

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050437.D
 Acq On : 14 Jul 2025 22:14
 Operator : RC/JU
 Sample : Q2489-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G2(6-12)

Quant Time: Jul 15 02:04:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 08 18:32:25 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

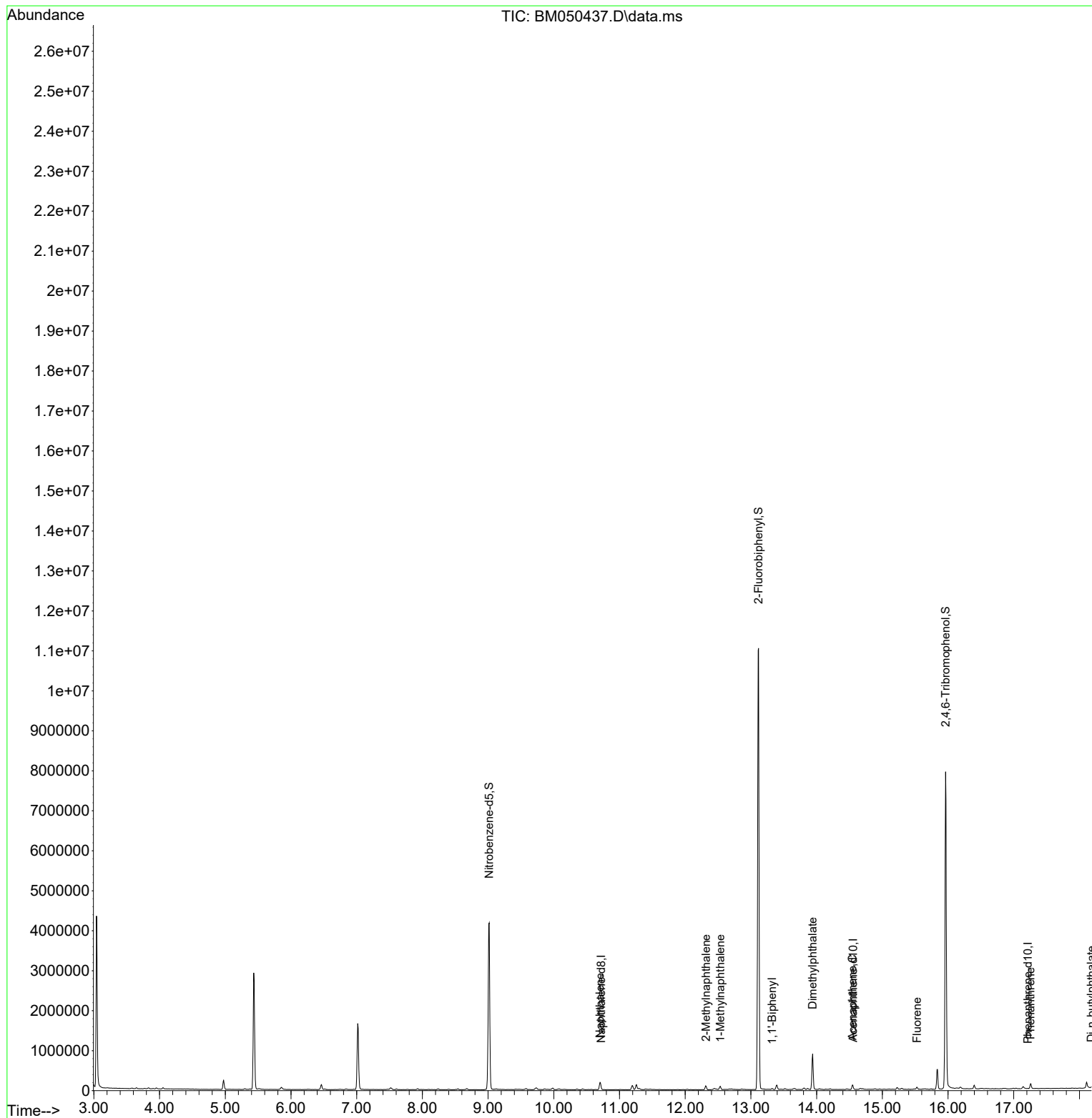
Internal Standards						
1) 1,4-Dichlorobenzene-d4	0.000	152	0	0.000	ng	-7.86
21) Naphthalene-d8	10.722	136	53	20.000	ng	# 0.06
39) Acenaphthene-d10	14.557	164	73	20.000	ng	# 0.06
64) Phenanthrene-d10	17.215	188	595	20.000	ng	#-0.01
76) Chrysene-d12	0.000	240	0m	20.000	ng	-21.46
86) Perylene-d12	24.444	264	69	20.000	ng	#-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1220295	0.000	ng	0.00
7) Phenol-d6	7.016	99	879374	0.000	ng	0.00
23) Nitrobenzene-d5	9.016	82	2318738	2232030.045	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	1451490	1636483.002	ng	-0.01
45) 2-Fluorobiphenyl	13.116	172	5822152	984926.479	ng	0.00
79) Terphenyl-d14	19.827	244	10281609	0.000	ng	-0.02
Target Compounds						
					Qvalue	
31) Naphthalene	10.704	128	148388	55121.405	ng	98
37) 2-Methylnaphthalene	12.316	142	44769	26338.511	ng	97
38) 1-Methylnaphthalene	12.533	142	35663	19825.377	ng	# 89
46) 1,1'-Biphenyl	13.322	154	9102	1680.539	ng	88
50) Dimethylphthalate	13.939	163	564587	112336.315	ng	99
52) Acenaphthene	14.545	154	39457	9509.771	ng	96
58) Fluorene	15.527	166	22346	4243.683	ng	94
71) Phenanthrene	17.257	178	74735	2219.733	ng	99
74) Di-n-butylphthalate	18.180	149	12453	375.194	ng	99
75) Fluoranthene	19.262	202	14900	407.088	ng	92

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

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