

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081524\
 Data File : BM047220.D
 Acq On : 15 Aug 2024 18:12
 Operator : RC/JU
 Sample : P3518-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OU4-TS-03-080724

Quant Time: Aug 16 01:19:23 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)

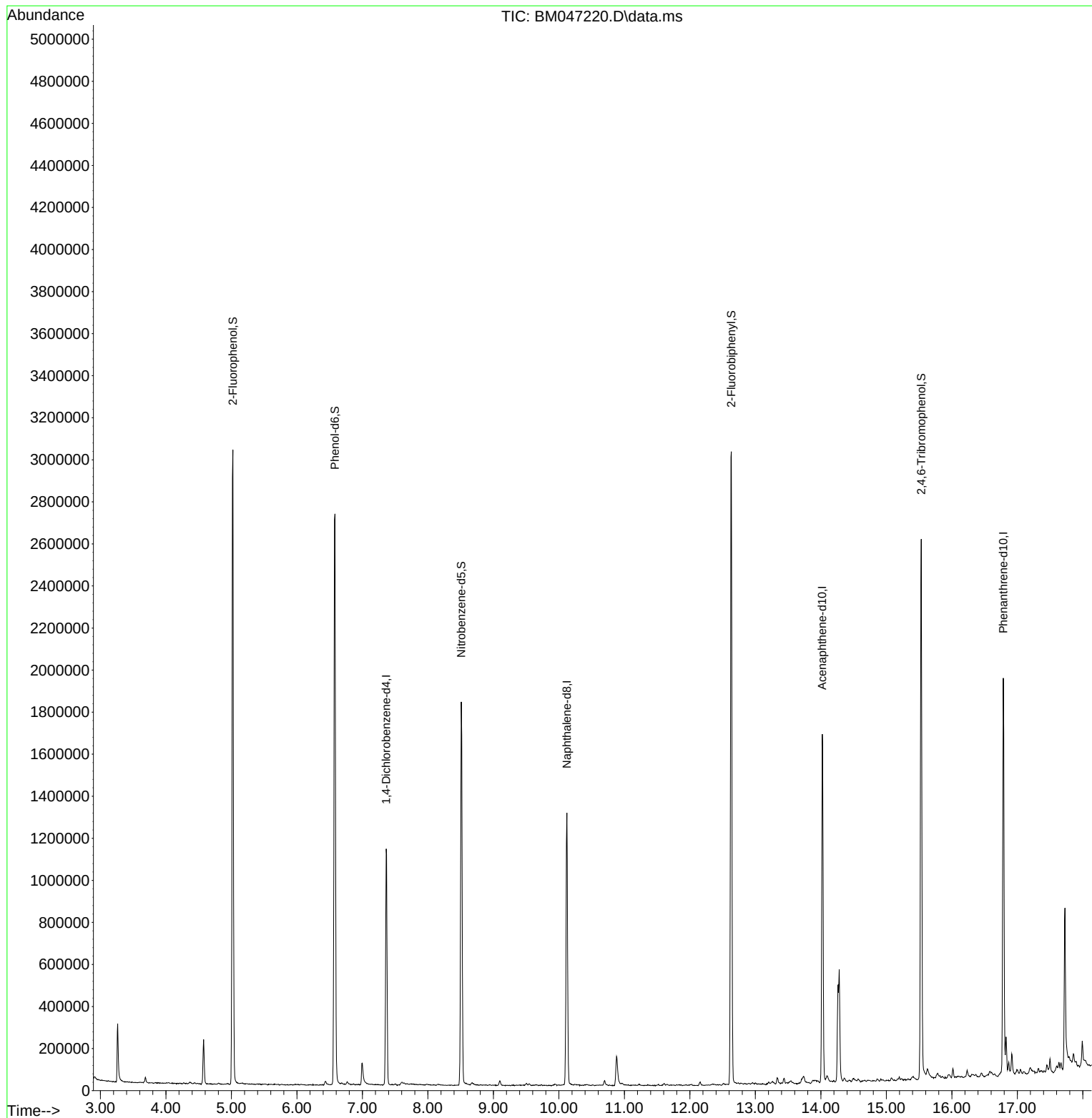
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	301755	20.000	ng	-0.01
21) Naphthalene-d8	10.122	136	1049790	20.000	ng	-0.01
39) Acenaphthene-d10	14.021	164	590390	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1113457	20.000	ng	0.00
76) Chrysene-d12	21.021	240	970223	20.000	ng	-0.01
86) Perylene-d12	23.721	264	1098695	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	1257620	74.470	ng	0.00
7) Phenol-d6	6.581	99	1593170	68.465	ng	0.00
23) Nitrobenzene-d5	8.510	82	1035399	49.663	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	482089	70.422	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	1611955	42.116	ng	-0.01
79) Terphenyl-d14	19.451	244	1869149	41.281	ng	-0.01
Target Compounds						
						Qvalue
75) Fluoranthene	18.862	202	217484	3.200	ng	99
78) Pyrene	19.227	202	207080	2.989	ng	99
83) Chrysene	21.062	228	131842	2.157	ng	97
88) Benzo(b)fluoranthene	22.880	252	200693	3.076	ng	# 95
90) Benzo(a)pyrene	23.591	252	140167	2.480	ng	# 92

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

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