

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091724\
 Data File : BM047539.D
 Acq On : 17 Sep 2024 09:27
 Operator : RC/JU
 Sample : PB163109BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB163109BL

Quant Time: Sep 17 11:08:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Sep 15 08:56:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	441001	20.000	ng	0.00
21) Naphthalene-d8	10.645	136	1746104	20.000	ng	0.00
39) Acenaphthene-d10	14.474	164	950653	20.000	ng	0.00
64) Phenanthrene-d10	17.209	188	1785434	20.000	ng	0.00
76) Chrysene-d12	21.439	240	1140059	20.000	ng	0.00
86) Perylene-d12	24.462	264	990611	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	4138782	149.429	ng	0.00
7) Phenol-d6	7.010	99	5232947	134.159	ng	0.00
23) Nitrobenzene-d5	9.004	82	3150090	90.287	ng	0.00
42) 2,4,6-Tribromophenol	15.962	330	1223566	133.157	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	6423567	93.079	ng	0.00
79) Terphenyl-d14	19.827	244	8633056	135.694	ng	0.00

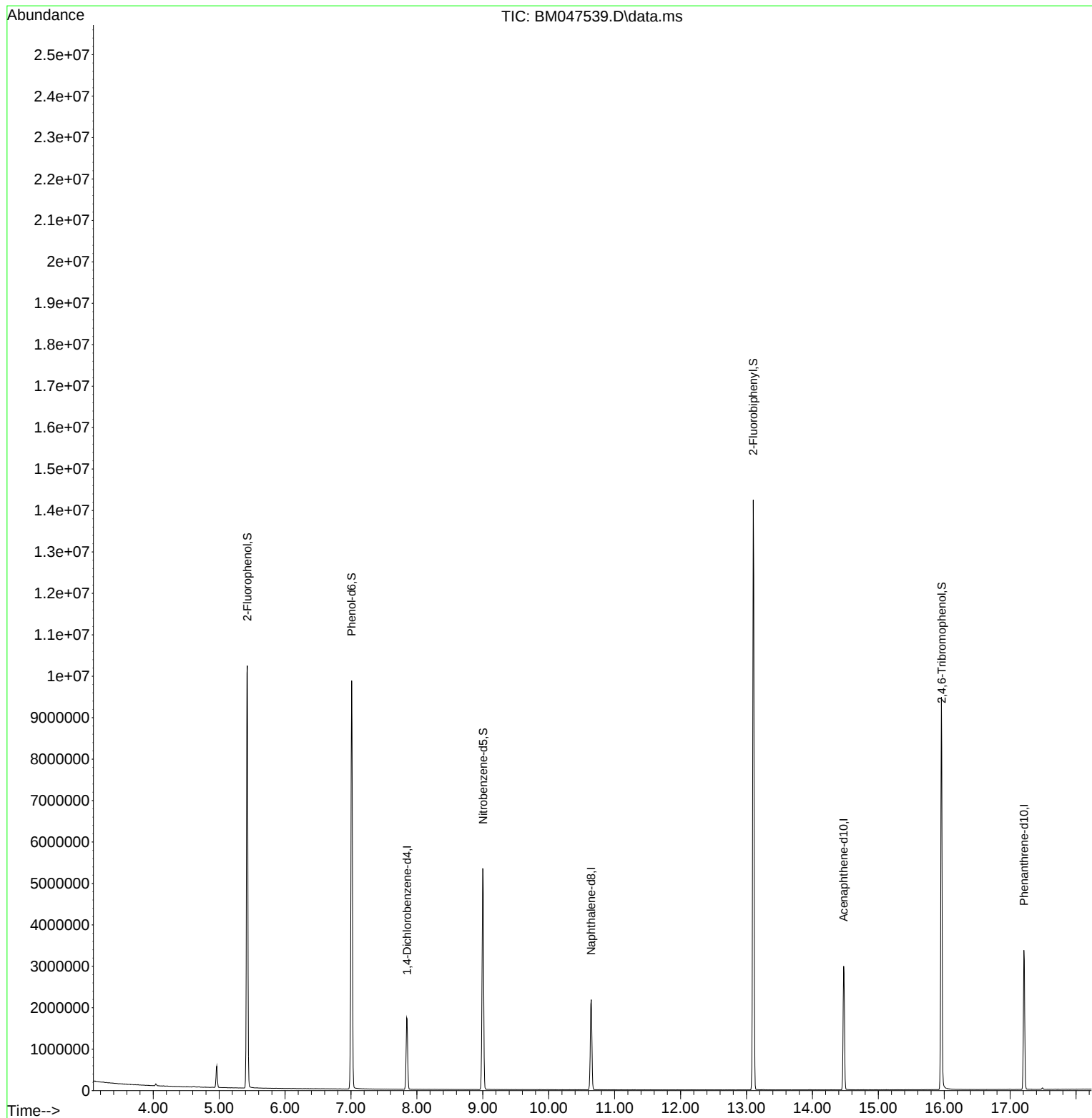
Target Compounds	Qvalue

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

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