

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041125\
 Data File : BM049899.D
 Acq On : 11 Apr 2025 13:38
 Operator : RC/JU
 Sample : Q1752-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-3

Quant Time: Apr 11 14:20:00 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 04:00:55 2025
 Response via : Initial Calibration

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

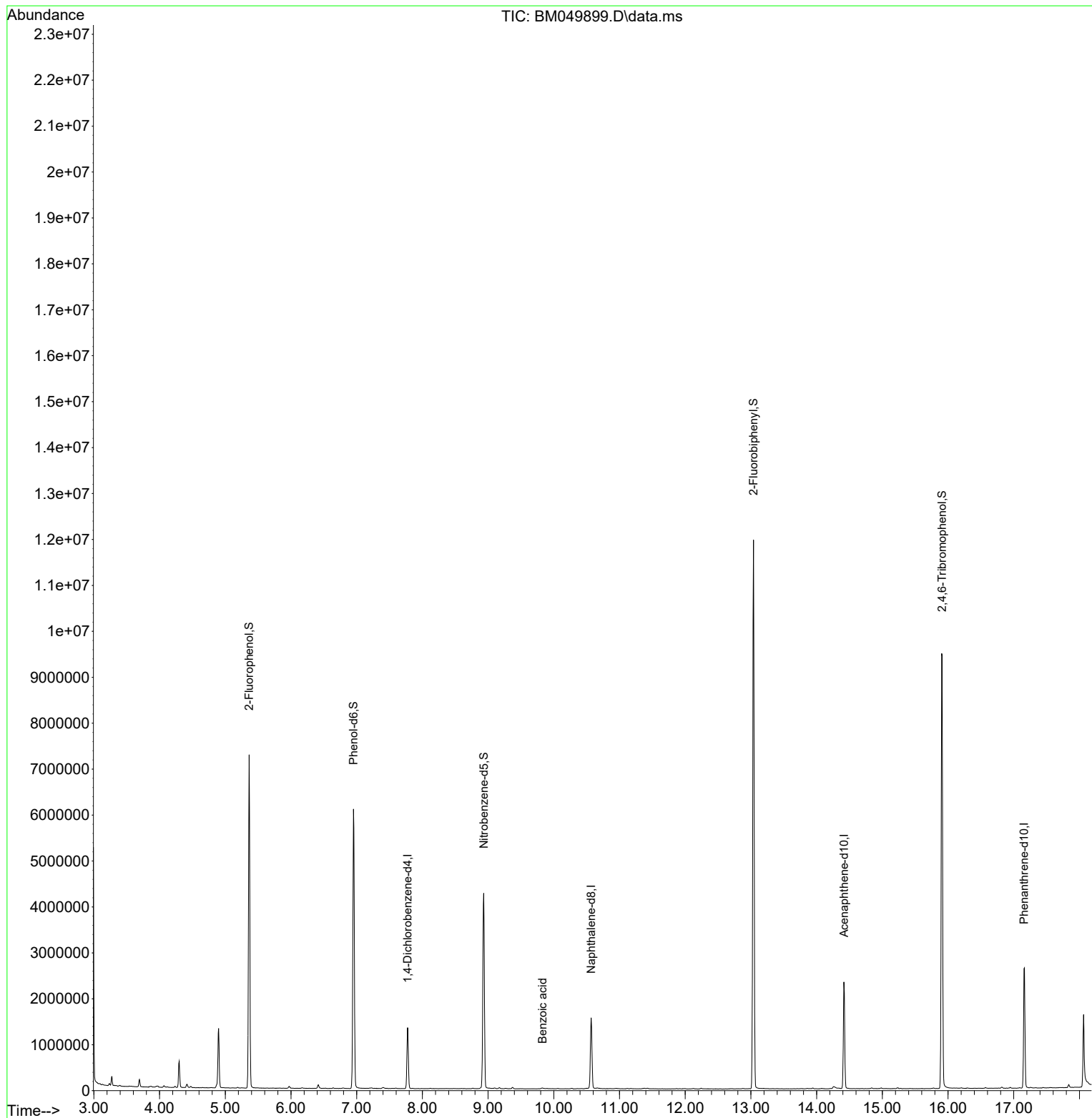
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	383103	20.000	ng	0.00
21) Naphthalene-d8	10.569	136	1310054	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	842074	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	1638534	20.000	ng	0.00
76) Chrysene-d12	21.397	240	1373545	20.000	ng	0.00
86) Perylene-d12	24.397	264	1040897	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	3174688	140.716	ng	0.00
7) Phenol-d6	6.951	99	3638660	129.628	ng	0.00
23) Nitrobenzene-d5	8.934	82	2477988	105.330	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	2114567	172.642	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	6478821	104.330	ng	0.00
79) Terphenyl-d14	19.786	244	9214121	125.716	ng	0.00
Target Compounds						
32) Benzoic acid	9.828	122	11502	9.099	ng	Qvalue 94

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

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