

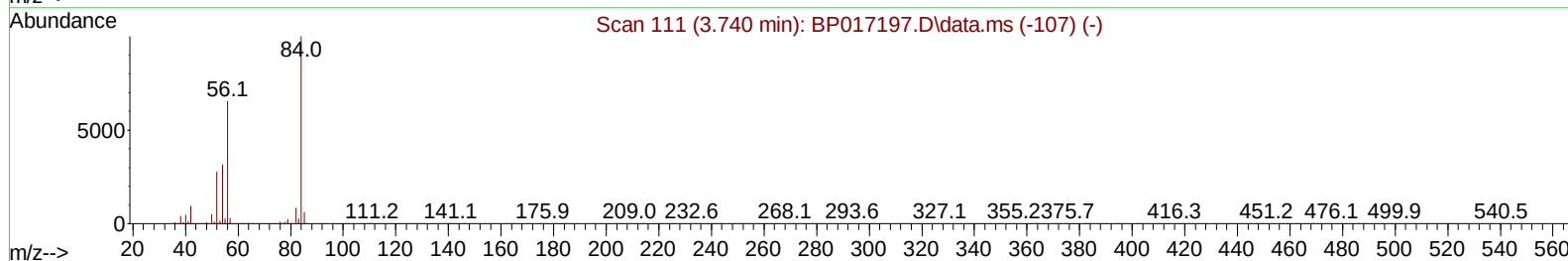
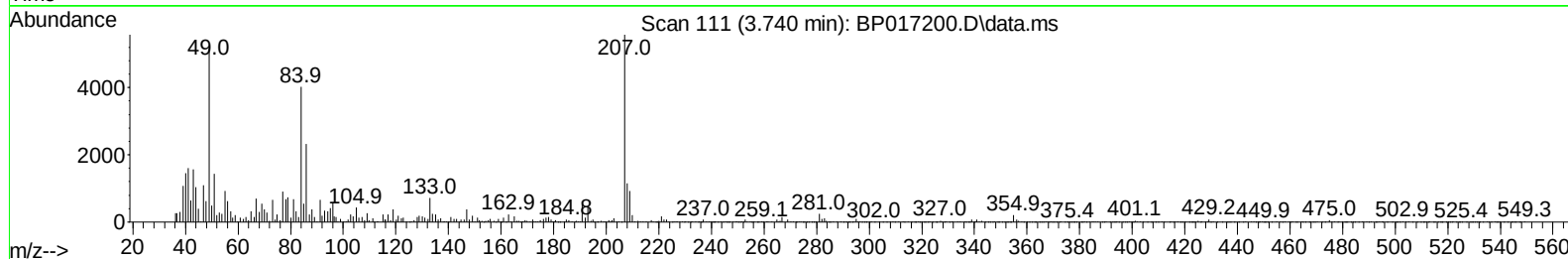
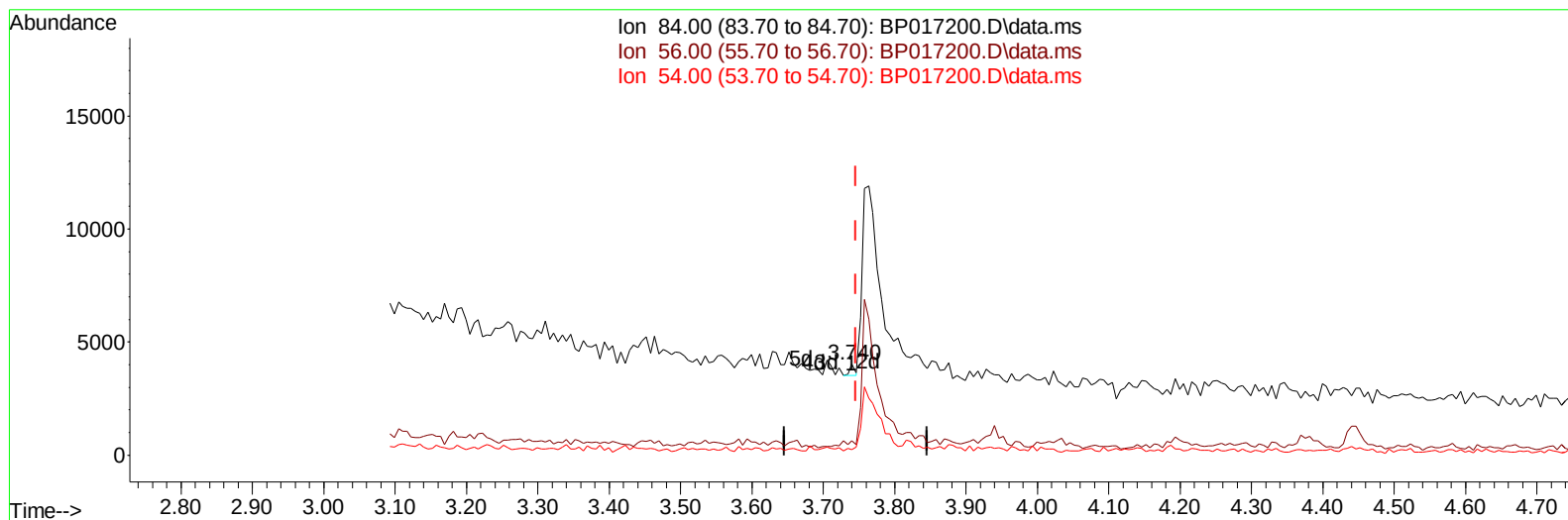
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017200.D
 Acq On : 18 Sep 2023 12:15
 Operator : MA/JU
 Sample : 04330-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYX8

