

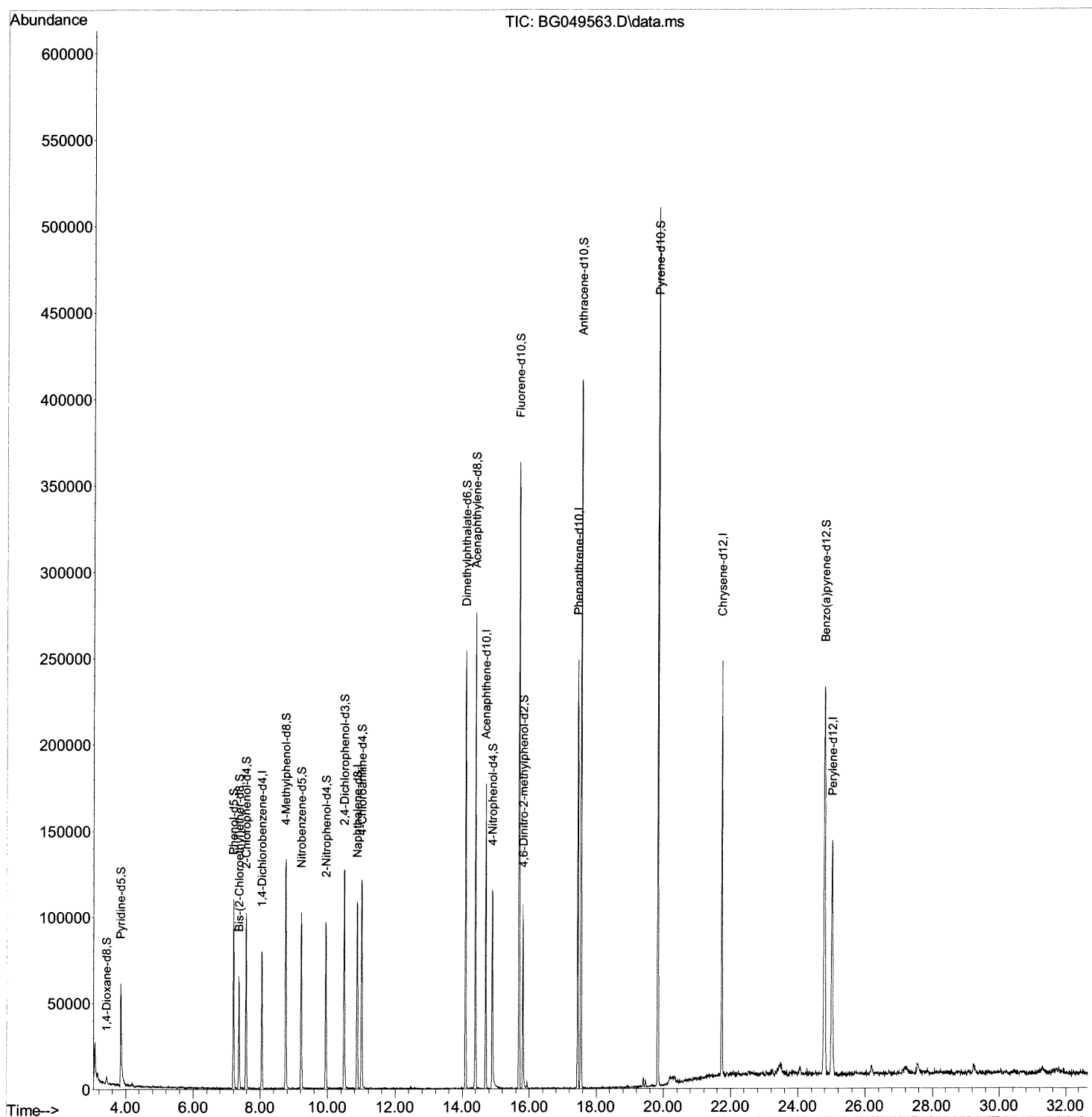
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
Data File : BG049563.D
Acq On : 29 Jul 2021 17:20
Operator : CG/JU
Sample : PB137967BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK967

Manual Integrations
APPROVED

mohammad
7/30/2021 12:00:10 PM

Quant Time: Jul 29 17:55:41 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 26 11:36:47 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

