

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050215.D
 Acq On : 06 Jun 2025 15:26
 Operator : RC/JU
 Sample : Q2194-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP-13

Manual Integrations
 APPROVED

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

Quant Time: Jun 06 16:25:04 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	294115	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1102386	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	616752	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1152106	20.000	ng	0.00
76) Chrysene-d12	21.380	240	1157843	20.000	ng	0.00
86) Perylene-d12	24.368	264	1249334	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2201110	124.841	ng	0.00
7) Phenol-d6	6.951	99	2570679	110.758	ng	0.00
23) Nitrobenzene-d5	8.934	82	1726328	81.444	ng	-0.01
42) 2,4,6-Tribromophenol	15.892	330	945363	134.761	ng	-0.01
45) 2-Fluorobiphenyl	13.039	172	3576220	78.503	ng	-0.01
79) Terphenyl-d14	19.780	244	4778745	78.213	ng	0.00
Target Compounds						
32) Benzoic acid	9.804	122	66073m	6.149	ng	Qvalue

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

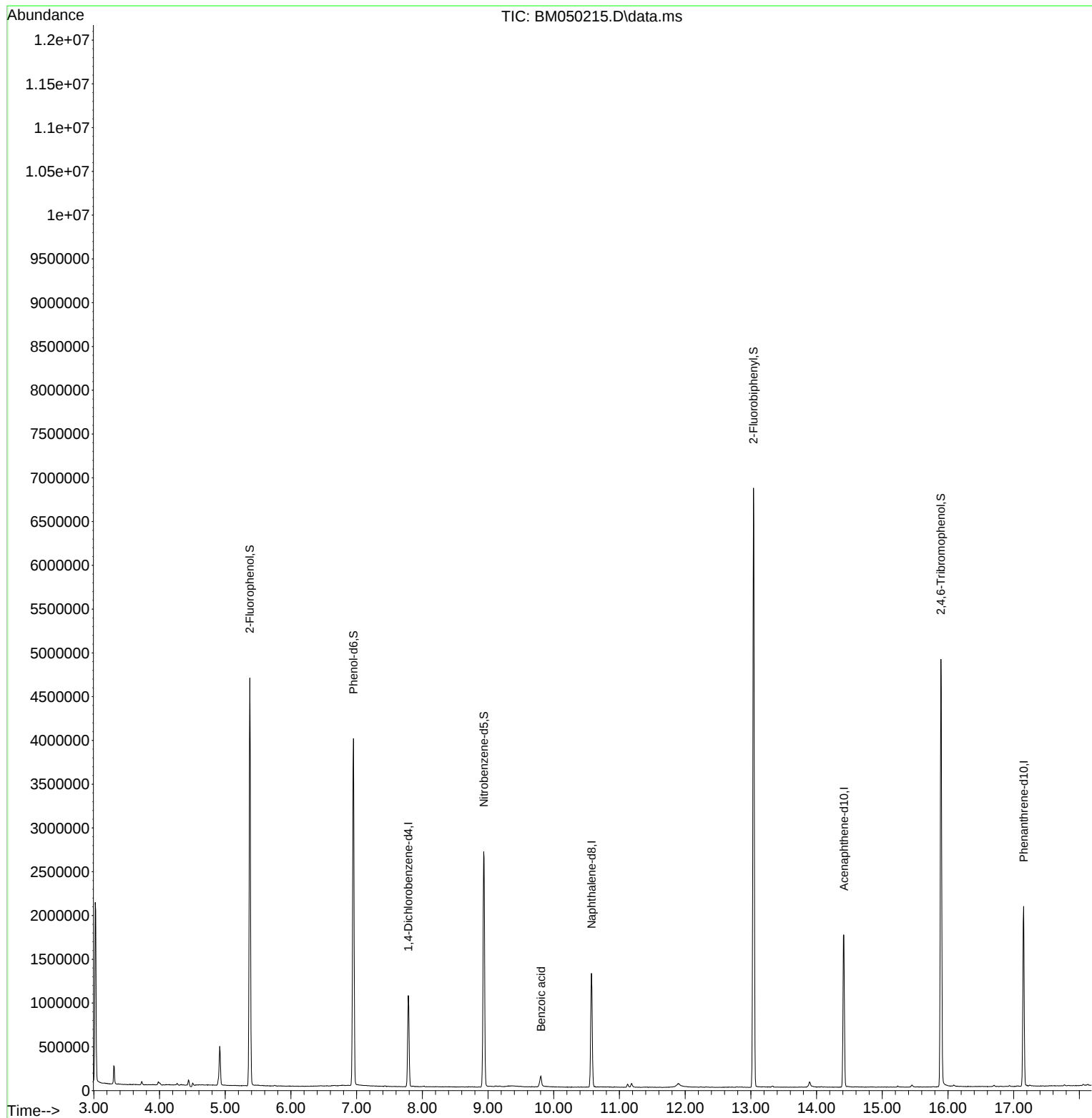
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