

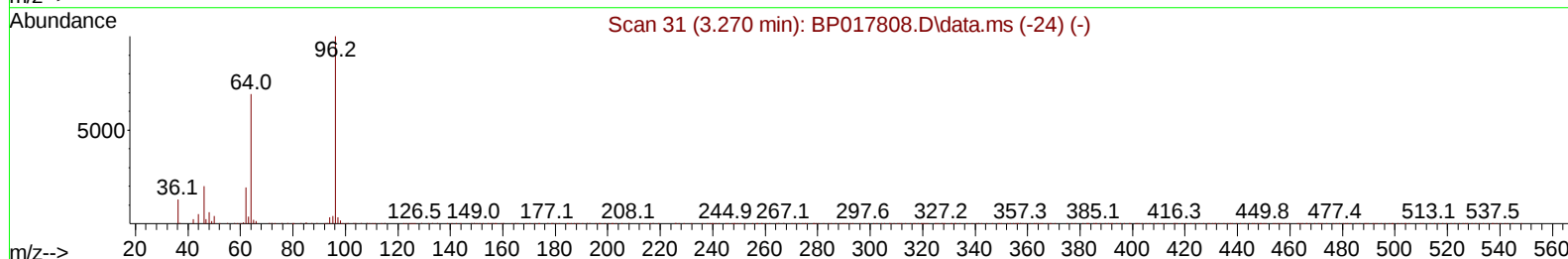
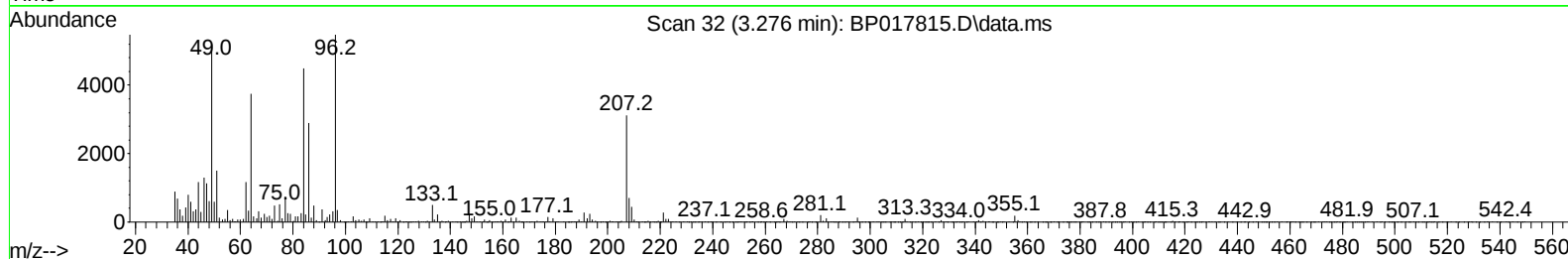
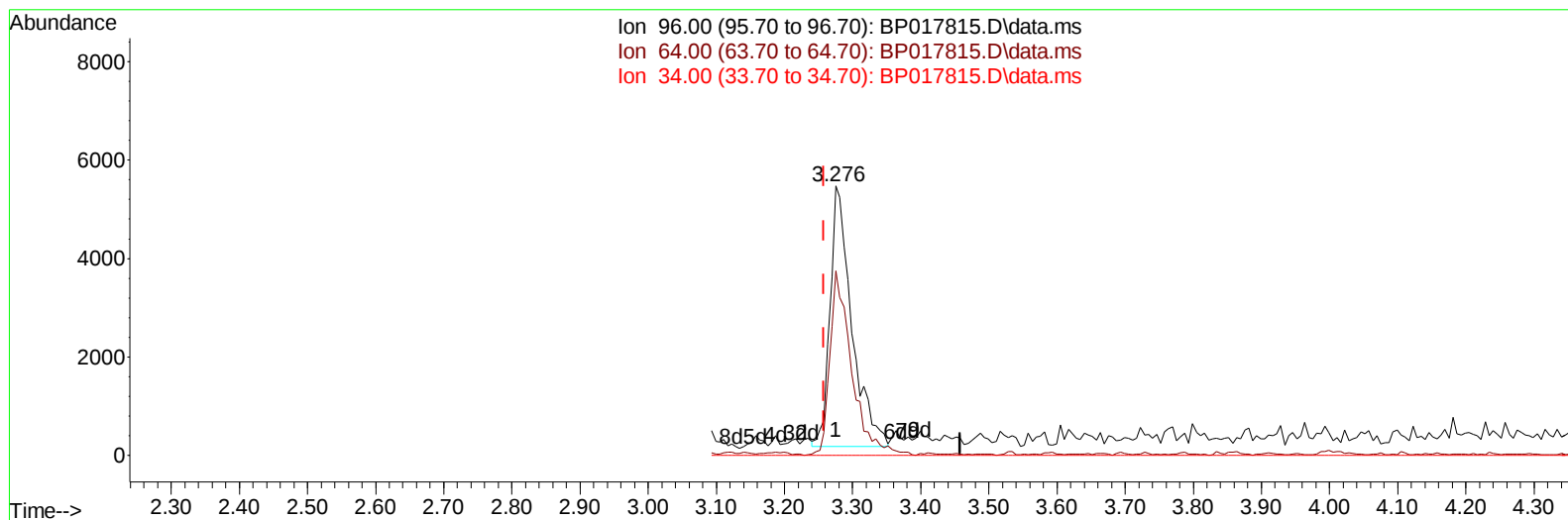
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102523\
 Data File : BP017815.D
 Acq On : 25 Oct 2023 22:40
 Operator : MA/JU
 Sample : 04953-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAS7

