

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050125\
 Data File : BM050072.D
 Acq On : 01 May 2025 20:00
 Operator : RC/JU
 Sample : Q1914-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SS-4

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 05/02/2025
 Supervised By :Jagrut Upadhyay 05/02/2025

Quant Time: May 02 01:12:18 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	266849	20.000	ng	-0.02
21) Naphthalene-d8	10.539	136	987063	20.000	ng	-0.02
39) Acenaphthene-d10	14.392	164	682416	20.000	ng	-0.02
64) Phenanthrene-d10	17.139	188	1413674	20.000	ng	-0.02
76) Chrysene-d12	21.380	240	1410216	20.000	ng	-0.02
86) Perylene-d12	24.368	264	1421527	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	1083091	71.073	ng	-0.02
7) Phenol-d6	6.934	99	1431584	74.742	ng	-0.02
23) Nitrobenzene-d5	8.904	82	834787	41.674	ng	-0.02
42) 2,4,6-Tribromophenol	15.892	330	719122	71.268	ng	-0.01
45) 2-Fluorobiphenyl	13.010	172	2208129	38.278	ng	-0.02
79) Terphenyl-d14	19.768	244	3399375	35.670	ng	-0.01
Target Compounds						Qvalue
71) Phenanthrene	17.186	178	614421	7.706	ng	100
75) Fluoranthene	19.203	202	1542165	16.474	ng	99
78) Pyrene	19.562	202	1336131	13.492	ng	100
81) Benzo(a)anthracene	21.362	228	806143	8.301	ng	98
83) Chrysene	21.427	228	729730	8.119	ng	98
87) Indeno(1,2,3-cd)pyrene	27.744	276	528099	4.986	ng	97
88) Benzo(b)fluoranthene	23.433	252	1089016	11.780	ng	97
89) Benzo(k)fluoranthene	23.486	252	326794m	3.495	ng	
90) Benzo(a)pyrene	24.227	252	738125	8.525	ng	98
92) Benzo(g,h,i)perylene	28.803	276	530889	6.422	ng	99

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

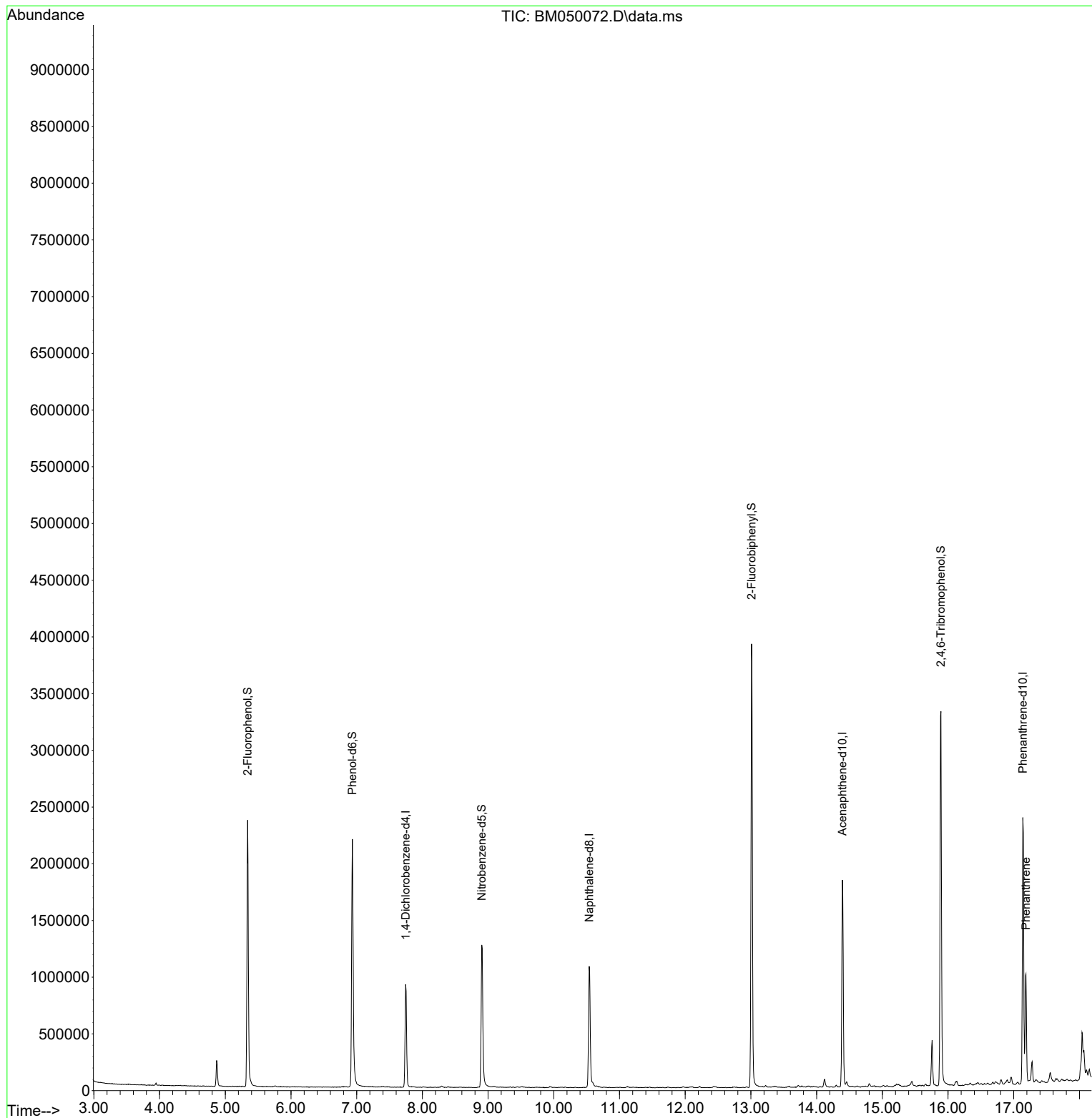
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