

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061125\
 Data File : BM050284.D
 Acq On : 11 Jun 2025 21:21
 Operator : RC/JU
 Sample : Q2267-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-20250605

Quant Time: Jun 12 02:56:54 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)

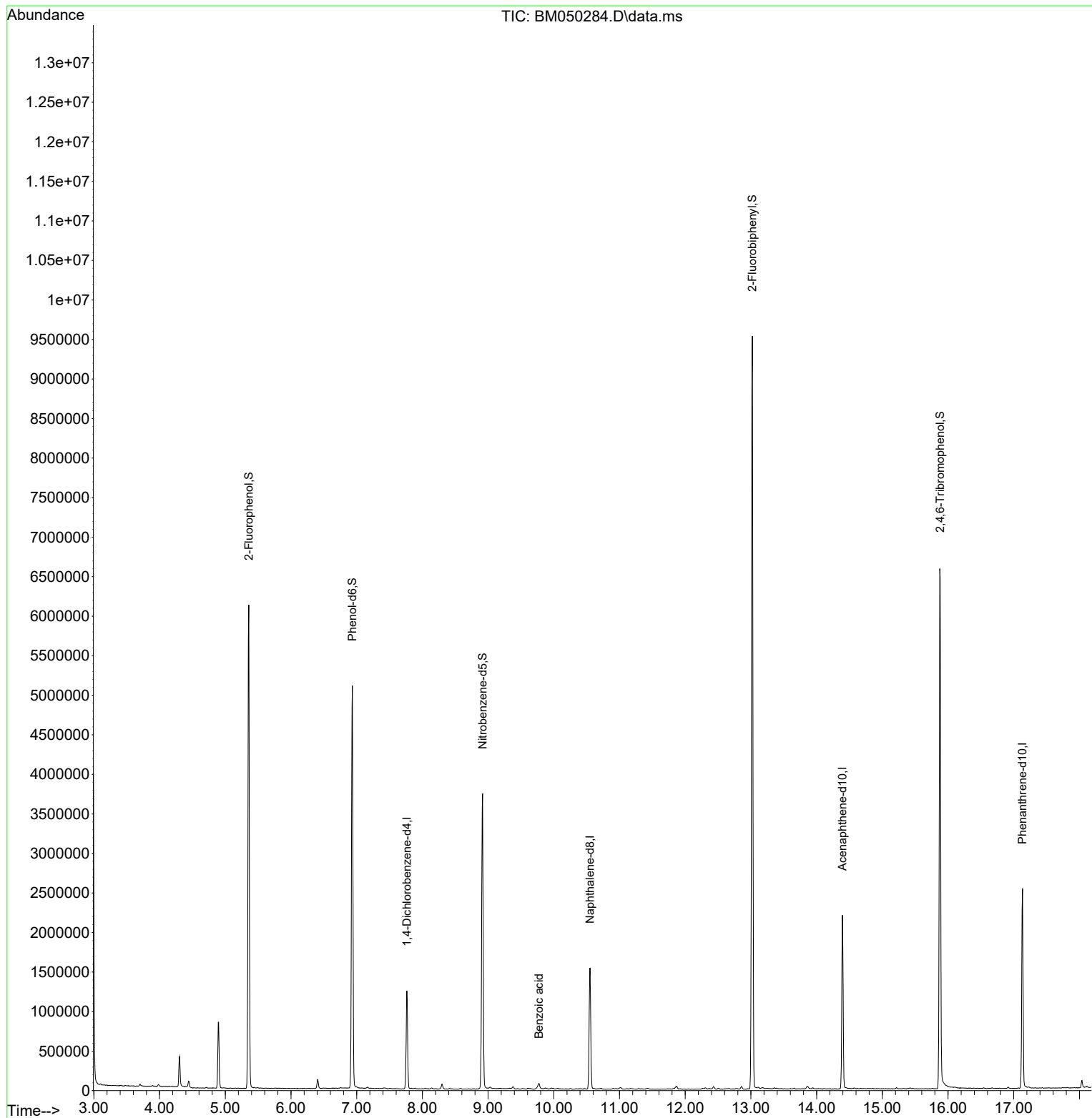
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	342435	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1297099	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	723910	20.000	ng	0.00
64) Phenanthrene-d10	17.133	188	1414148	20.000	ng	0.00
76) Chrysene-d12	21.356	240	1483946	20.000	ng	-0.02
86) Perylene-d12	24.333	264	828685	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2893731	140.966	ng	0.00
7) Phenol-d6	6.934	99	3268812	120.964	ng	0.00
23) Nitrobenzene-d5	8.916	82	2361569	94.689	ng	0.00
42) 2,4,6-Tribromophenol	15.880	330	1285779	156.156	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	4908961	91.807	ng	0.00
79) Terphenyl-d14	19.762	244	5362171	68.476	ng	0.00
Target Compounds						
32) Benzoic acid	9.775	122	35952	2.844	ng	Qvalue 94

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

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