

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081324\
 Data File : BM047169.D
 Acq On : 13 Aug 2024 10:09
 Operator : RC/JU
 Sample : PB162553BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162553BL

Quant Time: Aug 13 10:52:22 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	7.369	152	266916	20.000	ng	0.00
21) Naphthalene-d8	10.128	136	1056409	20.000	ng	0.00
39) Acenaphthene-d10	14.027	164	676276	20.000	ng	0.00
64) Phenanthrene-d10	16.792	188	1387815	20.000	ng	0.00
76) Chrysene-d12	21.027	240	1122401	20.000	ng	0.00
86) Perylene-d12	23.721	264	1330346	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	1625854	108.841	ng	0.00
7) Phenol-d6	6.581	99	2136786	103.813	ng	0.00
23) Nitrobenzene-d5	8.516	82	1404872	66.963	ng	0.00
42) 2,4,6-Tribromophenol	15.533	330	719760	91.788	ng	0.00
45) 2-Fluorobiphenyl	12.639	172	2720643	62.056	ng	0.00
79) Terphenyl-d14	19.456	244	3972397	75.837	ng	0.00

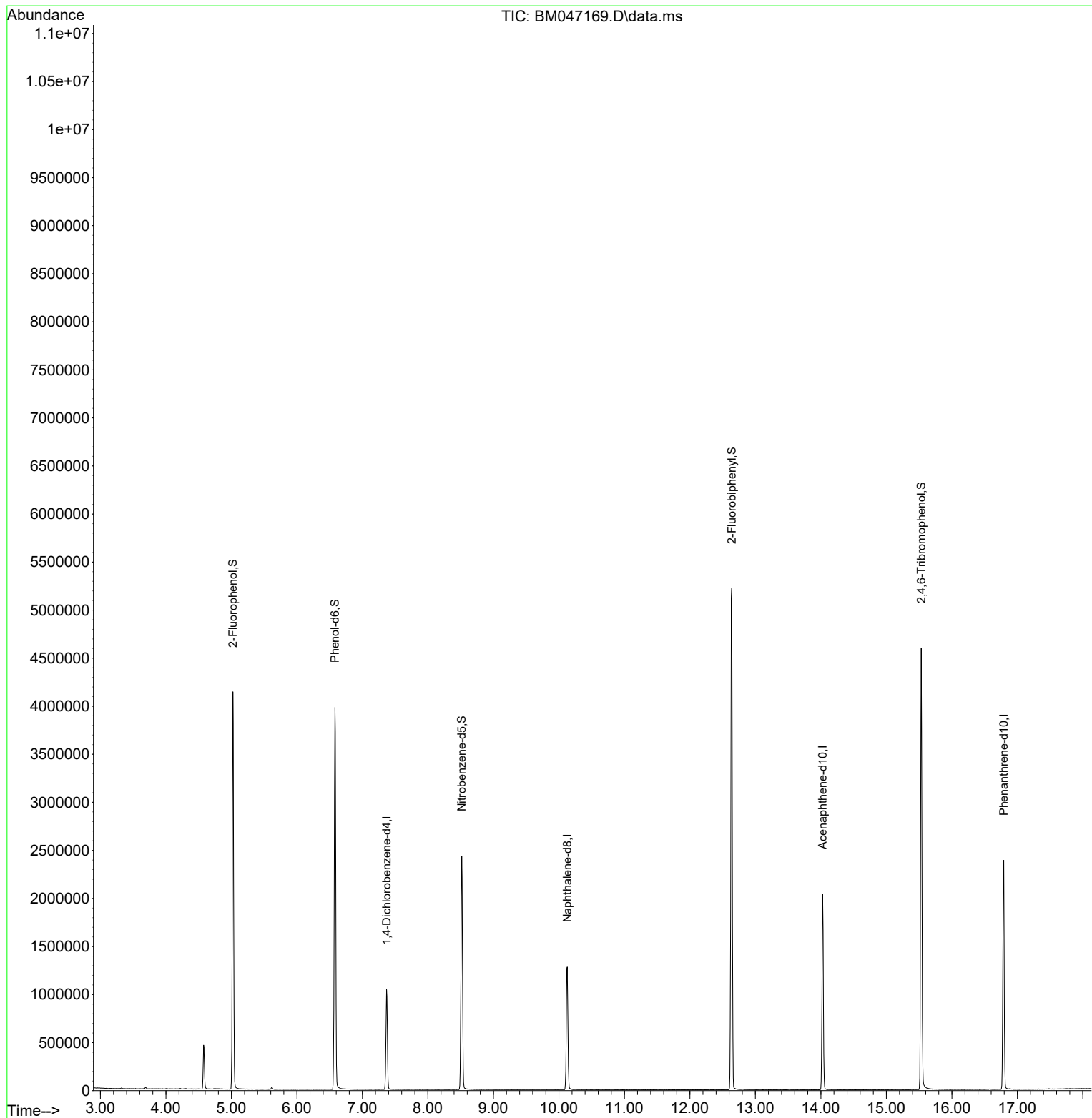
Target Compounds	Qvalue

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

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