

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071125\
 Data File : BM050415.D
 Acq On : 11 Jul 2025 16:24
 Operator : RC/JU
 Sample : PB168802BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168802BL

Quant Time: Jul 11 17:44:40 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	387286	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1348131	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	799830	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	1446767	20.000	ng	0.00
76) Chrysene-d12	21.450	240	1313023	20.000	ng	0.00
86) Perylene-d12	24.486	264	1392765	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3034104	134.853	ng	0.00
7) Phenol-d6	7.022	99	3688966	130.063	ng	0.00
23) Nitrobenzene-d5	9.016	82	2171017	82.159	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1211437	124.659	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	5368560	82.890	ng	0.00
79) Terphenyl-d14	19.839	244	7485657	100.308	ng	0.00

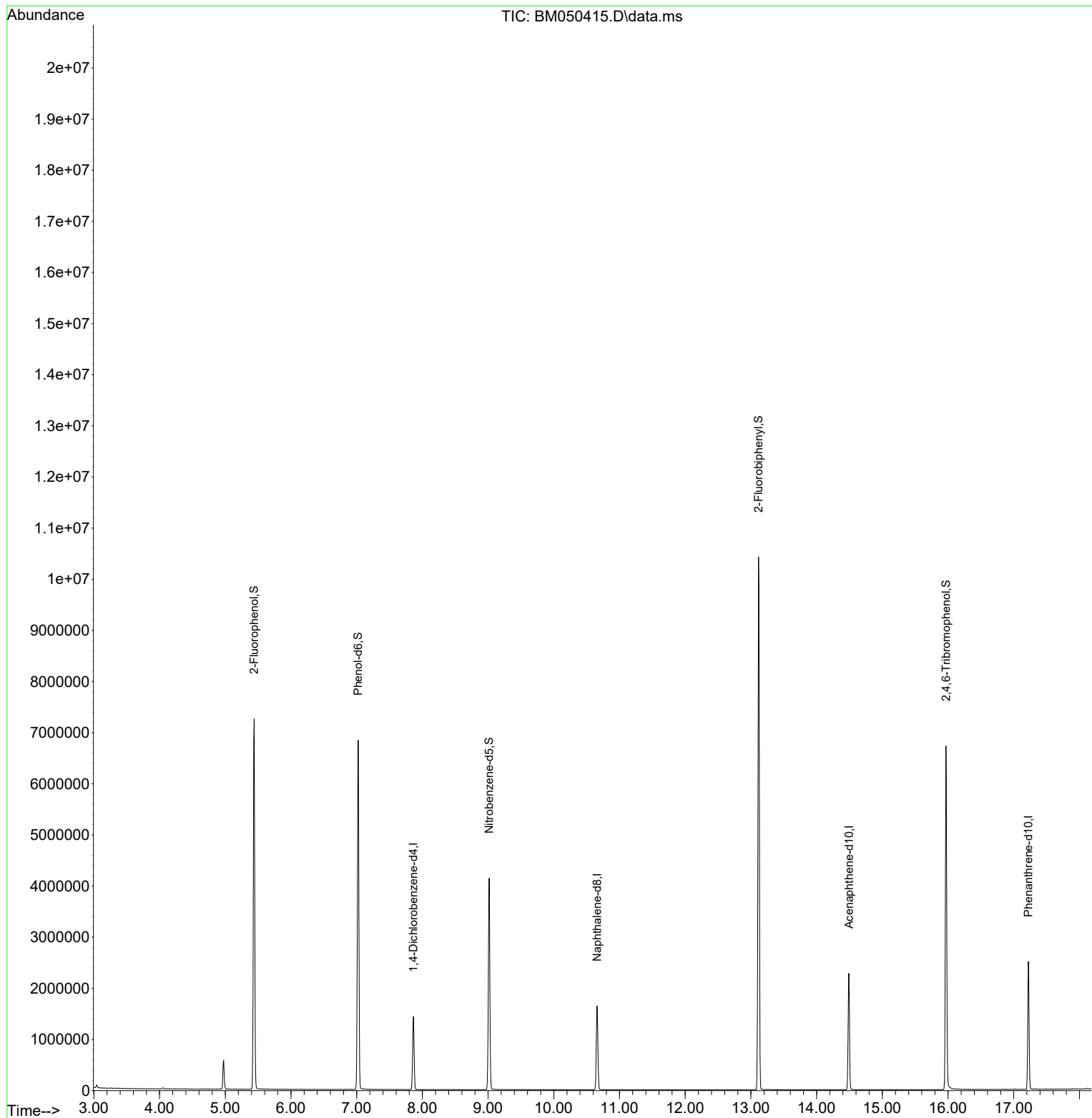
Target Compounds	Qvalue

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

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