

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092324\
 Data File : BM047605.D
 Acq On : 23 Sep 2024 15:44
 Operator : RC/JU
 Sample : P4139-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 350

Quant Time: Sep 23 19:32:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM092024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 21 02:22:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	298472	20.000	ng	-0.01
21) Naphthalene-d8	10.628	136	1177824	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	659267	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1442326	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1233781	20.000	ng	-0.01
86) Perylene-d12	24.450	264	1274216	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	1547509	79.454	ng	-0.01
7) Phenol-d6	6.999	99	2111688	82.284	ng	0.00
23) Nitrobenzene-d5	8.987	82	1312741	55.483	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	470839	77.128	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	2601154	55.828	ng	-0.01
79) Terphenyl-d14	19.821	244	4266602	62.049	ng	0.00

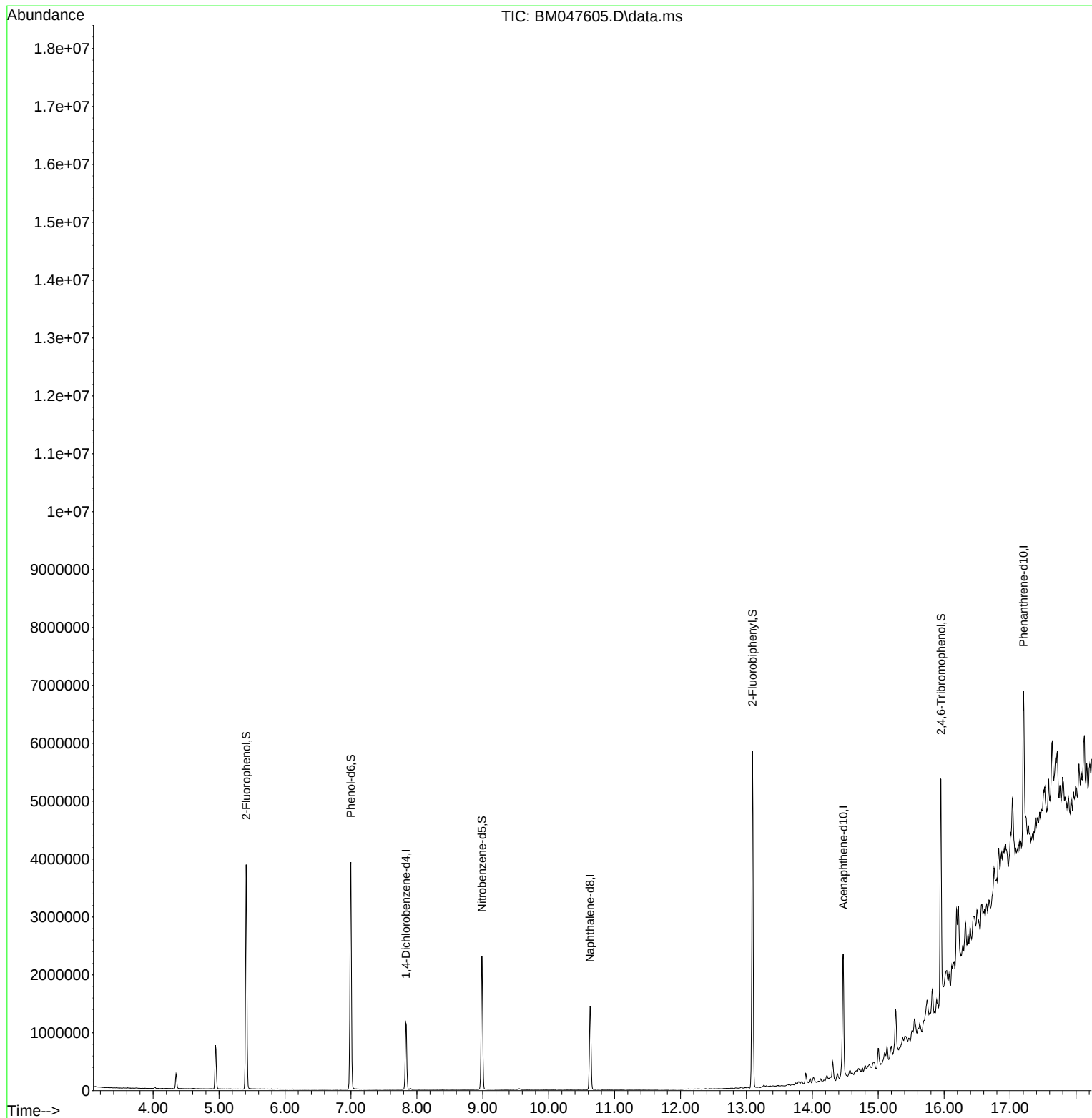
Target Compounds	Qvalue

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

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