Manual IntegrationsAPPROVED

Quant Time: Sep 18 22:31:03 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: BP017200.D\data.ms

(4) Pyridine-d5 (S)

3.740min (-0.006) 0.02 ng/uI

response 218

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.50	15.51#
54.00	29.40	6.27#
0.00	0.00	0.00

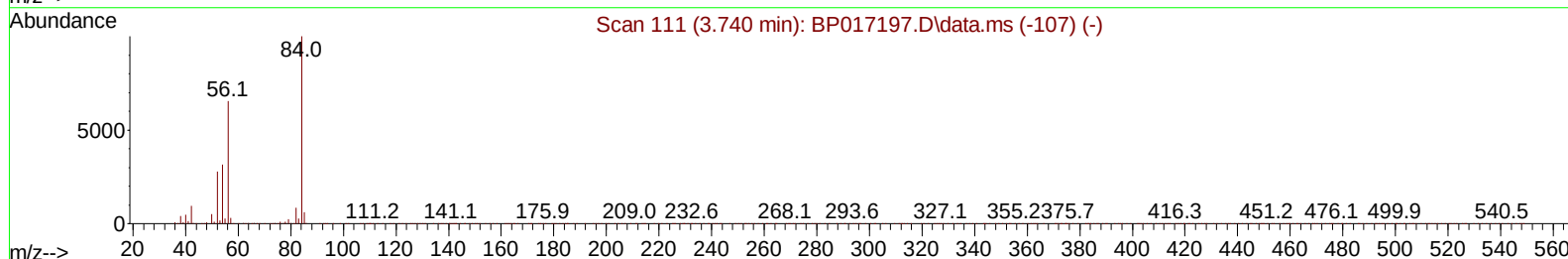
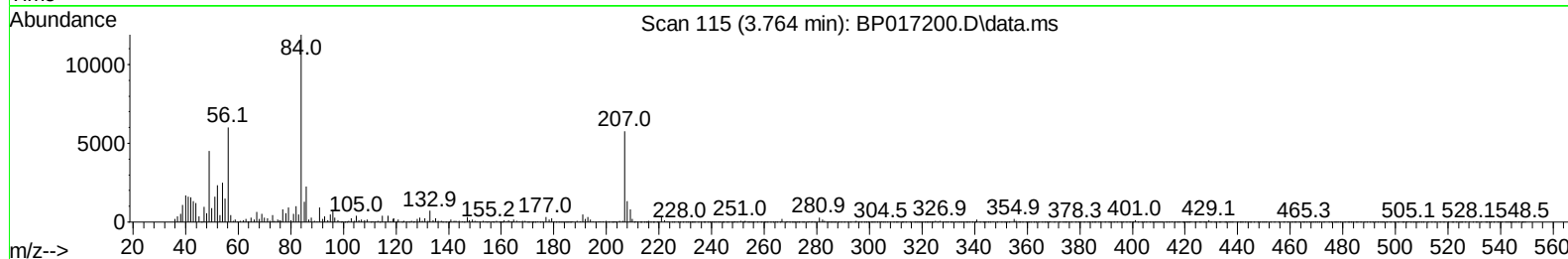
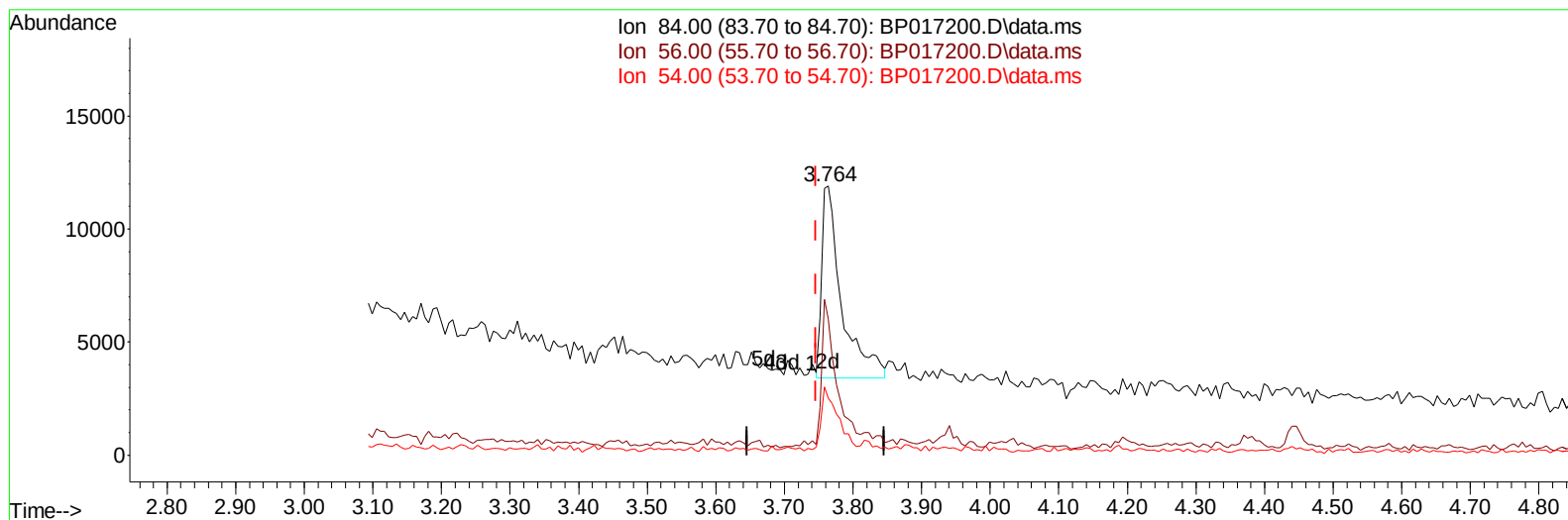
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(4) Pyridine-d5 (S)

