

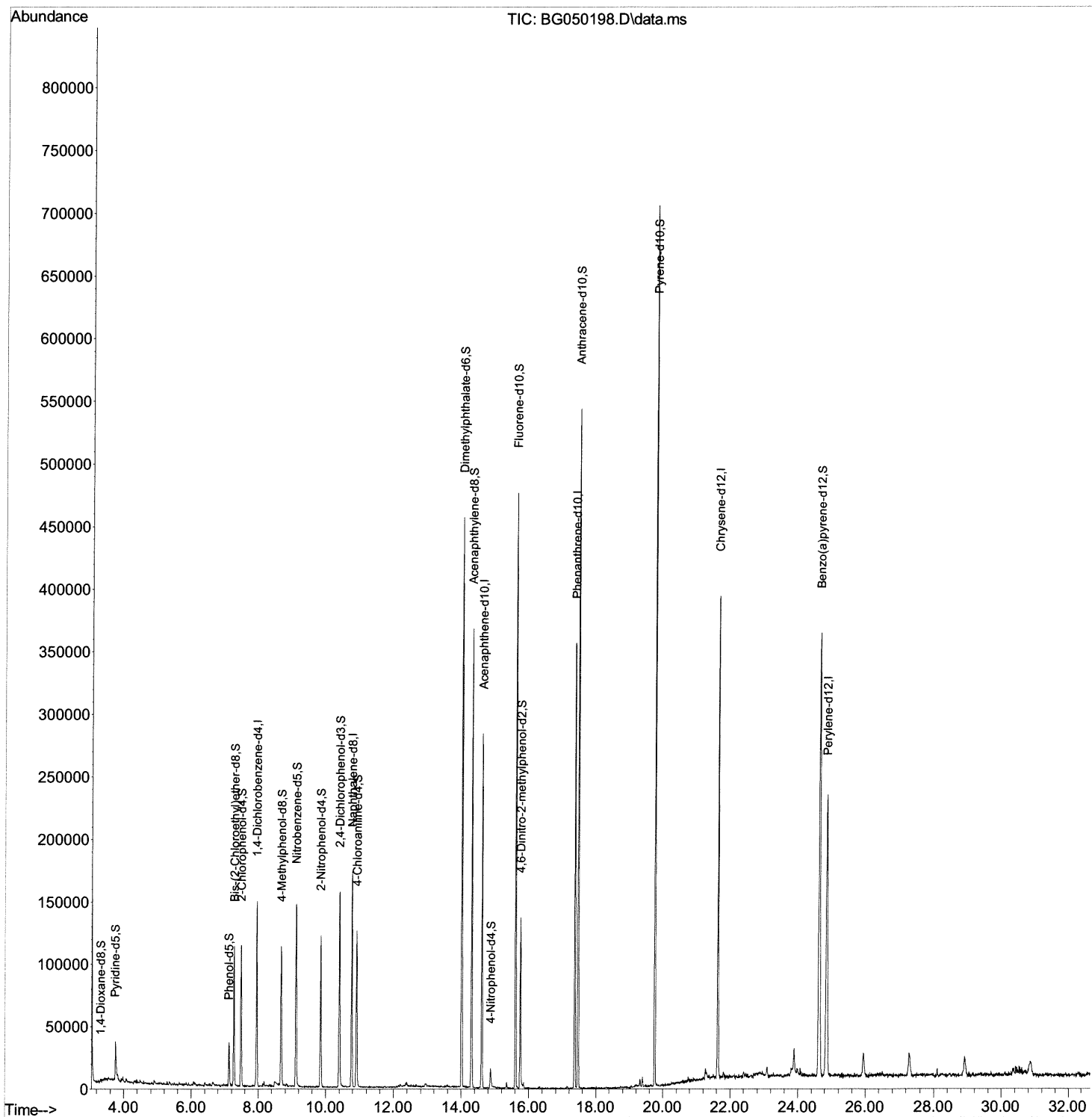
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050198.D
Acq On : 8 Sep 2021 13:50
Operator : CG/JU
Sample : M3566-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG7