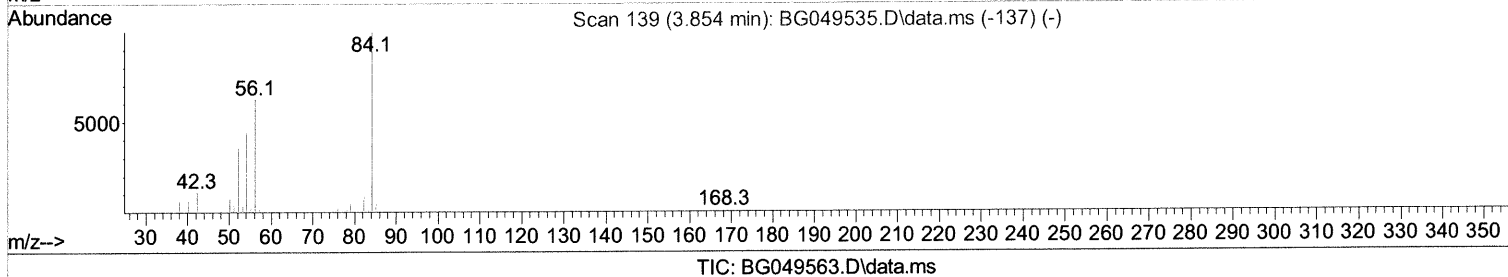
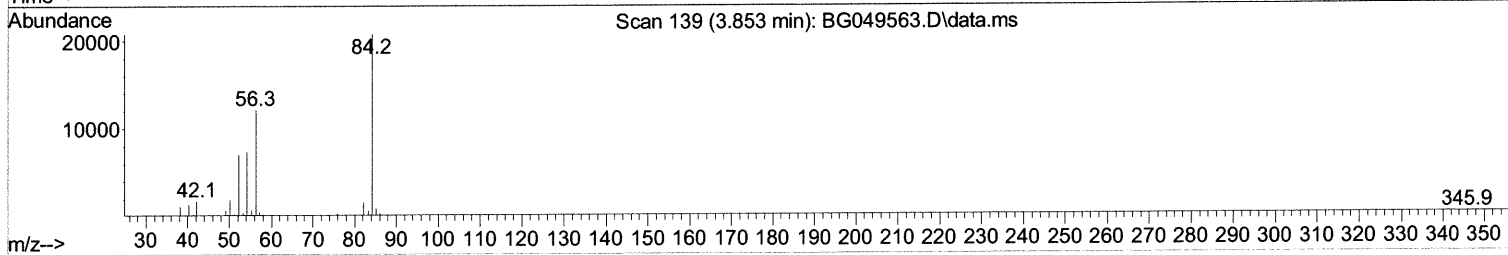
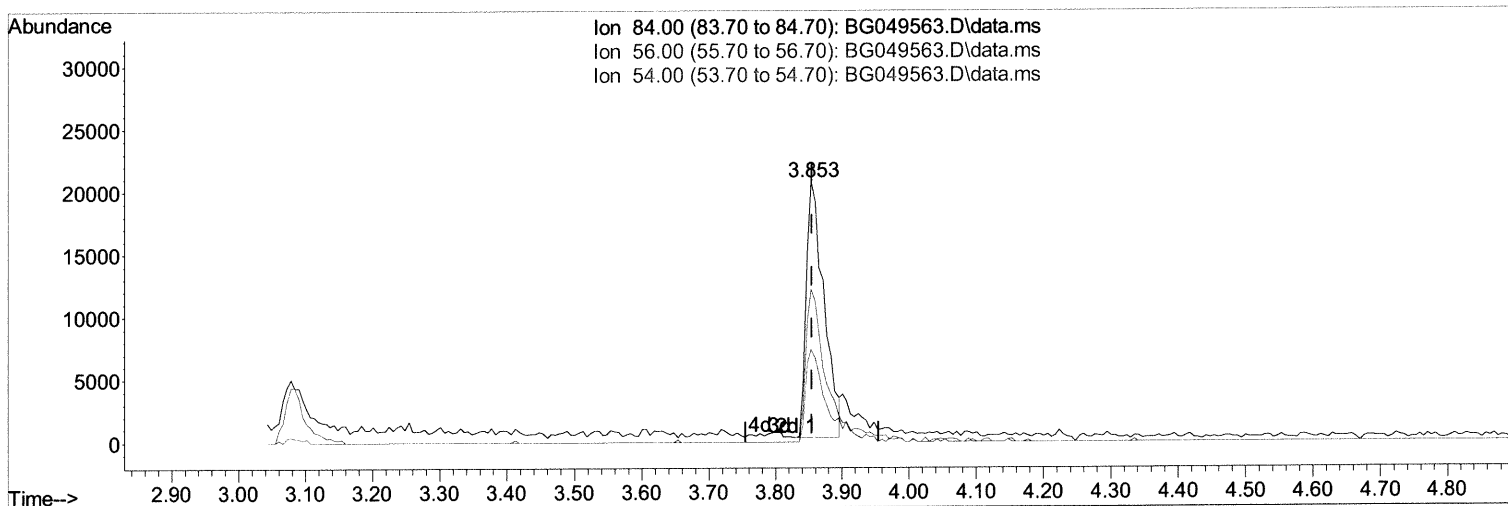
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(4) Pyridine-d5 (S)