Manual IntegrationsAPPROVED

Quant Time: Oct 26 04:58:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/27/2023
 Supervised By :mohammad ahmed 10/27/2023



TIC: BP017815.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.276min (+ 0.018) 3.29 ng/uL m

response 11725

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	71.80	68.57
34.00	0.00	0.00
0.00	0.00	0.00

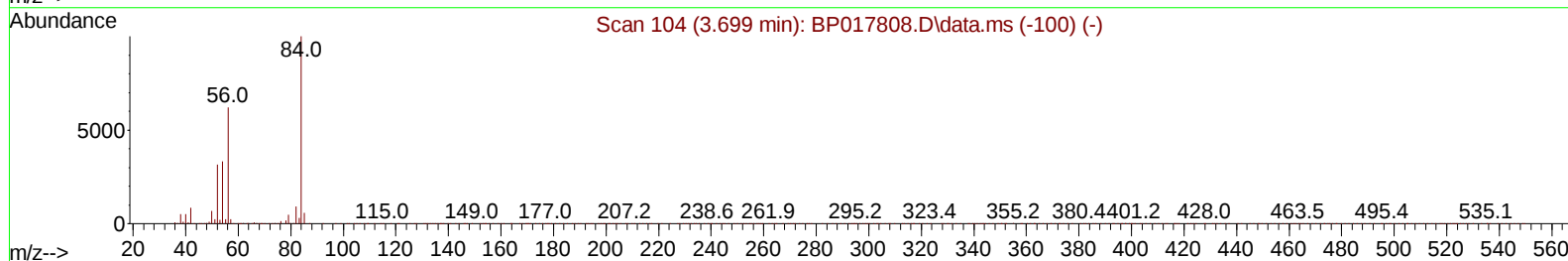
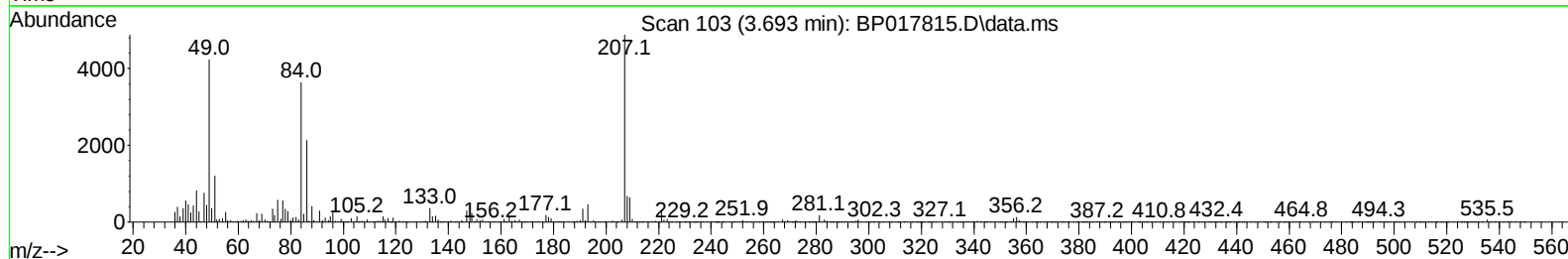
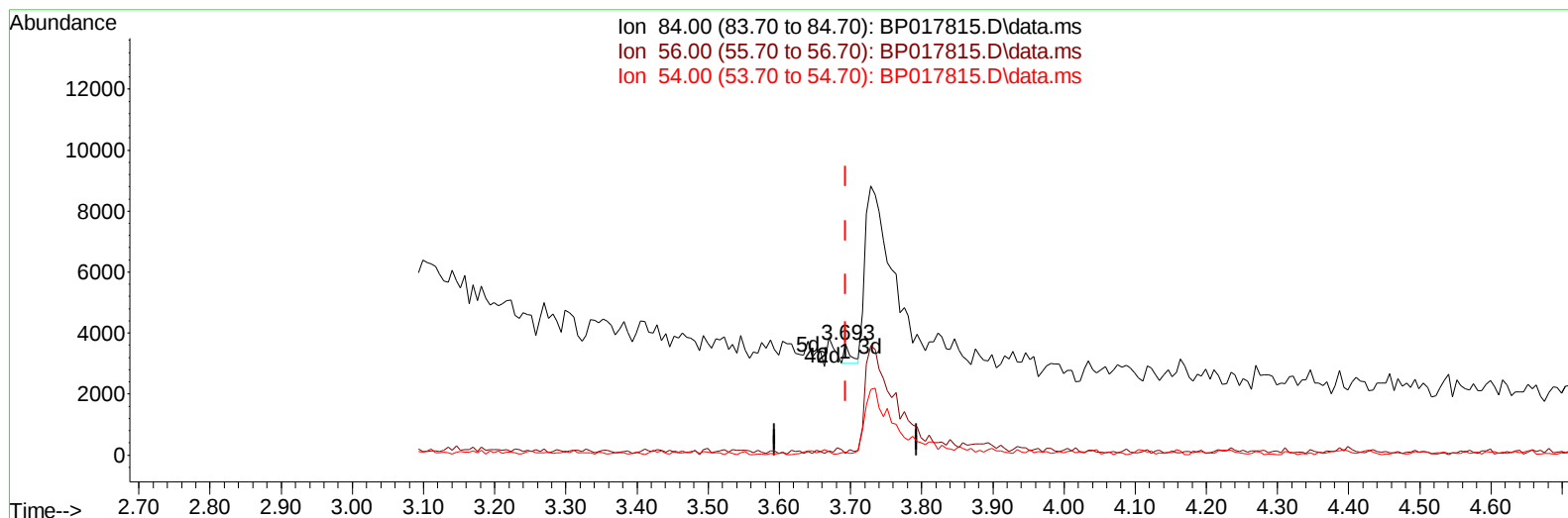
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(4) Pyridine-d5 (S)

