

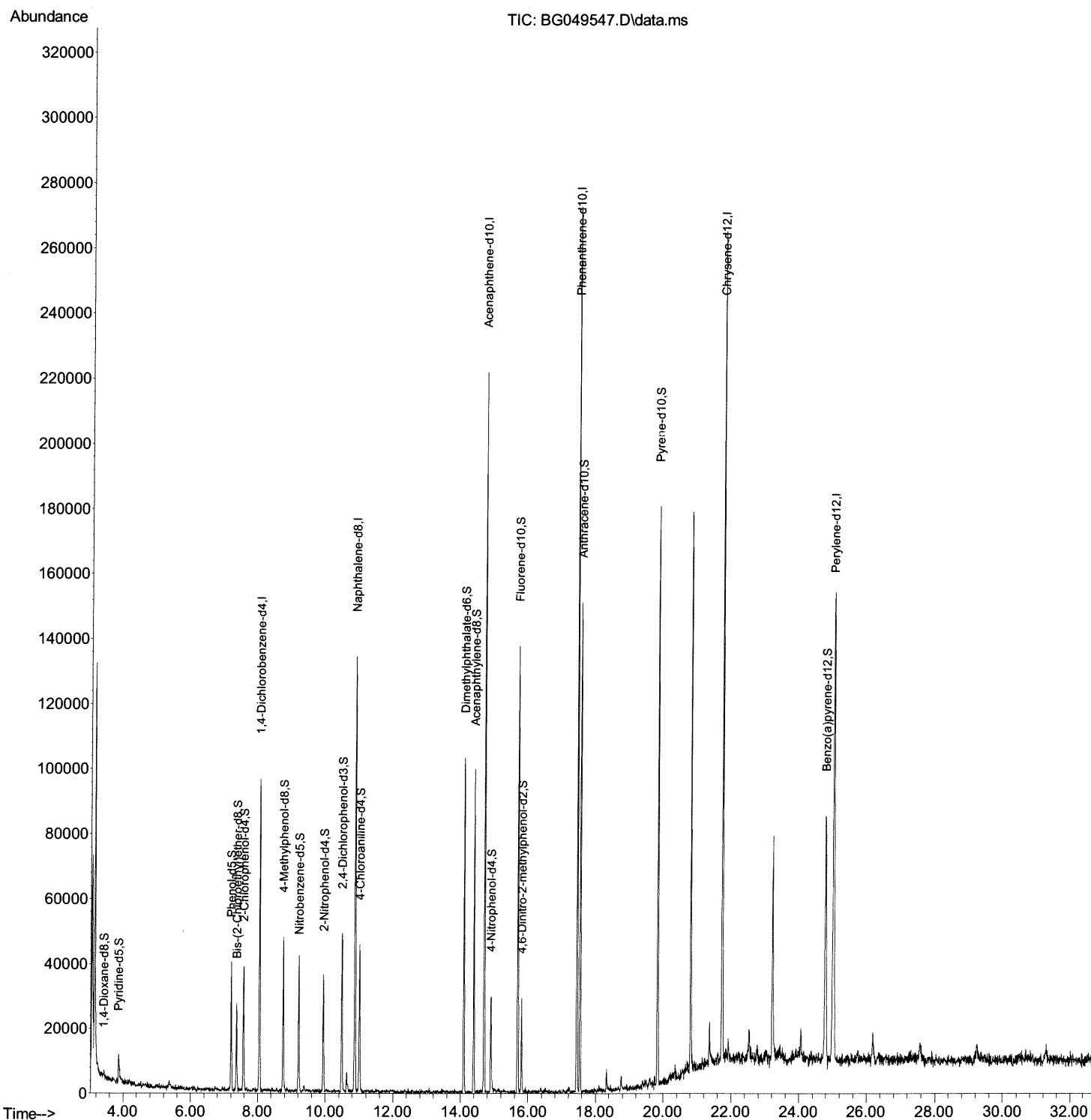
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072821\
Data File : BG049547.D
Acq On : 29 Jul 2021 00:37
Operator : CG/JU
Sample : M3106-07
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGCZ8

Manual Integrations
APPROVED

mohammad
7/29/2021 1:42:24 PM

Quant Time: Jul 29 04:43:38 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 29 04:40:56 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

