

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050274.D
 Acq On : 10 Jun 2025 12:38
 Operator : RC/JU
 Sample : Q2202-03DL 10X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-12-20250603DL

Quant Time: Jun 10 13:12:01 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)

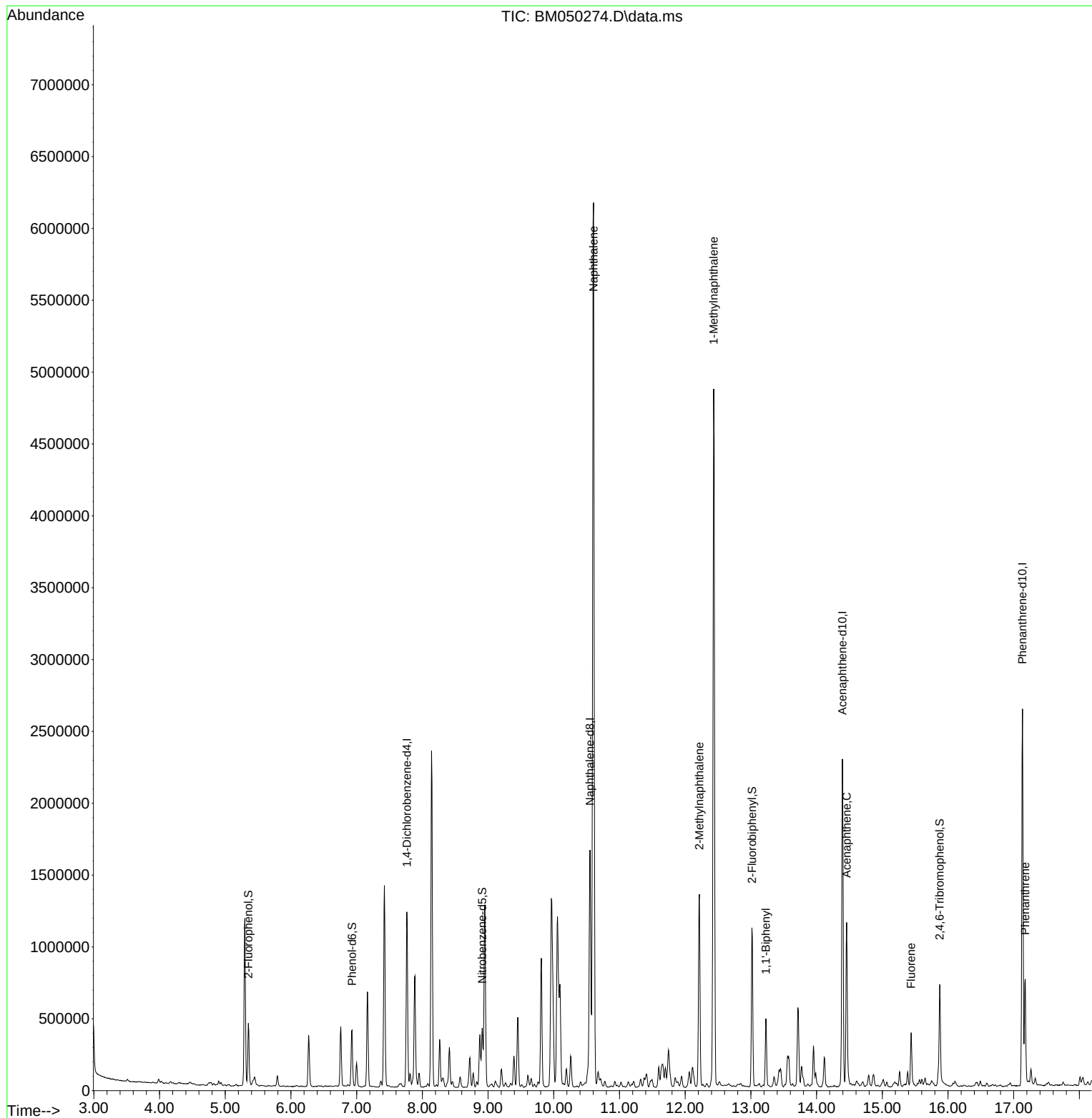
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	334632	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1355551	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	770915	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1509253	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1638874	20.000	ng	-0.01
86) Perylene-d12	24.339	264	1724623	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.352	112	202238	10.082	ng	0.00
7) Phenol-d6	6.928	99	150293	5.691	ng	0.00
23) Nitrobenzene-d5	8.910	82	242976	9.322	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	130318	14.862	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	562182	9.873	ng	-0.01
79) Terphenyl-d14	19.763	244	918863	10.625	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.604	128	5303692	75.675	ng	100
37) 2-Methylnaphthalene	12.216	142	652389	15.430	ng	100
38) 1-Methylnaphthalene	12.434	142	2300363	51.261	ng	99
46) 1,1'-Biphenyl	13.228	154	241213	4.179	ng	97
52) Acenaphthene	14.457	154	388430	8.555	ng	96
58) Fluorene	15.439	166	159079	3.127	ng	98
71) Phenanthrene	17.175	178	429996	4.987	ng	98

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

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