

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040125\
 Data File : BM049778.D
 Acq On : 01 Apr 2025 17:27
 Operator : RC/JU
 Sample : Q1632-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PIER-1-2

Quant Time: Apr 01 18:01:34 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 13 14:55:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	276000	20.000	ng	-0.01
21) Naphthalene-d8	10.586	136	944407	20.000	ng	-0.02
39) Acenaphthene-d10	14.433	164	612534	20.000	ng	-0.01
64) Phenanthrene-d10	17.174	188	1131243	20.000	ng	-0.01
76) Chrysene-d12	21.403	240	854565	20.000	ng	-0.02
86) Perylene-d12	24.409	264	783643	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	1446796	88.450	ng	-0.01
7) Phenol-d6	6.963	99	2219832	108.139	ng	-0.02
23) Nitrobenzene-d5	8.945	82	1405789	76.424	ng	-0.02
42) 2,4,6-Tribromophenol	15.921	330	532761	74.485	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	3909427	79.932	ng	-0.01
79) Terphenyl-d14	19.797	244	5293248	104.378	ng	-0.01

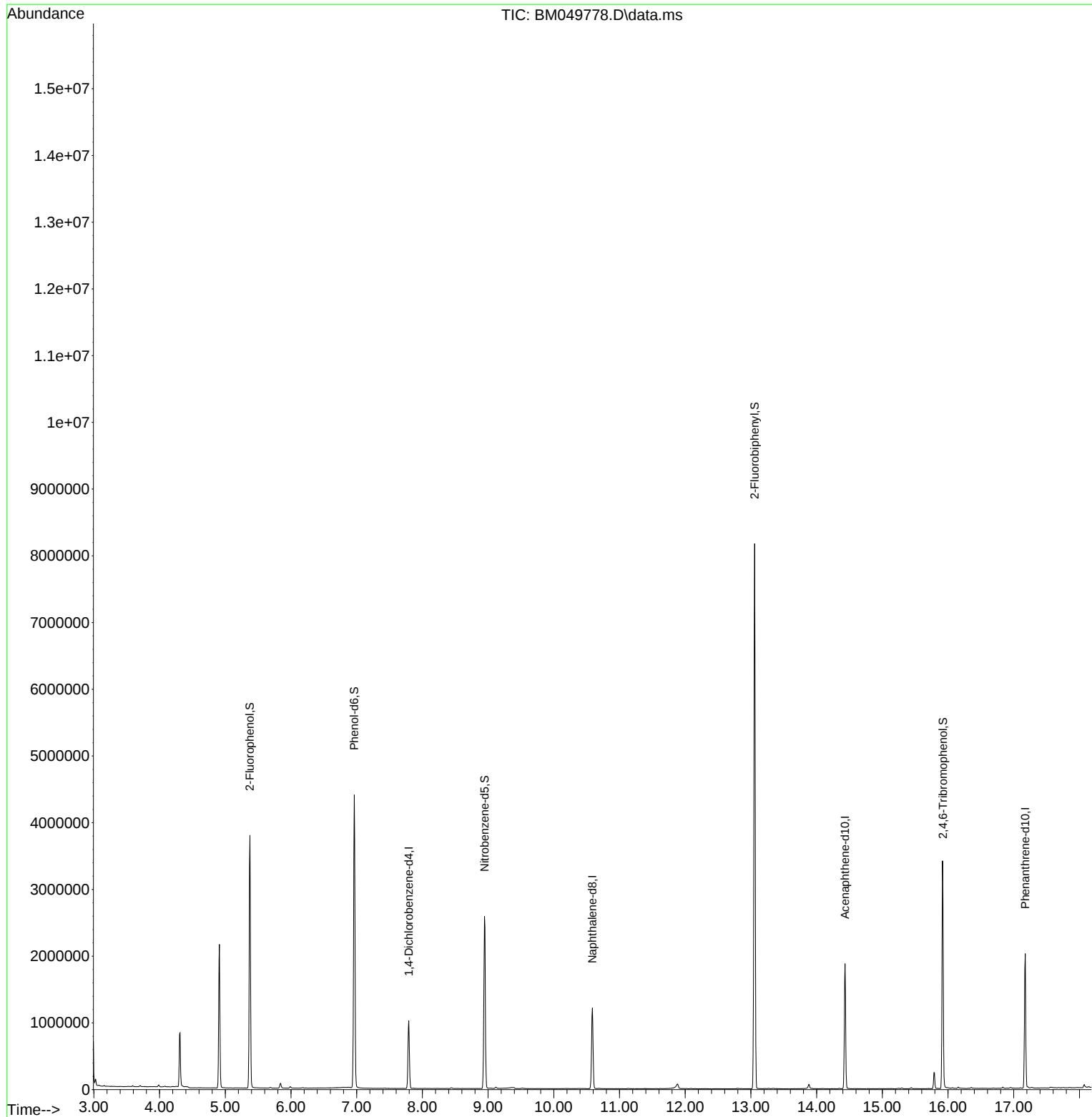
Target Compounds	Qvalue

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

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