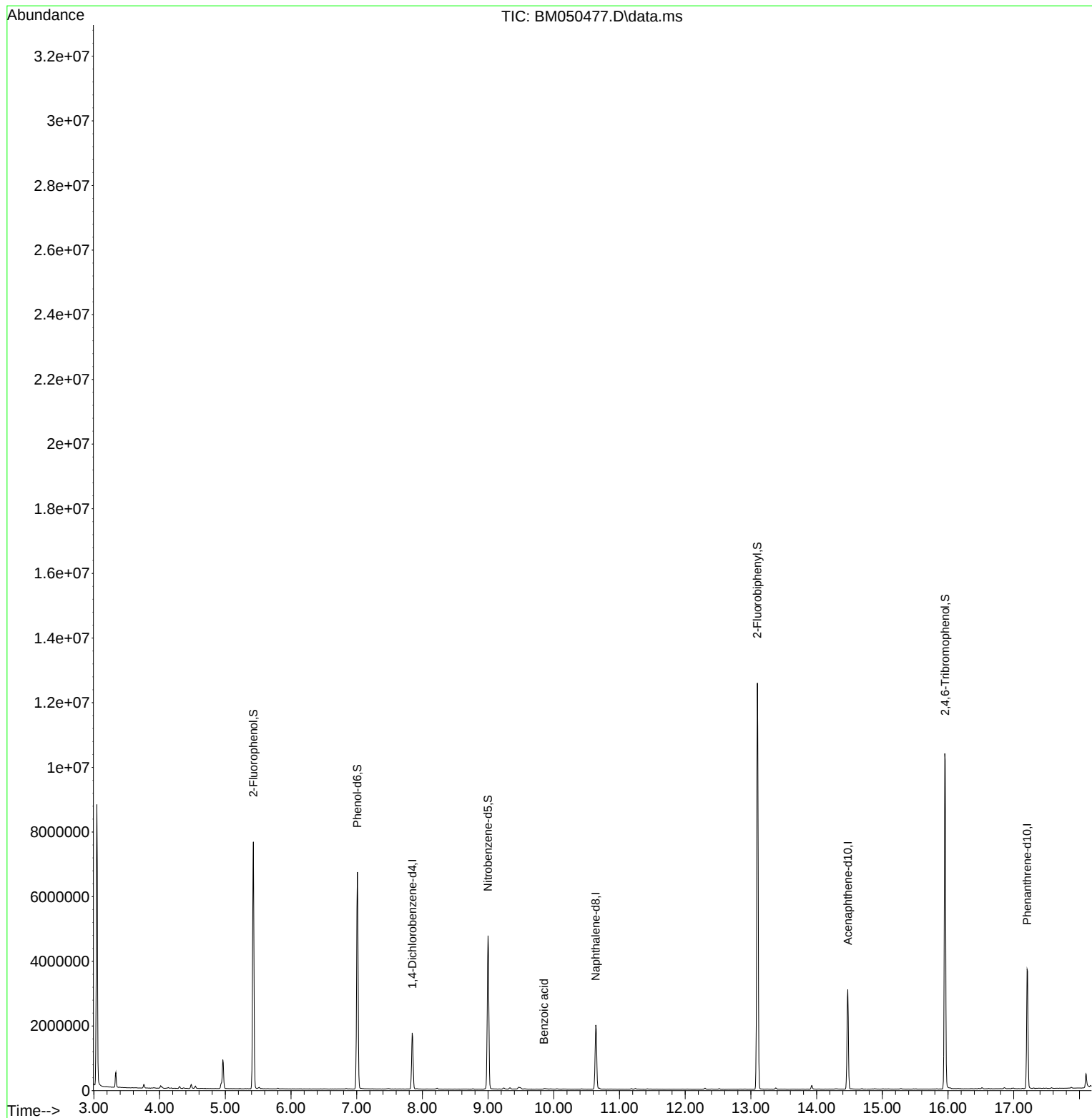
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