Manual Integrations
APPROVED

mohammad
9/9/2021 8:41:13 AM

Quant Time: Sep 08 14:35:24 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

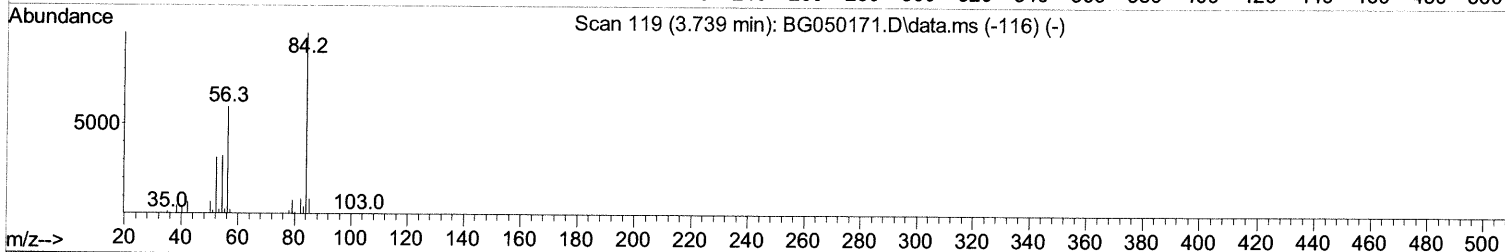
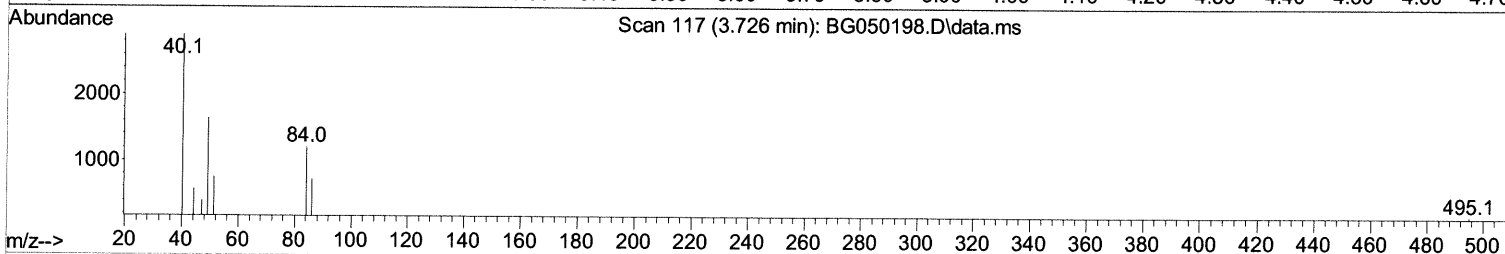
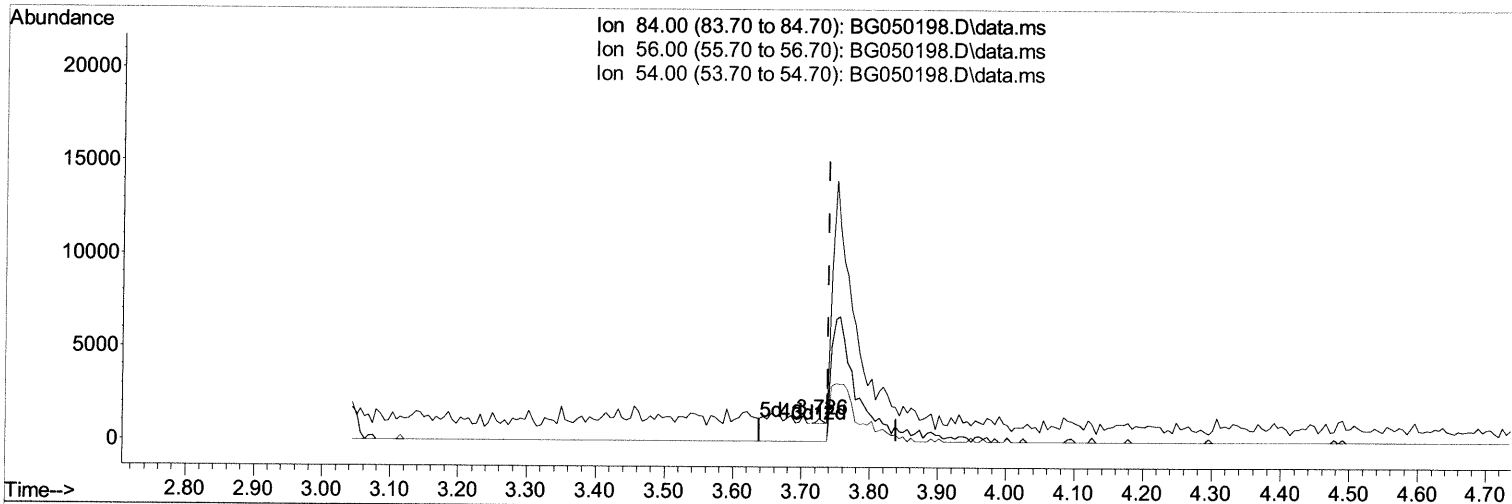
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TIC: BG050198.D\data.ms

