

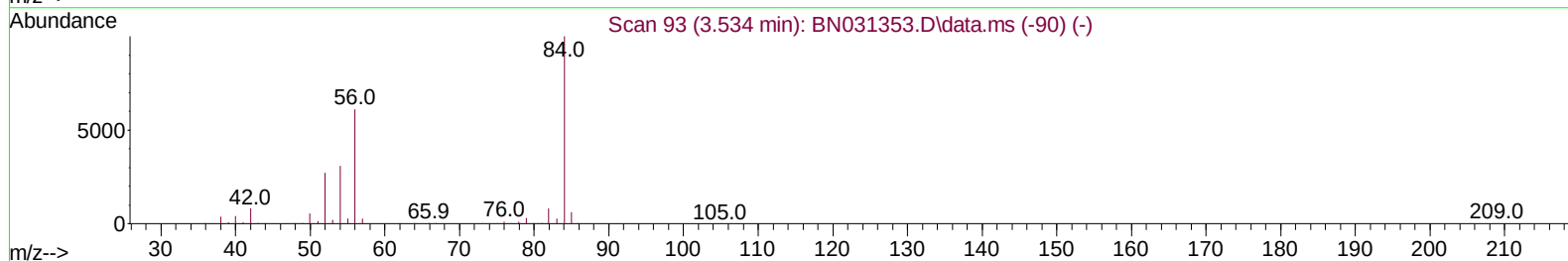
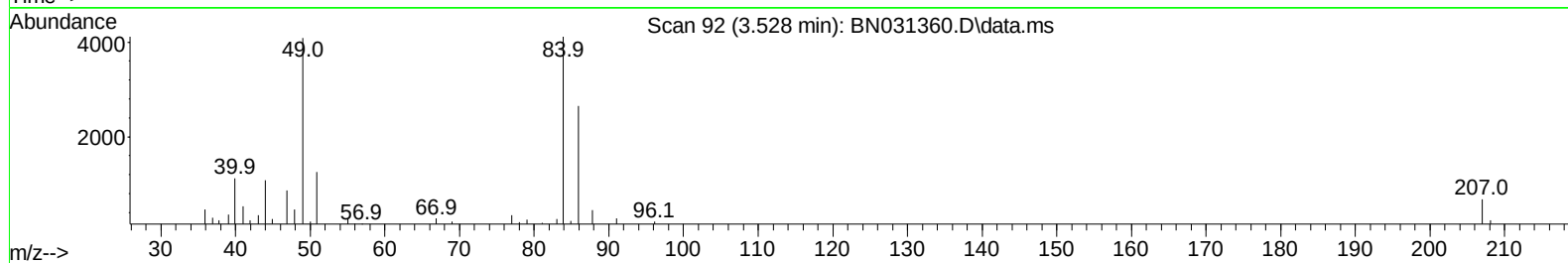
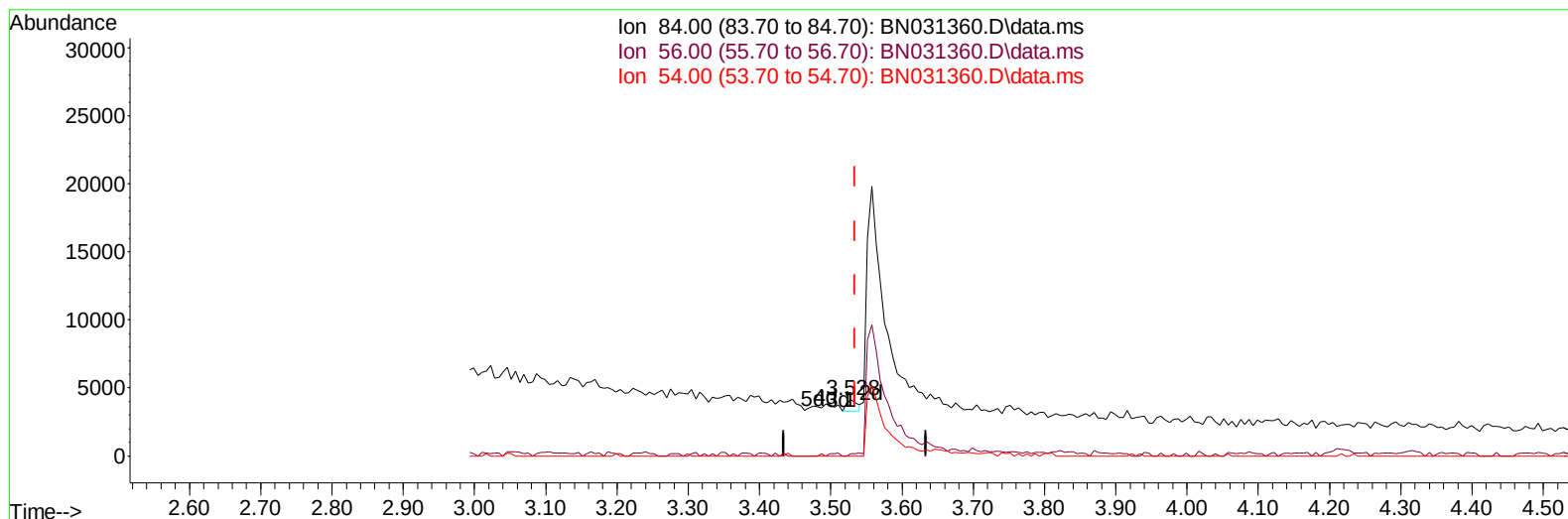
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050124\
Data File : BN031360.D
Acq On : 01 May 2024 23:11
Operator : MA/JU
Sample : P2104-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AA7

