

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050436.D
 Acq On : 14 Jul 2025 21:34
 Operator : RC/JU
 Sample : PB168728TB
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168728TB

Quant Time: Jul 15 02:04:45 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	354744	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1265655	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	770422	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	1489633	20.000	ng	-0.01
76) Chrysene-d12	21.439	240	1465640	20.000	ng	-0.02
86) Perylene-d12	24.468	264	1553412	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3173931	154.008	ng	0.00
7) Phenol-d6	7.022	99	3886308	149.590	ng	0.00
23) Nitrobenzene-d5	9.016	82	2309636	93.101	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1418630	151.552	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	5752719	92.212	ng	0.00
79) Terphenyl-d14	19.833	244	9302062	111.669	ng	-0.01

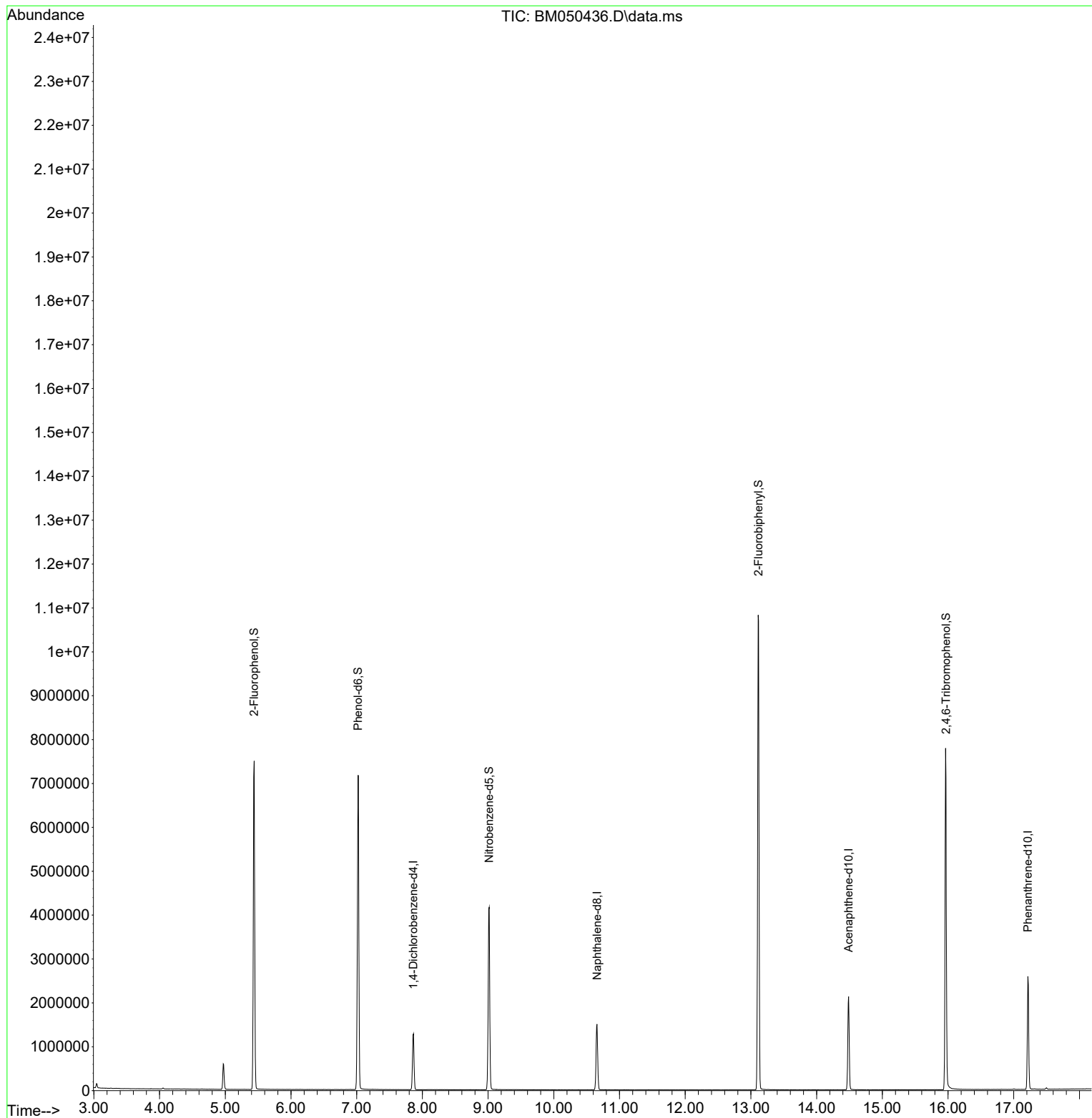
Target Compounds	Qvalue

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

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