

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071525\
 Data File : BM050448.D
 Acq On : 15 Jul 2025 10:10
 Operator : RC/JU
 Sample : PB168845BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168845BL

Quant Time: Jul 15 10:53:50 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 08 18:32:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	523100	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1859188	20.000	ng	-0.01
39) Acenaphthene-d10	14.480	164	1153910	20.000	ng	-0.01
64) Phenanthrene-d10	17.215	188	2194023	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	2232959	20.000	ng	-0.02
86) Perylene-d12	24.468	264	2515738	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3946471	129.863	ng	0.00
7) Phenol-d6	7.022	99	4832536	126.145	ng	0.00
23) Nitrobenzene-d5	9.010	82	2858462	78.439	ng	-0.01
42) 2,4,6-Tribromophenol	15.963	330	1972178	140.668	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	7350571	78.667	ng	-0.01
79) Terphenyl-d14	19.827	244	10843091	85.438	ng	-0.02

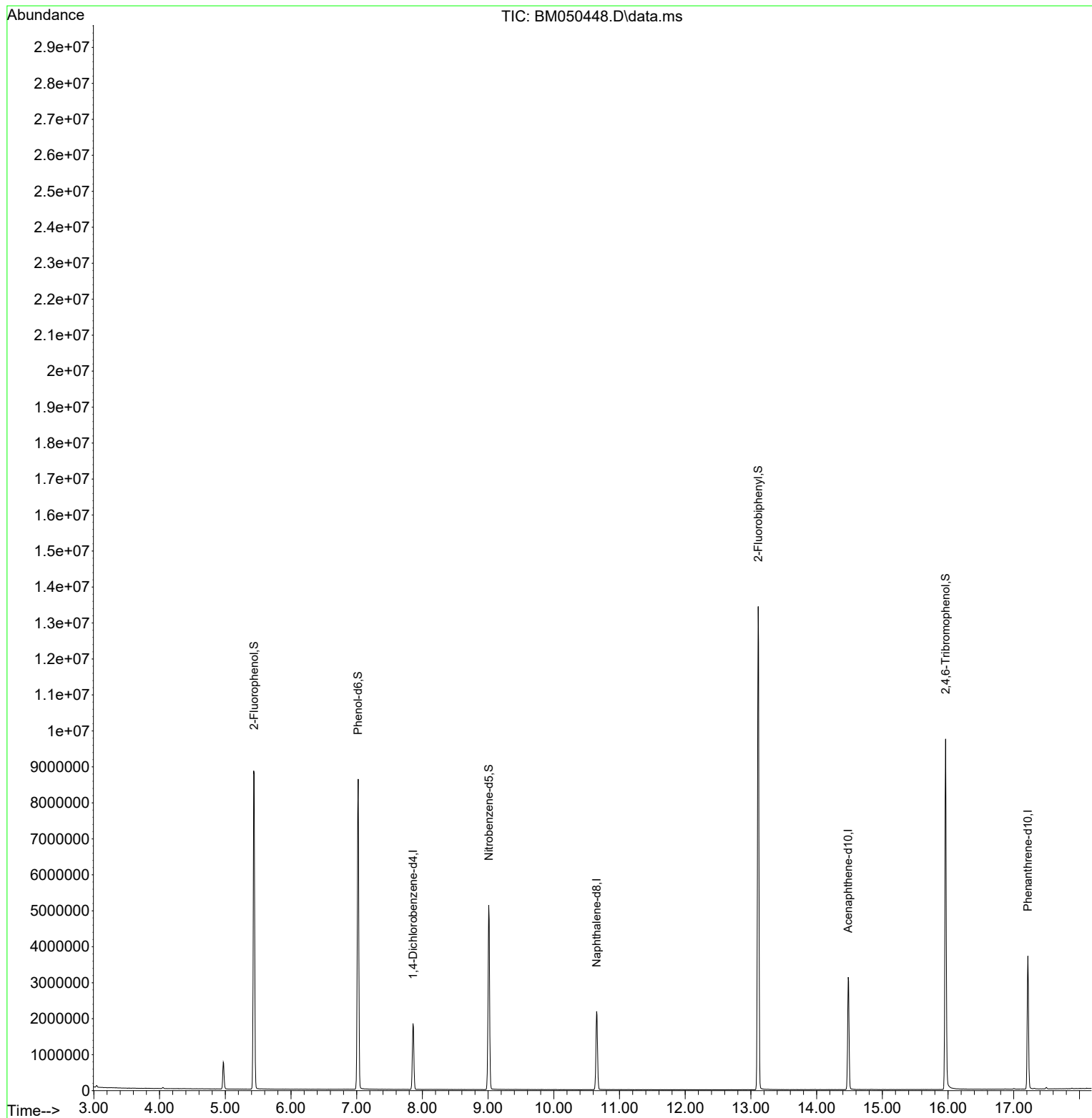
Target Compounds	Qvalue

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

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