

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102324\
 Data File : BM048209.D
 Acq On : 23 Oct 2024 18:25
 Operator : RC/JU
 Sample : PB164286BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB164286BL

Quant Time: Oct 24 01:01:28 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	7.722	152	158631	20.000	ng	0.00
21) Naphthalene-d8	10.516	136	584492	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	345617	20.000	ng	0.00
64) Phenanthrene-d10	17.110	188	674329	20.000	ng	0.00
76) Chrysene-d12	21.333	240	563717	20.000	ng	0.00
86) Perylene-d12	24.286	264	638791	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	1501590	159.117	ng	0.00
7) Phenol-d6	6.904	99	1770057	144.024	ng	0.00
23) Nitrobenzene-d5	8.881	82	1049093	101.344	ng	0.00
42) 2,4,6-Tribromophenol	15.857	330	602159	142.342	ng	0.00
45) 2-Fluorobiphenyl	12.986	172	2305028	102.340	ng	0.00
79) Terphenyl-d14	19.733	244	3044307	110.208	ng	0.00

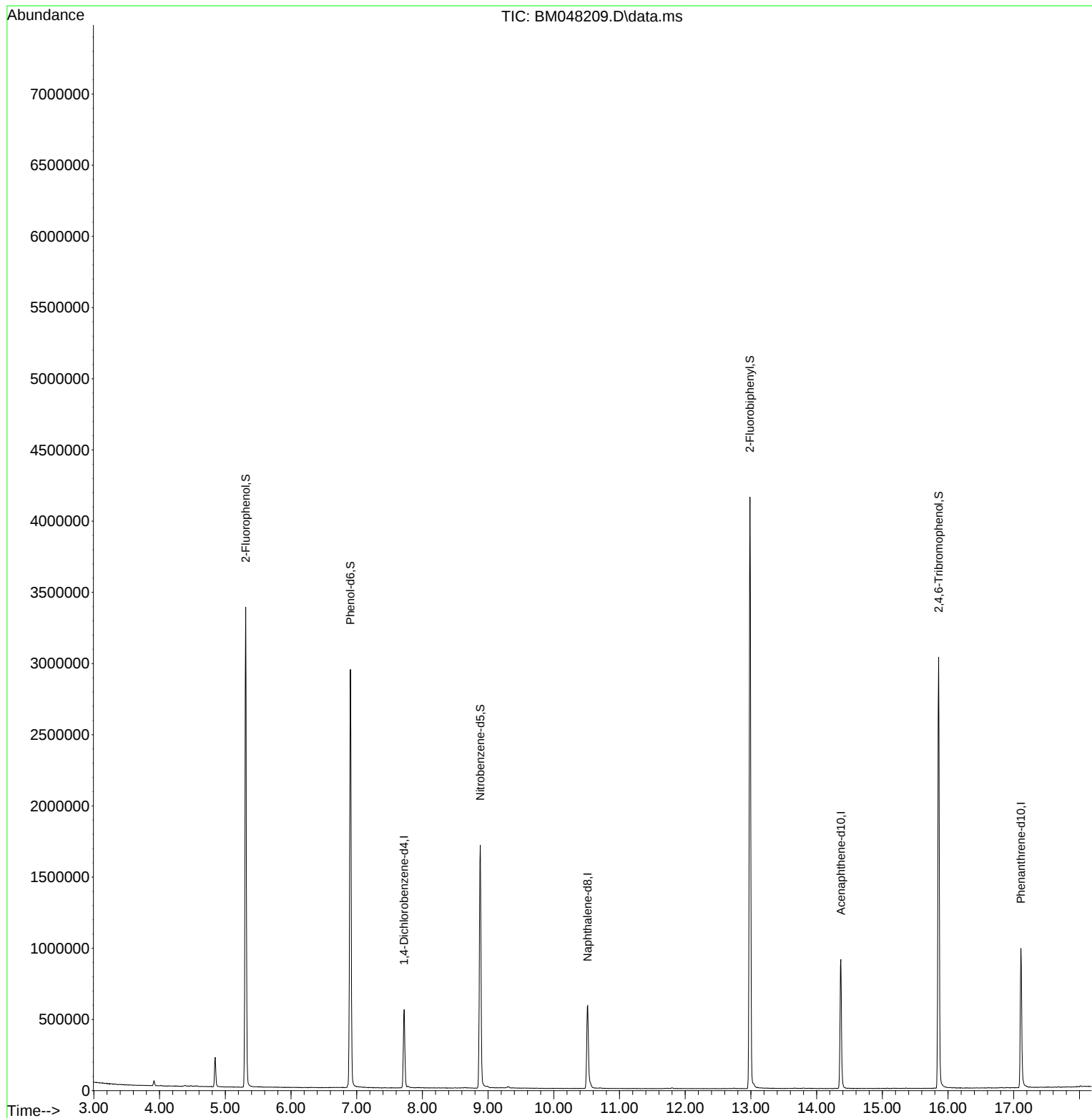
Target Compounds	Qvalue

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

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