Manual IntegrationsAPPROVED

Quant Time: May 02 00:09:20 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 02 00:07:34 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/02/2024
Supervised By :mohammad ahmed 05/02/2024



TIC: BN031360.D\data.ms

(4) Pyridine-d5 (S)

3.528min (-0.006) 0.08 ng/u1

response 863

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.50	4.18#
54.00	30.20	0.00#
0.00	0.00	0.00

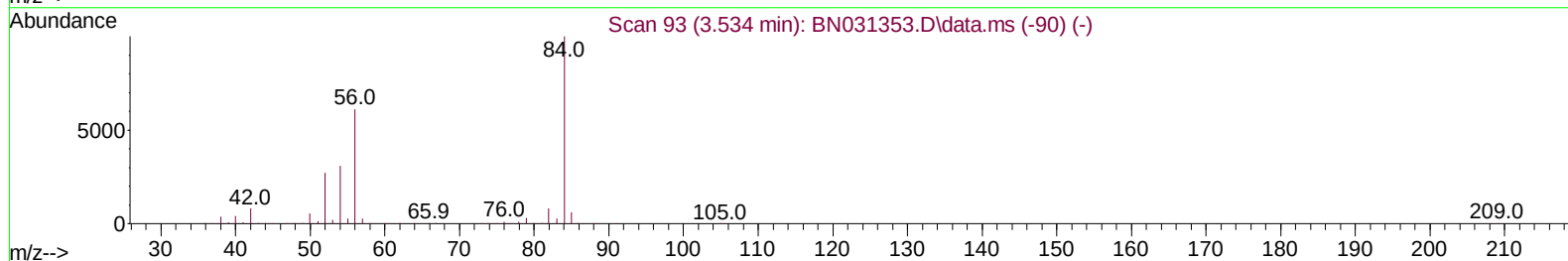
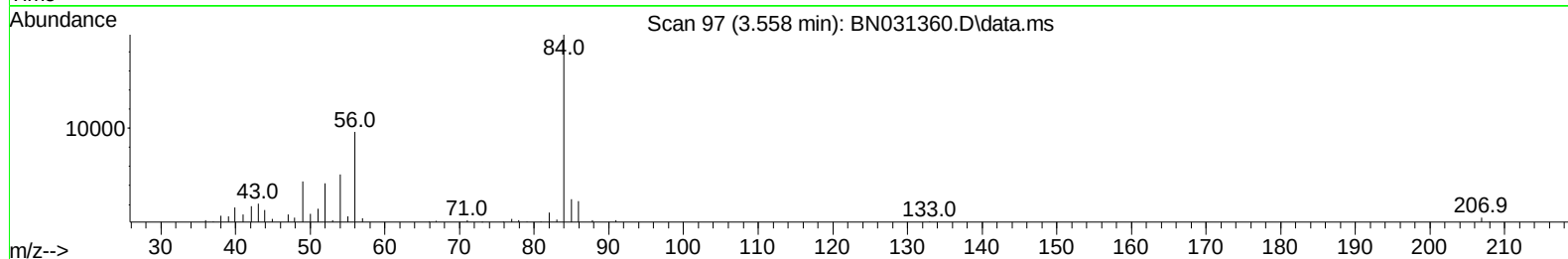
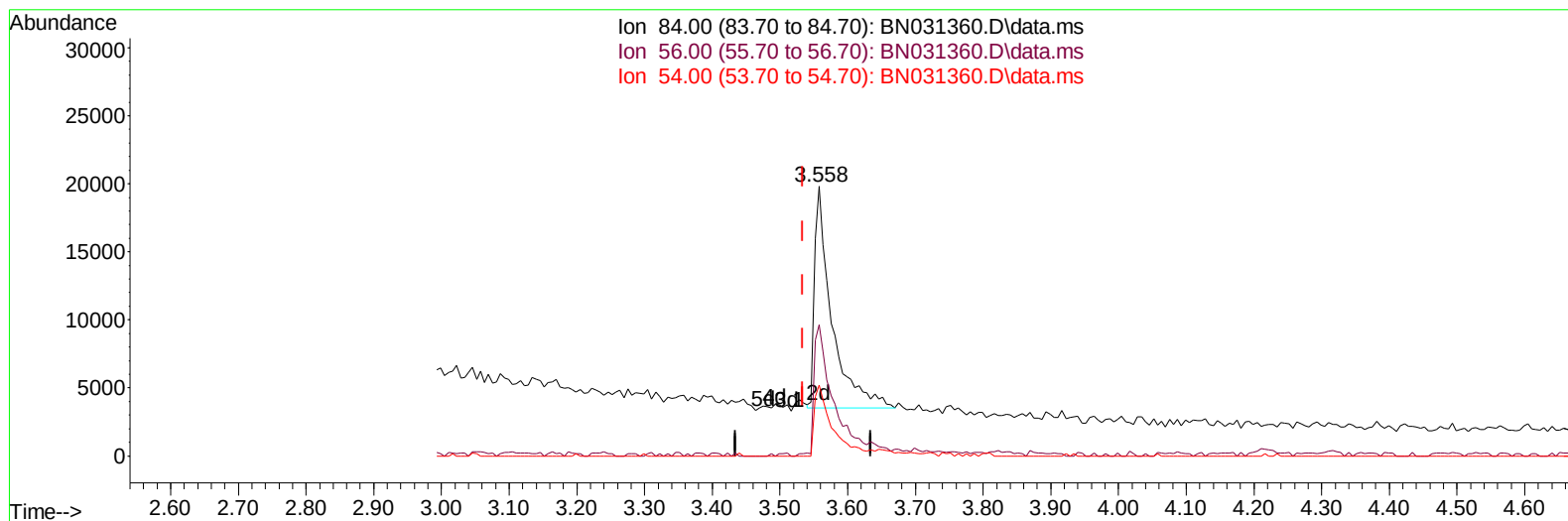
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(4) Pyridine-d5 (S)

3.558min (+ 0.024) 2.53 ng/ul m

response 29002

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.50	48.73#
54.00	30.20	26.22
0.00	0.00	0.00

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 Sample : P2104-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA7

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 05/02/2024
 Supervised By :mohammad ahmed 05/02/2024

Quant Time: May 02 00:30:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 02 00:07:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.610	152	186519	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	627409	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	454116	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.010	188	783596	20.000	ng/ul	0.00
79) Chrysene-d12	21.245	240	826911	20.000	ng/ul	-0.01
88) Perylene-d12	24.127	264	688583	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	13268	3.312	ng/uL	0.00
4) Pyridine-d5	3.558	84	29002m	2.528	ng/ul	0.02
7) Phenol-d5	6.793	99	114181	7.607	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.957	67	213545	24.300	ng/ul	0.00
11) 2-Chlorophenol-d4	7.146	132	110274	9.184	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	113795	9.149	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	156163	34.537	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	142124	29.205	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	73513	8.135	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	106259	8.114	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	921148	30.768	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	368560	10.610	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	48977	9.014	ng/ul	0.00
60) Fluorene-d10	15.257	176	340634	13.324	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	111219	27.600	ng/ul	0.00
73) Anthracene-d10	17.110	188	348629	10.780	ng/ul	0.00
81) Pyrene-d10	19.416	212	483997	9.611	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264	336880	10.568	ng/ul	-0.01

Target Compounds Qvalue

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

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