

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092324\
 Data File : BM047601.D
 Acq On : 23 Sep 2024 11:43
 Operator : RC/JU
 Sample : P3845-22
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RR-TRIAZINE-WP

Quant Time: Sep 23 12:31:02 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)

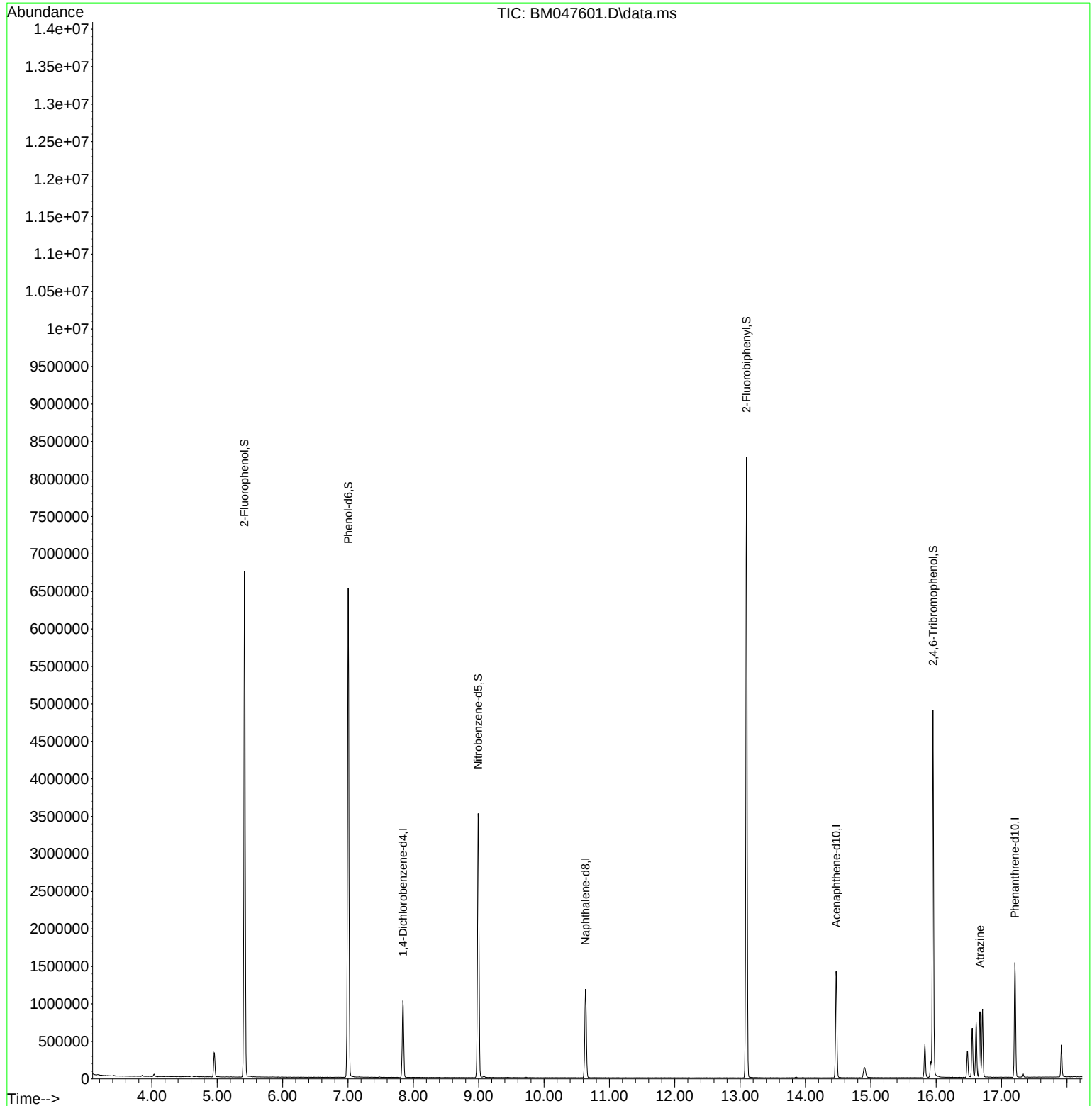
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	260539	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	961904	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	453032	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	806425	20.000	ng	0.00
76) Chrysene-d12	21.427	240	573772	20.000	ng	-0.01
86) Perylene-d12	24.450	264	644450	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	2674156	157.289	ng	0.00
7) Phenol-d6	7.004	99	3424961	152.888	ng	0.00
23) Nitrobenzene-d5	8.993	82	1998964	103.450	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	604557	144.116	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	3663225	114.415	ng	0.00
79) Terphenyl-d14	19.821	244	4350820	136.057	ng	0.00
Target Compounds						
69) Atrazine	16.668	200	123987	21.383	ng	Qvalue 97

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

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