

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050269.D
 Acq On : 10 Jun 2025 09:22
 Operator : RC/JU
 Sample : Q2266-08
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-5

Quant Time: Jun 10 10:40:30 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	7.763	152	330823	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1216311	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	298562	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	840437	20.000	ng	0.00
76) Chrysene-d12	21.362	240	156325	20.000	ng	-0.01
86) Perylene-d12	24.321	264	4747	20.000	ng	#-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2749596	138.646	ng	0.00
7) Phenol-d6	6.928	99	1664766	63.768	ng	0.00
23) Nitrobenzene-d5	8.916	82	2339637	100.040	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1067374	314.310	ng	0.00
45) 2-Fluorobiphenyl	13.021	172	4388234	198.988	ng	0.00
79) Terphenyl-d14	19.762	244	3607176	437.274	ng	0.00

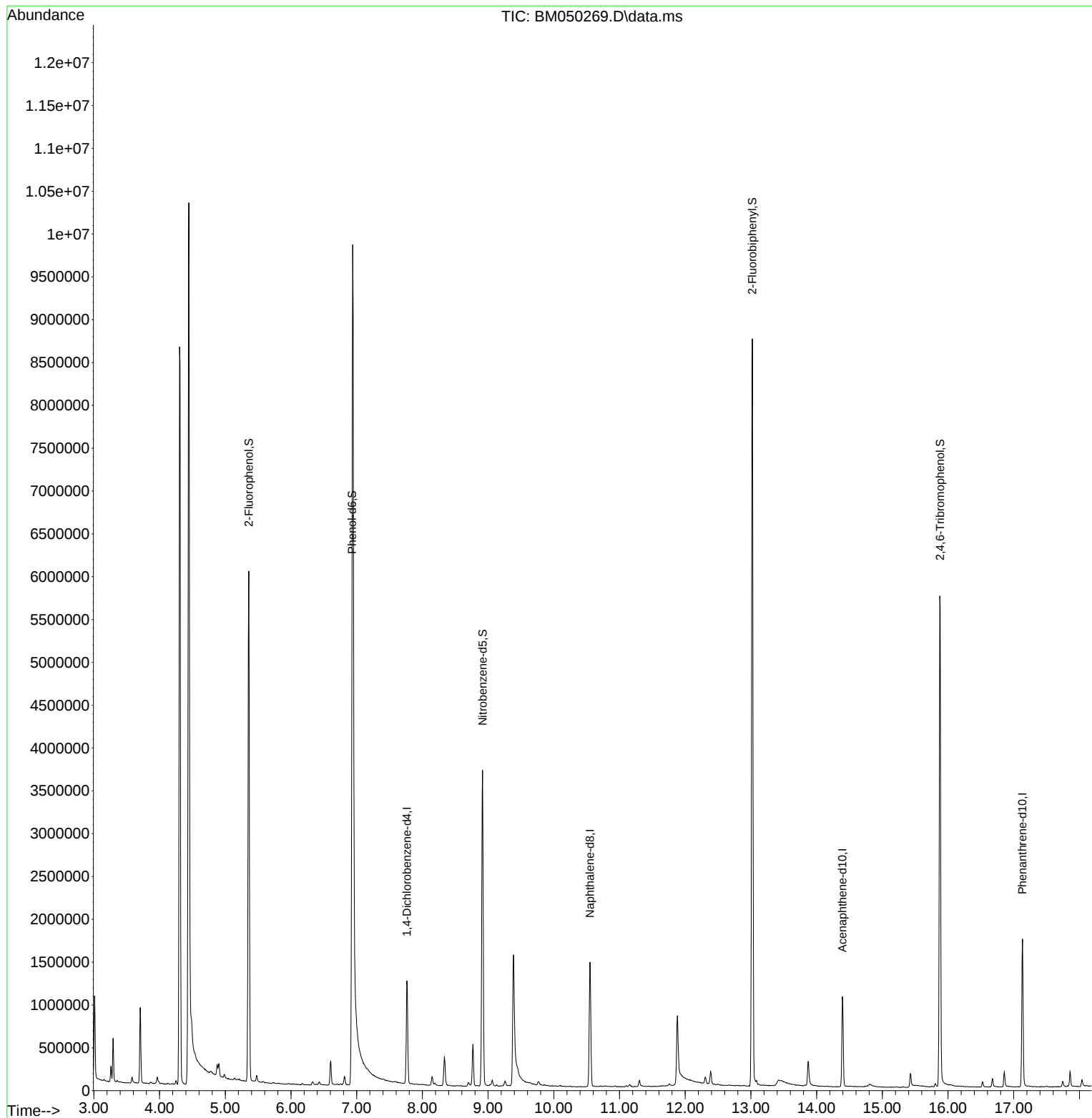
Target Compounds	Qvalue

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

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