

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090324\
 Data File : BM047424.D
 Acq On : 03 Sep 2024 14:05
 Operator : RC/JU
 Sample : P3788-01DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OR-02-082924DL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 09/04/2024
 Supervised By :mohammad ahmed 09/06/2024

Quant Time: Sep 03 14:43:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 26 18:11:40 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.340	152	182453	20.000	ng	-0.01
21) Naphthalene-d8	10.093	136	675436	20.000	ng	-0.01
39) Acenaphthene-d10	14.004	164	460896	20.000	ng	0.00
64) Phenanthrene-d10	16.769	188	945088	20.000	ng	0.00
76) Chrysene-d12	21.015	240	713107	20.000	ng	0.00
86) Perylene-d12	23.715	264	822417	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.011	112	97417	10.244	ng	0.00
7) Phenol-d6	6.581	99	115272	8.898	ng	0.01
23) Nitrobenzene-d5	8.504	82	72229	5.933	ng	0.00
42) 2,4,6-Tribromophenol	15.528	330	50274	9.041	ng	0.00
45) 2-Fluorobiphenyl	12.610	172	195270	6.510	ng	0.00
79) Terphenyl-d14	19.433	244	275383	6.874	ng	-0.01
Target Compounds						Qvalue
52) Acenaphthene	14.069	154	71875	3.089	ng	98
55) Dibenzofuran	14.416	168	82153	2.192	ng	# 96
58) Fluorene	15.069	166	117460	3.872	ng	96
71) Phenanthrene	16.816	178	1044111	22.749	ng	100
72) Anthracene	16.904	178	321177	7.178	ng	99
73) Carbazole	17.198	167	92027	2.143	ng	99
75) Fluoranthene	18.851	202	1567318	29.082	ng	100
78) Pyrene	19.216	202	1348417	24.654	ng	99
81) Benzo(a)anthracene	20.998	228	583942	12.557	ng	97
83) Chrysene	21.057	228	526769	12.042	ng	97
87) Indeno(1,2,3-cd)pyrene	26.721	276	377328	7.329	ng	95
88) Benzo(b)fluoranthene	22.874	252	691986	14.422	ng	97
89) Benzo(k)fluoranthene	22.921	252	233482m	5.133	ng	
90) Benzo(a)pyrene	23.592	252	527186	12.705	ng	97
91) Dibenzo(a,h)anthracene	26.739	278	95911	2.300	ng	# 84
92) Benzo(g,h,i)perylene	27.656	276	375995	8.672	ng	99

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

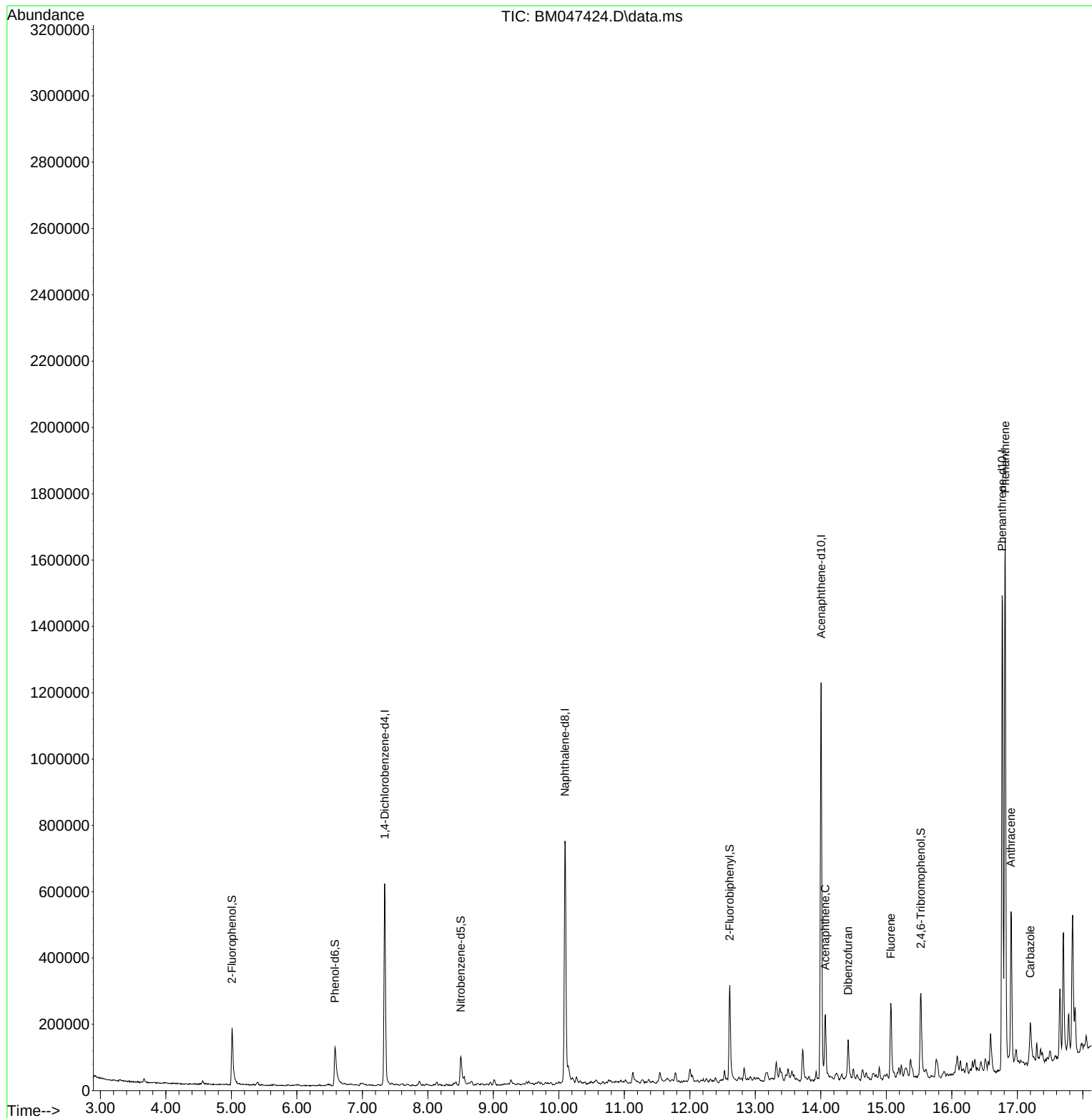
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