

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022825\
 Data File : BM049676.D
 Acq On : 28 Feb 2025 13:15
 Operator : RC/JU
 Sample : PB166930BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB166930BL

Quant Time: Feb 28 14:28:47 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 04:55:28 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	355589	20.000	ng	-0.02
21) Naphthalene-d8	10.604	136	1196115	20.000	ng	-0.02
39) Acenaphthene-d10	14.445	164	682743	20.000	ng	-0.02
64) Phenanthrene-d10	17.186	188	1208848	20.000	ng	-0.02
76) Chrysene-d12	21.427	240	924150	20.000	ng	-0.02
86) Perylene-d12	24.444	264	944069	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	3287031	148.650	ng	0.00
7) Phenol-d6	6.987	99	3916766	135.200	ng	0.00
23) Nitrobenzene-d5	8.963	82	2324979	99.983	ng	-0.02
42) 2,4,6-Tribromophenol	15.933	330	1272925	157.347	ng	-0.02
45) 2-Fluorobiphenyl	13.069	172	5612050	106.484	ng	-0.02
79) Terphenyl-d14	19.815	244	7773370	128.445	ng	-0.02

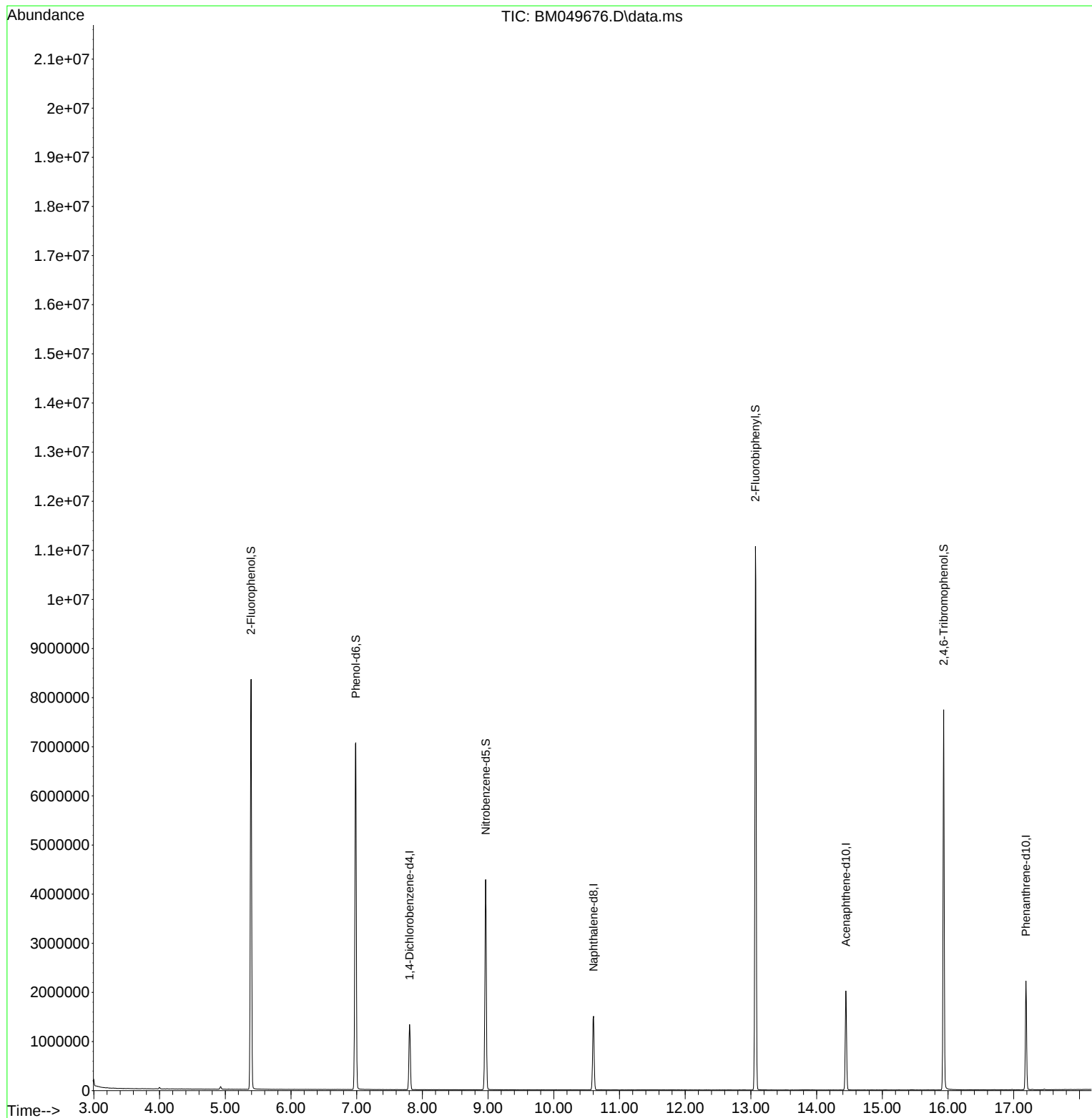
Target Compounds	Qvalue

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

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