

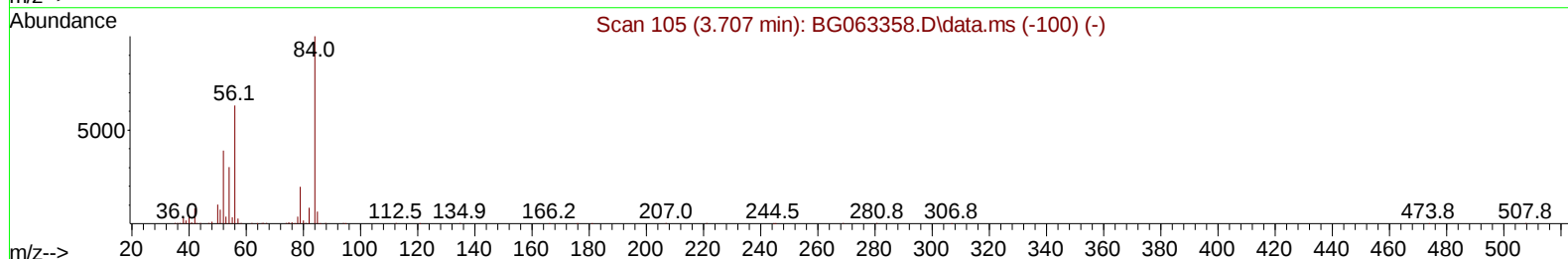
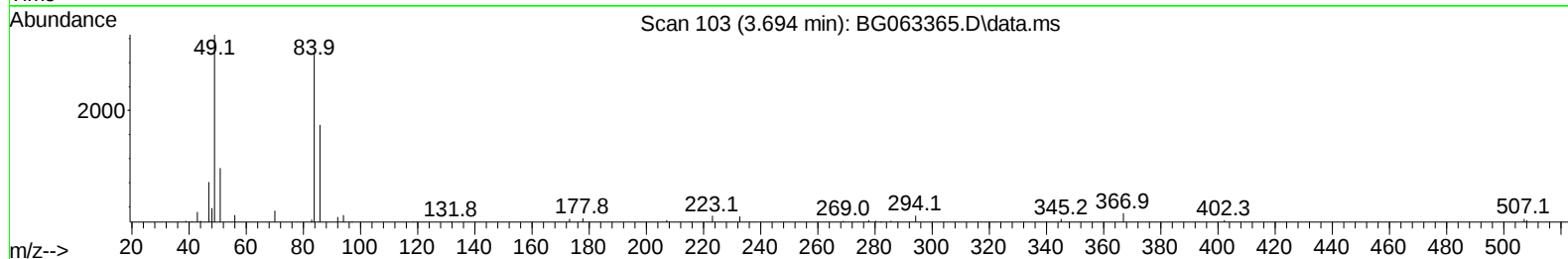
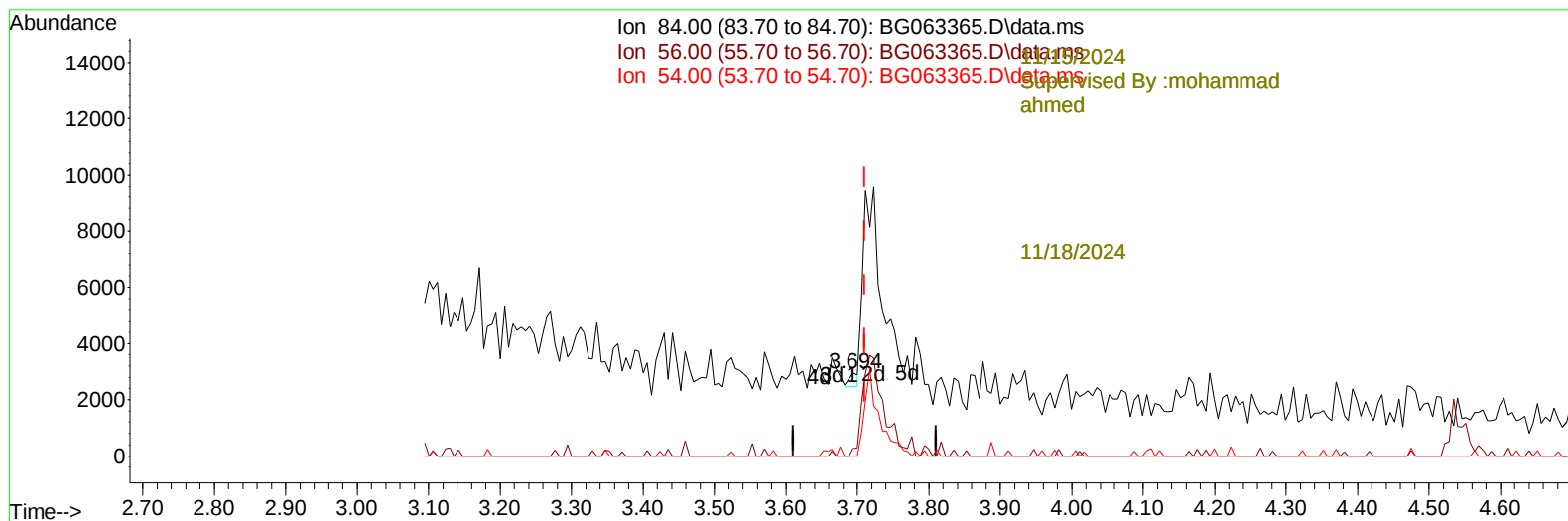
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063365.D
Acq On : 13 Nov 2024 21:52
Operator : RC/JU
Sample : P4801-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GCPB6

