

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091924\
 Data File : BM047577.D
 Acq On : 19 Sep 2024 19:12
 Operator : RC/JU
 Sample : P4103-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-14-7.5-10B

Quant Time: Sep 20 02:32:57 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)

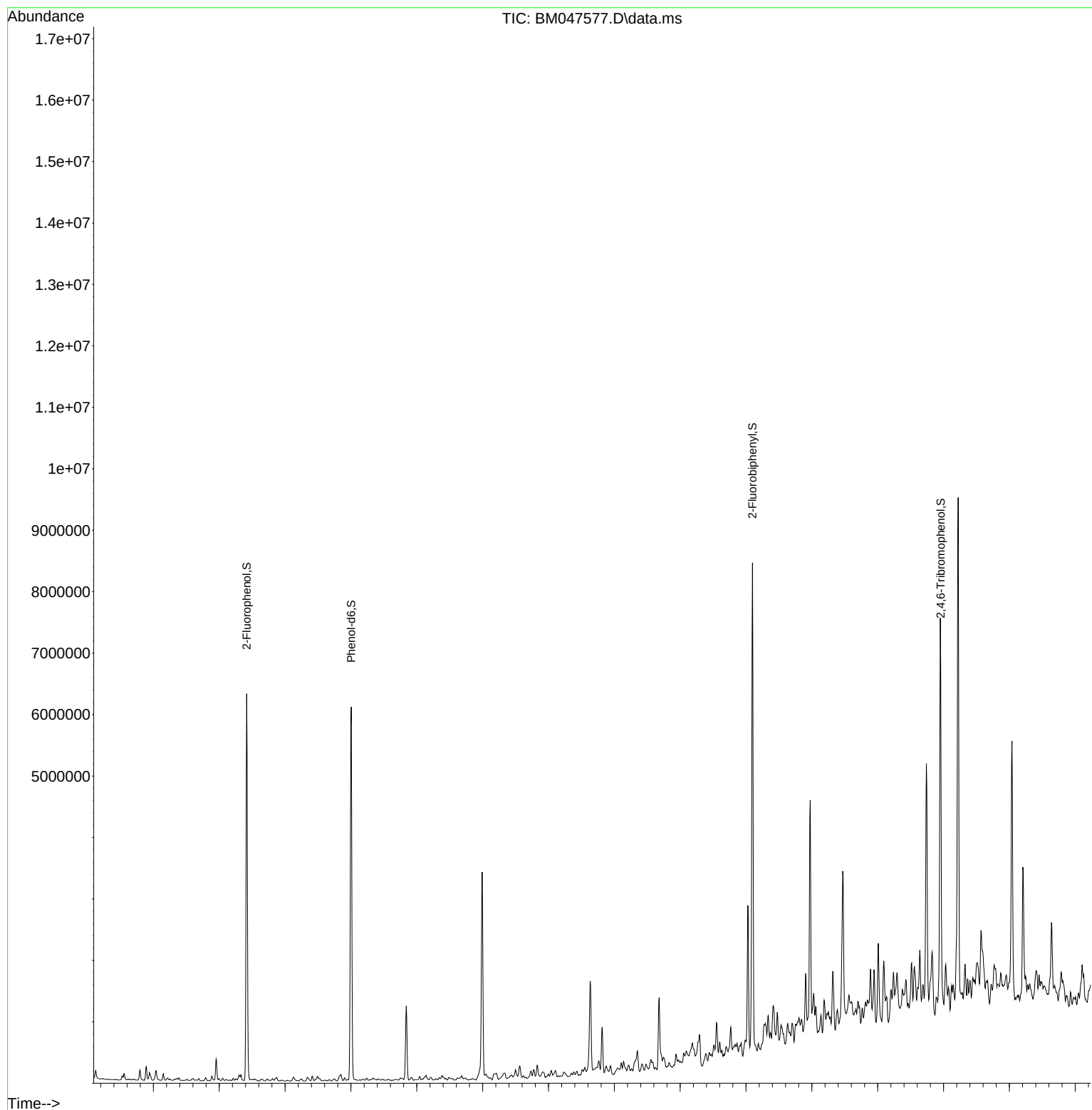
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	275927	20.000	ng	-0.01
21) Naphthalene-d8	10.634	136	1098397	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	619366	20.000	ng	-0.01
64) Phenanthrene-d10	17.204	188	1161651	20.000	ng	-0.01
76) Chrysene-d12	21.433	240	1002774	20.000	ng	-0.01
86) Perylene-d12	24.450	264	1123684	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	2468680	142.453	ng	0.00
7) Phenol-d6	7.004	99	3173715	130.043	ng	0.00
23) Nitrobenzene-d5	8.992	82	1878876	85.607	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	777214	129.823	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	3563836	79.263	ng	-0.01
79) Terphenyl-d14	19.821	244	5331522	95.273	ng	-0.01
Target Compounds						
58) Fluorene	15.516	166	109676	2.291	ng	Qvalue # 81

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

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