

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050249.D
 Acq On : 09 Jun 2025 19:40
 Operator : RC/JU
 Sample : Q2228-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP08-MHI-WC

Quant Time: Jun 10 01:26:17 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 09 11:54:38 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	344656	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	1293755	20.000	ng	0.00
39) Acenaphthene-d10	14.398	164	699023	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1337378	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1265522	20.000	ng	-0.01
86) Perylene-d12	24.344	264	635863	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2869963	138.907	ng	0.00
7) Phenol-d6	6.934	99	3323871	122.209	ng	0.00
23) Nitrobenzene-d5	8.916	82	2272103	91.337	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1180262	148.444	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	4569653	88.504	ng	0.00
79) Terphenyl-d14	19.762	244	5766516	86.349	ng	0.00

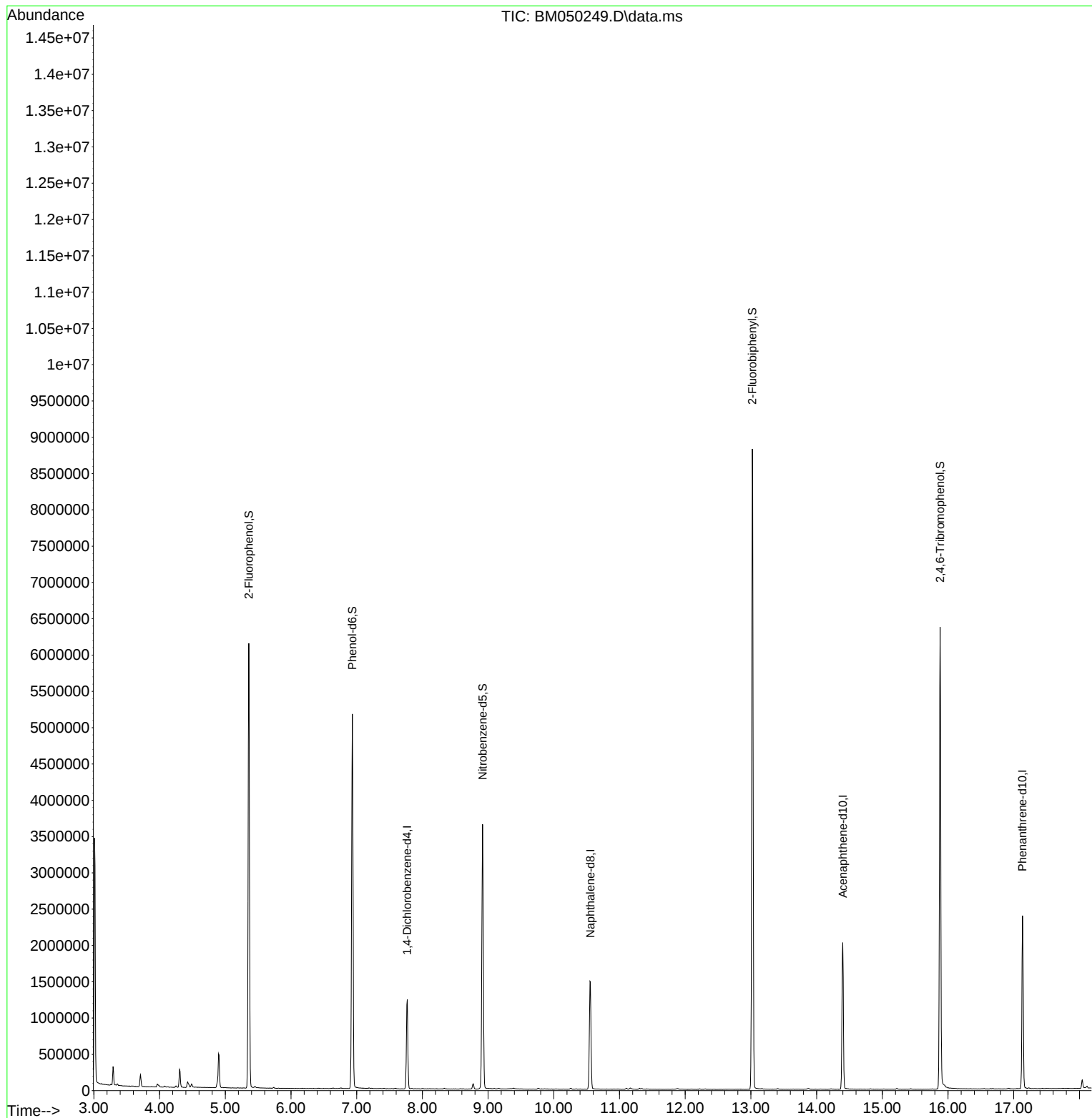
Target Compounds	Qvalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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