(4) Pyridine-d5 (S)

3.726min (-0.013) 0.05 ng/ul

response 167

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	0.00#
54.00	37.00	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

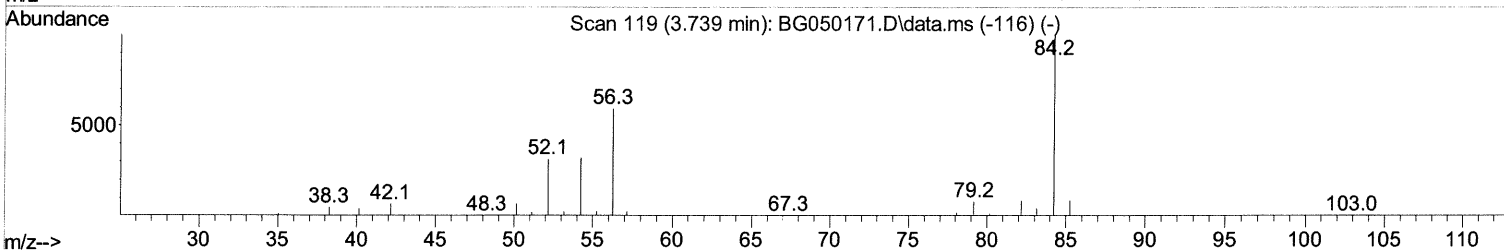
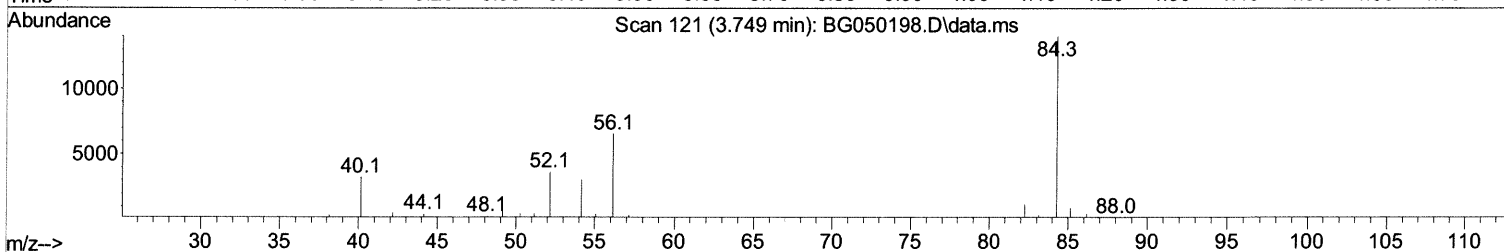
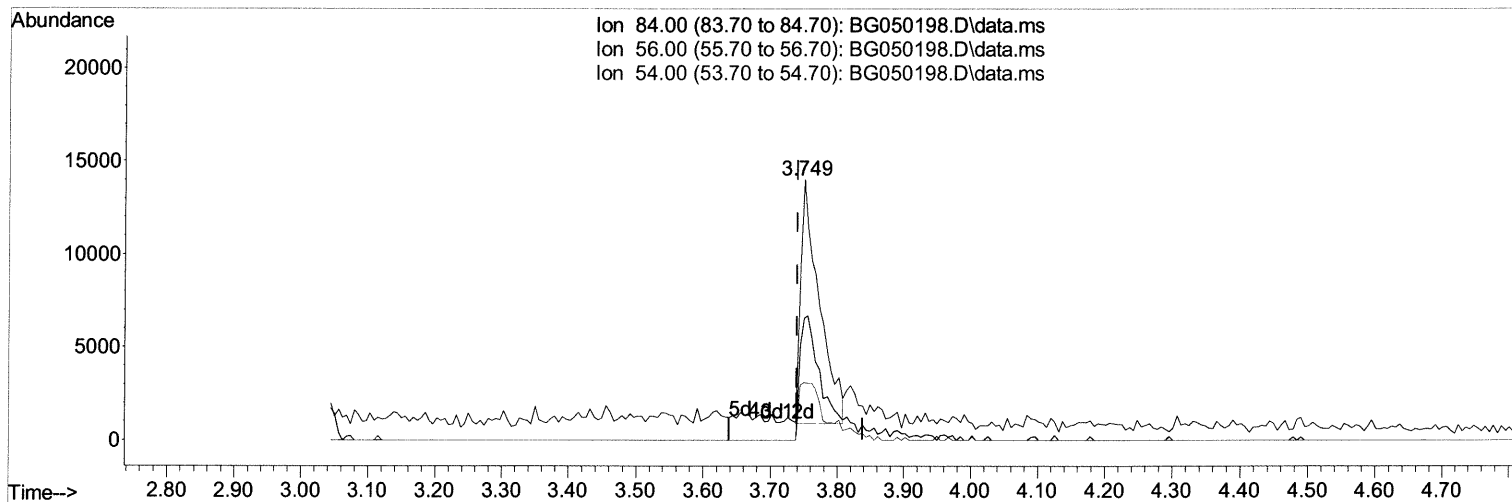
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TIC: BG050198.D\data.ms

