

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081524\
 Data File : BM047222.D
 Acq On : 15 Aug 2024 19:33
 Operator : RC/JU
 Sample : P3518-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OU4-TS-05-080724

Quant Time: Aug 16 01:19:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

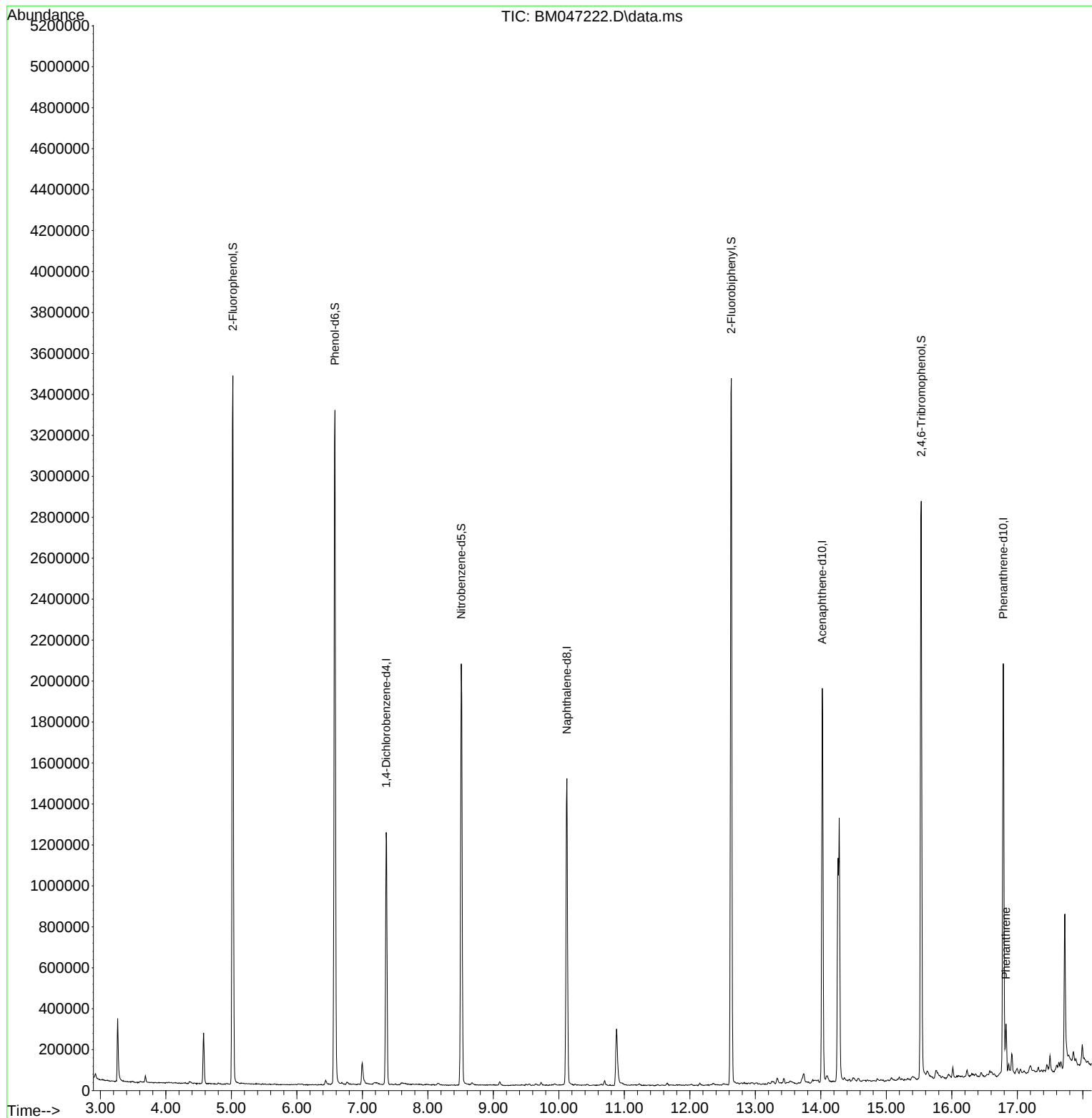
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	333710	20.000	ng	-0.01
21) Naphthalene-d8	10.122	136	1204983	20.000	ng	-0.01
39) Acenaphthene-d10	14.022	164	674607	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1184365	20.000	ng	0.00
76) Chrysene-d12	21.021	240	973369	20.000	ng	-0.01
86) Perylene-d12	23.715	264	1120925	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	1427988	76.461	ng	0.00
7) Phenol-d6	6.581	99	1848461	71.830	ng	0.00
23) Nitrobenzene-d5	8.510	82	1201558	50.210	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	534354	68.313	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	1862702	42.592	ng	-0.01
79) Terphenyl-d14	19.451	244	1923909	42.353	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	16.827	178	137967	2.321	ng	99
75) Fluoranthene	18.862	202	330584	4.573	ng	99
78) Pyrene	19.227	202	308272	4.435	ng	99
81) Benzo(a)anthracene	21.003	228	147929	2.289	ng	97
83) Chrysene	21.062	228	163086	2.660	ng	95
88) Benzo(b)fluoranthene	22.874	252	233684	3.510	ng	# 97
90) Benzo(a)pyrene	23.592	252	149319	2.589	ng	# 95

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

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