

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022825\
 Data File : BM049680.D
 Acq On : 28 Feb 2025 15:52
 Operator : RC/JU
 Sample : PB166933BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB166933BL

Quant Time: Feb 28 16:39:28 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 04:55:28 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	267882	20.000	ng	-0.02
21) Naphthalene-d8	10.599	136	881999	20.000	ng	-0.03
39) Acenaphthene-d10	14.439	164	497065	20.000	ng	-0.03
64) Phenanthrene-d10	17.180	188	897905	20.000	ng	-0.02
76) Chrysene-d12	21.415	240	738243	20.000	ng	-0.03
86) Perylene-d12	24.433	264	796061	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	1694874	101.743	ng	0.00
7) Phenol-d6	6.975	99	2016933	92.416	ng	0.00
23) Nitrobenzene-d5	8.957	82	1870101	109.062	ng	-0.02
42) 2,4,6-Tribromophenol	15.928	330	566821	96.238	ng	-0.02
45) 2-Fluorobiphenyl	13.069	172	4454073	116.082	ng	-0.02
79) Terphenyl-d14	19.810	244	6284003	129.983	ng	-0.02

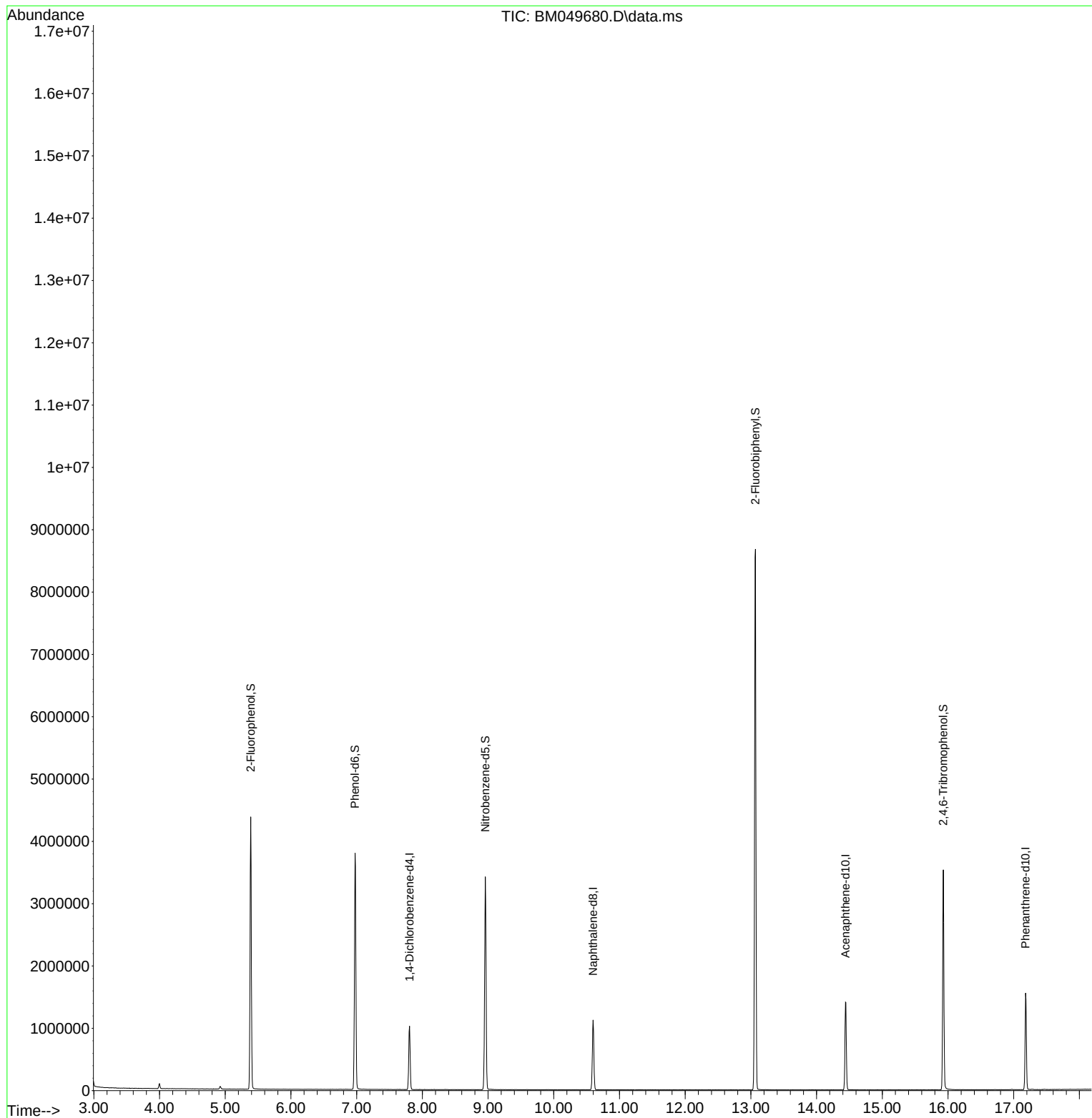
Target Compounds	Qvalue

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

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