(4) Pyridine-d5 (S)

3.749min (+ 0.011) 8.35 ng/ul m 09/10/21 JU

response 25765

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	46.82
54.00	37.00	21.95#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
 Data File : BG050198.D
 Acq On : 8 Sep 2021 13:50
 Operator : CG/JU
 Sample : M3566-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	42997	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	152241	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.611	164	101363	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.360	188	229391	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.637	240	245226	20.000	ng/ul	0.00
88) Perylene-d12	24.850	264	244731	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.309	96	2023	2.326	ng/uL	0.00
4) Pyridine-d5	3.749	84	25765m	8.346	ng/ul	> 0.01
7) Phenol-d5	7.121	99	22329	6.689	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.274	67	55110	28.266	ng/ul	0.00
11) 2-Chlorophenol-d4	7.485	132	58075	20.961	ng/ul	0.00
15) 4-Methylphenol-d8	8.672	113	46399	17.677	ng/ul	0.00
21) Nitrobenzene-d5	9.124	128	36255	34.695	ng/ul	0.00
24) 2-Nitrophenol-d4	9.847	143	37790	29.741	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	60175	24.813	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	71018	23.769	ng/ul	0.00
46) Dimethylphthalate-d6	14.012	166	317714	41.511	ng/ul	0.00
49) Acenaphthylene-d8	14.306	160	274590	30.130	ng/ul	0.00
54) 4-Nitrophenol-d4	14.864	143	4441	4.809	ng/ul	0.01
60) Fluorene-d10	15.604	176	194261	30.568	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	34044	24.603	ng/ul	0.00
73) Anthracene-d10	17.460	188	323958	31.459	ng/ul	0.00
81) Pyrene-d10	19.757	212	399386	32.697	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.627	264	392692	30.311	ng/ul	-0.01

Target Compounds Qvalue

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