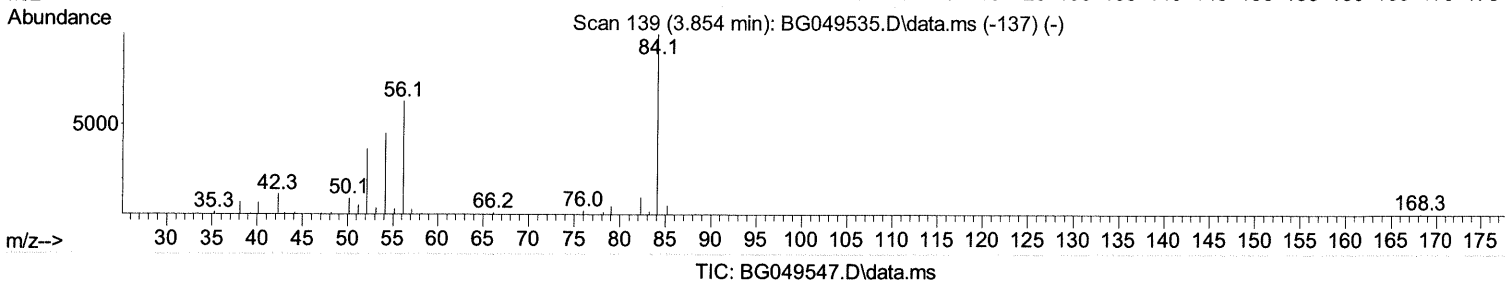
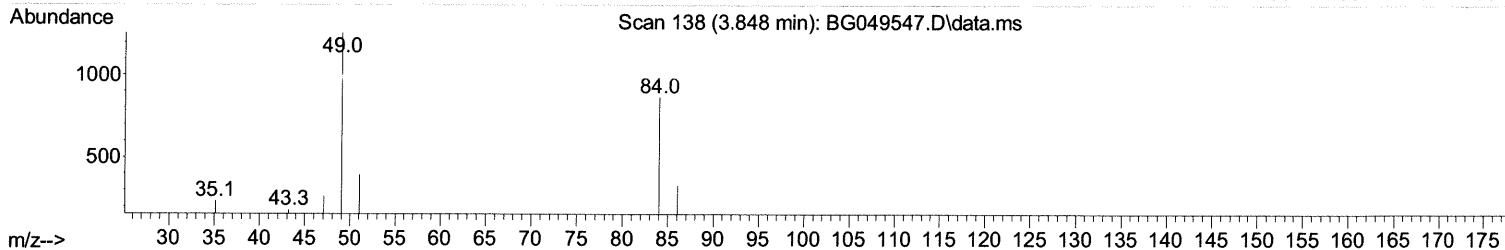
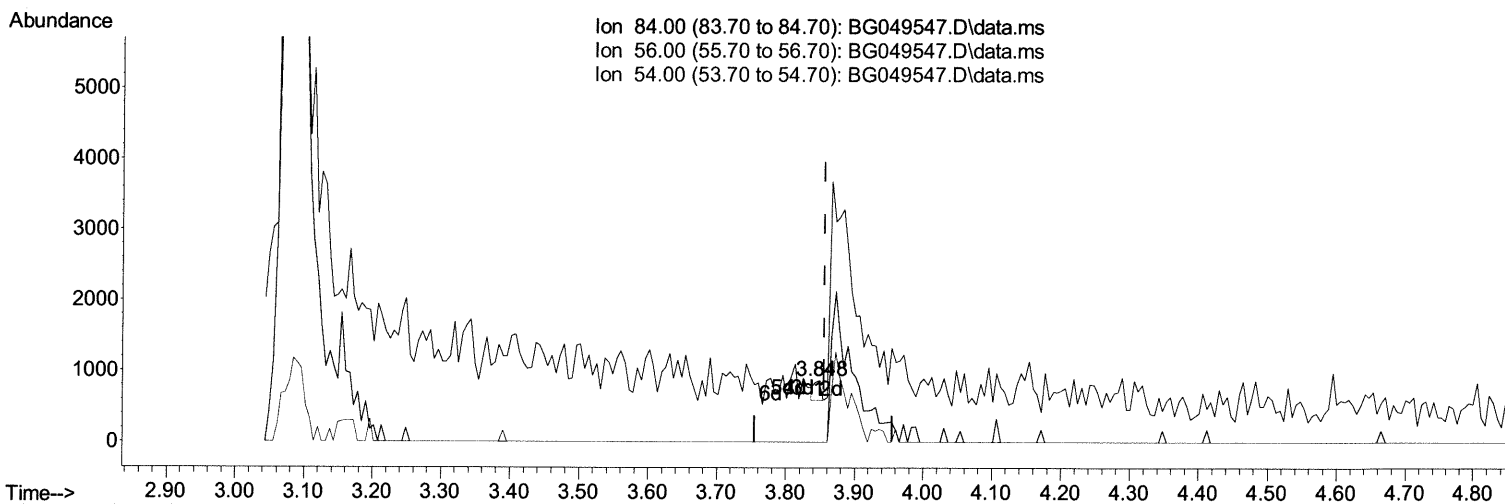
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(4) Pyridine-d5 (S)

