

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042125\
 Data File : BM049983.D
 Acq On : 21 Apr 2025 13:56
 Operator : RC/JU
 Sample : Q1832-36
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PL-714-COMP-6

Quant Time: Apr 21 14:32:19 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)

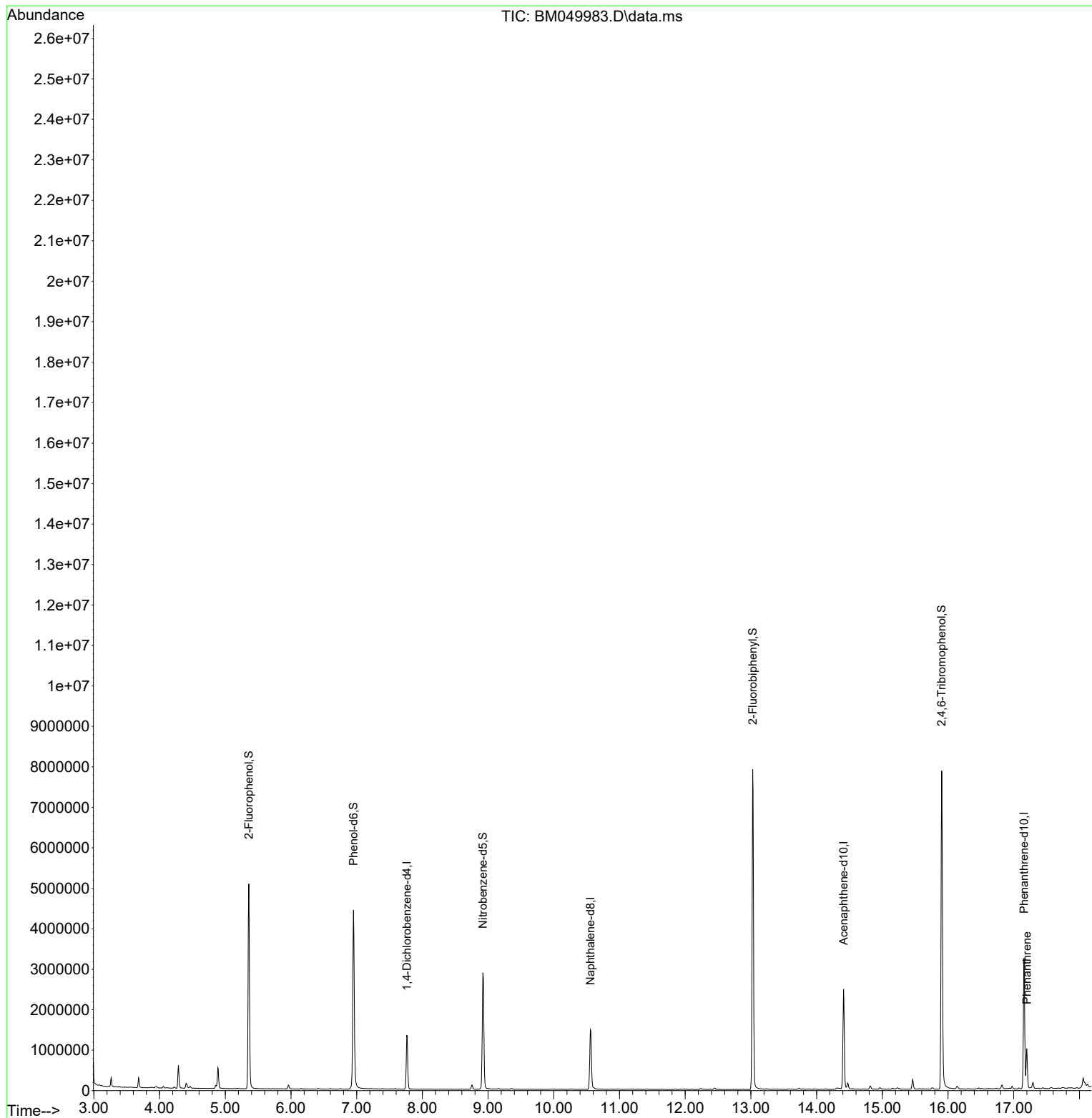
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	388091	20.000	ng	-0.02
21) Naphthalene-d8	10.557	136	1366819	20.000	ng	-0.02
39) Acenaphthene-d10	14.410	164	901913	20.000	ng	-0.01
64) Phenanthrene-d10	17.157	188	1970480	20.000	ng	0.00
76) Chrysene-d12	21.398	240	1933004	20.000	ng	0.00
86) Perylene-d12	24.403	264	523504	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2296804	100.496	ng	0.00
7) Phenol-d6	6.952	99	2790504	98.134	ng	0.00
23) Nitrobenzene-d5	8.922	82	1748303	71.227	ng	-0.01
42) 2,4,6-Tribromophenol	15.904	330	1789626	136.418	ng	0.00
45) 2-Fluorobiphenyl	13.028	172	4465796	67.142	ng	-0.02
79) Terphenyl-d14	19.786	244	10267242	99.541	ng	0.00
Target Compounds						
71) Phenanthrene	17.198	178	610224	5.649	ng	Qvalue 100

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

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