

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102424\
 Data File : BM048225.D
 Acq On : 24 Oct 2024 20:26
 Operator : RC/JU
 Sample : PB164366BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB164366BL

Quant Time: Oct 25 00:00:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 18:21:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	217658	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	789248	20.000	ng	0.00
39) Acenaphthene-d10	14.363	164	473893	20.000	ng	0.00
64) Phenanthrene-d10	17.110	188	888629	20.000	ng	0.00
76) Chrysene-d12	21.333	240	683470	20.000	ng	0.00
86) Perylene-d12	24.286	264	722400	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	1849767	142.855	ng	0.00
7) Phenol-d6	6.898	99	2235740	132.581	ng	0.00
23) Nitrobenzene-d5	8.875	82	1285890	91.993	ng	0.00
42) 2,4,6-Tribromophenol	15.857	330	724027	124.822	ng	0.00
45) 2-Fluorobiphenyl	12.980	172	2890105	93.583	ng	0.00
79) Terphenyl-d14	19.733	244	3440075	102.715	ng	0.00

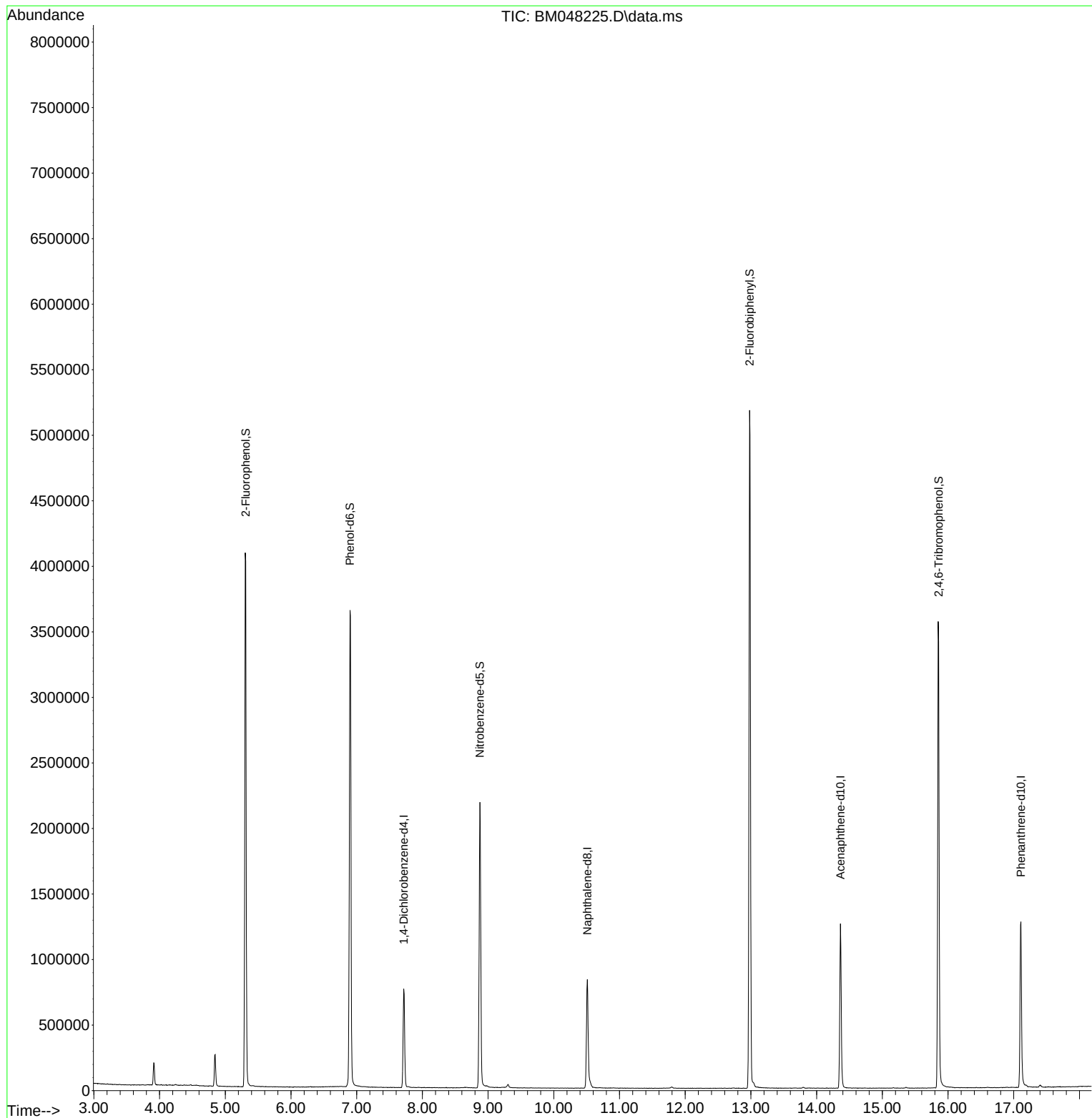
Target Compounds	Qvalue

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

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