

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050125\
 Data File : BM050065.D
 Acq On : 01 May 2025 15:26
 Operator : RC/JU
 Sample : Q1900-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-2

Quant Time: May 01 16:24:44 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 28 18:09:16 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

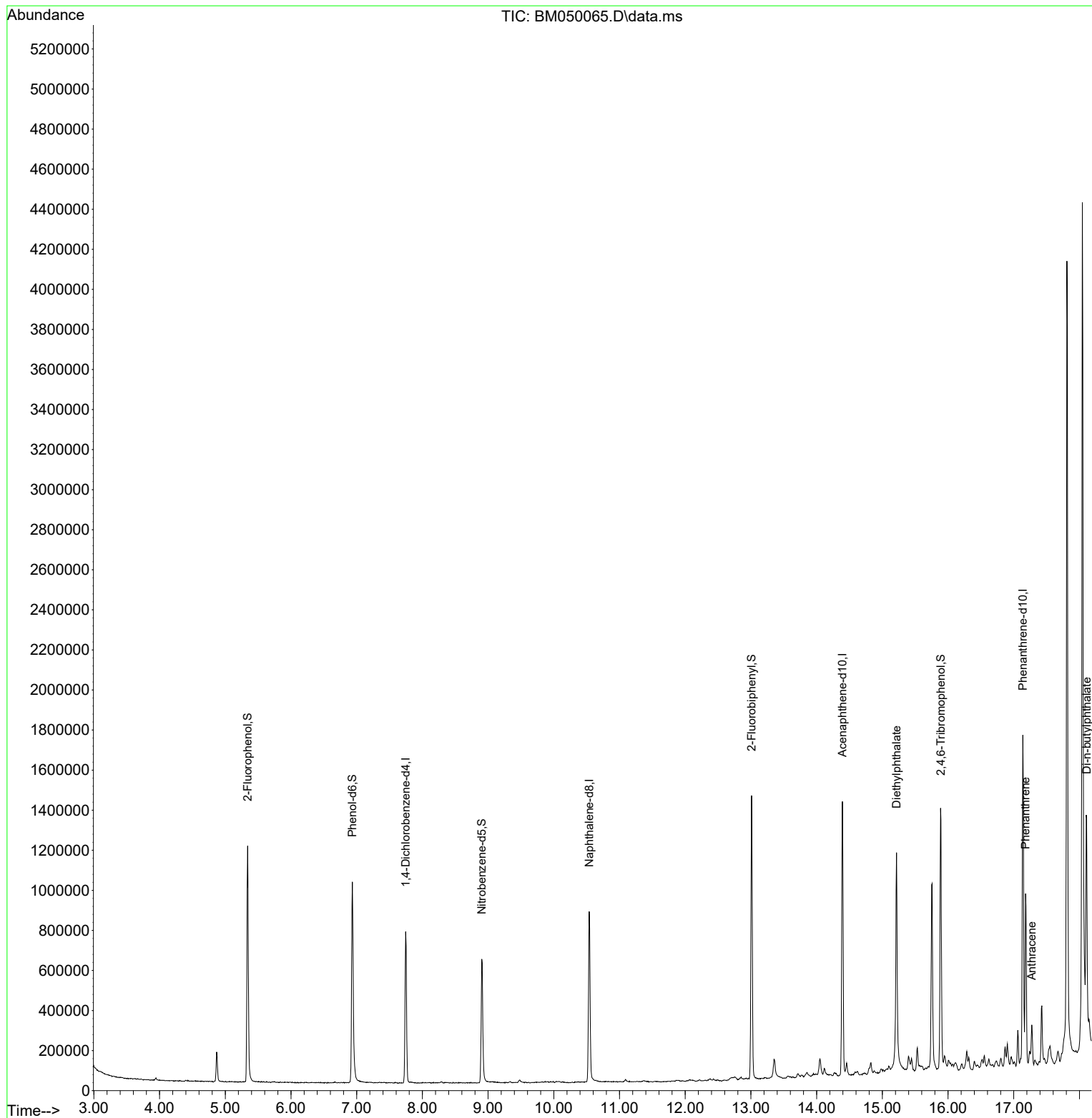
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.745	152	212536	20.000	ng	-0.02
21) Naphthalene-d8	10.539	136	734717	20.000	ng	-0.02
39) Acenaphthene-d10	14.392	164	479228	20.000	ng	-0.02
64) Phenanthrene-d10	17.139	188	957331	20.000	ng	-0.02
76) Chrysene-d12	21.380	240	1033347	20.000	ng	-0.02
86) Perylene-d12	24.368	264	1128226	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	532554	43.877	ng	-0.02
7) Phenol-d6	6.934	99	656138	43.011	ng	-0.02
23) Nitrobenzene-d5	8.904	82	389525	26.125	ng	-0.02
42) 2,4,6-Tribromophenol	15.892	330	256053	36.135	ng	-0.01
45) 2-Fluorobiphenyl	13.010	172	749520	18.502	ng	-0.02
79) Terphenyl-d14	19.768	244	1353156	19.377	ng	-0.01
Target Compounds						Qvalue
60) Diethylphthalate	15.215	149	882707	26.748	ng	98
71) Phenanthrene	17.180	178	541522	10.029	ng	100
72) Anthracene	17.274	178	129542	2.371	ng	99
74) Di-n-butylphthalate	18.109	149	942080	16.684	ng	100
75) Fluoranthene	19.203	202	1229086	19.388	ng	100
78) Pyrene	19.562	202	1033203	14.238	ng	99
81) Benzo(a)anthracene	21.362	228	578595	8.131	ng	98
83) Chrysene	21.427	228	501533	7.615	ng	98
87) Indeno(1,2,3-cd)pyrene	27.744	276	317240	3.774	ng	96
88) Benzo(b)fluoranthene	23.427	252	670304	9.136	ng	98
89) Benzo(k)fluoranthene	23.486	252	222812	3.002	ng	99
90) Benzo(a)pyrene	24.233	252	465357	6.772	ng	99
92) Benzo(g,h,i)perylene	28.797	276	300045	4.573	ng	100

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

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