

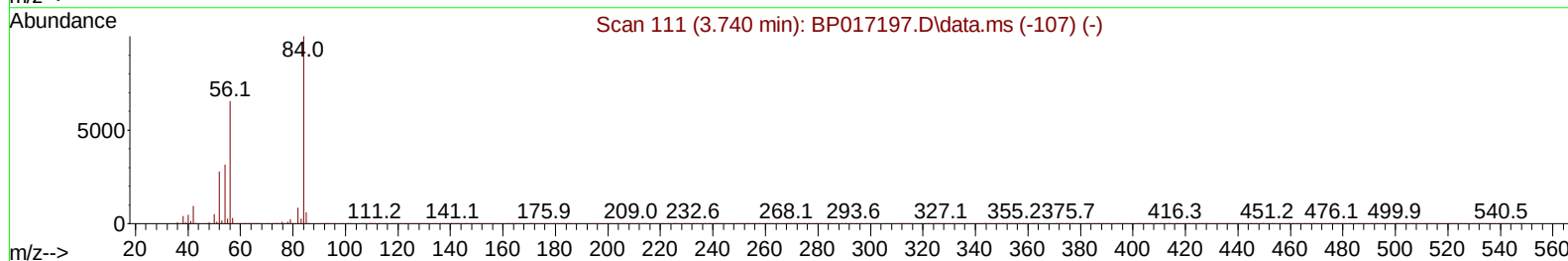
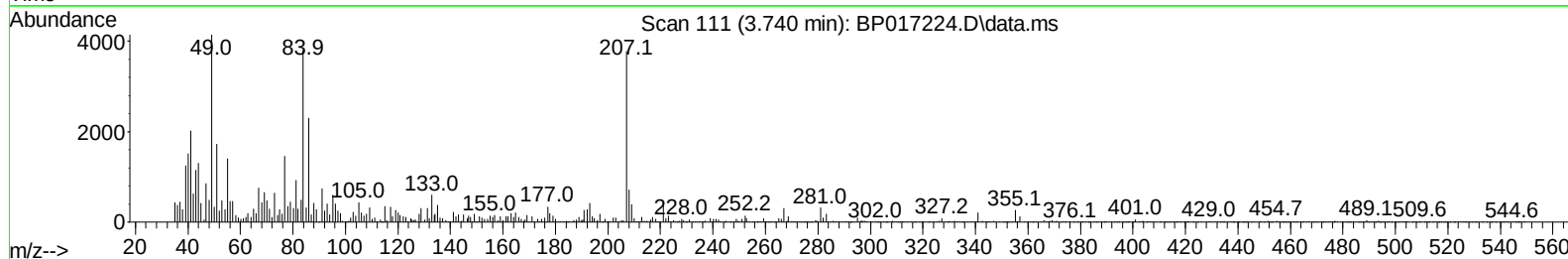
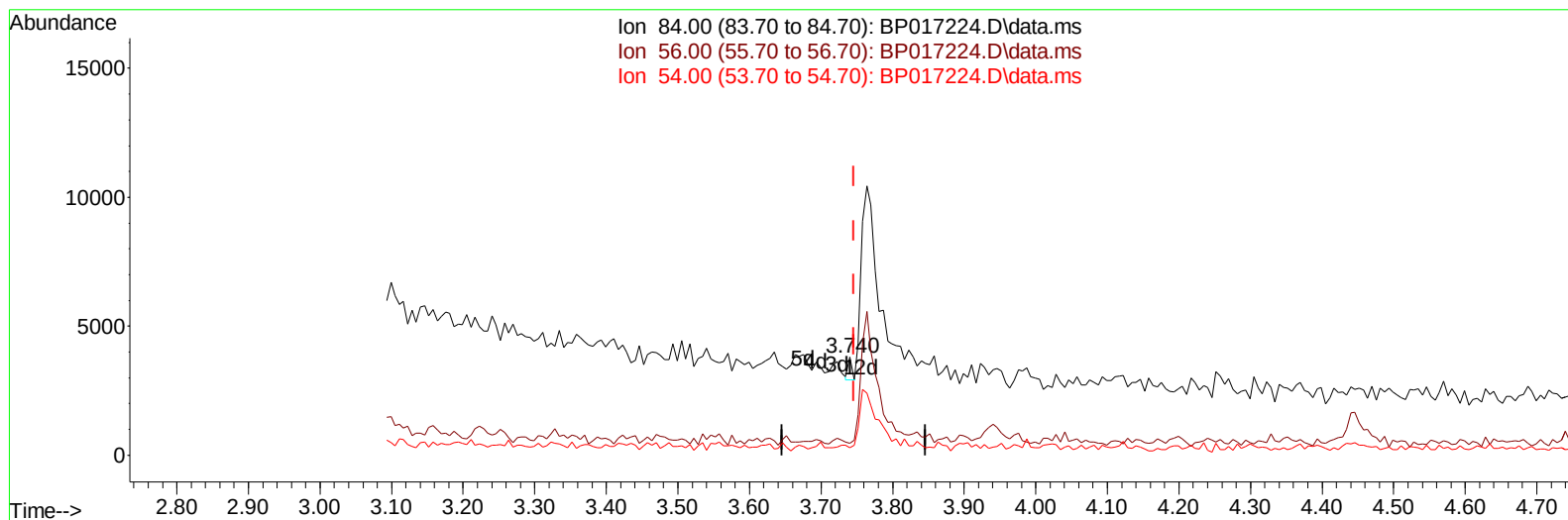
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017224.D
Acq On : 19 Sep 2023 01:58
Operator : MA/JU
Sample : 04332-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AC7