3.848min (-0.007) 0.11 ng/ul

response 187

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	0.00#
54.00	34.10	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

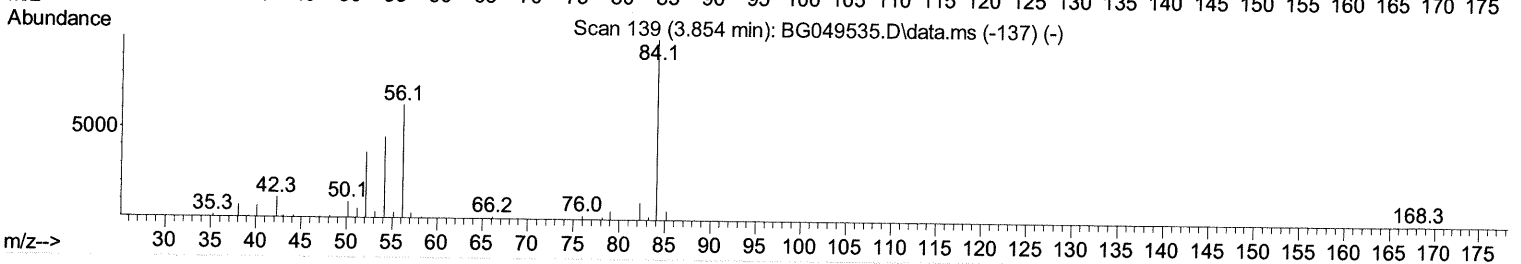
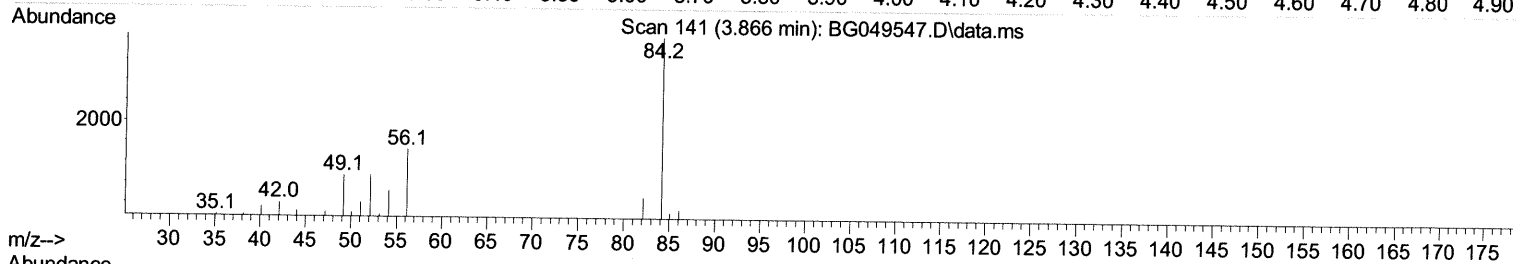
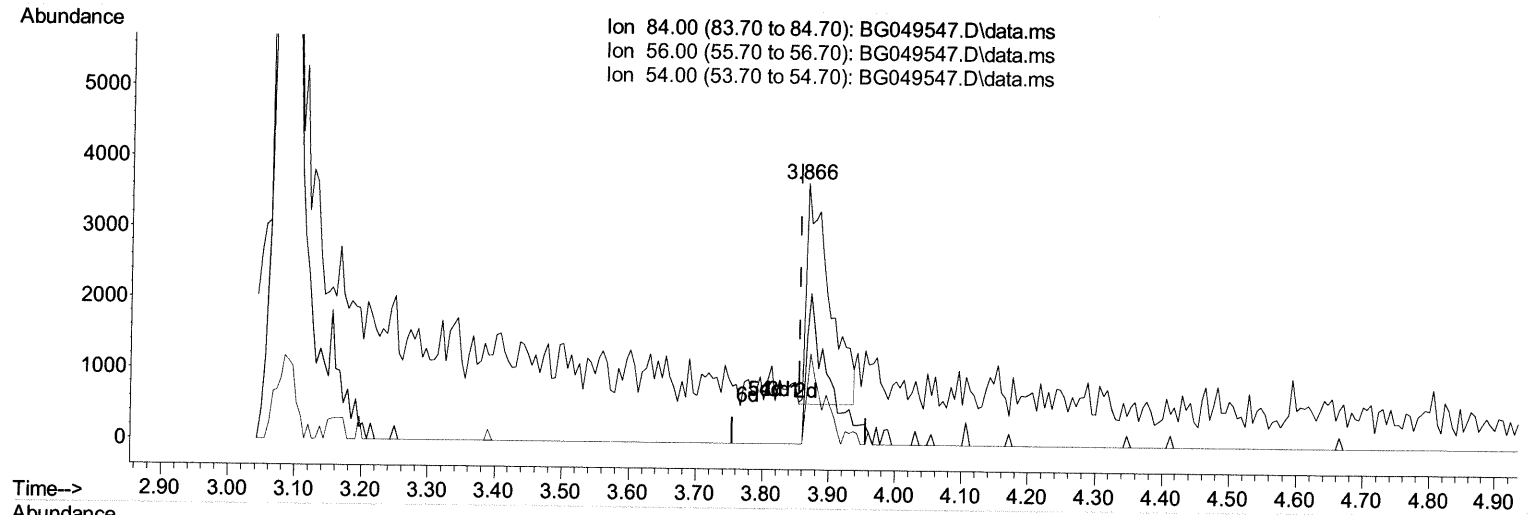
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TIC: BG049547.D\data.ms

