

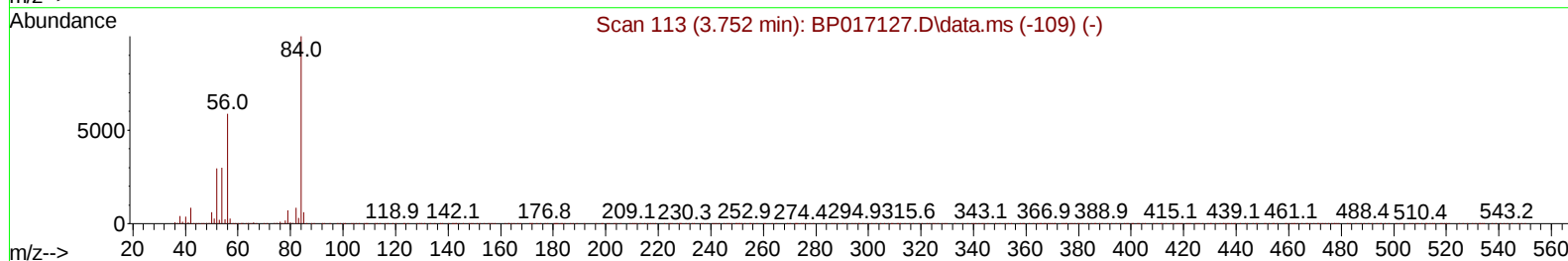
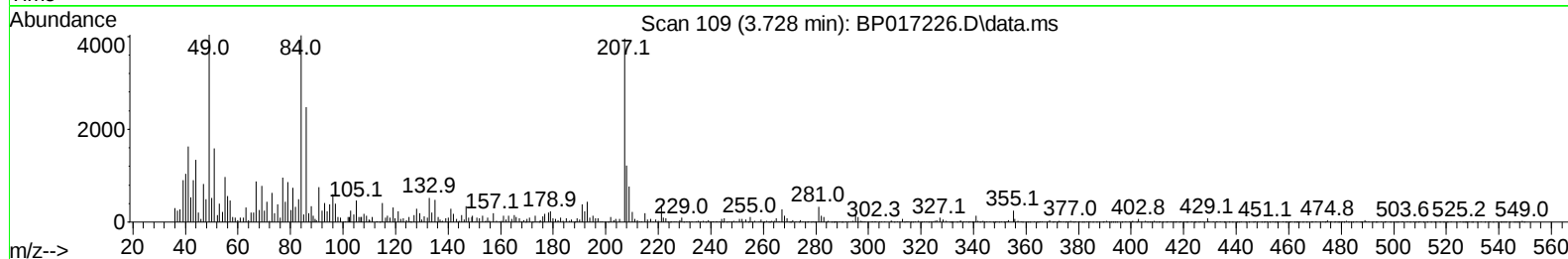
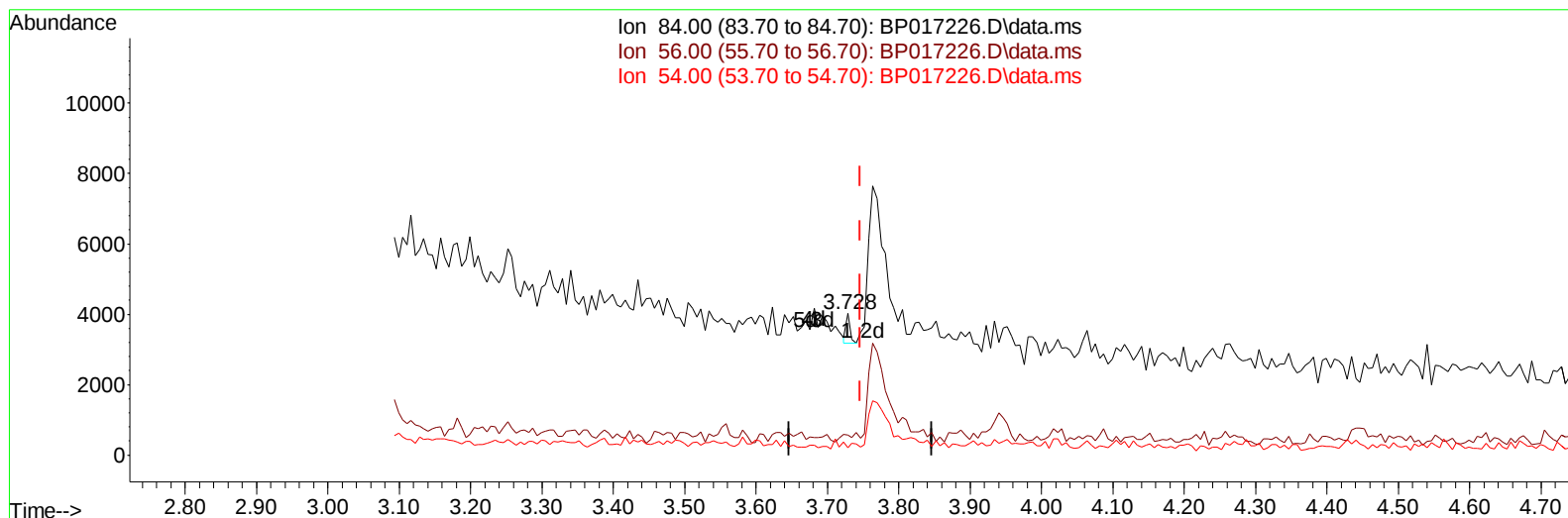
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017226.D
 Acq On : 19 Sep 2023 03:06
 Operator : MA/JU
 Sample : 04332-19
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AE6

