

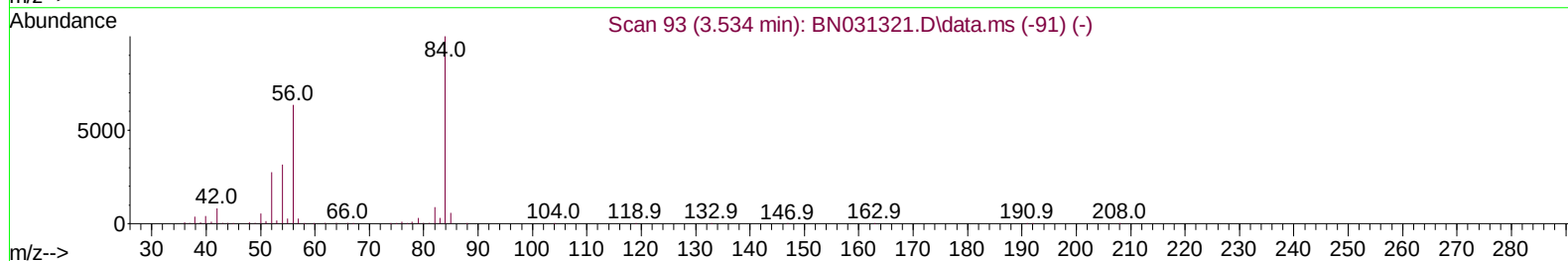
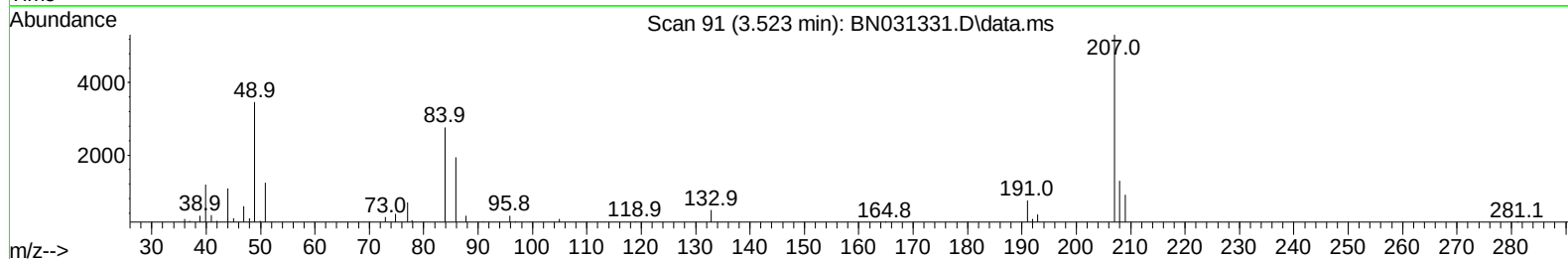
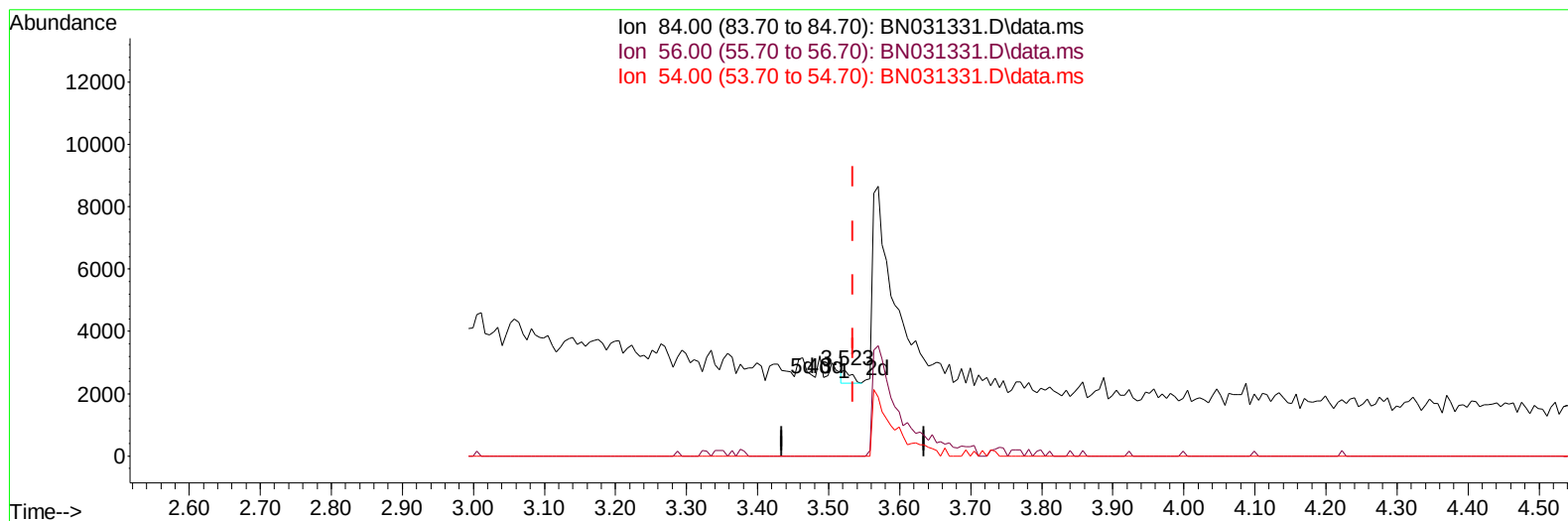
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN043024\
 Data File : BN031331.D
 Acq On : 30 Apr 2024 21:54
 Operator : MA/JU
 Sample : P2269-05DL 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0RQ2DL

