

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060925\
 Data File : BM050246.D
 Acq On : 09 Jun 2025 17:43
 Operator : RC/JU
 Sample : PB168311TB
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168311TB

Quant Time: Jun 10 01:25:35 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	7.769	152	348706	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1311818	20.000	ng	0.00
39) Acenaphthene-d10	14.398	164	755301	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1467647	20.000	ng	0.00
76) Chrysene-d12	21.368	240	1523440	20.000	ng	0.00
86) Perylene-d12	24.344	264	1647706	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2951685	141.203	ng	0.00
7) Phenol-d6	6.934	99	3637251	132.177	ng	0.00
23) Nitrobenzene-d5	8.916	82	2171348	86.085	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1216044	141.548	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	4669156	83.693	ng	0.00
79) Terphenyl-d14	19.768	244	6569575	81.720	ng	0.00

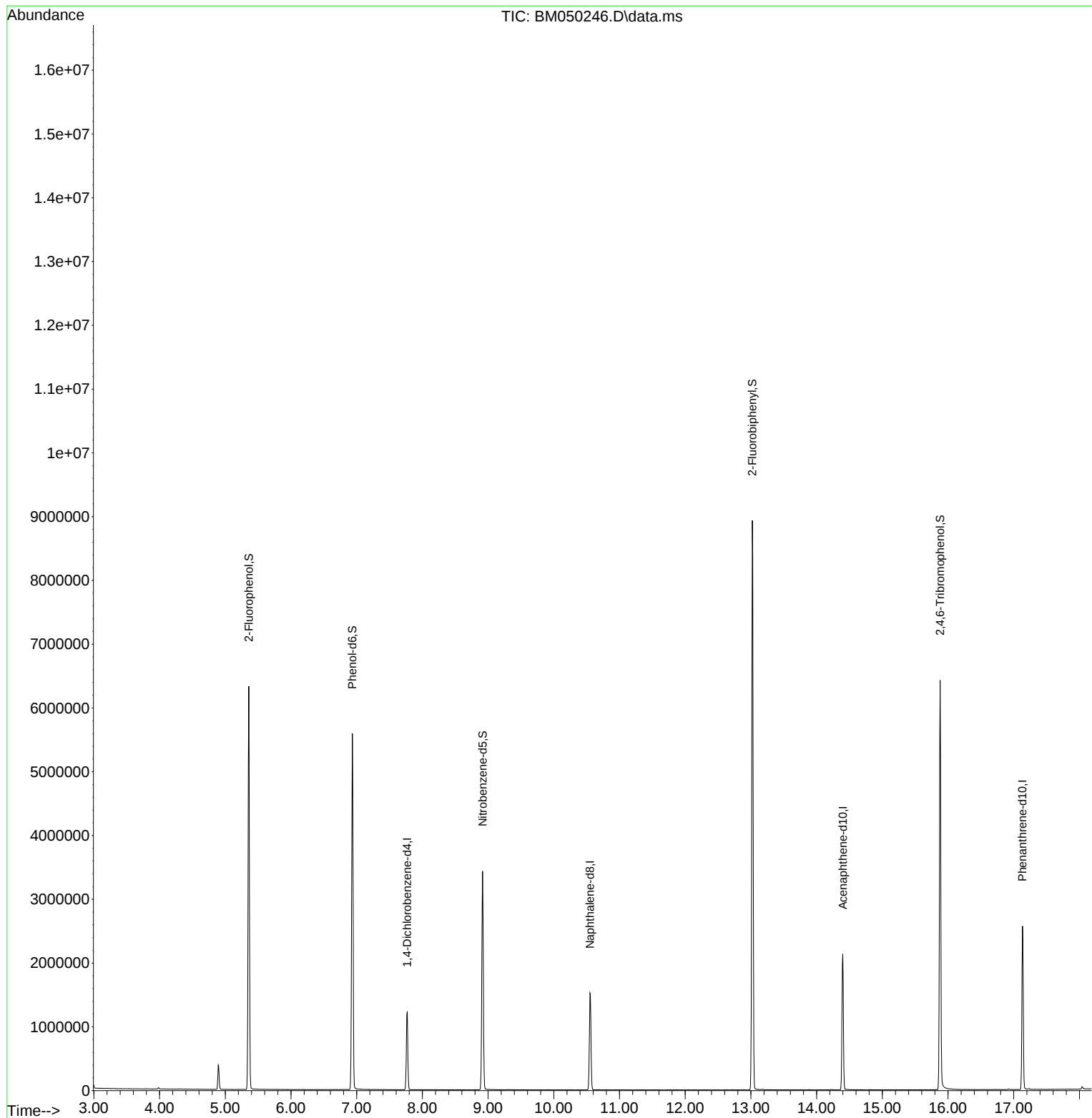
Target Compounds	Qvalue

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

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