Manual IntegrationsAPPROVED

Quant Time: Sep 19 03:36:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 16 03:18:48 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/19/2023
 Supervised By :mohammad ahmed 09/19/2023



TIC: BP017226.D\data.ms

(4) Pyridine-d5 (S)

3.728min (-0.018) 0.02 ng/u1

response 333

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.50	13.94#
54.00	29.40	5.39#
0.00	0.00	0.00

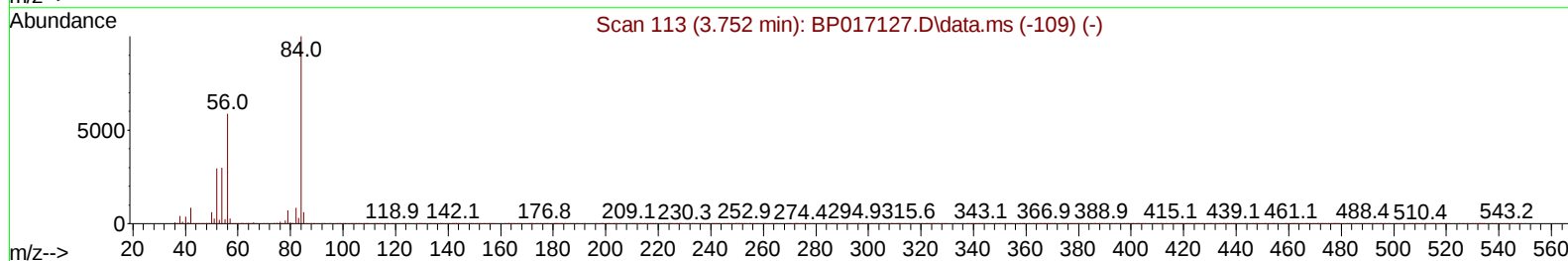
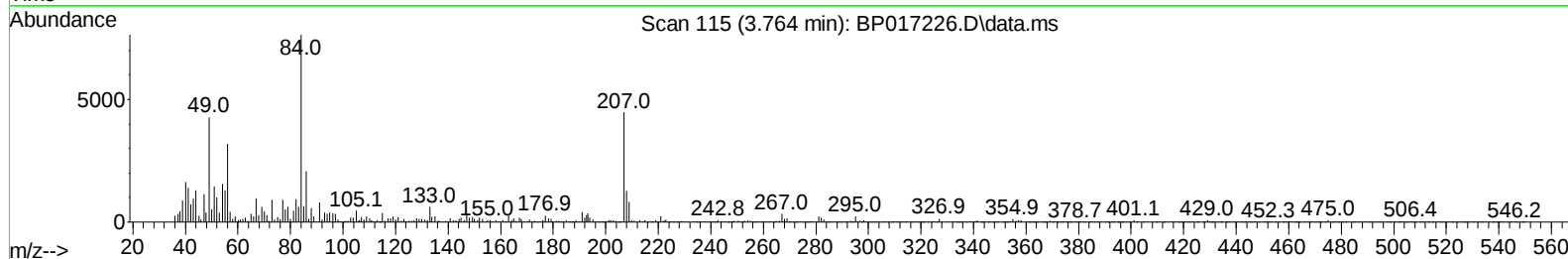
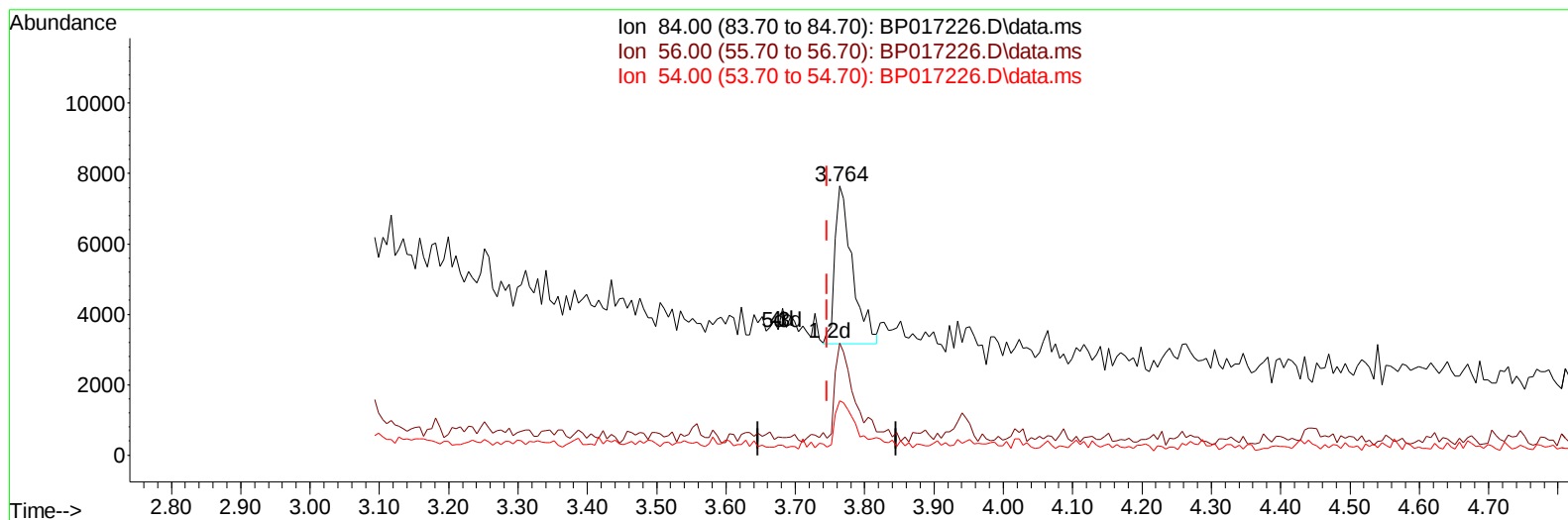
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TIC: BP017226.D\data.ms

(4) Pyridine-d5 (S)

3.764min (+ 0.018) 0.57 ng/ul m

response 7841

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.50	41.59#
54.00	29.40	20.30#
0.00	0.00	0.00

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 Sample : 04332-19
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AE6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/19/2023
 Supervised By :mohammad ahmed 09/19/2023

Quant Time: Sep 19 03:38:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.946	152		200988	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136		767667	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164		410727	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.375	188		845923	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240		840528	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264		1017618	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.311	96		14199	2.969	ng/uL	0.00
4) Pyridine-d5	3.764	84		7841m	0.573	ng/ul	0.02
7) Phenol-d5	7.099	99		247559	15.148	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67		165963	16.247	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132		214368	16.442	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113		168758	12.460	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128		98340	18.294	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143		102250	17.818	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165		175575	15.820	ng/ul	0.00
31) 4-Chloroaniline-d4	10.934	131		142127	8.496	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166		487069	16.664	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160		565726	17.058	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143		47667	8.438	ng/ul	0.00
60) Fluorene-d10	15.598	176		414774	16.412	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200		35003	7.406	ng/ul	0.00
73) Anthracene-d10	17.475	188		607431	16.901	ng/ul	0.00
81) Pyrene-d10	19.716	212		793974	18.012	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264		837813	17.444	ng/ul	0.00
Target Compounds							
16) Acetophenone	8.793	105		50203	2.368	ng/ul	Qvalue 97

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

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