

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071025\
 Data File : BM050389.D
 Acq On : 10 Jul 2025 14:33
 Operator : RC/JU
 Sample : PB168689TB
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168689TB

Quant Time: Jul 10 15:09:37 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 08 18:32:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	428549	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1515361	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	896439	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	1634836	20.000	ng	0.00
76) Chrysene-d12	21.450	240	1402149	20.000	ng	0.00
86) Perylene-d12	24.485	264	1470081	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3495276	140.392	ng	0.00
7) Phenol-d6	7.022	99	4263708	135.853	ng	0.00
23) Nitrobenzene-d5	9.016	82	2456670	82.709	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	1470709	135.029	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	6111858	84.197	ng	0.00
79) Terphenyl-d14	19.839	244	8274820	103.835	ng	0.00

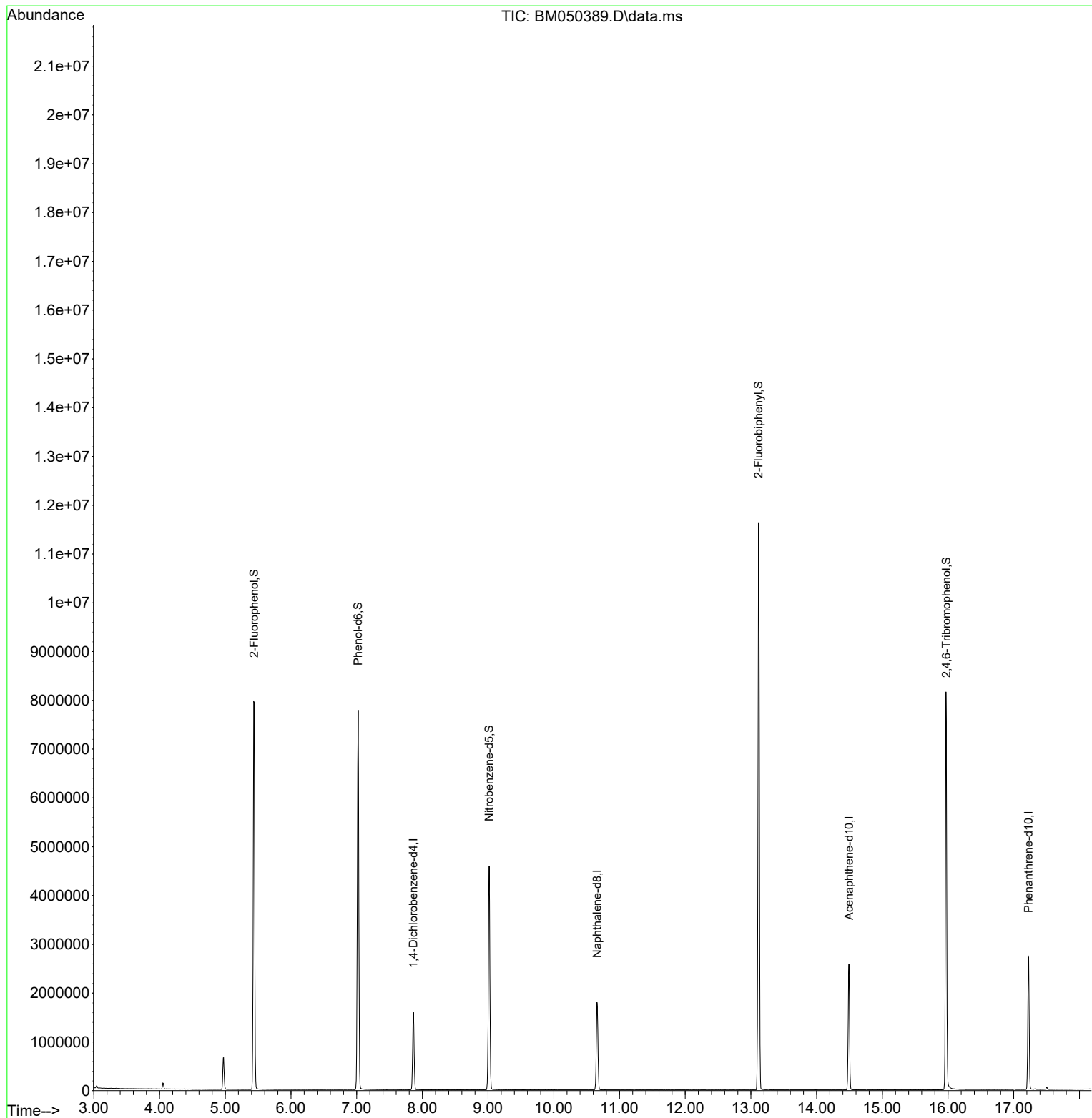
Target Compounds	Qvalue

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

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