

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050273.D
 Acq On : 10 Jun 2025 11:59
 Operator : RC/JU
 Sample : Q2231-02DL 10X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-14-20250604DL

Quant Time: Jun 10 12:31:06 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)

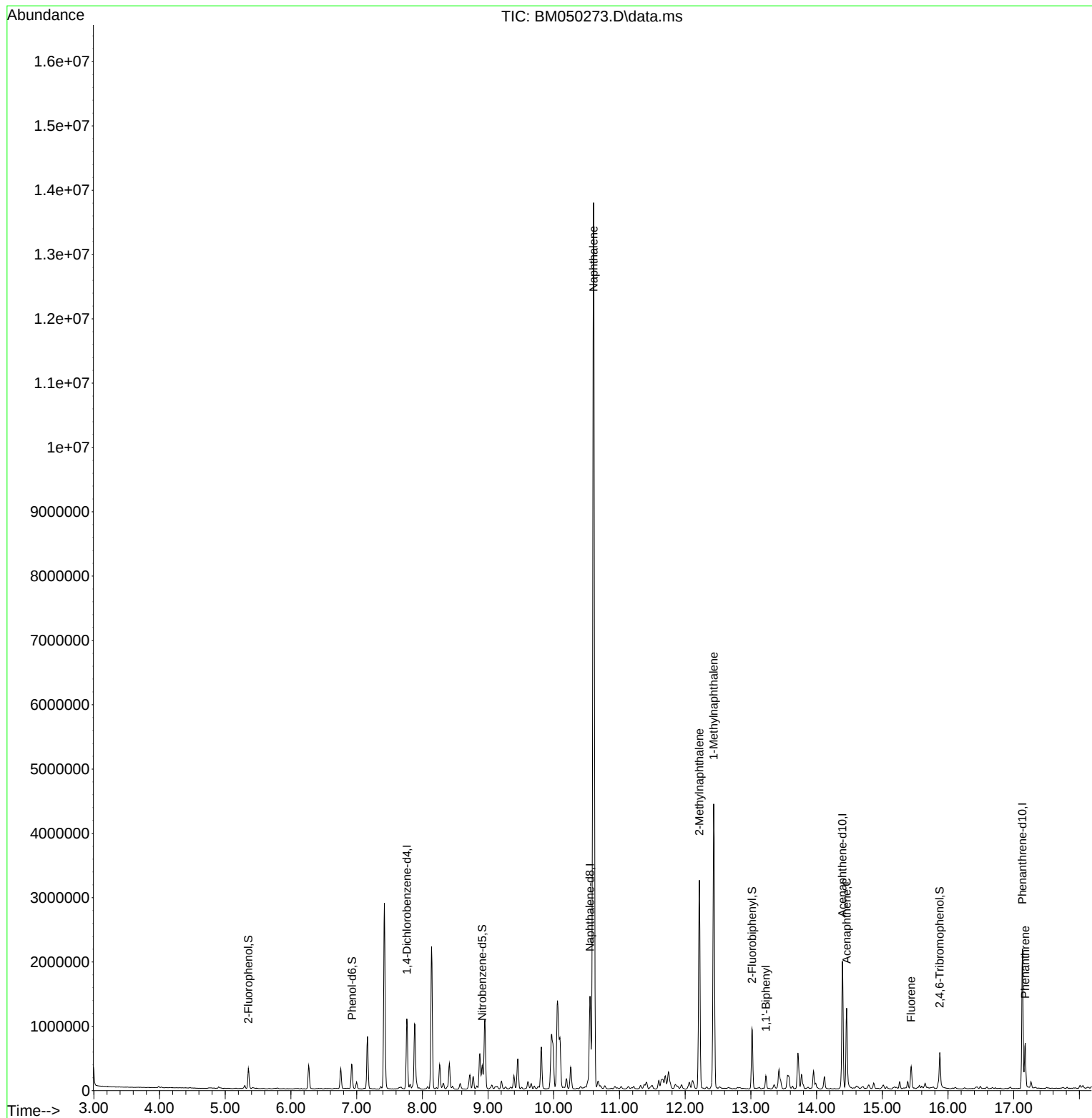
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	306070	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1203869	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	655906	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1230960	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1295114	20.000	ng	-0.01
86) Perylene-d12	24.339	264	1344466	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.352	112	143963	7.846	ng	0.00
7) Phenol-d6	6.928	99	104618	4.331	ng	0.00
23) Nitrobenzene-d5	8.910	82	213430	9.220	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	108039	14.482	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	487246	10.057	ng	-0.01
79) Terphenyl-d14	19.757	244	751319	10.993	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	10.604	128	12286529	197.397	ng	99
37) 2-Methylnaphthalene	12.216	142	1528618	40.709	ng	100
38) 1-Methylnaphthalene	12.434	142	2058141	51.642	ng	99
46) 1,1'-Biphenyl	13.228	154	100305	2.042	ng	98
52) Acenaphthene	14.457	154	421719	10.917	ng	99
58) Fluorene	15.439	166	143496	3.315	ng	98
71) Phenanthrene	17.175	178	419160	5.961	ng	98

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

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