

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040125\
 Data File : BM049779.D
 Acq On : 01 Apr 2025 18:06
 Operator : RC/JU
 Sample : Q1662-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-6

Quant Time: Apr 01 19:19:37 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	310337	20.000	ng	-0.01
21) Naphthalene-d8	10.586	136	1060057	20.000	ng	-0.02
39) Acenaphthene-d10	14.433	164	672580	20.000	ng	-0.01
64) Phenanthrene-d10	17.174	188	1210736	20.000	ng	-0.01
76) Chrysene-d12	21.403	240	835223	20.000	ng	-0.02
86) Perylene-d12	24.409	264	746205	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	3592376	195.320	ng	0.00
7) Phenol-d6	6.969	99	4463605	193.386	ng	-0.01
23) Nitrobenzene-d5	8.951	82	2674812	129.549	ng	-0.01
42) 2,4,6-Tribromophenol	15.921	330	1836095	233.787	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	7058484	131.433	ng	-0.01
79) Terphenyl-d14	19.797	244	8720938	175.951	ng	-0.01

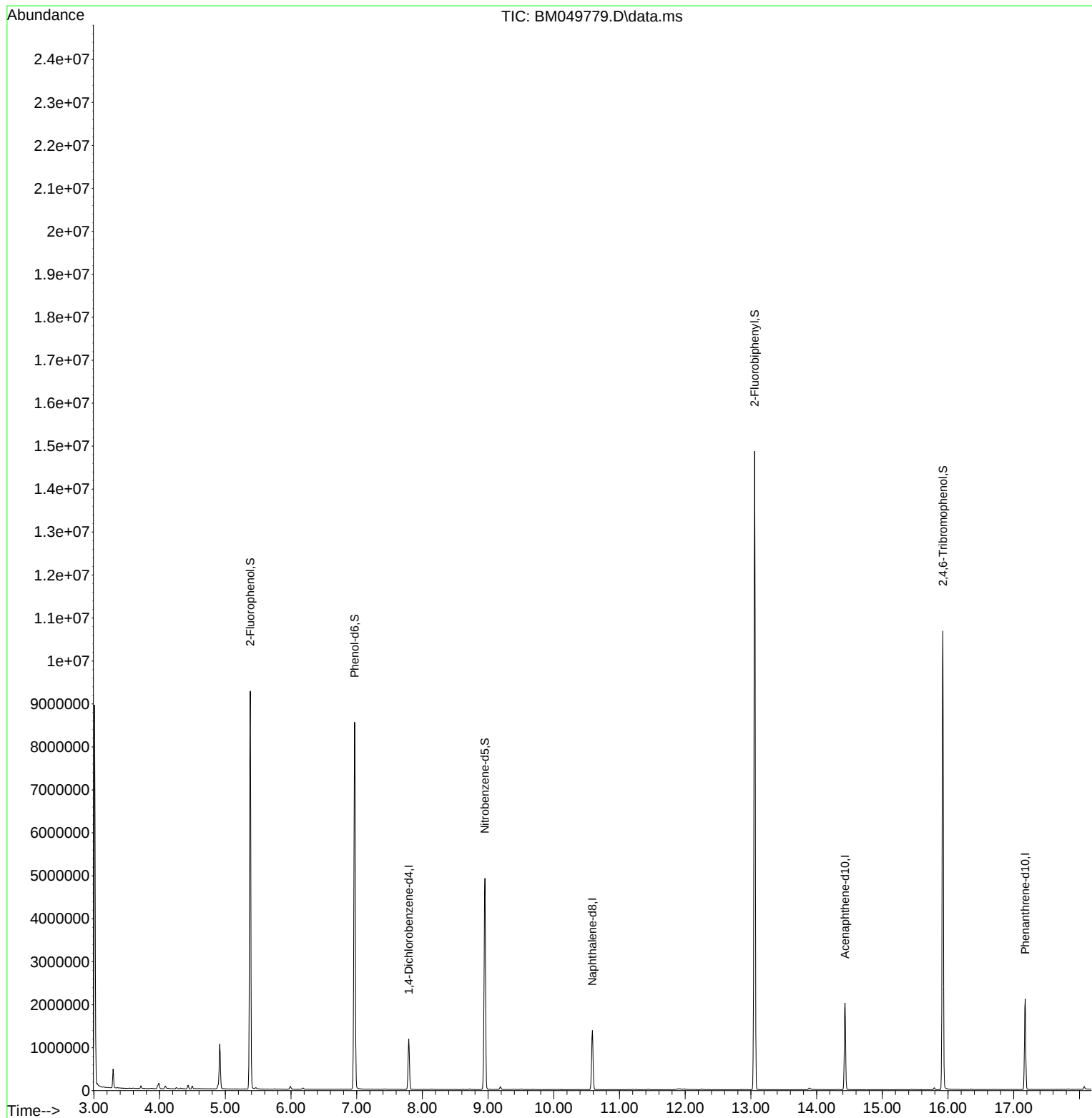
Target Compounds	Qvalue

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

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