

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082024\
 Data File : BM047281.D
 Acq On : 20 Aug 2024 22:14
 Operator : RC/JU
 Sample : P3645-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 916-J-WS-081524

Quant Time: Aug 21 02:50:40 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	7.357	152	312772	20.000	ng	-0.02
21) Naphthalene-d8	10.110	136	1142660	20.000	ng	-0.02
39) Acenaphthene-d10	14.016	164	745797	20.000	ng	-0.02
64) Phenanthrene-d10	16.780	188	1530970	20.000	ng	-0.01
76) Chrysene-d12	21.021	240	1309543	20.000	ng	-0.01
86) Perylene-d12	23.715	264	1291407	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.016	112	1270694	72.594	ng	0.00
7) Phenol-d6	6.569	99	1090156	45.199	ng	-0.02
23) Nitrobenzene-d5	8.504	82	1744202	76.861	ng	-0.02
42) 2,4,6-Tribromophenol	15.527	330	1147929	132.745	ng	-0.01
45) 2-Fluorobiphenyl	12.628	172	3947185	81.639	ng	-0.02
79) Terphenyl-d14	19.451	244	5714616	93.506	ng	-0.01

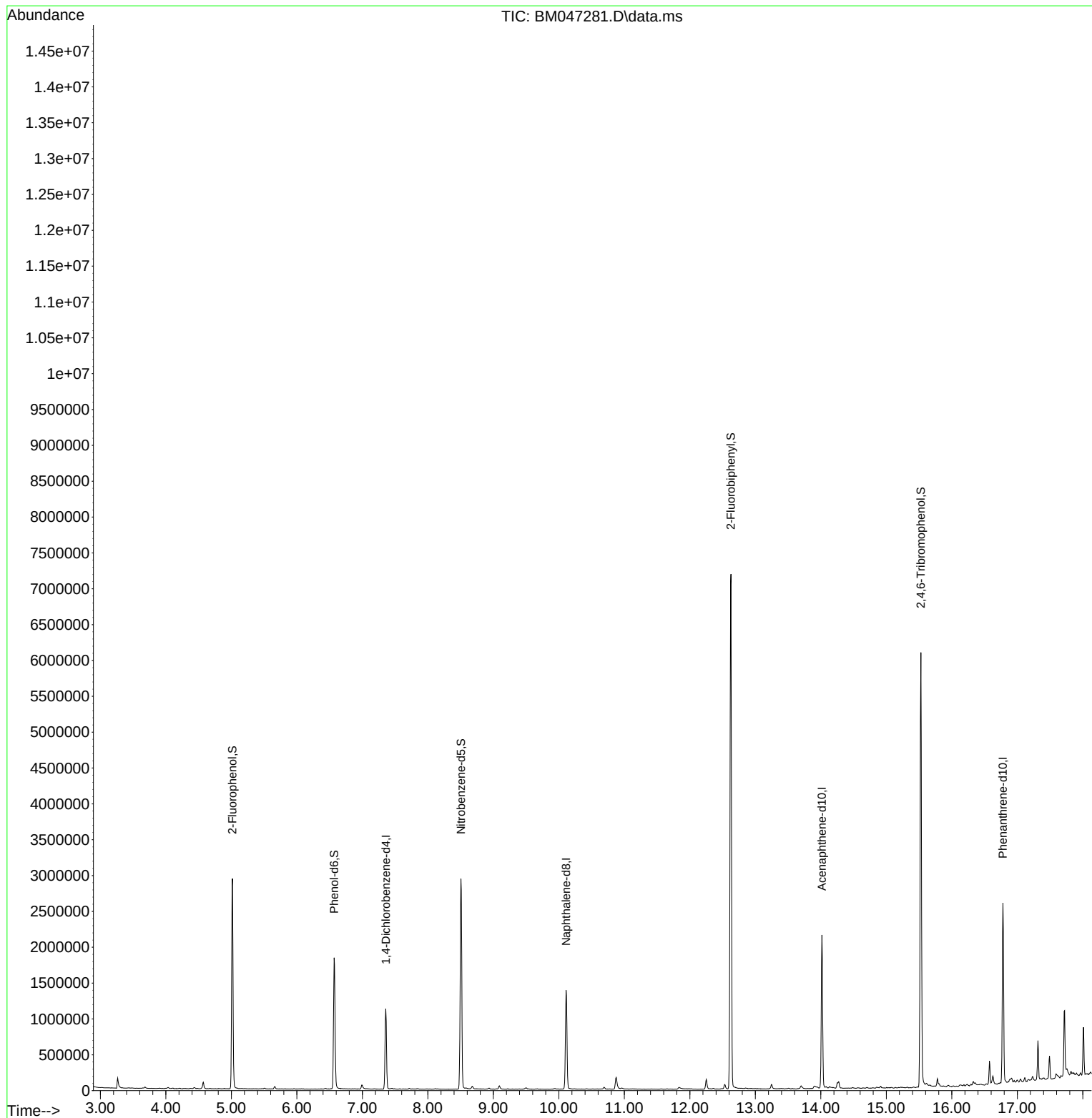
Target Compounds	Qvalue

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

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