

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050217.D
 Acq On : 06 Jun 2025 16:45
 Operator : RC/JU
 Sample : Q2207-18
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BU-703-COMP-02

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 06/09/2025
 Supervised By :Jagrut Upadhyay 06/09/2025

Quant Time: Jun 06 17:25:20 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	292016	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1122506	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	609641	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1145449	20.000	ng	0.00
76) Chrysene-d12	21.380	240	1073649	20.000	ng	0.00
86) Perylene-d12	24.368	264	686399	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2097030	119.793	ng	0.00
7) Phenol-d6	6.951	99	2443008	106.014	ng	0.00
23) Nitrobenzene-d5	8.934	82	1711052	79.277	ng	-0.01
42) 2,4,6-Tribromophenol	15.892	330	886301	127.815	ng	-0.01
45) 2-Fluorobiphenyl	13.039	172	3469477	77.048	ng	-0.01
79) Terphenyl-d14	19.780	244	4338429	76.575	ng	0.00
Target Compounds						
32) Benzoic acid	9.792	122	23822m	2.177	ng	Qvalue

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

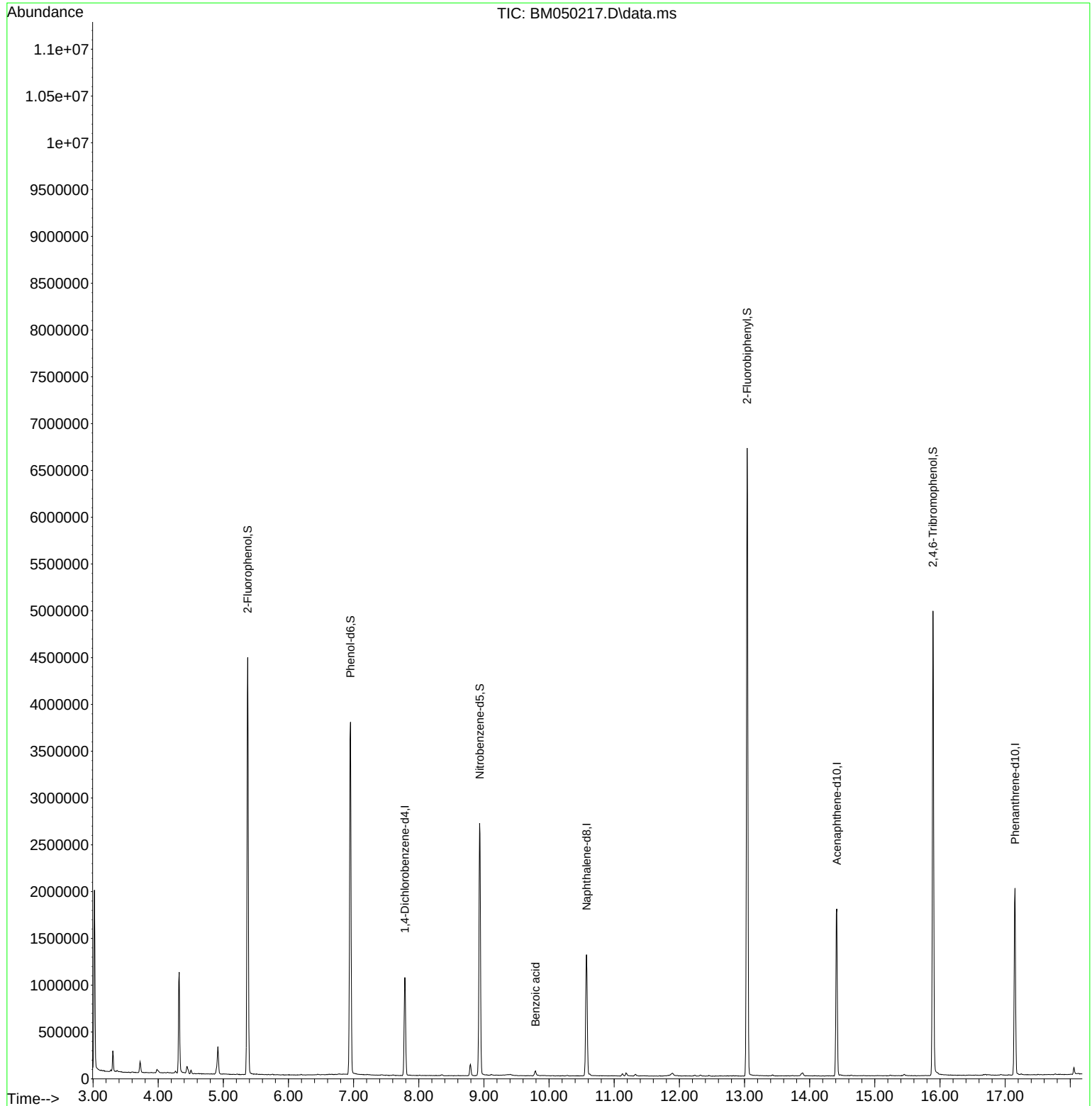
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