Manual IntegrationsAPPROVED

Quant Time: Sep 19 02:34:12 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: BP017224.D\data.ms

(4) Pyridine-d5 (S)

3.740min (-0.006) 0.02 ng/uI

response 309

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.50	12.16#
54.00	29.40	7.53#
0.00	0.00	0.00

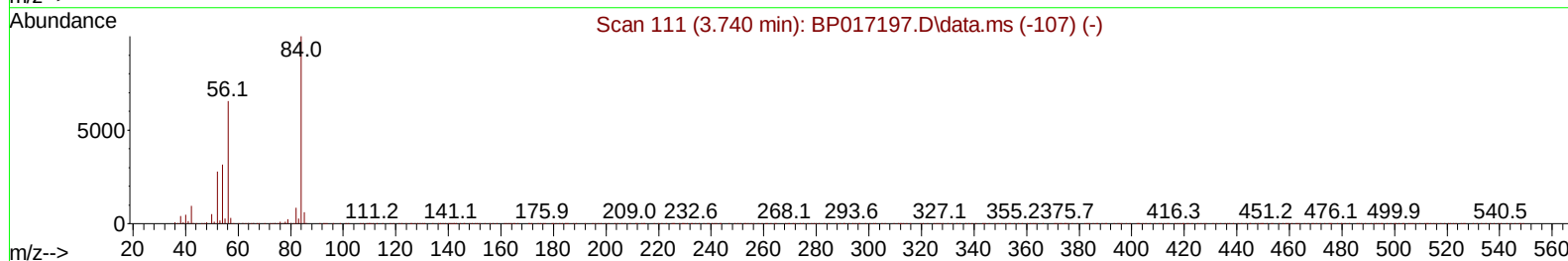
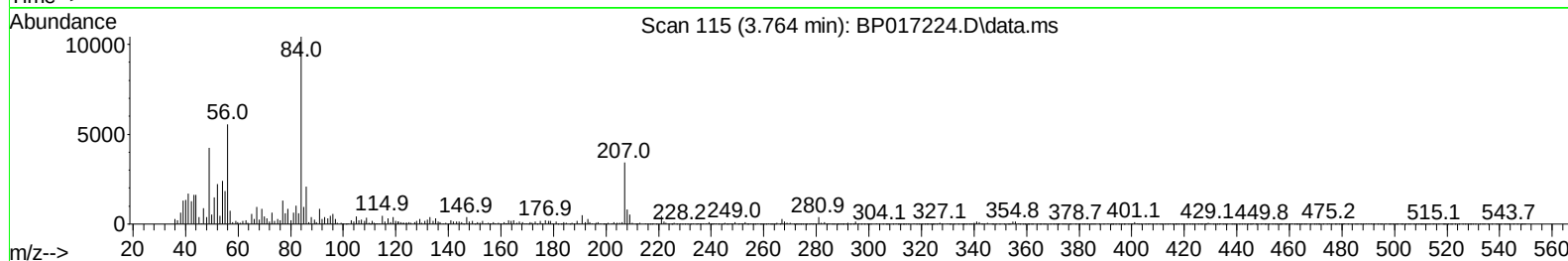
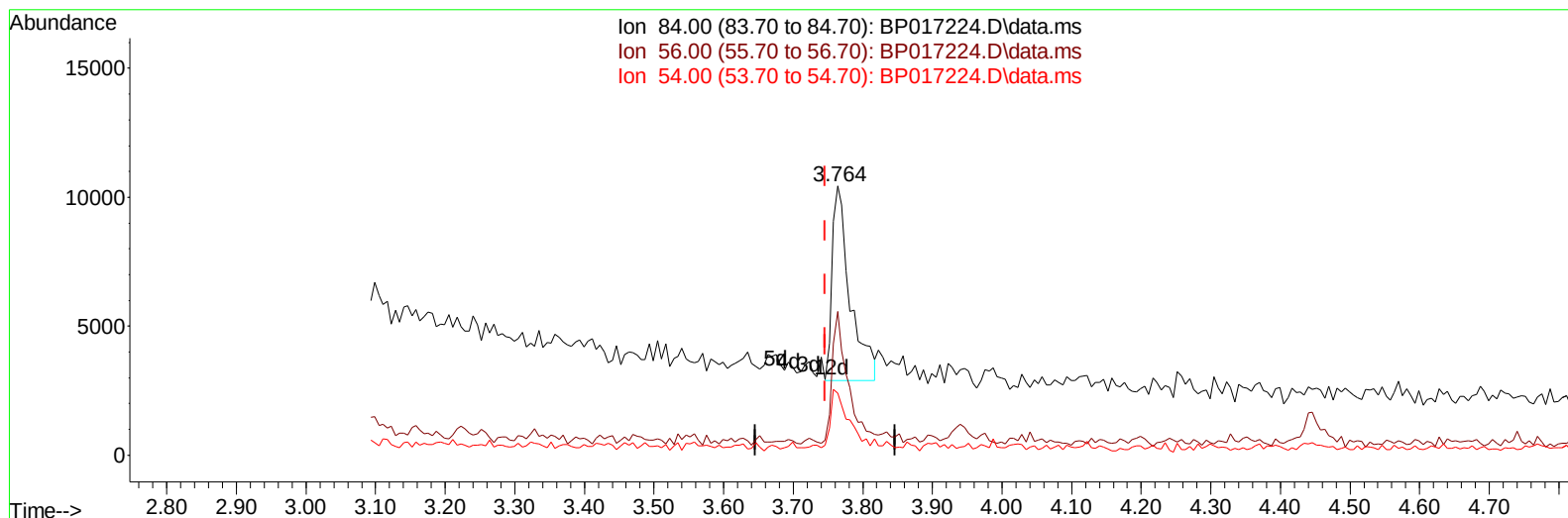
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TIC: BP017224.D\data.ms