3.853min (-0.002) 28.59 ng/ul

response 37553

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	58.57
54.00	34.10	35.41
0.00	0.00	0.00

Quantitation Report (Qedit)

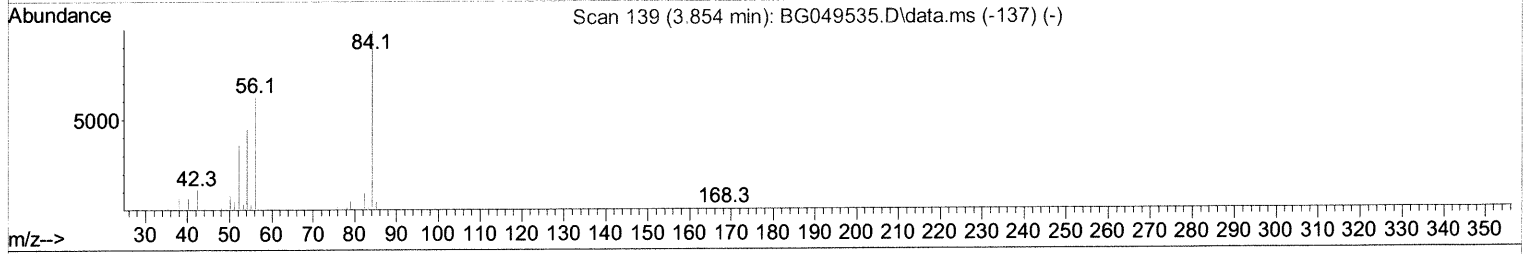
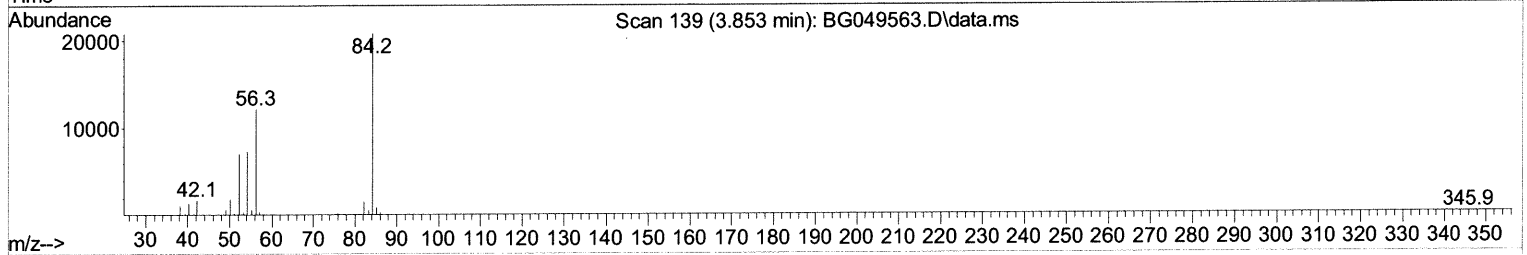
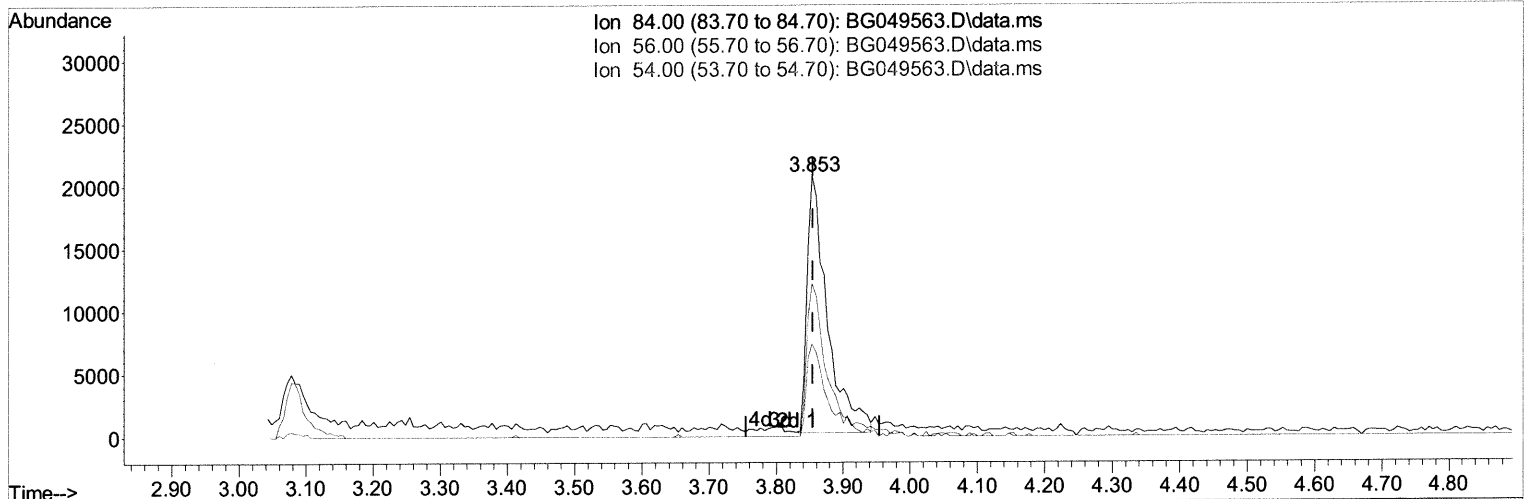
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 Response via : Initial Calibration



