

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081924\
 Data File : BM047249.D
 Acq On : 19 Aug 2024 12:21
 Operator : RC/JU
 Sample : PB162757BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162757BL

Quant Time: Aug 19 13:02:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	138664	20.000	ng	-0.01
21) Naphthalene-d8	10.122	136	543168	20.000	ng	-0.01
39) Acenaphthene-d10	14.027	164	380393	20.000	ng	0.00
64) Phenanthrene-d10	16.792	188	888654	20.000	ng	0.00
76) Chrysene-d12	21.033	240	1101328	20.000	ng	0.00
86) Perylene-d12	23.738	264	1446229	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	981765	126.511	ng	0.00
7) Phenol-d6	6.581	99	1282289	119.918	ng	0.00
23) Nitrobenzene-d5	8.510	82	820053	76.021	ng	-0.01
42) 2,4,6-Tribromophenol	15.539	330	636024	144.200	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	1970497	79.905	ng	-0.01
79) Terphenyl-d14	19.462	244	4117692	80.115	ng	0.00

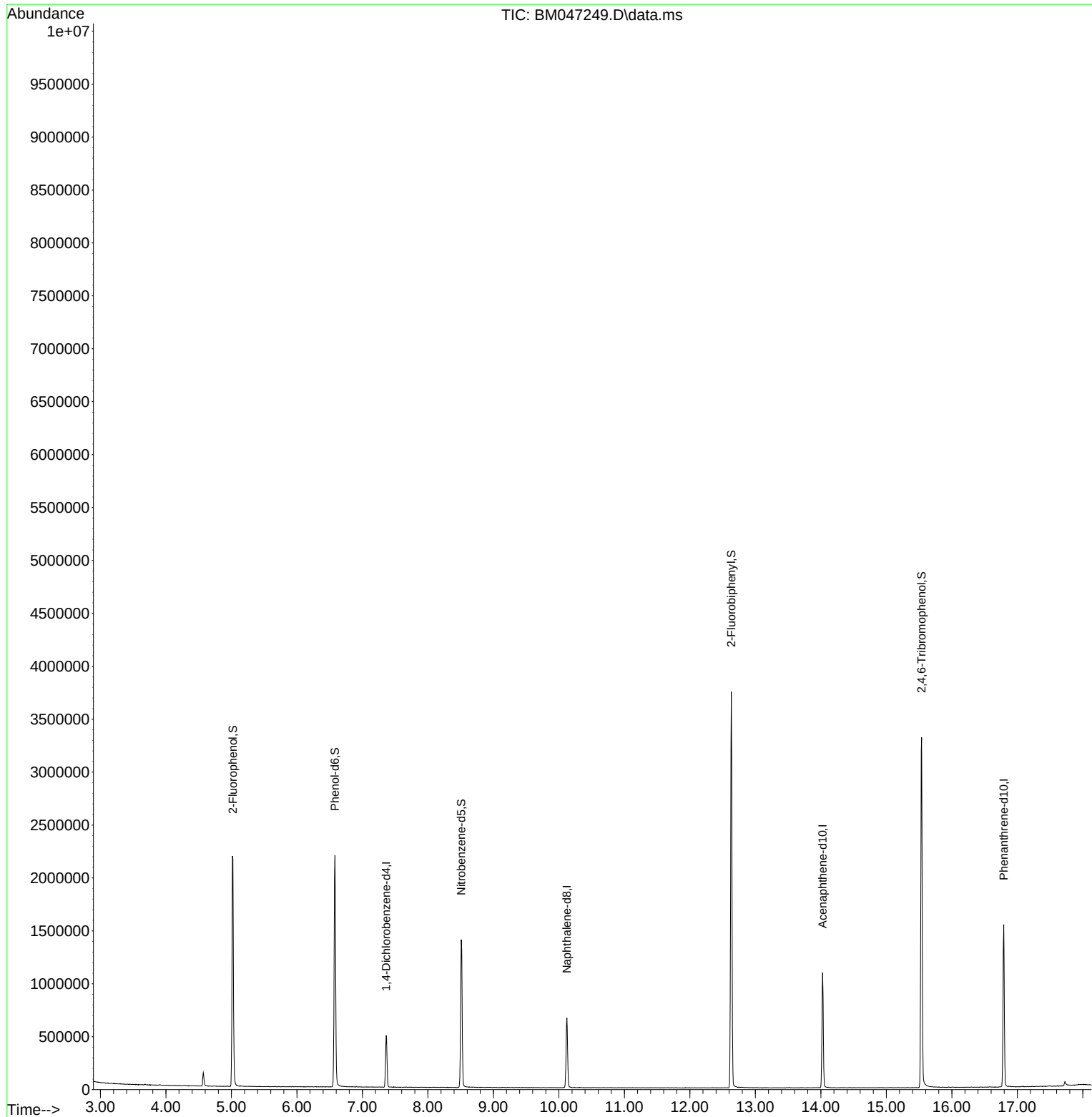
Target Compounds	Qvalue

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

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