(4) Pyridine-d5 (S)

3.866min (+ 0.011) 4.40 ng/ul m 07/30/21JU

response 7446

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	39.69#
54.00	34.10	18.20#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072821\
 Data File : BG049547.D
 Acq On : 29 Jul 2021 00:37
 Operator : CG/JU
 Sample : M3106-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGCZ8

Manual Integrations
 APPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.048	152	26946	20.000	ng/ul	0.00
20) Naphthalene-d8	10.868	136	107028	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.698	164	77601	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.447	188	165663	20.000	ng/ul	0.00
79) Chrysene-d12	21.730	240	152692	20.000	ng/ul	0.00
88) Perylene-d12	25.002	264	157487	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.431	96	779	1.323	ng/uL	0.00
4) Pyridine-d5	3.866	84	7446m	4.401	ng/ul	0.01
7) Phenol-d5	7.202	99	21155	8.874	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.367	67	12420	9.198	ng/ul	0.00
11) 2-Chlorophenol-d4	7.578	132	17045	9.974	ng/ul	0.00
15) 4-Methylphenol-d8	8.753	113	17663	9.829	ng/ul	0.00
21) Nitrobenzene-d5	9.211	128	7838	9.525	ng/ul	0.00
24) 2-Nitrophenol-d4	9.946	143	10705	11.102	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.486	165	17811	10.156	ng/ul	0.00
31) 4-Chloroaniline-d4	11.003	131	23466	9.601	ng/ul	0.00
46) Dimethylphthalate-d6	14.093	166	63582	10.696	ng/ul	0.00
49) Acenaphthylene-d8	14.393	160	74835	10.839	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	7204	7.359	ng/ul	0.01
60) Fluorene-d10	15.691	176	55154	10.770	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.808	200	6411	6.245	ng/ul	0.00
73) Anthracene-d10	17.547	188	89952	11.683	ng/ul	0.00
81) Pyrene-d10	19.832	212	99298	12.718	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.773	264	80997	9.785	ng/ul	0.00

Target Compounds Qvalue

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