Manual IntegrationsAPPROVED

Quant Time: Nov 14 03:32:30 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 15:45:52 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/15/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BG063365.D\data.ms

(4) Pyridine-d5 (S)

3.694min (-0.017) 0.06 ng/u1

response 418

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.80	8.96#
54.00	35.10	0.00#
0.00	0.00	0.00

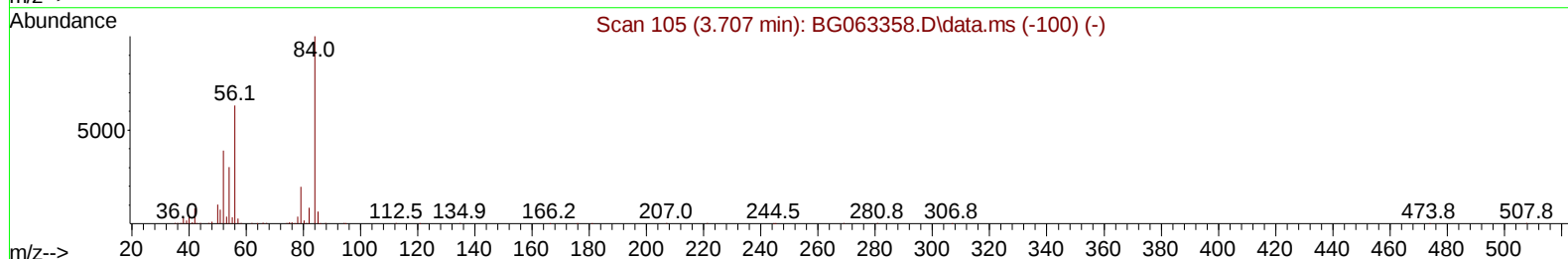
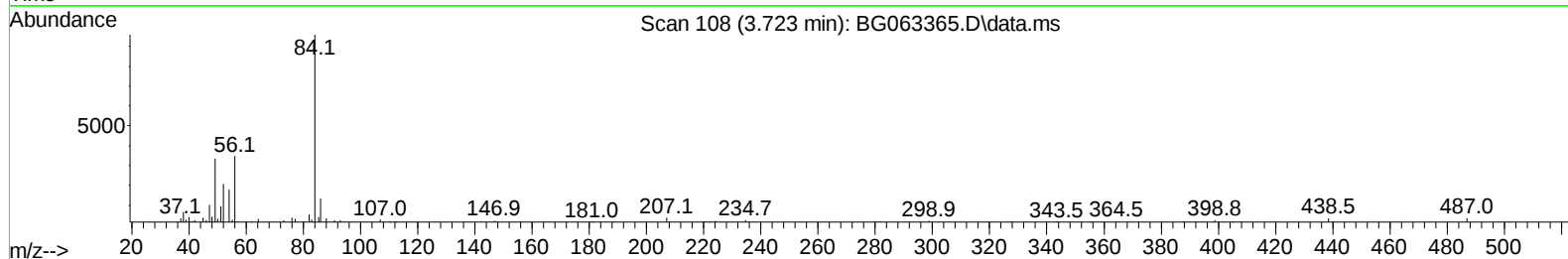
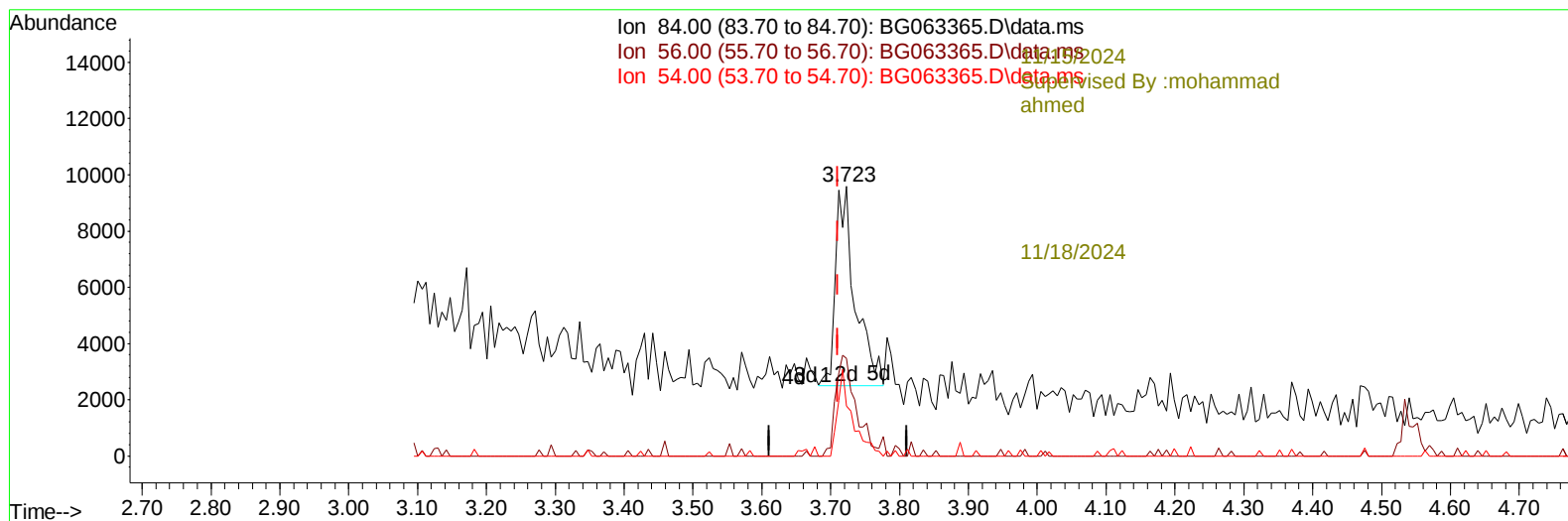
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(4) Pyridine-d5 (S)

