

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081524\
 Data File : BM047221.D
 Acq On : 15 Aug 2024 18:53
 Operator : RC/JU
 Sample : P3518-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OU4-TS-04-080724

Quant Time: Aug 16 01:19:45 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)

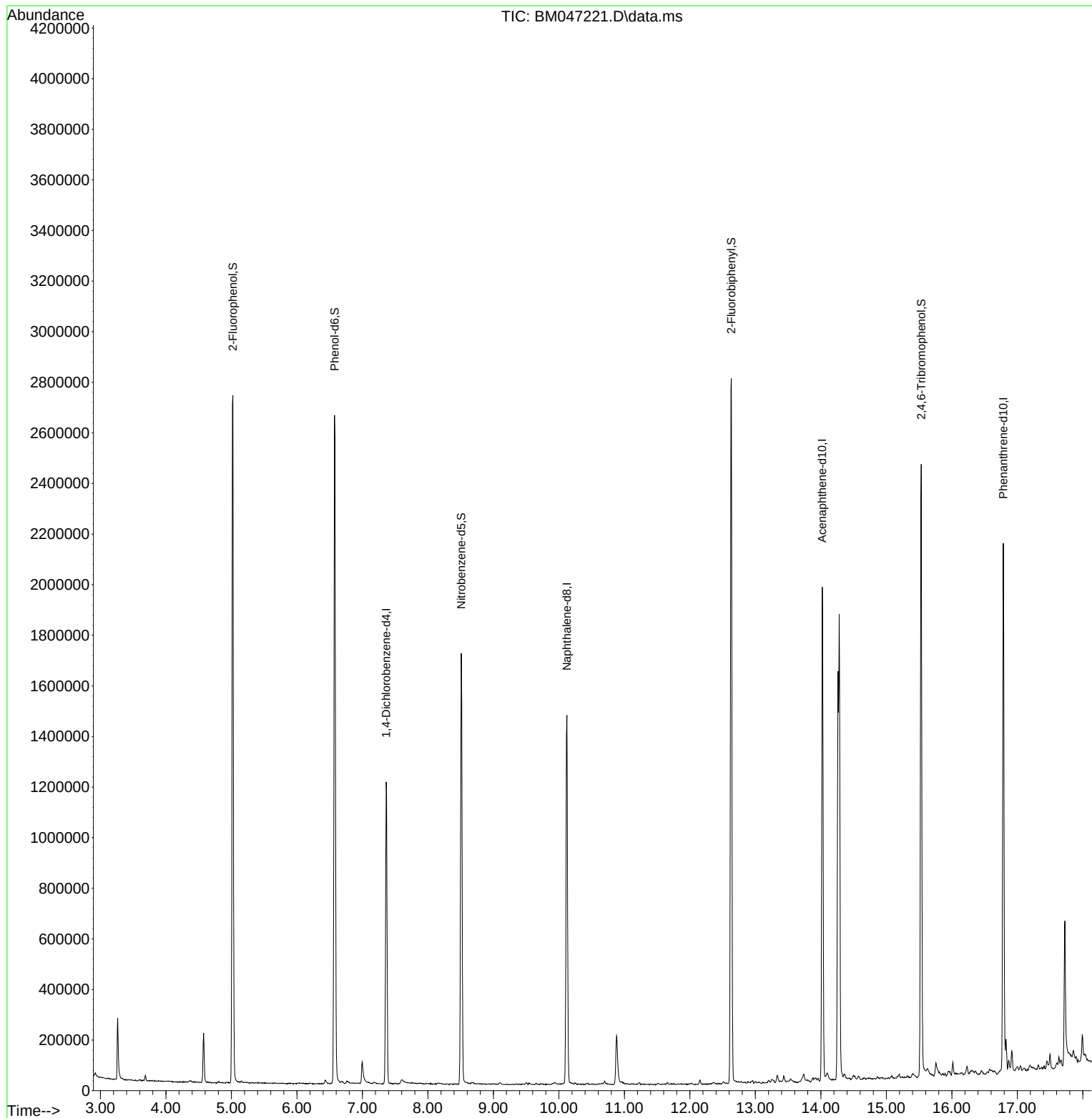
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	326789	20.000	ng	-0.01
21) Naphthalene-d8	10.122	136	1192472	20.000	ng	-0.01
39) Acenaphthene-d10	14.021	164	681602	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1195572	20.000	ng	0.00
76) Chrysene-d12	21.021	240	953020	20.000	ng	-0.01
86) Perylene-d12	23.721	264	1083116	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	1155066	63.158	ng	0.00
7) Phenol-d6	6.575	99	1510761	59.950	ng	-0.01
23) Nitrobenzene-d5	8.510	82	971221	41.011	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	437371	55.340	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	1534382	34.724	ng	-0.01
79) Terphenyl-d14	19.451	244	1559342	35.060	ng	-0.01
Target Compounds						
75) Fluoranthene	18.862	202	161031	2.206	ng	99
78) Pyrene	19.227	202	146360	2.151	ng	99

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

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