

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081624\
 Data File : BM047233.D
 Acq On : 16 Aug 2024 13:51
 Operator : RC/JU
 Sample : P3478-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TS07

Quant Time: Aug 16 14:28:22 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)

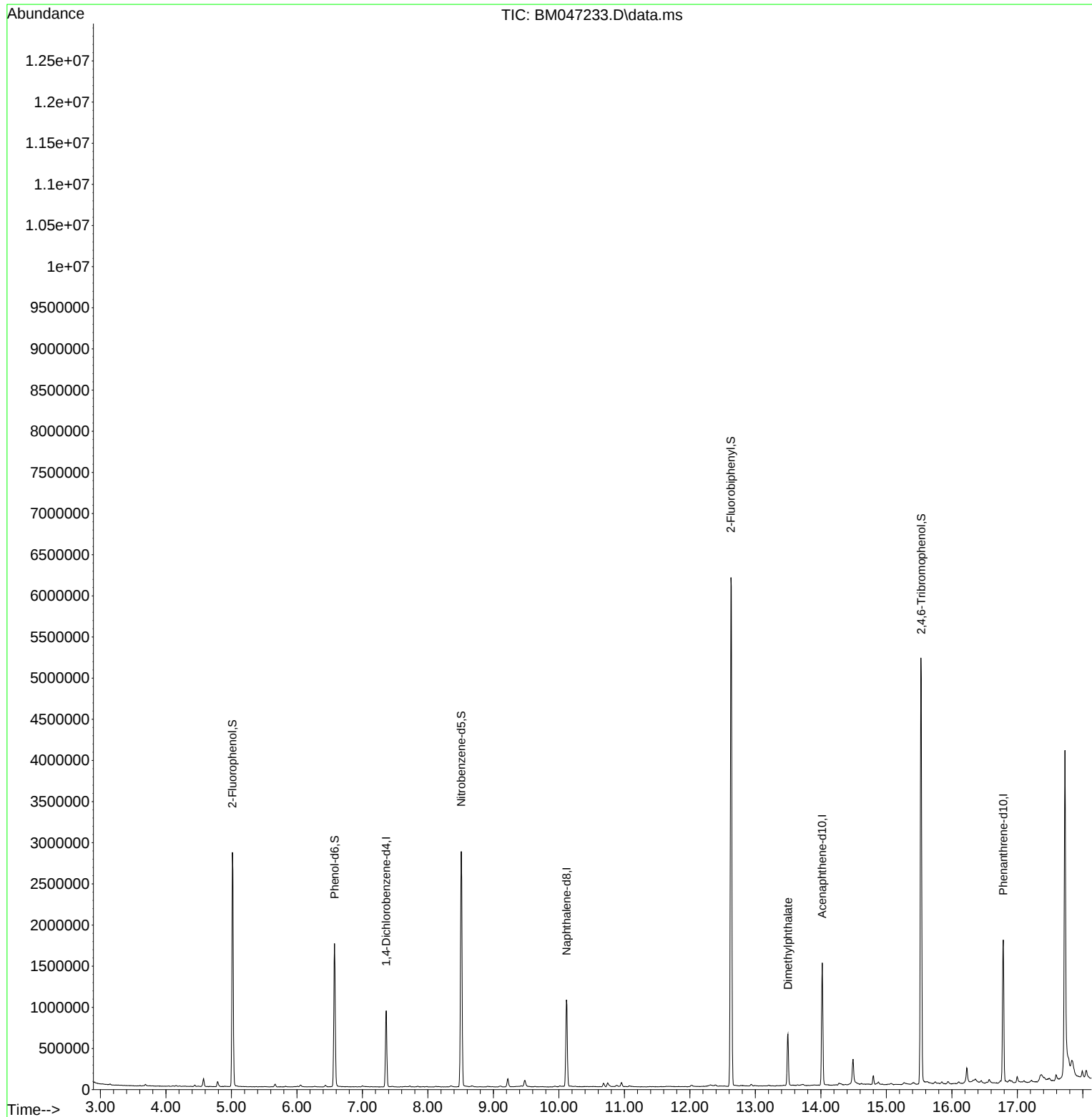
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	237293	20.000	ng	-0.01
21) Naphthalene-d8	10.116	136	855543	20.000	ng	-0.02
39) Acenaphthene-d10	14.021	164	507793	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1034778	20.000	ng	0.00
76) Chrysene-d12	21.021	240	1022079	20.000	ng	-0.01
86) Perylene-d12	23.715	264	1190964	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.016	112	1175995	88.554	ng	0.00
7) Phenol-d6	6.575	99	971600	53.097	ng	-0.01
23) Nitrobenzene-d5	8.510	82	1666200	98.065	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	936165	158.997	ng	0.00
45) 2-Fluorobiphenyl	12.627	172	3347270	101.680	ng	-0.02
79) Terphenyl-d14	19.450	244	5019509	105.233	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.498	163	400424	11.291	ng	Qvalue 100

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

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