

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050248.D
 Acq On : 09 Jun 2025 19:01
 Operator : RC/JU
 Sample : Q2236-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-A2-04-C

Quant Time: Jun 10 01:26:00 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)

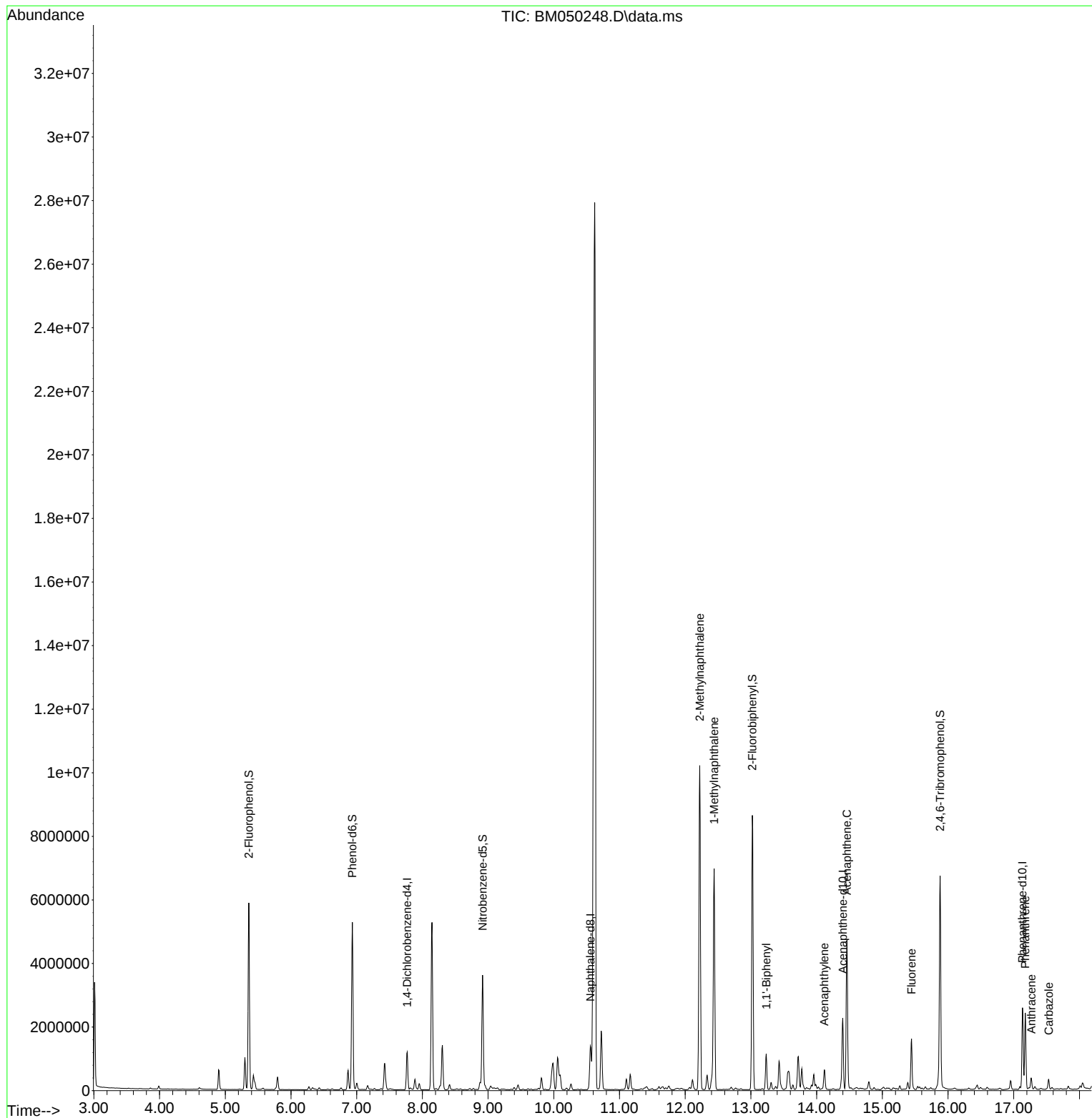
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	334436	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	1312450	20.000	ng	0.00
39) Acenaphthene-d10	14.398	164	761229	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1459849	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1563302	20.000	ng	-0.01
86) Perylene-d12	24.344	264	1685599	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2752148	137.275	ng	0.00
7) Phenol-d6	6.934	99	3189329	120.845	ng	0.00
23) Nitrobenzene-d5	8.916	82	2263913	89.712	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1287864	148.741	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	4606639	81.929	ng	0.00
79) Terphenyl-d14	19.762	244	6889305	83.512	ng	0.00
Target Compounds						
						Qvalue
37) 2-Methylnaphthalene	12.222	142	4908486	119.906	ng	98
38) 1-Methylnaphthalene	12.439	142	3260277	75.038	ng	99
46) 1,1'-Biphenyl	13.233	154	580651	10.187	ng	99
49) Acenaphthylene	14.116	152	258538	3.607	ng	# 96
52) Acenaphthene	14.463	154	1714018	38.233	ng	98
58) Fluorene	15.445	166	693734	13.810	ng	99
71) Phenanthrene	17.174	178	1412684	16.939	ng	99
72) Anthracene	17.268	178	225734	2.706	ng	99
73) Carbazole	17.533	167	198686	2.614	ng	100

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

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