

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061125\
 Data File : BM050278.D
 Acq On : 11 Jun 2025 17:24
 Operator : RC/JU
 Sample : PB168422BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168422BL

Quant Time: Jun 11 18:24:46 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	379949	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1397818	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	753306	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1427852	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1514576	20.000	ng	-0.01
86) Perylene-d12	24.338	264	1590017	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	3541836	155.502	ng	0.00
7) Phenol-d6	6.934	99	4283310	142.856	ng	0.00
23) Nitrobenzene-d5	8.916	82	2535242	94.328	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1329100	155.118	ng	0.00
45) 2-Fluorobiphenyl	13.021	172	5339966	95.971	ng	0.00
79) Terphenyl-d14	19.762	244	7459640	93.334	ng	0.00

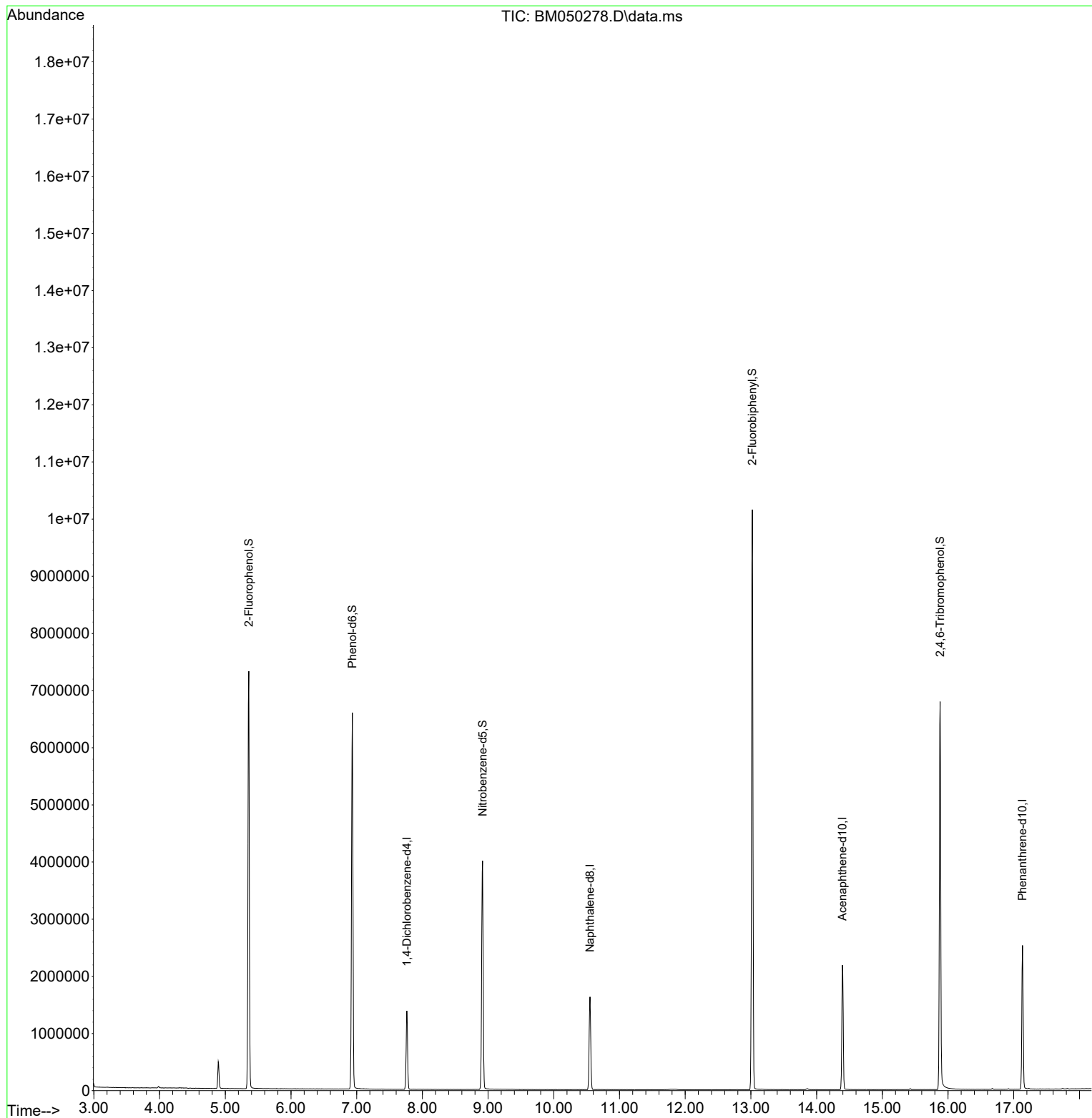
Target Compounds	Qvalue

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

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