Manual IntegrationsAPPROVED

Quant Time: Apr 30 23:33:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 30 23:30:44 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/01/2024
 Supervised By :mohammad ahmed 05/02/2024



TIC: BN031331.D\data.ms

(4) Pyridine-d5 (S)

3.523min (-0.012) 0.04 ng/u1

response 353

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

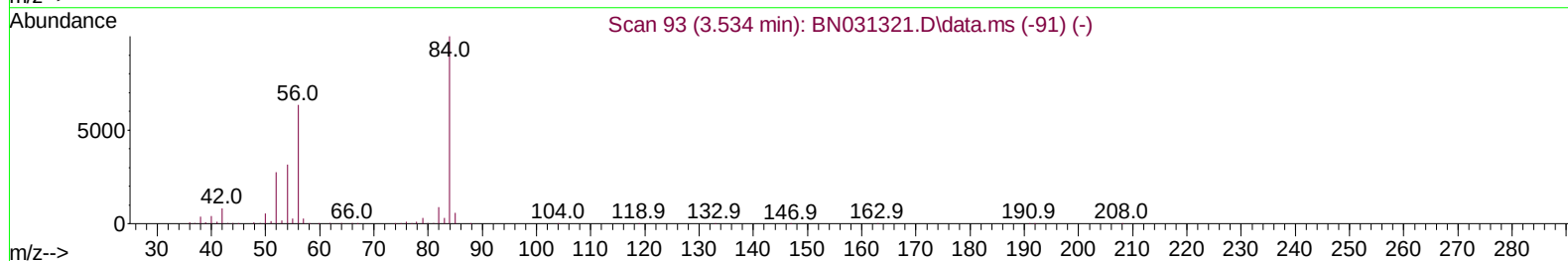
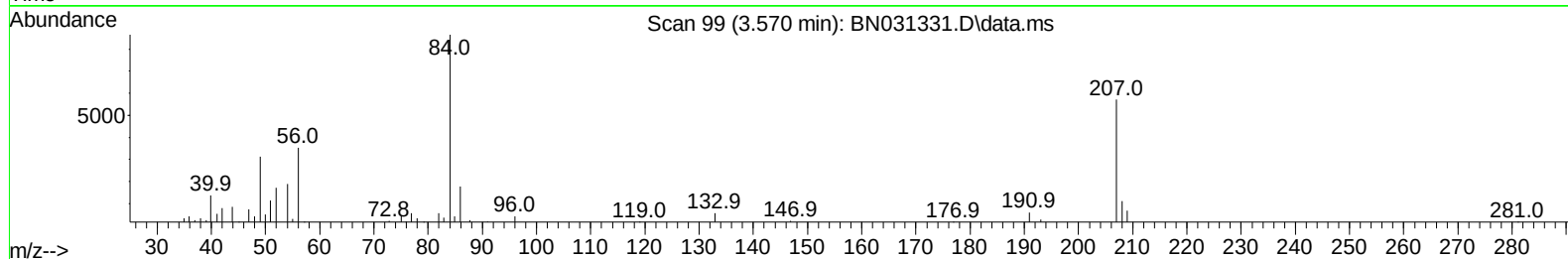
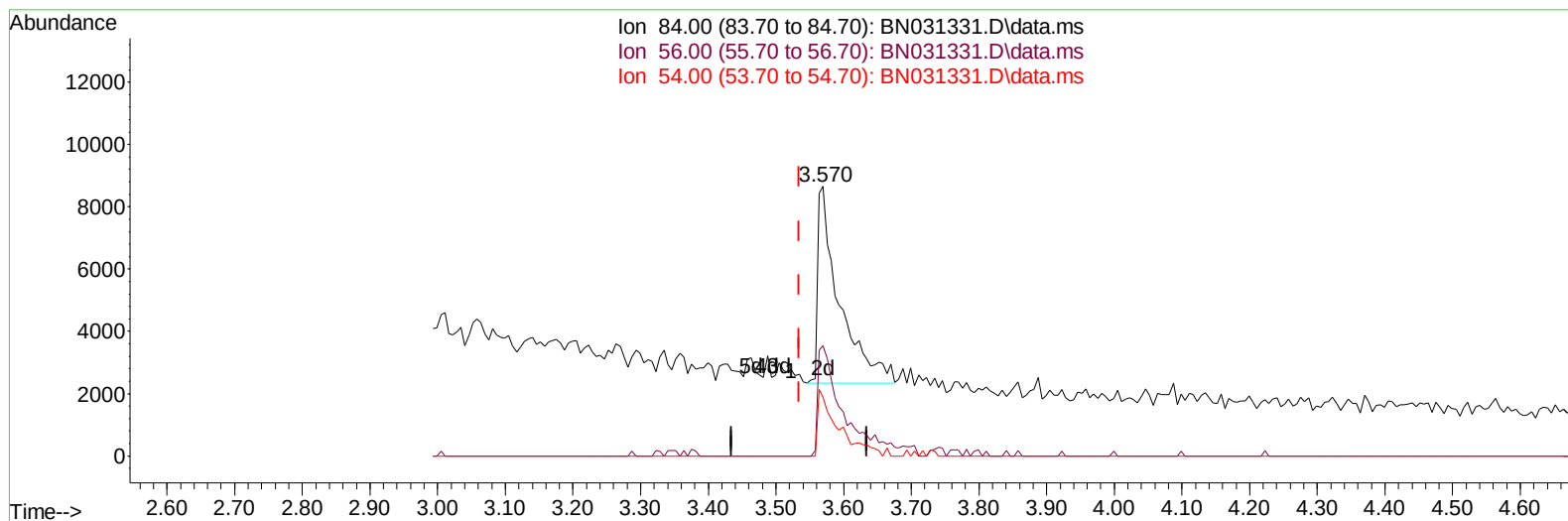
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(4) Pyridine-d5 (S)