3.764min (+ 0.018) 1.27 ng/ul m

response 17100

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.50	50.51
54.00	29.40	21.05#
0.00	0.00	0.00

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Quant Time: Sep 18 22:46:50 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	197075	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	841423	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	510970	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.375	188	1072965	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	729111	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264	837118	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	14350	3.060	ng/uL	0.00
4) Pyridine-d5	3.764	84	17100m	1.275	ng/ul	0.02
7) Phenol-d5	7.093	99	227134	14.174	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	166069	16.581	ng/ul	0.00
11) 2-Chlorophenol-d4	7.464	132	194063	15.180	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	62533	4.709	ng/ul	0.00
21) Nitrobenzene-d5	9.146	128	91333	15.502	ng/ul	0.00
24) 2-Nitrophenol-d4	9.858	143	95945	15.253	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	161600	13.284	ng/ul	0.00
31) 4-Chloroaniline-d4	10.934	131	104967	5.725	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	622410	17.117	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	681389	16.515	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	37513	5.338	ng/ul	0.00
60) Fluorene-d10	15.598	176	568303	18.076	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	31751	5.297	ng/ul	0.00
73) Anthracene-d10	17.475	188	800793	17.566	ng/ul	0.00
81) Pyrene-d10	19.716	212	897560	23.473	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	692169	17.519	ng/ul	-0.01
Target Compounds						
						Qvalue
16) Acetophenone	8.793	105	63835	3.071	ng/ul	98
36) 2-Methylnaphthalene	12.428	142	28781	1.002	ng/ul	91
72) Phenanthrene	17.416	178	60042	1.114	ng/ul	98
80) Fluoranthene	19.386	202	45993	1.029	ng/ul	98

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

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