

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061325\
 Data File : BM050317.D
 Acq On : 14 Jun 2025 03:00
 Operator : RC/JU
 Sample : PB168428TB
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168428TB

Quant Time: Jun 14 04:28:02 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	427909	20.000	ng	0.00
21) Naphthalene-d8	10.545	136	1575829	20.000	ng	-0.01
39) Acenaphthene-d10	14.392	164	855467	20.000	ng	0.00
64) Phenanthrene-d10	17.127	188	1557192	20.000	ng	-0.01
76) Chrysene-d12	21.356	240	1581816	20.000	ng	-0.02
86) Perylene-d12	24.321	264	1677028	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	3698811	144.193	ng	0.00
7) Phenol-d6	6.934	99	4492289	133.033	ng	0.00
23) Nitrobenzene-d5	8.910	82	2676296	88.328	ng	0.00
42) 2,4,6-Tribromophenol	15.874	330	1383440	142.178	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	5595242	88.550	ng	-0.01
79) Terphenyl-d14	19.757	244	7293777	87.380	ng	-0.01

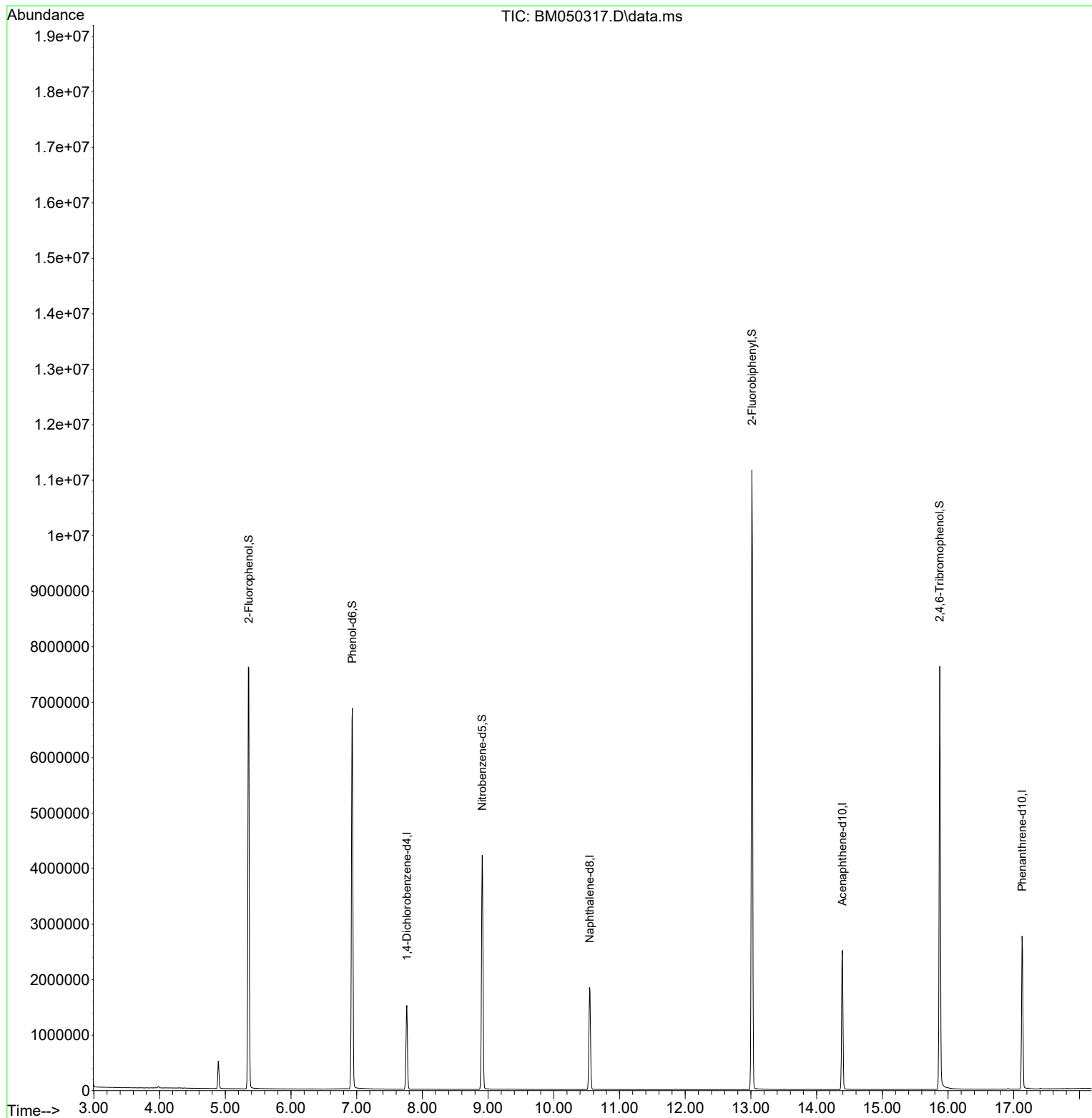
Target Compounds	Qvalue

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

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