

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050216.D
 Acq On : 06 Jun 2025 16:06
 Operator : RC/JU
 Sample : Q2207-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BU-703-COMP-01

Quant Time: Jun 06 16:54:05 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 05 16:20:25 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Rahul
 Chavli

06/09/2025
 Supervised By :Jagrut
 Upadhyay

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	289452	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1124813	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	638368	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1162706	20.000	ng	0.00
76) Chrysene-d12	21.380	240	1061848	20.000	ng	0.00
86) Perylene-d12	24.362	264	1128394	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2181497	125.722	ng	0.00
7) Phenol-d6	6.951	99	2599499	113.804	ng	0.00
23) Nitrobenzene-d5	8.934	82	1742254	80.557	ng	-0.01
42) 2,4,6-Tribromophenol	15.892	330	925629	127.480	ng	-0.01
45) 2-Fluorobiphenyl	13.039	172	3623948	76.857	ng	-0.01
79) Terphenyl-d14	19.774	244	4374976	78.078	ng	-0.01
Target Compounds						
32) Benzoic acid	9.792	122	27885m	2.543	ng	Qvalue

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

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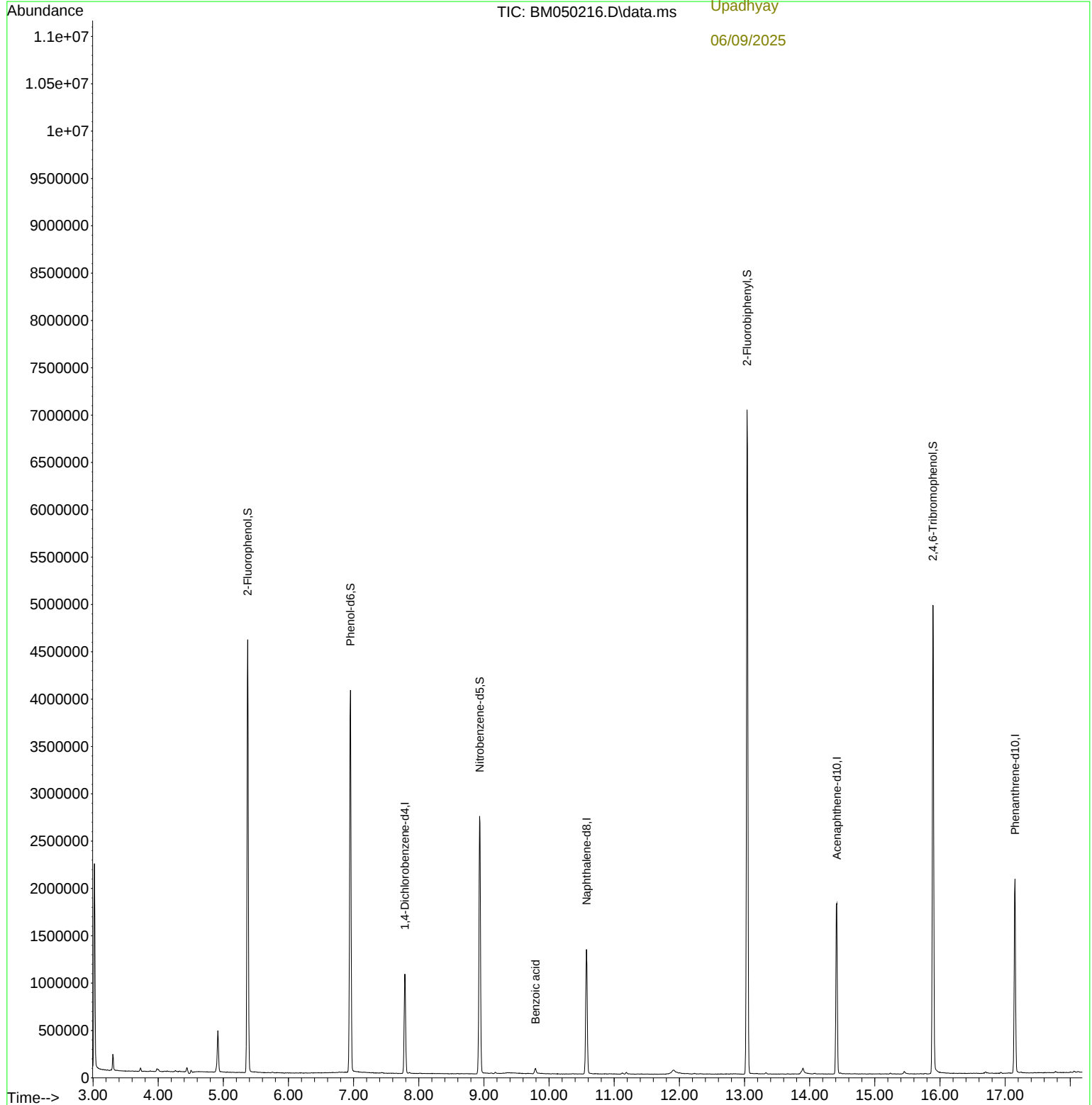
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