

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050525\
 Data File : BM050108.D
 Acq On : 05 May 2025 20:55
 Operator : RC/JU
 Sample : Q1922-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MH-R

Quant Time: May 06 05:20:27 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 28 18:09:16 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.745	152	415751	20.000	ng	-0.02
21) Naphthalene-d8	10.539	136	1427299	20.000	ng	-0.02
39) Acenaphthene-d10	14.392	164	905134	20.000	ng	-0.02
64) Phenanthrene-d10	17.139	188	1708333	20.000	ng	-0.02
76) Chrysene-d12	21.380	240	1558229	20.000	ng	-0.02
86) Perylene-d12	24.374	264	1604513	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	1623497	68.379	ng	-0.02
7) Phenol-d6	6.934	99	2131315	71.421	ng	-0.02
23) Nitrobenzene-d5	8.904	82	1141357	39.404	ng	-0.02
42) 2,4,6-Tribromophenol	15.892	330	865810	64.692	ng	-0.01
45) 2-Fluorobiphenyl	13.010	172	2712924	35.457	ng	-0.02
79) Terphenyl-d14	19.768	244	4655192	44.207	ng	-0.01

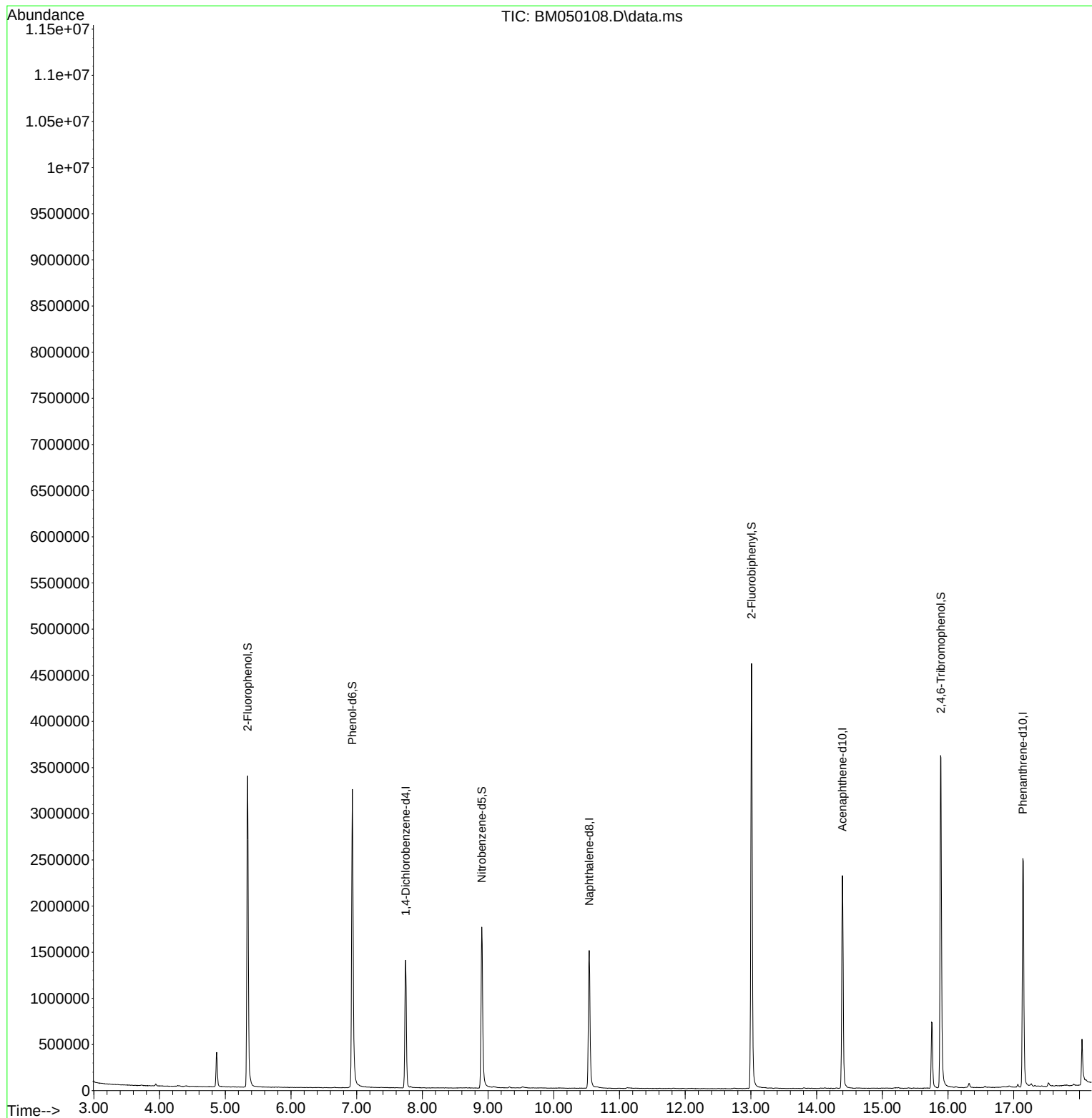
Target Compounds	Qvalue

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

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