

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041125\
 Data File : BM049901.D
 Acq On : 11 Apr 2025 14:57
 Operator : RC/JU
 Sample : Q1753-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-2

Quant Time: Apr 11 15:35:21 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)

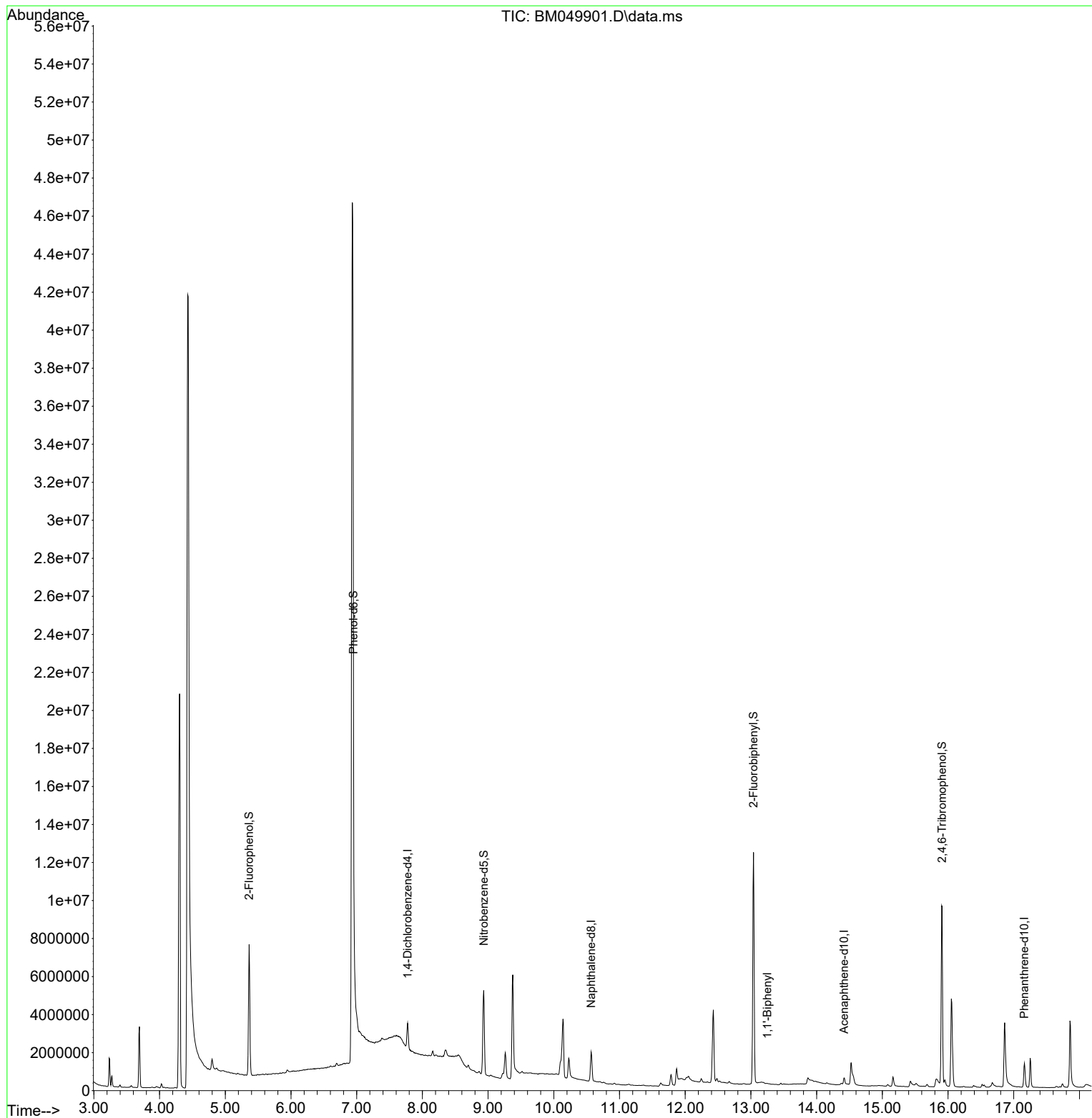
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	378429	20.000	ng	0.00
21) Naphthalene-d8	10.569	136	1169151	20.000	ng	# 0.00
39) Acenaphthene-d10	14.416	164	68395	20.000	ng	# 0.00
64) Phenanthrene-d10	17.163	188	521187	20.000	ng	0.00
76) Chrysene-d12	21.403	240	2630	20.000	ng	# 0.00
86) Perylene-d12	24.415	264	301	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	2953160	132.513	ng	0.00
7) Phenol-d6	6.951	99	384070	13.852	ng	0.00
23) Nitrobenzene-d5	8.934	82	2546713	121.297	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	1982601	1992.904	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	6645650	1317.578	ng	0.00
79) Terphenyl-d14	19.786	244	2165574	15431.153	ng	0.00
Target Compounds						
46) 1,1'-Biphenyl	13.251	154	13818	2.718	ng	Qvalue 92

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

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