

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040925\
 Data File : BM049871.D
 Acq On : 09 Apr 2025 17:03
 Operator : RC/JU
 Sample : Q1712-01DL2 50X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 Z-05ADL2

Quant Time: Apr 09 17:42:34 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 04:00:55 2025
 Response via : Initial Calibration

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

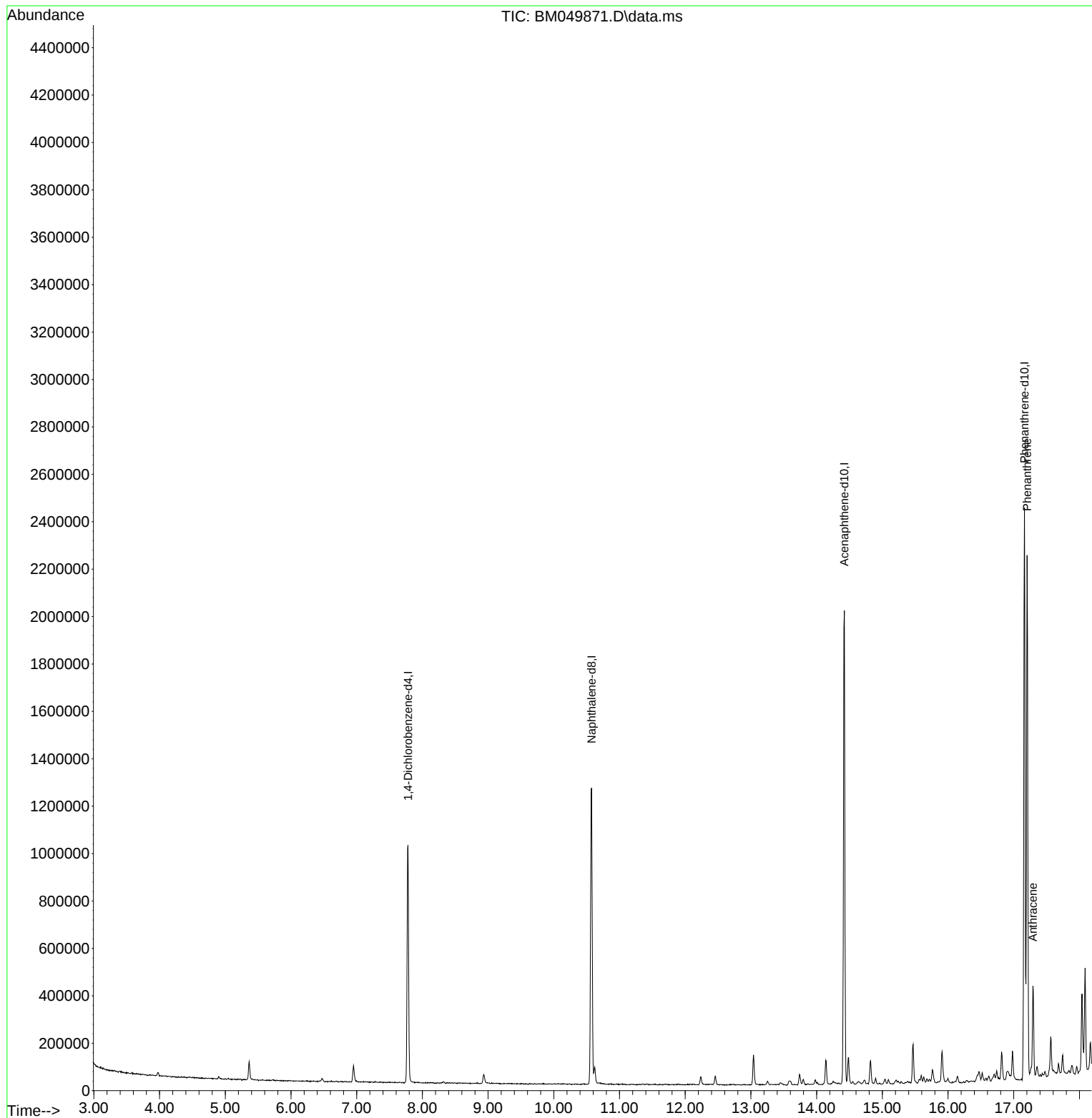
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	308569	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1104161	20.000	ng	0.00
39) Acenaphthene-d10	14.421	164	735135	20.000	ng	0.00
64) Phenanthrene-d10	17.162	188	1424172	20.000	ng	0.00
76) Chrysene-d12	21.397	240	1366865	20.000	ng	0.00
86) Perylene-d12	24.397	264	1468519	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	34800	1.915	ng	0.00
7) Phenol-d6	6.951	99	42034	1.859	ng	0.00
23) Nitrobenzene-d5	8.934	82	25482	1.285	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	18096	1.692	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	67361	1.243	ng	0.00
79) Terphenyl-d14	19.786	244	91784	1.258	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.204	178	1334833	17.096	ng	100
72) Anthracene	17.292	178	250896	3.261	ng	99
75) Fluoranthene	19.221	202	1711967	17.833	ng	99
78) Pyrene	19.580	202	1566993	16.634	ng	100
81) Benzo(a)anthracene	21.380	228	834995	9.192	ng	99
83) Chrysene	21.444	228	831753	9.631	ng	98
87) Indeno(1,2,3-cd)pyrene	27.779	276	391187	3.574	ng	98
88) Benzo(b)fluoranthene	23.456	252	972601	10.370	ng	99
89) Benzo(k)fluoranthene	23.509	252	533000m	5.864	ng	
90) Benzo(a)pyrene	24.256	252	604097	7.330	ng	98
92) Benzo(g,h,i)perylene	28.832	276	363735	3.948	ng	95

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

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