

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040225\
 Data File : BM049789.D
 Acq On : 02 Apr 2025 14:07
 Operator : RC/JU
 Sample : Q1680-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-4

Quant Time: Apr 02 14:45:07 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 13 14:55:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	283645	20.000	ng	-0.02
21) Naphthalene-d8	10.580	136	957401	20.000	ng	-0.02
39) Acenaphthene-d10	14.427	164	601649	20.000	ng	-0.02
64) Phenanthrene-d10	17.168	188	1212495	20.000	ng	-0.02
76) Chrysene-d12	21.403	240	1401093	20.000	ng	-0.02
86) Perylene-d12	24.397	264	1440719	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	1725938	102.671	ng	-0.02
7) Phenol-d6	6.957	99	2157044	102.248	ng	-0.02
23) Nitrobenzene-d5	8.939	82	1282496	68.775	ng	-0.02
42) 2,4,6-Tribromophenol	15.915	330	855826	121.818	ng	-0.02
45) 2-Fluorobiphenyl	13.051	172	3108434	64.704	ng	-0.02
79) Terphenyl-d14	19.792	244	5297131	63.710	ng	-0.02

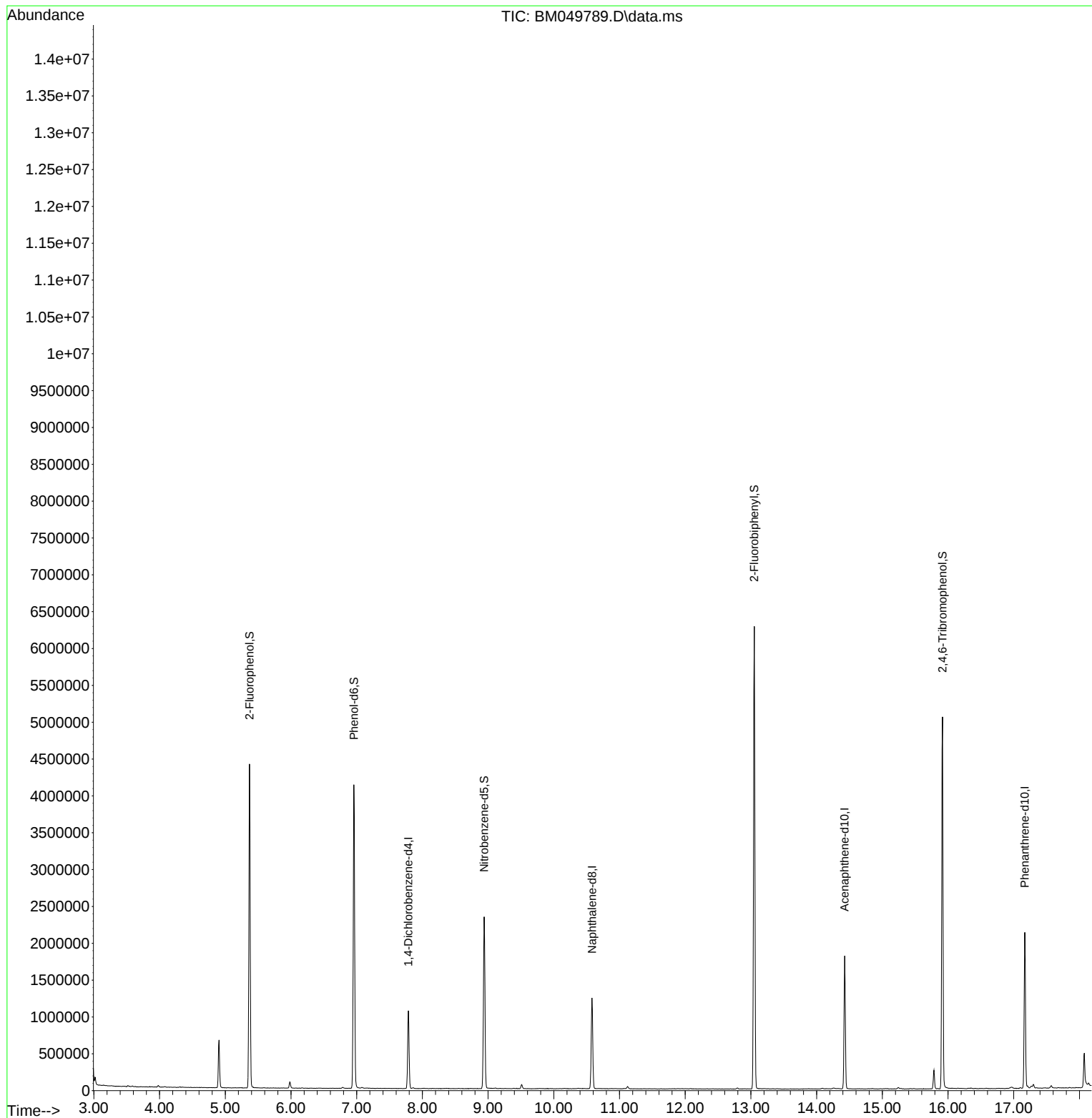
Target Compounds	Qvalue

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

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