

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072225\
 Data File : BM050481.D
 Acq On : 22 Jul 2025 11:29
 Operator : RC/JU
 Sample : PB168929BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168929BL

Quant Time: Jul 22 19:41:19 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 17 16:20:15 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	552021	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	1914515	20.000	ng	0.00
39) Acenaphthene-d10	14.468	164	1205881	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	2563541	20.000	ng	0.00
76) Chrysene-d12	21.427	240	2673293	20.000	ng	0.00
86) Perylene-d12	24.450	264	2939248	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	4104774	127.996	ng	0.00
7) Phenol-d6	7.010	99	5060084	125.165	ng	0.00
23) Nitrobenzene-d5	8.998	82	2995161	79.815	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	2243185	153.102	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	7772805	79.601	ng	0.00
79) Terphenyl-d14	19.815	244	12429504	81.806	ng	0.00

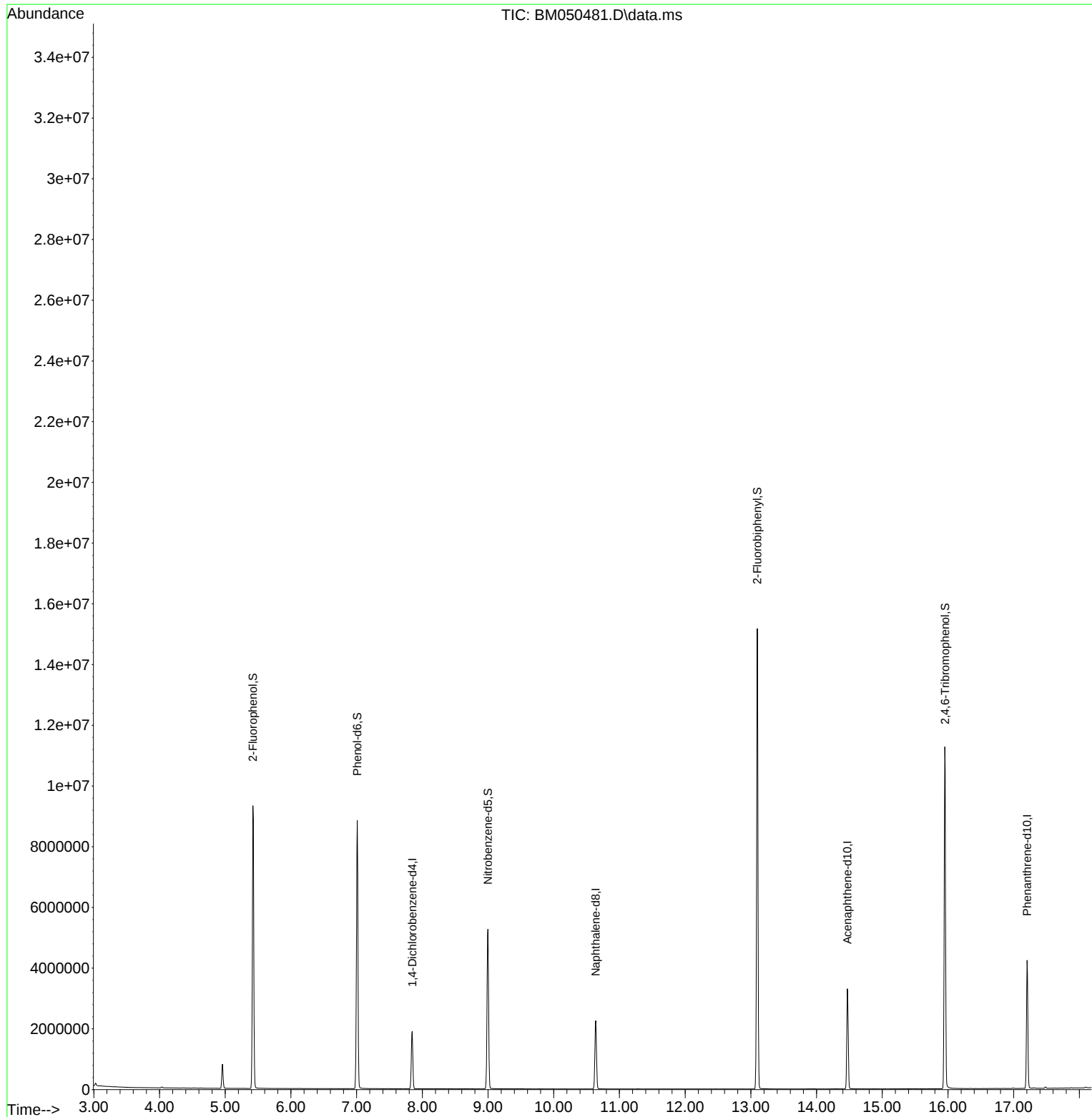
Target Compounds	Qvalue

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

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