

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061325\
 Data File : BM050301.D
 Acq On : 13 Jun 2025 14:35
 Operator : RC/JU
 Sample : PB168376BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168376BL

Quant Time: Jun 13 15:06:31 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 09 11:54:38 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	440199	20.000	ng	0.00
21) Naphthalene-d8	10.545	136	1615194	20.000	ng	-0.01
39) Acenaphthene-d10	14.392	164	895324	20.000	ng	0.00
64) Phenanthrene-d10	17.127	188	1712472	20.000	ng	-0.01
76) Chrysene-d12	21.356	240	1763403	20.000	ng	-0.02
86) Perylene-d12	24.327	264	1843126	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	3659530	138.679	ng	0.00
7) Phenol-d6	6.934	99	4521078	130.148	ng	0.00
23) Nitrobenzene-d5	8.910	82	2678000	86.230	ng	0.00
42) 2,4,6-Tribromophenol	15.874	330	1402951	137.765	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	5629702	85.129	ng	-0.01
79) Terphenyl-d14	19.756	244	7704026	82.790	ng	-0.01

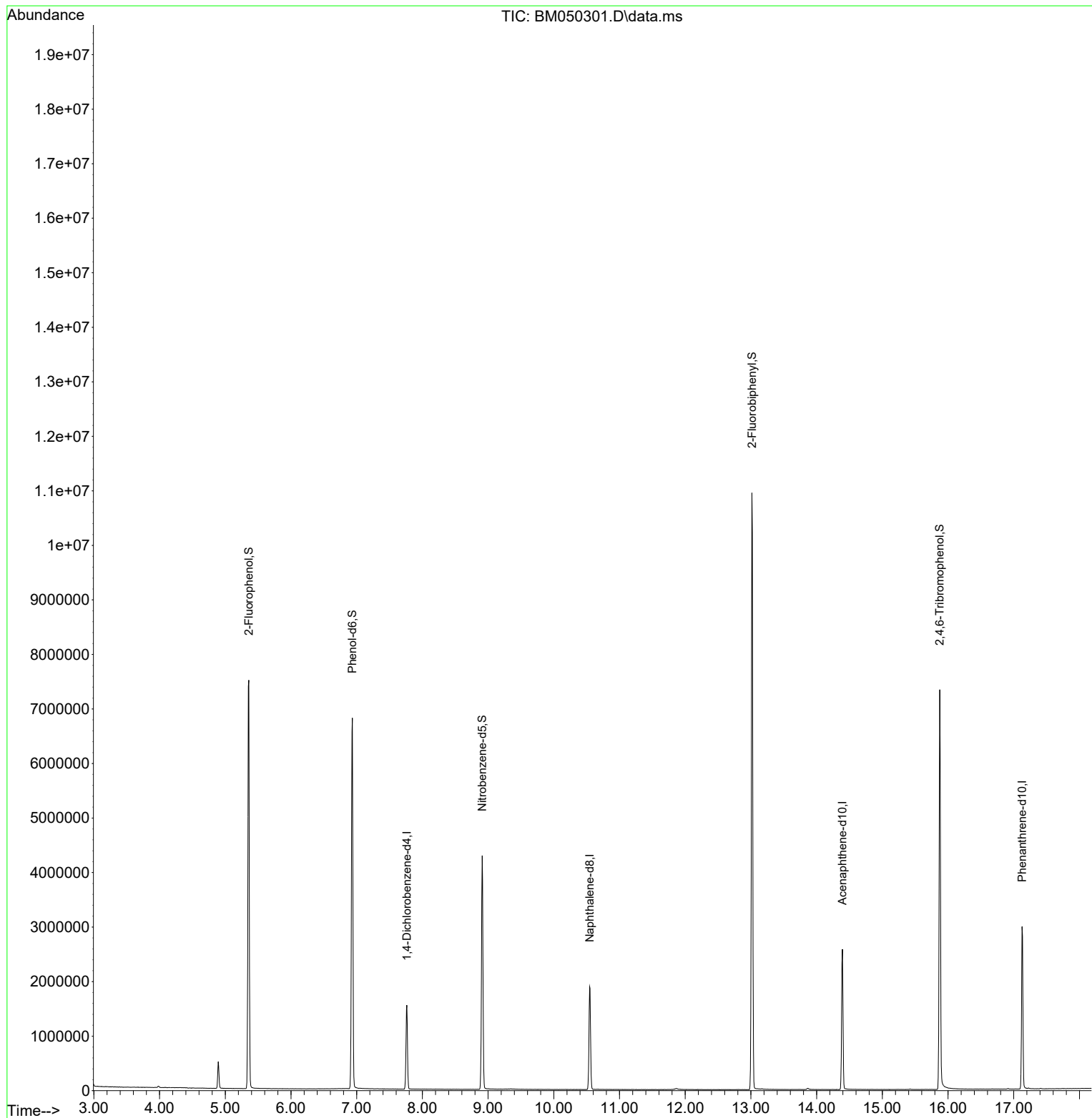
Target Compounds	Qvalue

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

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