

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091924\
 Data File : BM047573.D
 Acq On : 19 Sep 2024 16:33
 Operator : RC/JU
 Sample : PB163344BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB163344BL

Quant Time: Sep 19 17:20:29 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	7.839	152	319571	20.000	ng	-0.01
21) Naphthalene-d8	10.639	136	1227567	20.000	ng	0.00
39) Acenaphthene-d10	14.468	164	572655	20.000	ng	-0.01
64) Phenanthrene-d10	17.204	188	1053943	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	713655	20.000	ng	-0.01
86) Perylene-d12	24.450	264	703338	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	1995844	99.440	ng	0.00
7) Phenol-d6	7.004	99	2564878	90.743	ng	0.00
23) Nitrobenzene-d5	8.992	82	2413561	98.397	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	408784	73.852	ng	-0.01
45) 2-Fluorobiphenyl	13.098	172	4300764	103.455	ng	-0.01
79) Terphenyl-d14	19.821	244	5168848	129.786	ng	-0.01

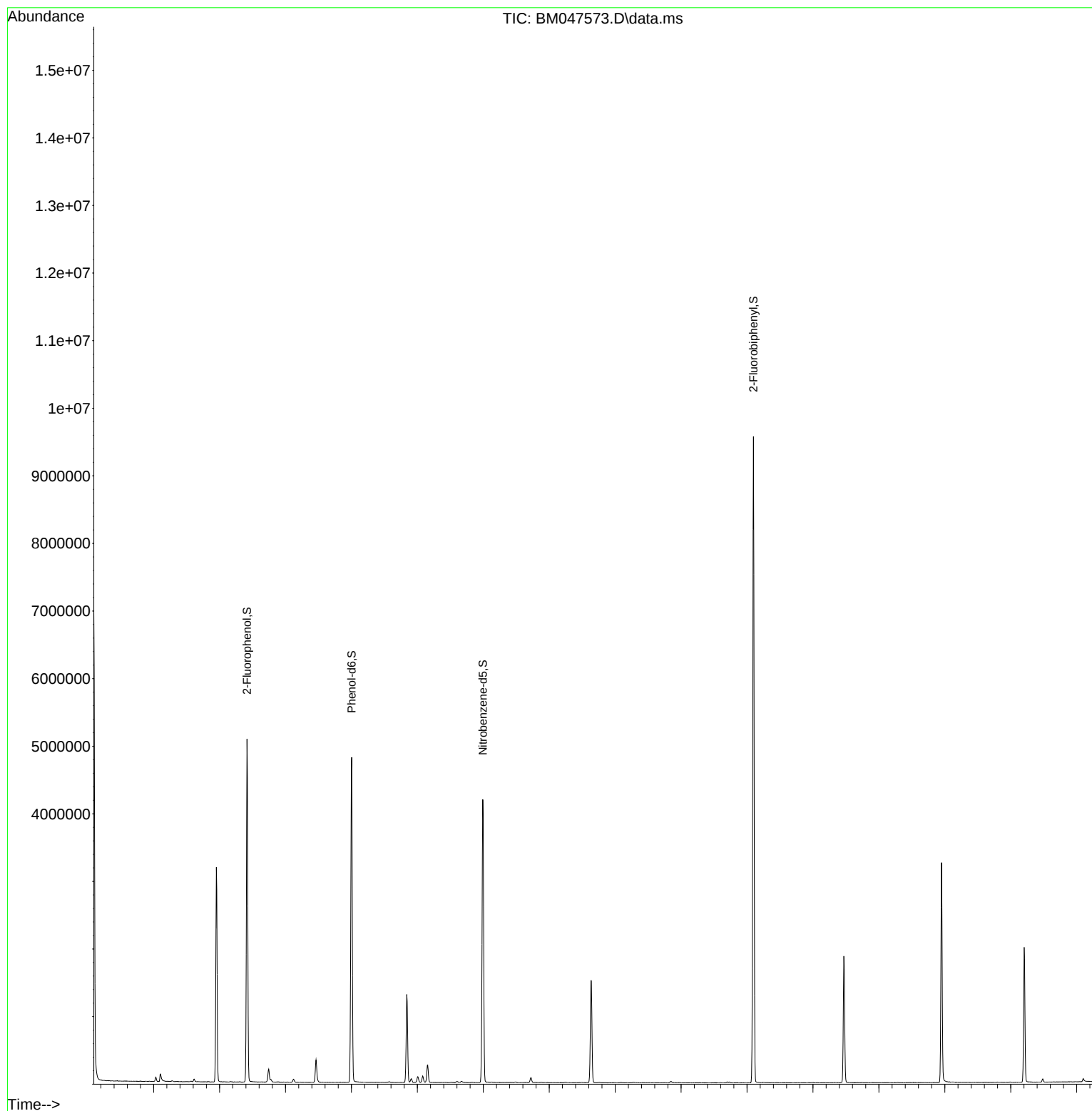
Target Compounds Qvalue

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

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