

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050226.D
 Acq On : 06 Jun 2025 23:22
 Operator : RC/JU
 Sample : PB168271TB
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168271TB

Quant Time: Jun 07 01:22:04 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	7.787	152	358609	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1359938	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	776933	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1434193	20.000	ng	0.00
76) Chrysene-d12	21.374	240	1254575	20.000	ng	-0.01
86) Perylene-d12	24.362	264	1261316	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2842427	132.221	ng	0.00
7) Phenol-d6	6.951	99	3513251	124.146	ng	0.00
23) Nitrobenzene-d5	8.934	82	2064741	78.962	ng	-0.01
42) 2,4,6-Tribromophenol	15.892	330	1128266	127.674	ng	-0.01
45) 2-Fluorobiphenyl	13.039	172	4591404	80.008	ng	-0.01
79) Terphenyl-d14	19.774	244	5446768	82.273	ng	-0.01

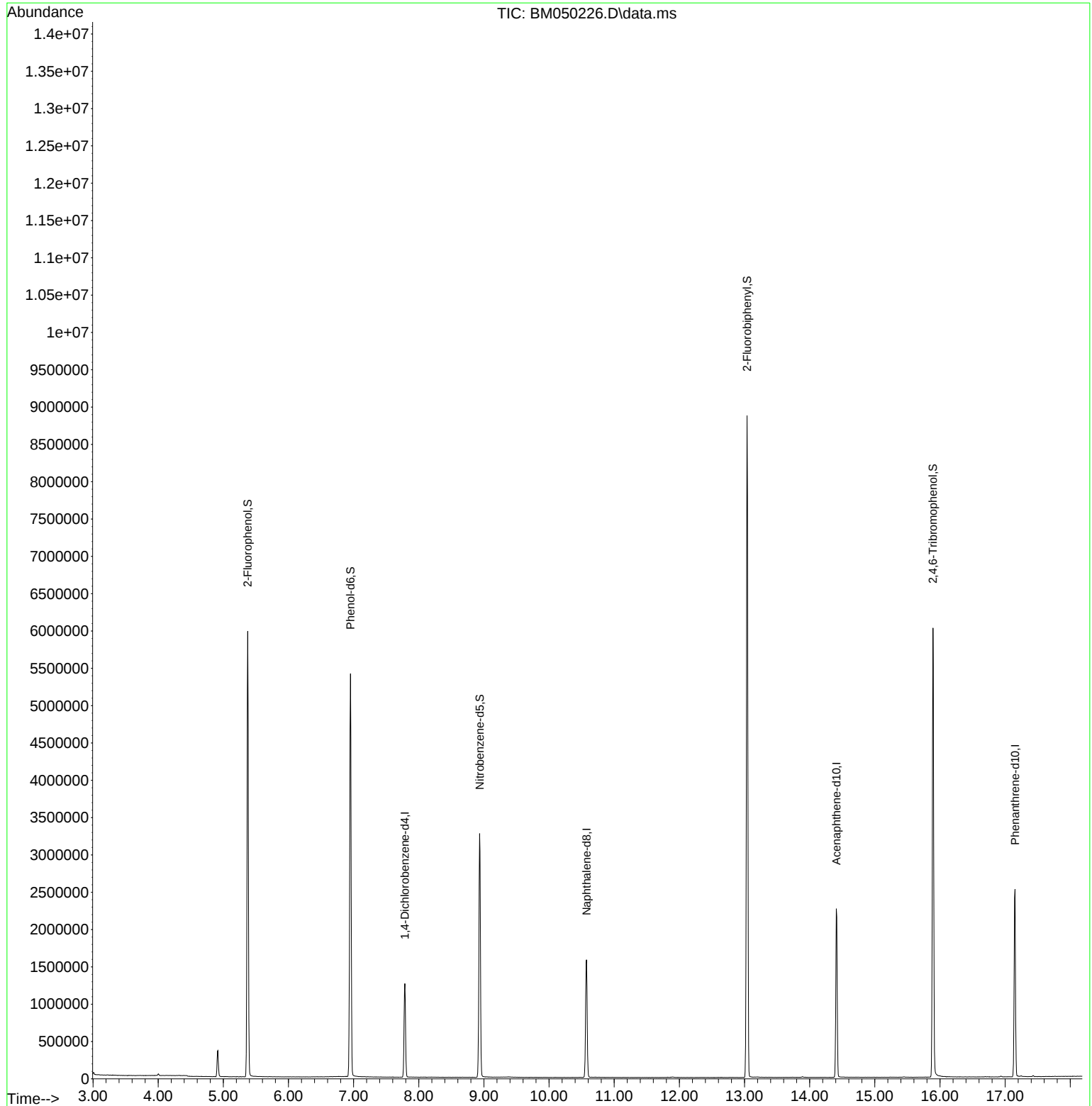
Target Compounds	Qvalue

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

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