3.723min (+ 0.013) 2.00 ng/u1 m

response 13821

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.80	36.18#
54.00	35.10	18.53#
0.00	0.00	0.00

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 Sample : P4801-17
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
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Quant Time: Nov 14 03:55:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 15:45:52 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/15/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	94753	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.621	136	391481	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.469	164	334759	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.219	188	831815	20.000	ng/ul	0.00
79) Chrysene-d12	21.473	240	886277	20.000	ng/ul	# 0.00
88) Perylene-d12	24.505	264	951834	20.000	ng/ul	-0.01
-----11/18/2024						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	9881	4.715	ng/uL	0.00
4) Pyridine-d5	3.723	84	13821m	1.996	ng/ul	0.01
7) Phenol-d5	7.002	99	103396	12.093	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.160	67	124436	24.812	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.366	132	136997	24.686	ng/ul	-0.01
15) 4-Methylphenol-d8	8.541	113	149718	20.659	ng/ul	0.00
21) Nitrobenzene-d5	8.982	128	75177	27.315	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.704	143	95124	26.742	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.251	165	178575	25.581	ng/ul	0.00
31) 4-Chloroaniline-d4	10.756	131	34413	3.785	ng/ul	0.00
46) Dimethylphthalate-d6	13.876	166	710248	27.086	ng/ul	0.00
49) Acenaphthylene-d8	14.158	160	645079	25.134	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	34041	9.482	ng/ul	0.00
60) Fluorene-d10	15.462	176	657788	29.723	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.592	200	160637	30.728	ng/ul	0.00
73) Anthracene-d10	17.319	188	1227570	34.337	ng/ul	0.00
81) Pyrene-d10	19.616	212	1637097	36.842	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.299	264	1560932	33.966	ng/ul	0.00

Target Compounds Qvalue

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

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