

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040425\
 Data File : BM049818.D
 Acq On : 04 Apr 2025 11:17
 Operator : RC/JU
 Sample : PB167436BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB167436BL

Quant Time: Apr 04 12:01:27 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	411389	20.000	ng	-0.02
21) Naphthalene-d8	10.580	136	1389933	20.000	ng	-0.02
39) Acenaphthene-d10	14.427	164	844008	20.000	ng	-0.02
64) Phenanthrene-d10	17.168	188	1547901	20.000	ng	-0.02
76) Chrysene-d12	21.403	240	1260667	20.000	ng	-0.02
86) Perylene-d12	24.403	264	1277571	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	3778442	154.974	ng	-0.01
7) Phenol-d6	6.963	99	4545239	148.552	ng	-0.02
23) Nitrobenzene-d5	8.945	82	2727991	100.767	ng	-0.02
42) 2,4,6-Tribromophenol	15.915	330	1955686	198.436	ng	-0.02
45) 2-Fluorobiphenyl	13.051	172	6997158	103.827	ng	-0.02
79) Terphenyl-d14	19.797	244	9561220	127.804	ng	-0.01

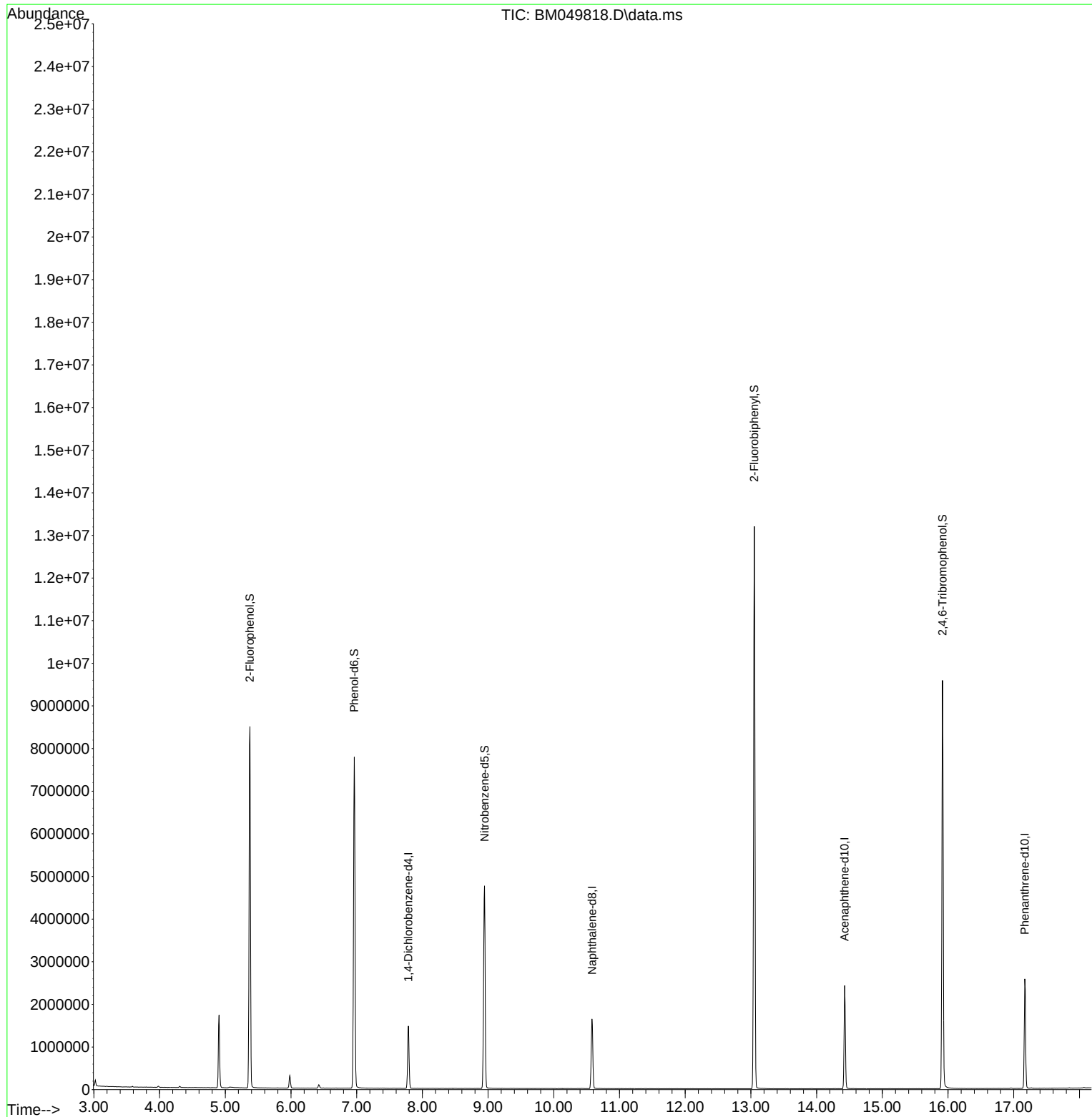
Target Compounds	Qvalue

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

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