3.693min (+ 0.000) 0.05 ng/u1

response 431

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	55.10	1.23#
54.00	30.40	2.63#
0.00	0.00	0.00

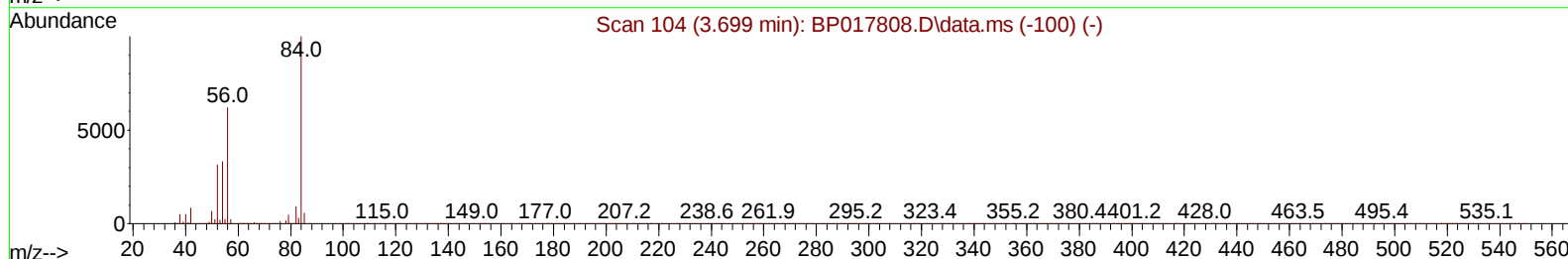
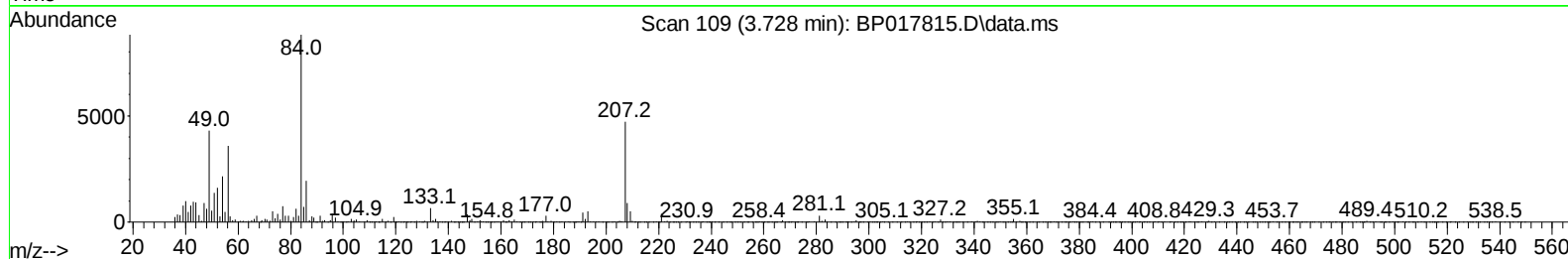
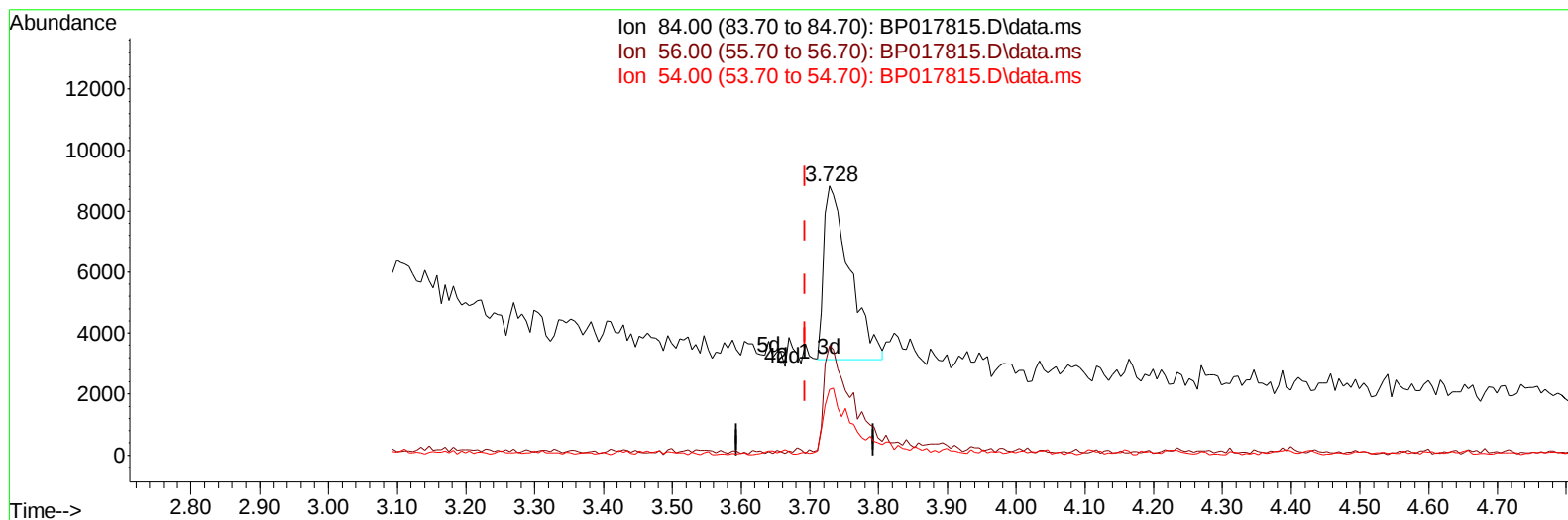
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TIC: BP017815.D\data.ms

(4) Pyridine-d5 (S)

3.728min (+ 0.035) 1.59 ng/u1 m

response 14804

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	55.10	40.70#
54.00	30.40	24.44
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/27/2023
 Supervised By :mohammad ahmed 10/27/2023

Quant Time: Oct 26 05:53:28 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	132684	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	491934	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.575	164	275520	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.369	188	553291	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	482493	20.000	ng/ul	0.00
88) Perylene-d12	24.063	264	506951	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	11725m	3.287	ng/uL	0.02
4) Pyridine-d5	3.728	84	14804m	1.585	ng/ul	0.04
7) Phenol-d5	7.046	99	173950	15.755	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	114833	16.841	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	145615	16.170	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	120872	13.940	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	64932	16.289	ng/ul	0.00
24) 2-Nitrophenol-d4	9.805	143	69137	15.697	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	124963	14.464	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	126197	11.161	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	382872	16.554	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	413770	16.041	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	28875	8.762	ng/ul	0.00
60) Fluorene-d10	15.581	176	322034	16.107	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.734	200	32020	8.786	ng/ul	0.00
73) Anthracene-d10	17.469	188	472446	16.709	ng/ul	0.00
81) Pyrene-d10	19.728	212	575752	19.033	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.892	264	502373	17.792	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.746	105	29999	2.201	ng/ul	91

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

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