(4) Pyridine-d5 (S)

3.764min (+ 0.018) 0.95 ng/ul m

response 13468

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.50	53.31
54.00	29.40	23.13#
0.00	0.00	0.00

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Quant Time: Sep 19 02:39:57 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	209323	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	793718	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	449147	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.375	188	990363	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	976601	20.000	ng/ul	0.00
88) Perylene-d12	24.004	264	1170506	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	15965	3.205	ng/uL	0.00
4) Pyridine-d5	3.764	84	13468m	0.945	ng/ul	0.02
7) Phenol-d5	7.099	99	238058	13.986	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	178302	16.760	ng/ul	0.00
11) 2-Chlorophenol-d4	7.464	132	222905	16.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	88477	6.272	ng/ul	0.00
21) Nitrobenzene-d5	9.146	128	109713	19.740	ng/ul	0.00
24) 2-Nitrophenol-d4	9.858	143	115574	19.478	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	200131	17.441	ng/ul	0.00
31) 4-Chloroaniline-d4	10.934	131	65316	3.776	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	597434	18.692	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	686523	18.930	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	81476	13.189	ng/ul	0.00
60) Fluorene-d10	15.598	176	517777	18.735	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	50403	9.109	ng/ul	0.00
73) Anthracene-d10	17.475	188	833872	19.817	ng/ul	0.00
81) Pyrene-d10	19.716	212	1070109	20.893	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	1082844	19.600	ng/ul	0.00
Target Compounds						
						Qvalue
6) Benzaldehyde	7.105	77	134836	14.229	ng/ul	98
8) Phenol	7.128	94	18323	1.021	ng/ul	96
16) Acetophenone	8.799	105	85844	3.888	ng/ul	98

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

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