

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050925\
 Data File : BM050159.D
 Acq On : 09 May 2025 14:48
 Operator : RC/JU
 Sample : Q1964-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MH-LL

Quant Time: May 09 15:24:35 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 08 12:03:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	349102	20.000	ng	-0.03
21) Naphthalene-d8	10.534	136	1219858	20.000	ng	-0.03
39) Acenaphthene-d10	14.392	164	784394	20.000	ng	-0.02
64) Phenanthrene-d10	17.139	188	1450953	20.000	ng	-0.02
76) Chrysene-d12	21.386	240	1072762	20.000	ng	-0.01
86) Perylene-d12	24.386	264	831862	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	2589598	129.893	ng	-0.02
7) Phenol-d6	6.934	99	2932105	117.015	ng	-0.02
23) Nitrobenzene-d5	8.904	82	2034319	82.177	ng	-0.02
42) 2,4,6-Tribromophenol	15.892	330	1613952	139.154	ng	-0.01
45) 2-Fluorobiphenyl	13.004	172	5258391	79.304	ng	-0.03
79) Terphenyl-d14	19.768	244	7123815	98.264	ng	-0.01

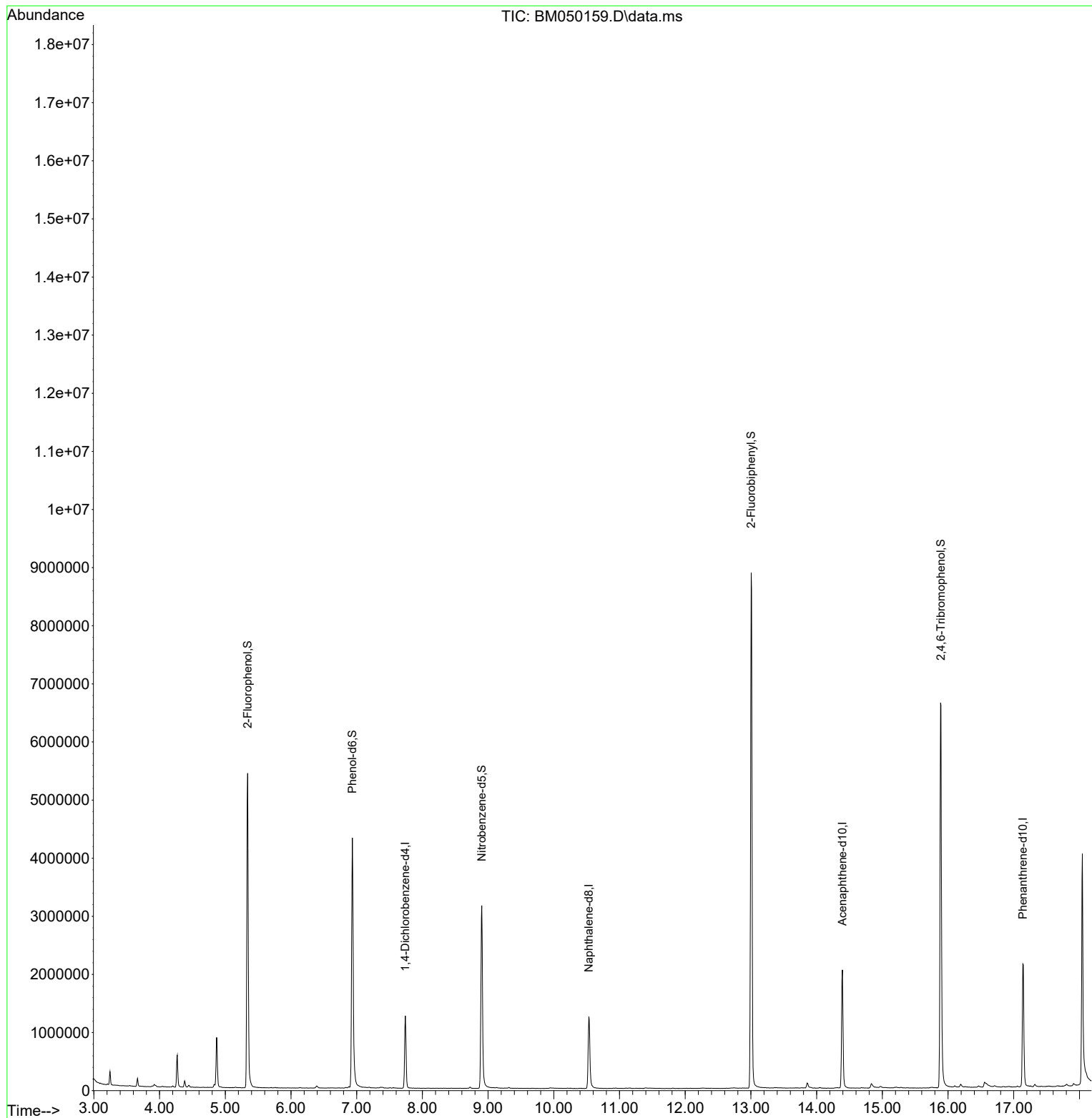
Target Compounds	Qvalue

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

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