

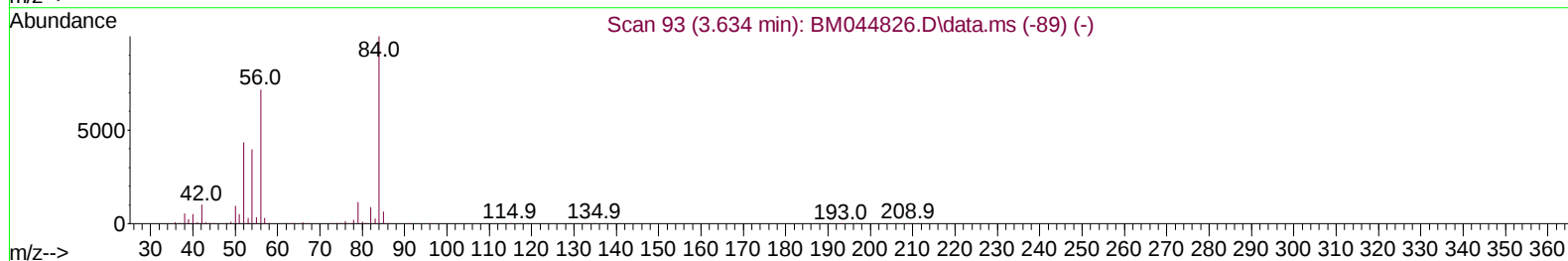
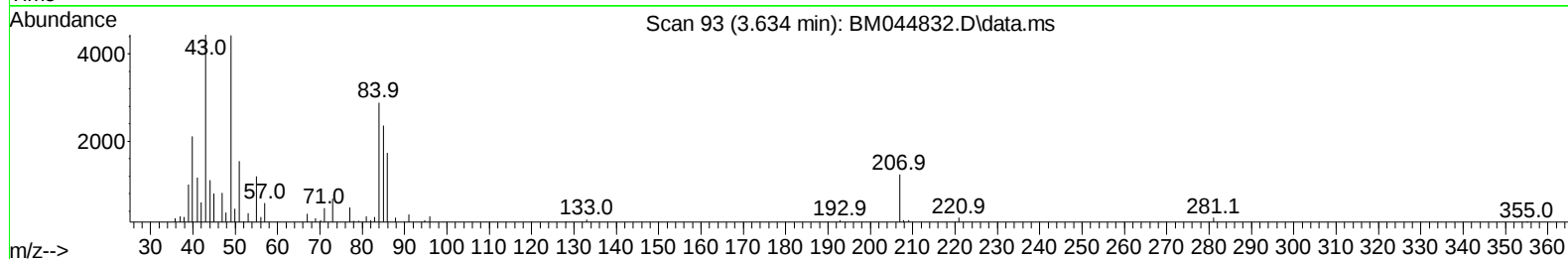
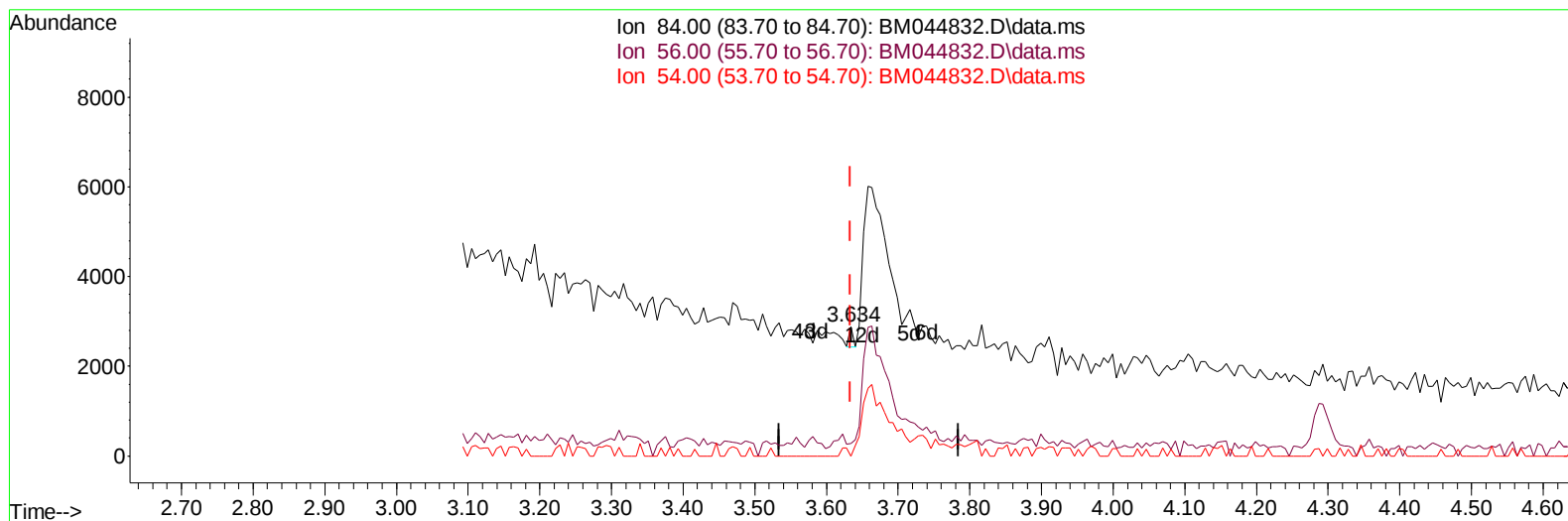
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM033024\
Data File : BM044832.D
Acq On : 30 Mar 2024 15:03
Operator : MA/JU
Sample : P1887-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0798

