

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050224.D
 Acq On : 06 Jun 2025 22:03
 Operator : RC/JU
 Sample : PB168314BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168314BL

Quant Time: Jun 07 01:21:07 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	7.786	152	315983	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1181137	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	658679	20.000	ng	0.00
64) Phenanthrene-d10	17.145	188	1198309	20.000	ng	-0.01
76) Chrysene-d12	21.374	240	1134031	20.000	ng	-0.01
86) Perylene-d12	24.356	264	1193795	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2551038	134.675	ng	0.00
7) Phenol-d6	6.951	99	3124058	125.285	ng	0.00
23) Nitrobenzene-d5	8.933	82	1813937	79.872	ng	-0.01
42) 2,4,6-Tribromophenol	15.892	330	935918	124.922	ng	-0.01
45) 2-Fluorobiphenyl	13.039	172	3959669	81.387	ng	-0.01
79) Terphenyl-d14	19.774	244	4894065	81.782	ng	-0.01

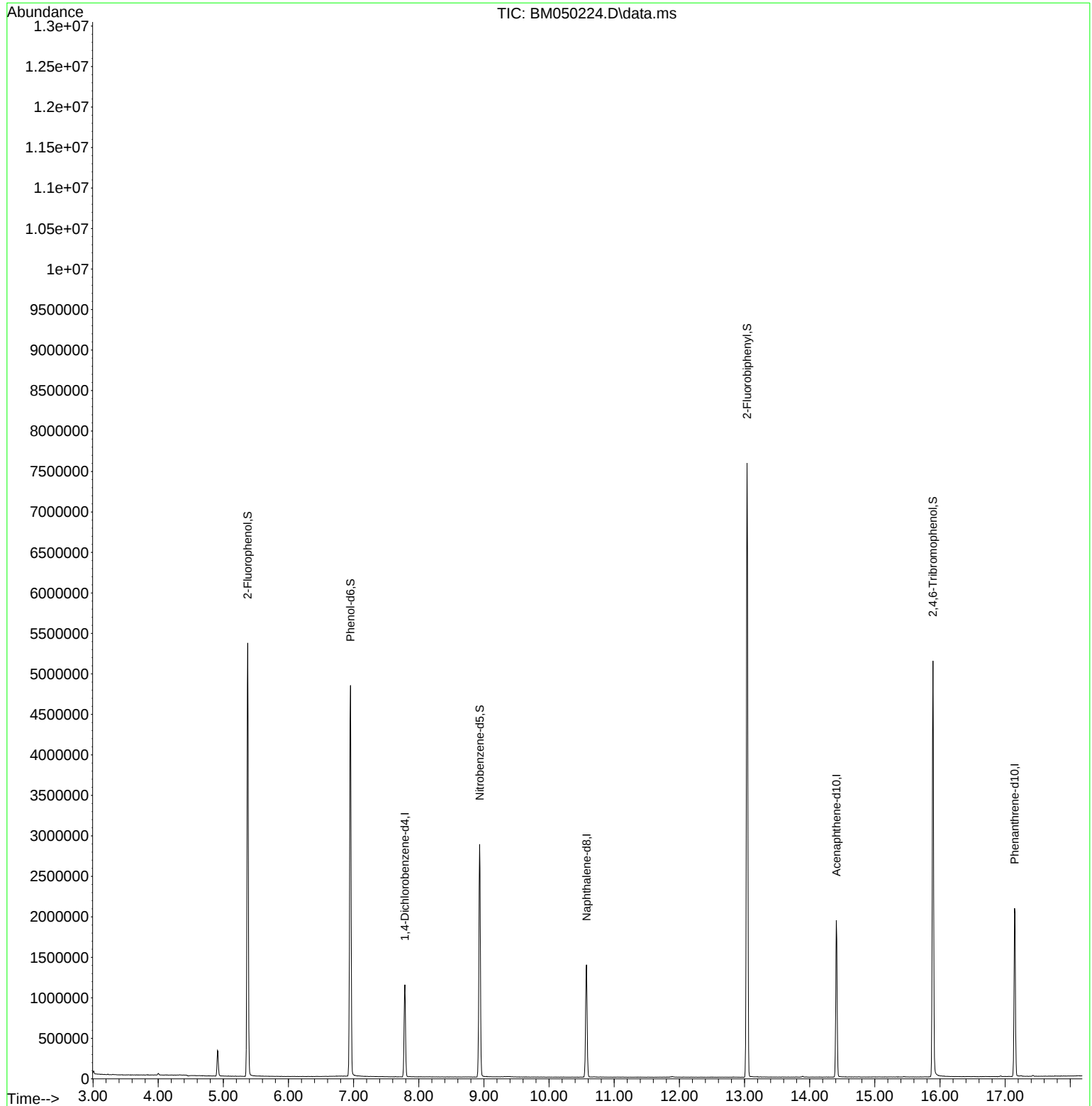
Target Compounds	Qvalue

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

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