

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061325\
 Data File : BM050310.D
 Acq On : 13 Jun 2025 21:44
 Operator : RC/JU
 Sample : Q2296-16
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-4

Quant Time: Jun 13 23:21:03 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	380880	20.000	ng	0.00
21) Naphthalene-d8	10.546	136	1422513	20.000	ng	#-0.01
39) Acenaphthene-d10	14.392	164	742819	20.000	ng	0.00
64) Phenanthrene-d10	17.128	188	1363568	20.000	ng	-0.01
76) Chrysene-d12	21.357	240	1452631	20.000	ng	-0.02
86) Perylene-d12	24.327	264	1211865	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	3353808	146.887	ng	0.00
7) Phenol-d6	6.928	99	3827381	127.338	ng	0.00
23) Nitrobenzene-d5	8.910	82	2674521	97.783	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	1238230	146.553	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	5038782	91.836	ng	-0.01
79) Terphenyl-d14	19.757	244	6645330	86.691	ng	-0.01

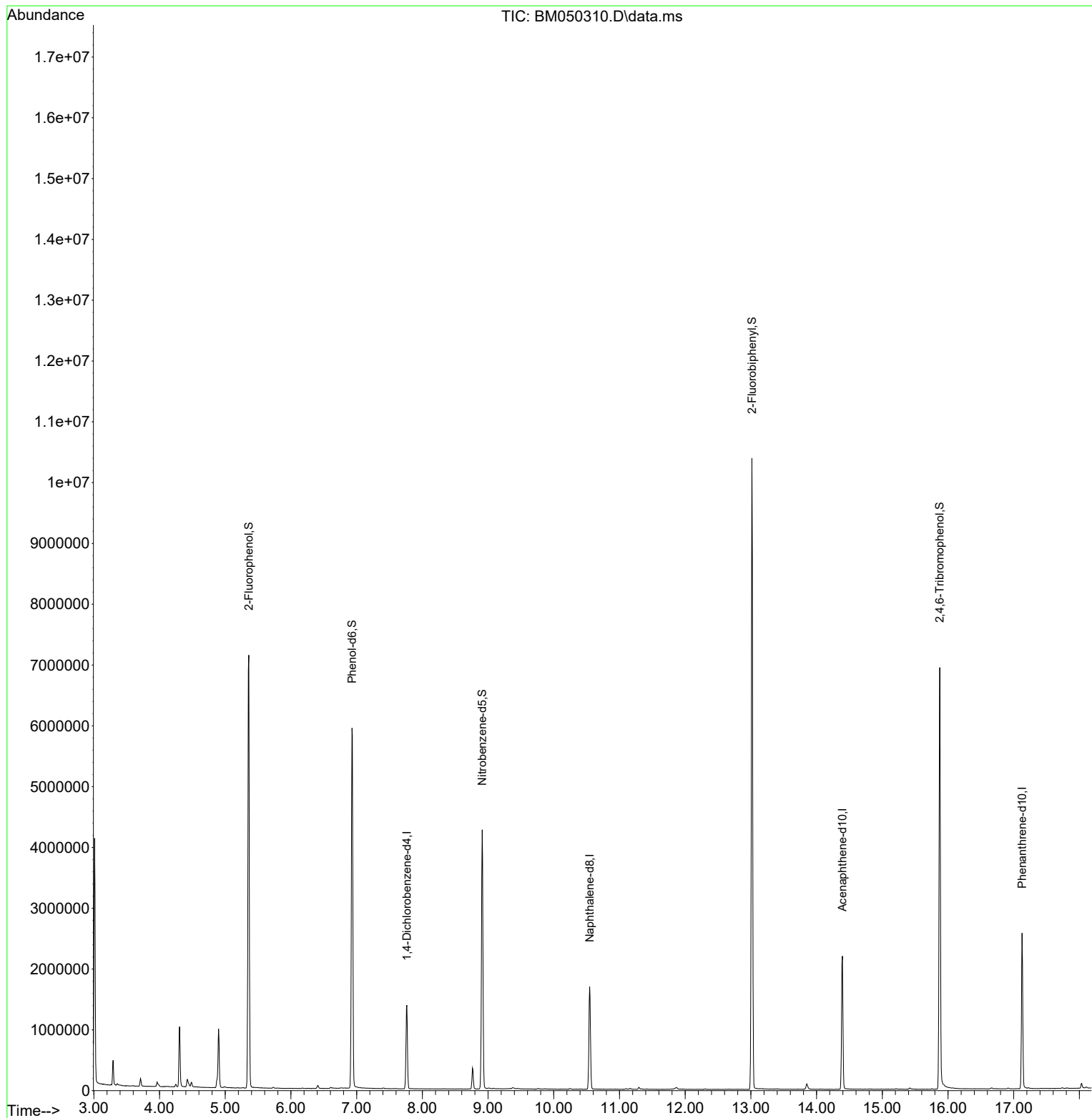
Target Compounds	Qvalue

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

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