Manual IntegrationsAPPROVED

Quant Time: Mar 31 15:18:55 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Mar 31 13:54:26 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/01/2024
Supervised By :mohammad ahmed 04/02/2024



TIC: BM044832.D\data.ms

(4) Pyridine-d5 (S)

3.634min (+ 0.000) 0.02 ng/u1

response 162

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	9.53#
54.00	29.60	0.00#
0.00	0.00	0.00

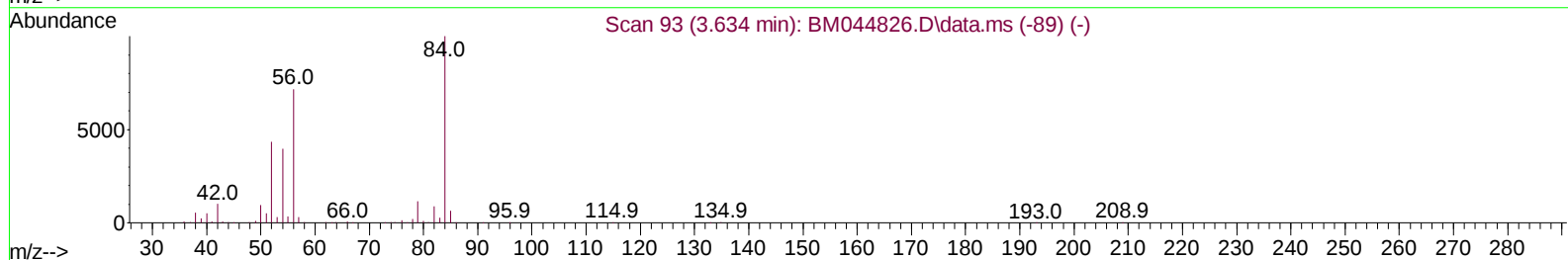
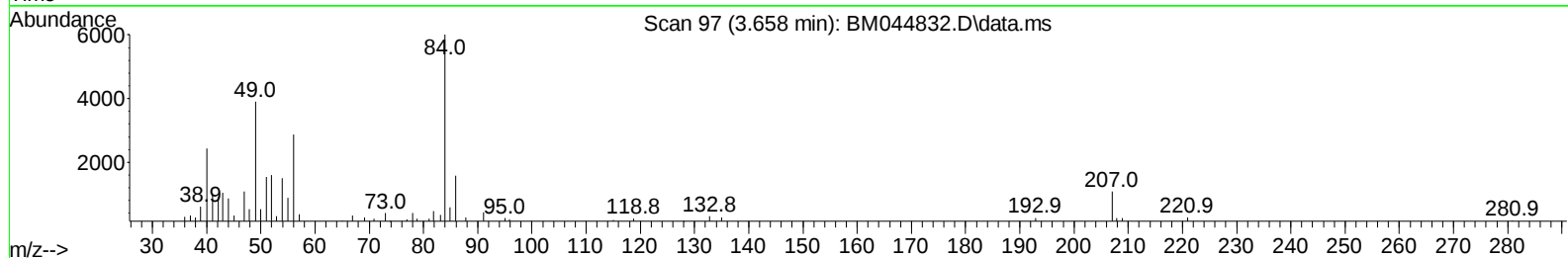
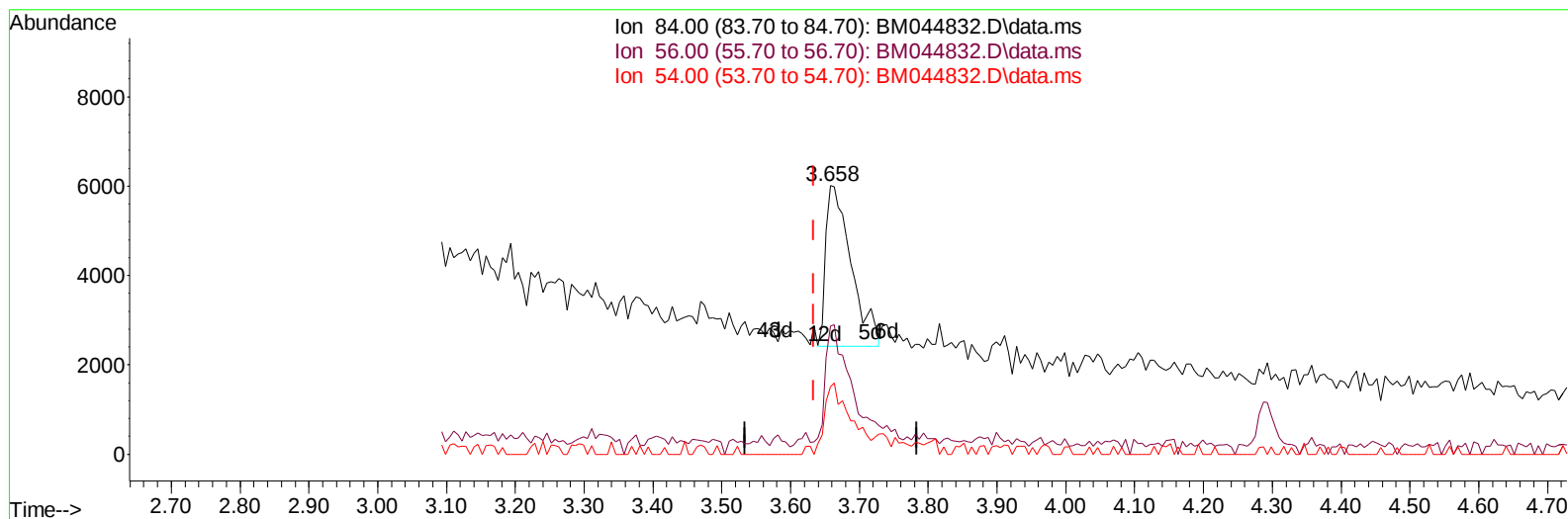
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(4) Pyridine-d5 (S)

3.658min (+ 0.024) 1.22 ng/ul m

response 9177

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	47.64#
54.00	29.60	25.05
0.00	0.00	0.00

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 QLast Update : Sun Mar 31 13:54:26 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.669	152	86447	20.000	ng/ul	0.00
20) Naphthalene-d8	10.445	136	347361	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	199747	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.068	188	388515	20.000	ng/ul	0.00
79) Chrysene-d12	21.268	240	358628	20.000	ng/ul	0.00
88) Perylene-d12	23.503	264	409527	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	5014	1.967	ng/uL	0.00
4) Pyridine-d5	3.658	84	9177m	1.223	ng/ul	0.02
7) Phenol-d5	6.846	99	72823	8.194	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	49151	8.953	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132	59993	9.185	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113	42538	6.094	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128	29499	10.085	ng/ul	0.00
24) 2-Nitrophenol-d4	9.540	143	32382	10.269	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	48611	9.058	ng/ul	0.01
31) 4-Chloroaniline-d4	10.616	131	2466	0.277	ng/ul	0.02
46) Dimethylphthalate-d6	13.733	166	158399	10.109	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	186381	10.401	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	22044	6.784	ng/ul	0.02
60) Fluorene-d10	15.316	176	143323	10.885	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	14839	5.884	ng/ul	0.00
73) Anthracene-d10	17.168	188	229354	12.350	ng/ul	0.00
81) Pyrene-d10	19.468	212	282073	12.824	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.356	264	278676	12.946	ng/ul	0.00

Target Compounds Qvalue

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

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