3.570min (+ 0.035) 1.63 ng/ul m

response 14108

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.50	40.87#
54.00	30.20	21.90#
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024
 Supervised By :mohammad ahmed 05/02/2024

Quant Time: May 01 00:13:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 30 23:30:44 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	140364	20.000	ng/uL	0.01
20) Naphthalene-d8	10.399	136	656648	20.000	ng/uL	0.00
38) Acenaphthene-d10	14.263	164	451477	20.000	ng/uL	0.00
64) Phenanthrene-d10	17.010	188	1015930	20.000	ng/uL	0.00
79) Chrysene-d12	21.251	240	803881	20.000	ng/uL	0.00
88) Perylene-d12	24.133	264	729725	20.000	ng/uL	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	2774	0.920	ng/uL	0.01
4) Pyridine-d5	3.570	84	14108m	1.634	ng/uL	0.04
7) Phenol-d5	6.793	99	28542	2.527	ng/uL	0.00
9) Bis-(2-Chloroethyl)eth...	6.964	67	80674	12.199	ng/uL	0.00
11) 2-Chlorophenol-d4	7.152	132	84036	9.300	ng/uL	0.00
15) 4-Methylphenol-d8	8.322	113	54584	5.832	ng/uL	0.00
21) Nitrobenzene-d5	8.769	128	49847	10.533	ng/uL	0.00
24) 2-Nitrophenol-d4	9.487	143	49917	9.801	ng/uL	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	94409	9.982	ng/uL	0.00
31) 4-Chloroaniline-d4	10.540	131	84278	6.149	ng/uL	0.00
46) Dimethylphthalate-d6	13.675	166	367450	12.345	ng/uL	0.00
49) Acenaphthylene-d8	13.951	160	393480	11.393	ng/uL	0.00
54) 4-Nitrophenol-d4	14.487	143	8054	1.491	ng/uL	0.00
60) Fluorene-d10	15.257	176	327225	12.874	ng/uL	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	33159	6.347	ng/uL	0.00
73) Anthracene-d10	17.110	188	520450	12.413	ng/uL	0.00
81) Pyrene-d10	19.416	212	593801	12.129	ng/uL	0.00
92) Benzo(a)pyrene-d12	23.939	264	410025	12.137	ng/uL	-0.01
Target Compounds						
					Qvalue	
22) Nitrobenzene	8.811	77	309869	27.905	ng/uL	99
39) 1,2,4,5-Tetrachloroben...	12.434	216	239918	20.589	ng/uL	96
75) 1,2,3,4-Tetrachloroben...	13.046	216	510370	39.661	ng/uL	96
76) Pentachlorobenzene	14.575	250	14867	1.178	ng/uL	98

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

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