

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM123124\
 Data File : BM049311.D
 Acq On : 31 Dec 2024 19:06
 Operator : RC/JU
 Sample : PB165894BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB165894BL

Quant Time: Dec 31 23:21:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM123124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 23:06:33 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.528	152	159339	20.000	ng	0.00
21) Naphthalene-d8	10.292	136	552549	20.000	ng	0.00
39) Acenaphthene-d10	14.169	164	333009	20.000	ng	0.00
64) Phenanthrene-d10	16.921	188	628110	20.000	ng	0.00
76) Chrysene-d12	21.150	240	535240	20.000	ng	0.00
86) Perylene-d12	23.933	264	541026	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.169	112	1815204	178.058	ng	0.00
7) Phenol-d6	6.728	99	2211391	165.590	ng	0.00
23) Nitrobenzene-d5	8.675	82	1328650	113.779	ng	0.00
42) 2,4,6-Tribromophenol	15.668	330	766532	172.375	ng	0.00
45) 2-Fluorobiphenyl	12.780	172	3016909	118.053	ng	0.00
79) Terphenyl-d14	19.574	244	4478342	138.481	ng	0.00

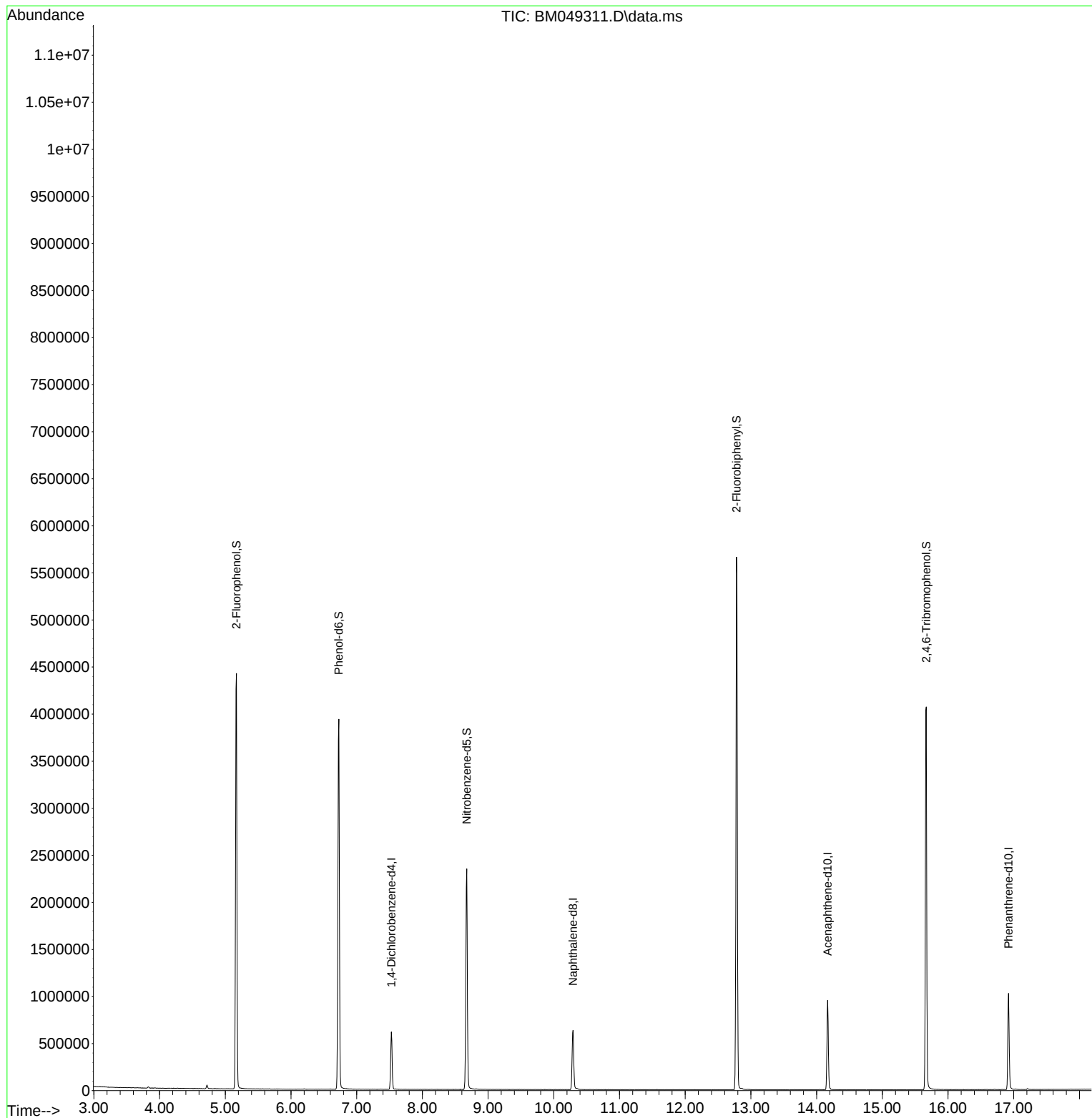
Target Compounds	Qvalue

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

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