

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092324\
 Data File : BM047606.D
 Acq On : 23 Sep 2024 16:23
 Operator : RC/JU
 Sample : P4139-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 371

Quant Time: Sep 23 17:02:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM092024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 21 02:22:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	298364	20.000	ng	-0.01
21) Naphthalene-d8	10.633	136	1138621	20.000	ng	0.00
39) Acenaphthene-d10	14.468	164	601031	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1171927	20.000	ng	0.00
76) Chrysene-d12	21.427	240	993688	20.000	ng	-0.01
86) Perylene-d12	24.444	264	1073973	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	1628488	83.642	ng	-0.01
7) Phenol-d6	6.998	99	2176748	84.850	ng	0.00
23) Nitrobenzene-d5	8.986	82	1323908	57.881	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	434723	78.112	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	2523385	59.406	ng	-0.01
79) Terphenyl-d14	19.821	244	3754272	67.790	ng	0.00

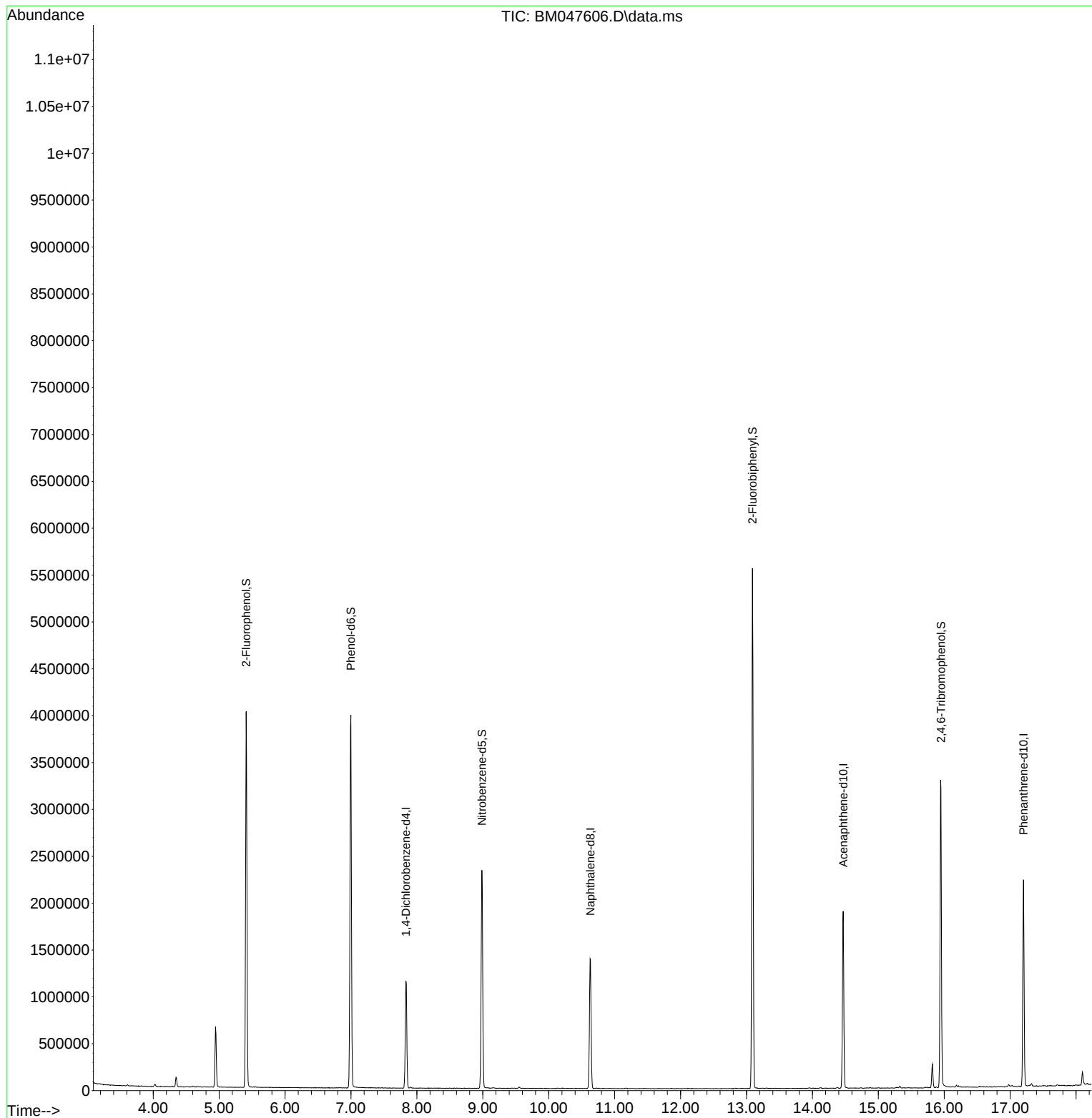
Target Compounds	Qvalue

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

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