TIC: BG049563.D\data.ms

(4) Pyridine-d5 (S)

3.853min (-0.002) 32.93 ng/ul m 08/02/21 JU

response 43244

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	58.50	58.57
54.00	34.10	35.41
0.00	0.00	0.00

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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	8.048	152	20920	20.000	ng/ul	0.00
20) Naphthalene-d8	10.868	136	89476	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.698	164	65372	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.447	188	153543	20.000	ng/ul	0.00
79) Chrysene-d12	21.729	240	149286	20.000	ng/ul	0.00
88) Perylene-d12	25.013	264	151878	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.425	96	2622	5.734	ng/ul	-0.01
4) Pyridine-d5	3.853	84	43244m	32.926	ng/ul	0.00
7) Phenol-d5	7.202	99	58052	31.364	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.372	67	32859	31.346	ng/ul	0.00
11) 2-Chlorophenol-d4	7.578	132	42666	32.158	ng/ul	0.00
15) 4-Methylphenol-d8	8.753	113	45734	32.781	ng/ul	0.00
21) Nitrobenzene-d5	9.217	128	22781	33.116	ng/ul	0.00
24) 2-Nitrophenol-d4	9.945	143	25612	31.772	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.492	165	44752	30.524	ng/ul	0.00
31) 4-Chloroaniline-d4	11.003	131	63205	30.933	ng/ul	0.00
46) Dimethylphthalate-d6	14.098	166	162726	32.496	ng/ul	0.00
49) Acenaphthylene-d8	14.392	160	197881	34.022	ng/ul	0.00
54) 4-Nitrophenol-d4	14.891	143	25455	30.868	ng/ul	0.00
60) Fluorene-d10	15.690	176	144733	33.549	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.814	200	25156	26.439	ng/ul	0.01
73) Anthracene-d10	17.547	188	246124	34.490	ng/ul	0.00
81) Pyrene-d10	19.838	212	277189	36.312	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.784	264	258006	32.321	ng/ul	0.02

Target Compounds Qvalue

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