

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092324\
 Data File : BM047600.D
 Acq On : 23 Sep 2024 11:04
 Operator : RC/JU
 Sample : PB163344BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB163344BL

Quant Time: Sep 23 18:57:30 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	7.845	152	304832	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	1161519	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	523773	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	914542	20.000	ng	0.00
76) Chrysene-d12	21.427	240	616013	20.000	ng	-0.01
86) Perylene-d12	24.450	264	584879	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	1920175	96.531	ng	0.00
7) Phenol-d6	7.004	99	2440958	93.130	ng	0.00
23) Nitrobenzene-d5	8.992	82	2245447	96.235	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	338589	69.812	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	4038464	109.099	ng	0.00
79) Terphenyl-d14	19.821	244	4365635	127.158	ng	0.00

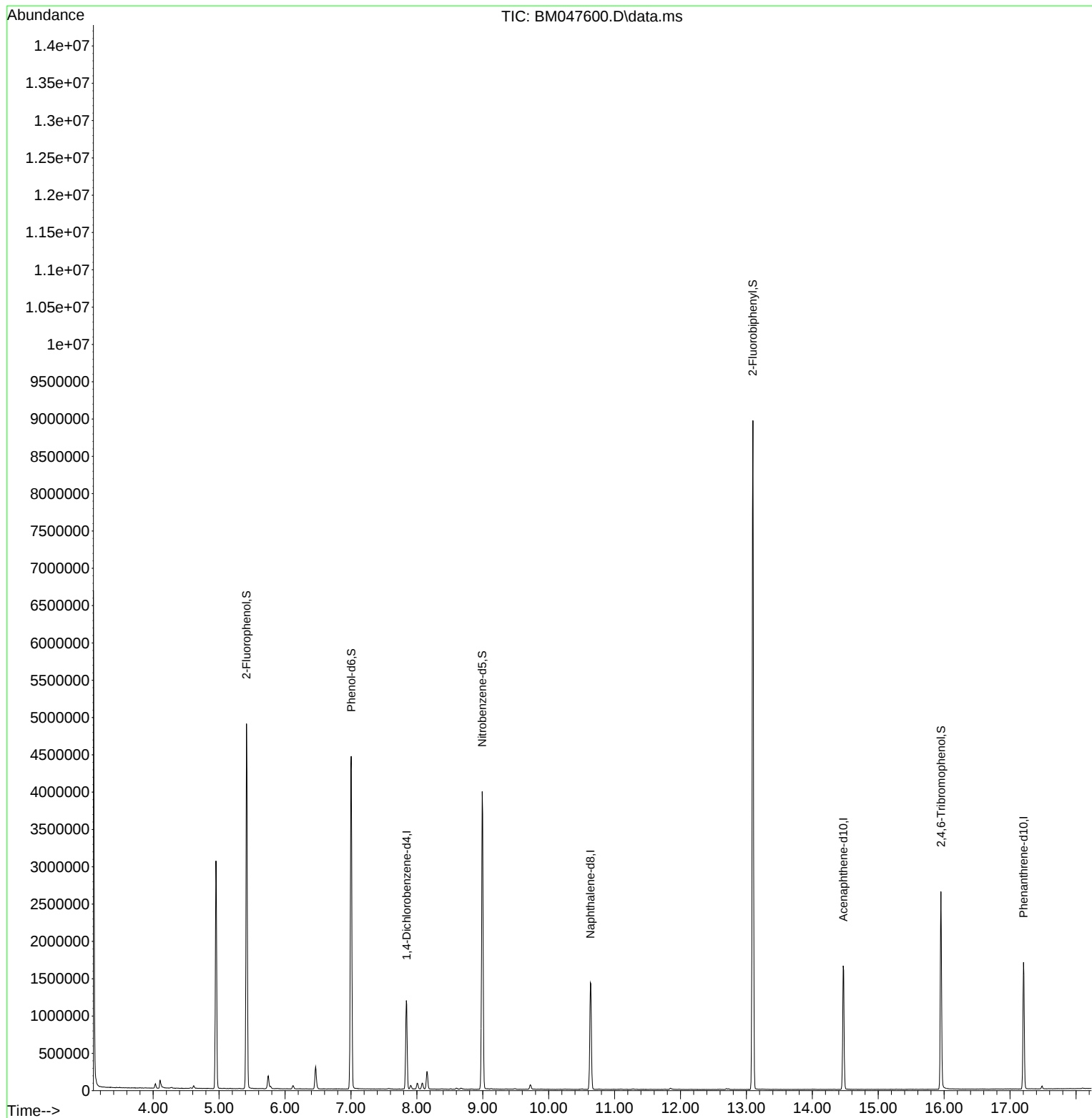
Target Compounds	Qvalue

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

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