

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041625\
 Data File : BM049948.D
 Acq On : 16 Apr 2025 12:05
 Operator : RC/JU
 Sample : PB167587TB
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB167587TB

Quant Time: Apr 16 12:44:18 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 04:00:55 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	340959	20.000	ng	-0.01
21) Naphthalene-d8	10.563	136	1139922	20.000	ng	-0.01
39) Acenaphthene-d10	14.416	164	732723	20.000	ng	0.00
64) Phenanthrene-d10	17.162	188	1442241	20.000	ng	0.00
76) Chrysene-d12	21.403	240	1379497	20.000	ng	0.00
86) Perylene-d12	24.409	264	1521991	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2587891	128.885	ng	0.00
7) Phenol-d6	6.951	99	3096568	123.951	ng	0.00
23) Nitrobenzene-d5	8.928	82	1879437	91.811	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	1513490	142.009	ng	0.00
45) 2-Fluorobiphenyl	13.033	172	4915373	90.966	ng	-0.01
79) Terphenyl-d14	19.786	244	8274746	112.413	ng	0.00

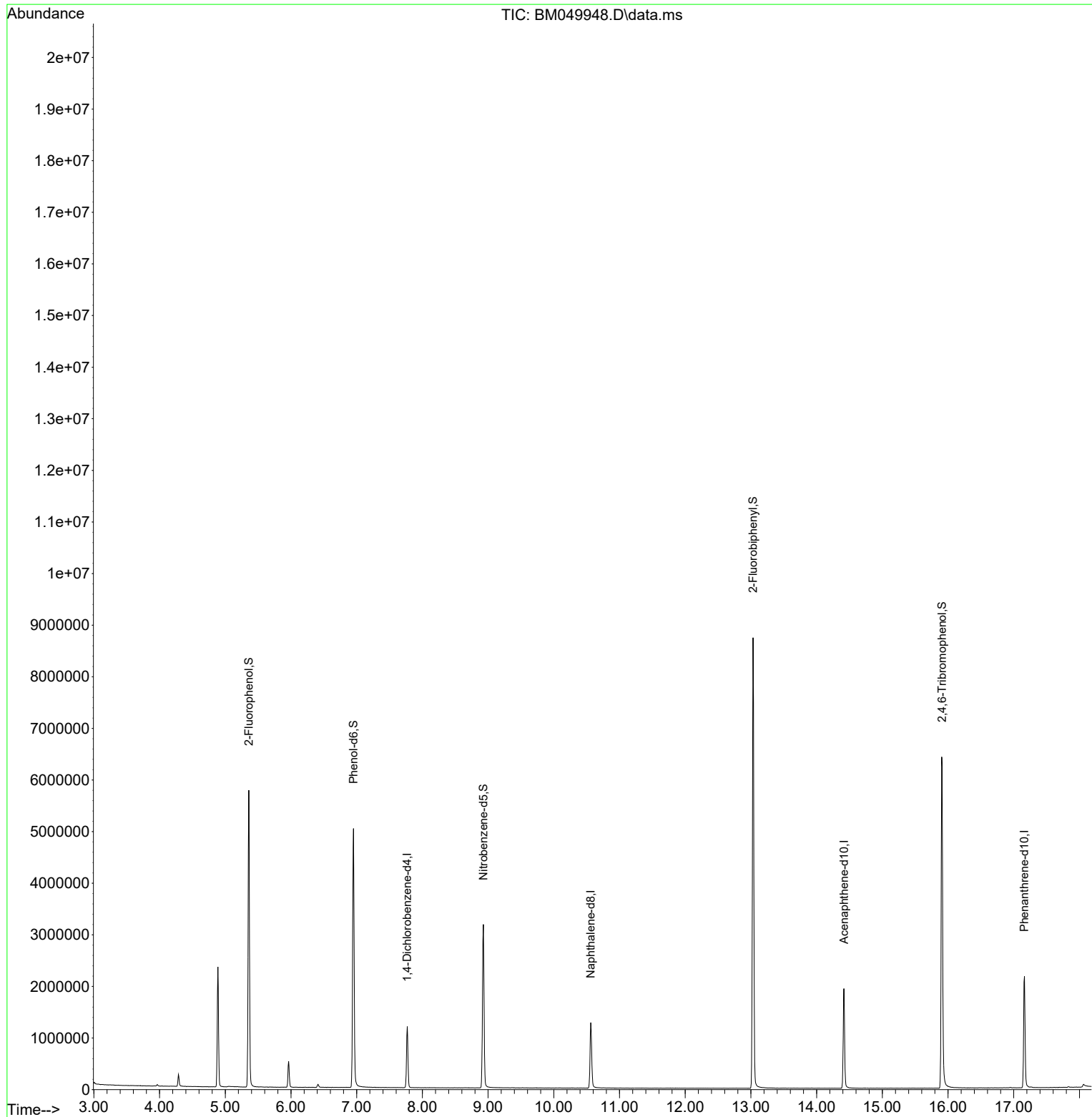
Target Compounds	Qvalue

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

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