

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050207.D
 Acq On : 06 Jun 2025 10:08
 Operator : RC/JU
 Sample : PB168269BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168269BL

Quant Time: Jun 06 10:53:13 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 05 16:20:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	307416	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1142669	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	634427	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1133741	20.000	ng	0.00
76) Chrysene-d12	21.380	240	982238	20.000	ng	0.00
86) Perylene-d12	24.368	264	1022855	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2527466	137.149	ng	0.00
7) Phenol-d6	6.951	99	3136982	129.309	ng	0.00
23) Nitrobenzene-d5	8.934	82	1832446	83.403	ng	-0.01
42) 2,4,6-Tribromophenol	15.892	330	906461	125.615	ng	-0.01
45) 2-Fluorobiphenyl	13.039	172	3884866	82.902	ng	-0.01
79) Terphenyl-d14	19.780	244	4336326	83.660	ng	0.00

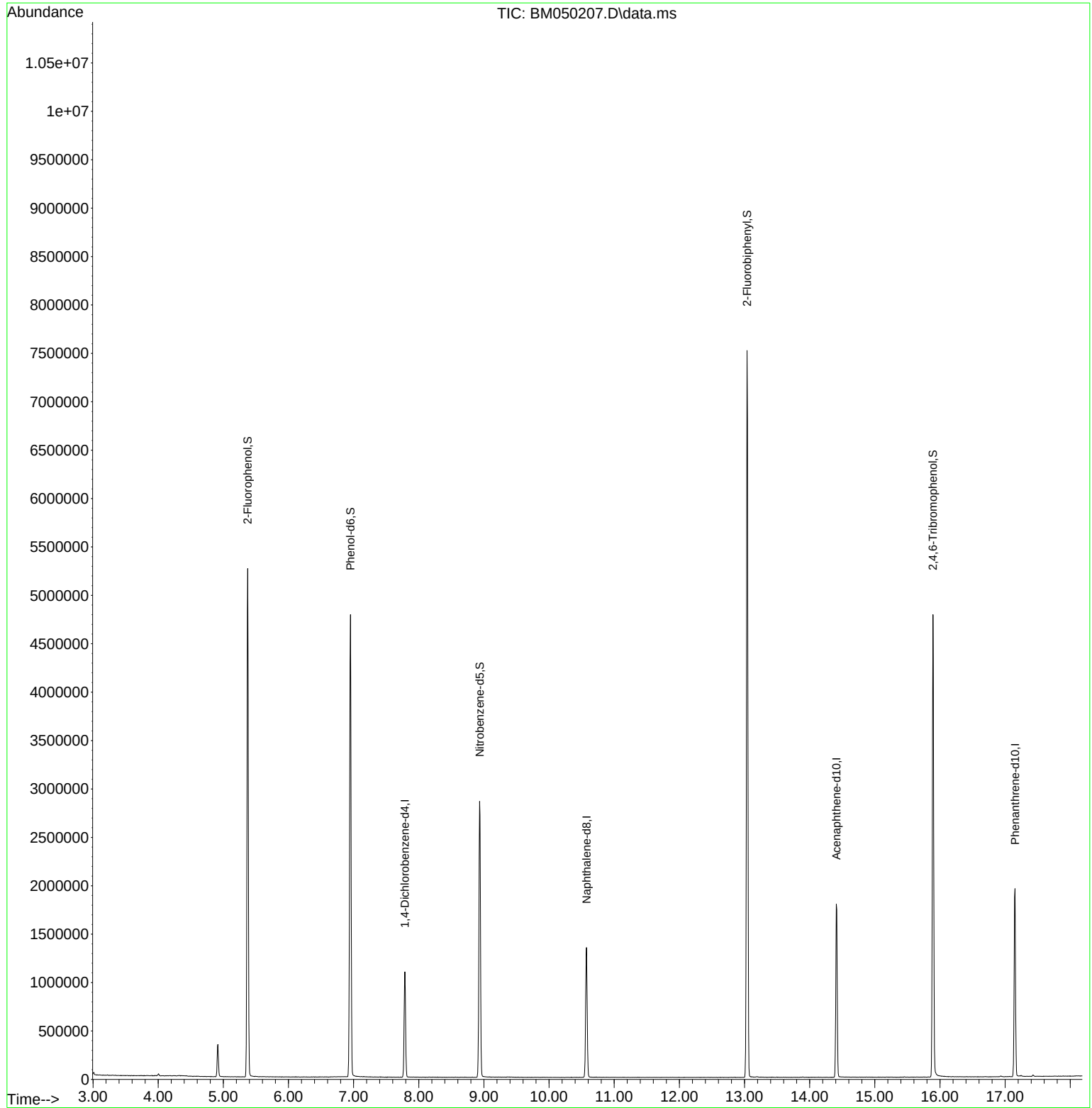
Target Compounds	Qvalue

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

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