

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041125\
 Data File : BM049900.D
 Acq On : 11 Apr 2025 14:18
 Operator : RC/JU
 Sample : Q1753-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-1

Quant Time: Apr 11 14:58:50 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	372488	20.000	ng	0.00
21) Naphthalene-d8	10.569	136	1259224	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	792777	20.000	ng	0.00
64) Phenanthrene-d10	17.163	188	1485571	20.000	ng	0.00
76) Chrysene-d12	21.398	240	1295856	20.000	ng	0.00
86) Perylene-d12	24.397	264	1363805	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	3164020	144.239	ng	0.00
7) Phenol-d6	6.951	99	3562757	130.541	ng	0.00
23) Nitrobenzene-d5	8.934	82	2487274	109.992	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	1967380	170.613	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	6387041	109.248	ng	0.00
79) Terphenyl-d14	19.786	244	9025998	130.533	ng	0.00

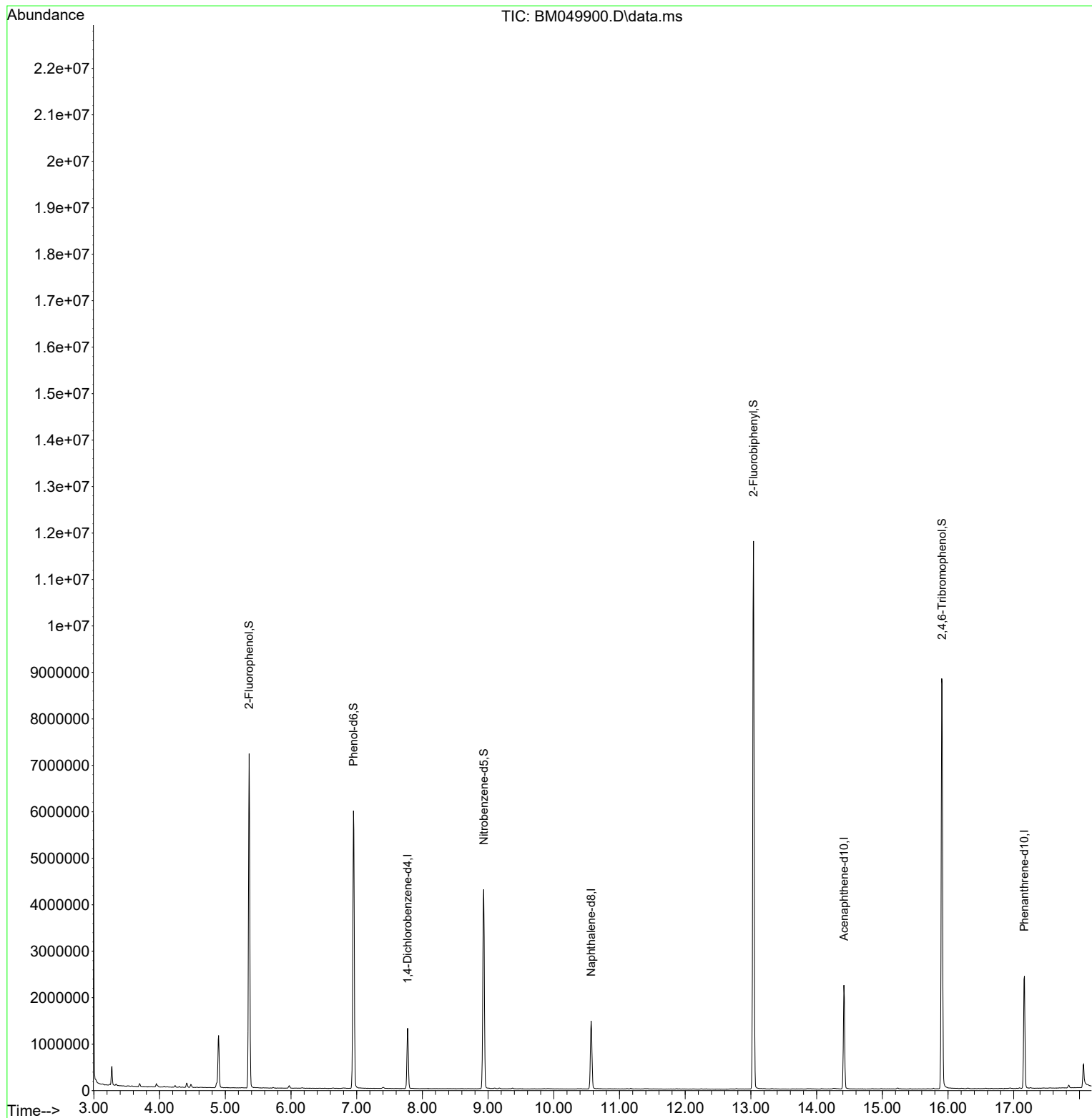
Target Compounds	Qvalue

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

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